

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
 Data File : BN034340.D  
 Acq On : 02 Oct 2024 16:57  
 Operator : RC/JU  
 Sample : P4016-18  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 DDC23

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
 Title : SVOA CALIBRATION

Signal : TIC: BN034340.D\data.ms

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 4.593       | 282           | 290         | 297          | rVB      | 3750078        | 6051127       | 11.80%          | 0.413%        |
| 2         | 5.463       | 430           | 438         | 445          | rVB2     | 13194572       | 26836374      | 52.32%          | 1.830%        |
| 3         | 5.704       | 472           | 479         | 486          | rBV2     | 3573797        | 7448924       | 14.52%          | 0.508%        |
| 4         | 5.834       | 486           | 501         | 506          | rVB      | 6688652        | 20984512      | 40.91%          | 1.431%        |
| 5         | 5.916       | 509           | 515         | 523          | rBV3     | 4250657        | 8950849       | 17.45%          | 0.610%        |
| 6         | 5.993       | 523           | 528         | 532          | rBV      | 6006522        | 9189079       | 17.91%          | 0.627%        |
| 7         | 6.057       | 532           | 539         | 543          | rBV3     | 4673207        | 10925568      | 21.30%          | 0.745%        |
| 8         | 6.098       | 543           | 546         | 551          | rVB      | 4403783        | 6499505       | 12.67%          | 0.443%        |
| 9         | 6.322       | 576           | 584         | 589          | rBV4     | 3174713        | 7962342       | 15.52%          | 0.543%        |
| 10        | 6.387       | 589           | 595         | 598          | rBV2     | 5186383        | 10559778      | 20.59%          | 0.720%        |
| 11        | 6.422       | 598           | 601         | 606          | rVB      | 7237174        | 10941058      | 21.33%          | 0.746%        |
| 12        | 6.481       | 606           | 611         | 614          | rBV      | 9336184        | 17136354      | 33.41%          | 1.168%        |
| 13        | 6.510       | 614           | 616         | 620          | rVB      | 6947850        | 8241203       | 16.07%          | 0.562%        |
| 14        | 6.587       | 620           | 629         | 634          | rVB3     | 9206762        | 23442104      | 45.70%          | 1.598%        |
| 15        | 6.645       | 634           | 639         | 644          | rVB      | 6594591        | 10424782      | 20.32%          | 0.711%        |
| 16        | 6.798       | 659           | 665         | 669          | rBV      | 7499080        | 12195146      | 23.78%          | 0.832%        |
| 17        | 6.857       | 669           | 675         | 679          | rBV2     | 7448505        | 12274140      | 23.93%          | 0.837%        |
| 18        | 6.945       | 685           | 690         | 701          | rVB4     | 3567263        | 11307040      | 22.04%          | 0.771%        |
| 19        | 7.057       | 701           | 709         | 717          | rBV3     | 16950865       | 46871835      | 91.38%          | 3.196%        |
| 20        | 7.245       | 733           | 741         | 744          | rBV3     | 2530357        | 6684187       | 13.03%          | 0.456%        |
| 21        | 7.398       | 760           | 767         | 771          | rVB      | 10073633       | 18114594      | 35.32%          | 1.235%        |
| 22        | 7.498       | 777           | 784         | 795          | rVB4     | 10756023       | 29546770      | 57.60%          | 2.015%        |
| 23        | 7.628       | 796           | 806         | 810          | rBV6     | 5169601        | 15262959      | 29.76%          | 1.041%        |
| 24        | 7.681       | 810           | 815         | 823          | rVB7     | 5376903        | 15742924      | 30.69%          | 1.073%        |
| 25        | 7.834       | 836           | 841         | 846          | rBV2     | 4295937        | 7480652       | 14.58%          | 0.510%        |
| 26        | 7.887       | 846           | 850         | 854          | rBV2     | 5313610        | 8310124       | 16.20%          | 0.567%        |
| 27        | 7.945       | 854           | 860         | 865          | rVB2     | 11918343       | 22055153      | 43.00%          | 1.504%        |
| 28        | 7.998       | 865           | 869         | 872          | rBV      | 7329582        | 11457576      | 22.34%          | 0.781%        |
| 29        | 8.039       | 872           | 876         | 879          | rBV2     | 7033591        | 10585625      | 20.64%          | 0.722%        |
| 30        | 8.075       | 879           | 882         | 893          | rVB2     | 10039222       | 17081499      | 33.30%          | 1.165%        |
| 31        | 8.175       | 893           | 899         | 902          | rBV      | 9302274        | 18037676      | 35.17%          | 1.230%        |
| 32        | 8.345       | 923           | 928         | 932          | rBV      | 5839946        | 9435421       | 18.40%          | 0.643%        |
| 33        | 8.398       | 932           | 937         | 943          | rVB      | 6909382        | 11128992      | 21.70%          | 0.759%        |
| 34        | 8.498       | 943           | 954         | 963          | rBV2     | 8767235        | 23724927      | 46.25%          | 1.618%        |
| 35        | 8.651       | 971           | 980         | 984          | rVB      | 15289518       | 42131687      | 82.14%          | 2.873%        |
| 36        | 8.745       | 984           | 996         | 1002         | rBV7     | 6528310        | 18927040      | 36.90%          | 1.291%        |

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Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Title : SVOA CALIBRATION

|    |        |      |      |      |      |          |          |         |        |
|----|--------|------|------|------|------|----------|----------|---------|--------|
| 37 | 8.828  | 1002 | 1010 | 1013 | rBV3 | 3992526  | 8124835  | 15.84%  | 0.554% |
| 38 | 9.075  | 1048 | 1052 | 1063 | rVB3 | 5381257  | 13015264 | 25.37%  | 0.887% |
| 39 | 9.186  | 1063 | 1071 | 1077 | rVB2 | 4610847  | 9020802  | 17.59%  | 0.615% |
| 40 | 9.269  | 1077 | 1085 | 1088 | rBV4 | 3426682  | 9345852  | 18.22%  | 0.637% |
| 41 | 9.328  | 1088 | 1095 | 1099 | rBV4 | 4237574  | 9487102  | 18.50%  | 0.647% |
| 42 | 9.475  | 1110 | 1120 | 1125 | rVB2 | 6168906  | 14837421 | 28.93%  | 1.012% |
| 43 | 9.534  | 1125 | 1130 | 1134 | rBV2 | 4724657  | 8392611  | 16.36%  | 0.572% |
| 44 | 9.616  | 1140 | 1144 | 1147 | rBV  | 4942438  | 7532902  | 14.69%  | 0.514% |
| 45 | 9.716  | 1153 | 1161 | 1166 | rBV4 | 5967070  | 12185556 | 23.76%  | 0.831% |
| 46 | 9.869  | 1181 | 1187 | 1192 | rBV2 | 6634540  | 10718345 | 20.90%  | 0.731% |
| 47 | 10.175 | 1226 | 1239 | 1248 | rBV6 | 13186377 | 51293244 | 100.00% | 3.497% |
| 48 | 10.304 | 1254 | 1261 | 1265 | rBV3 | 9958960  | 17461841 | 34.04%  | 1.191% |
| 49 | 10.345 | 1265 | 1268 | 1273 | rVB  | 4660949  | 6447513  | 12.57%  | 0.440% |
| 50 | 10.563 | 1298 | 1305 | 1316 | rVB  | 14278906 | 28182960 | 54.94%  | 1.922% |
| 51 | 10.692 | 1317 | 1327 | 1332 | rBV  | 7921266  | 16997738 | 33.14%  | 1.159% |
| 52 | 11.210 | 1408 | 1415 | 1420 | rBV2 | 10363926 | 20680736 | 40.32%  | 1.410% |
| 53 | 11.275 | 1420 | 1426 | 1432 | rVB2 | 9778236  | 19513142 | 38.04%  | 1.331% |
| 54 | 11.451 | 1446 | 1456 | 1461 | rBV  | 9404287  | 17447122 | 34.01%  | 1.190% |
| 55 | 11.628 | 1475 | 1486 | 1489 | rBV  | 12572085 | 29795485 | 58.09%  | 2.032% |
| 56 | 11.657 | 1489 | 1491 | 1495 | rVV  | 9497236  | 11914673 | 23.23%  | 0.812% |
| 57 | 11.833 | 1516 | 1521 | 1524 | rVV  | 7627764  | 13836012 | 26.97%  | 0.943% |
| 58 | 11.875 | 1524 | 1528 | 1533 | rVB  | 11458862 | 18413831 | 35.90%  | 1.256% |
| 59 | 12.033 | 1549 | 1555 | 1564 | rBV3 | 5865889  | 16728138 | 32.61%  | 1.141% |
| 60 | 12.280 | 1591 | 1597 | 1600 | rBV2 | 3683278  | 7066661  | 13.78%  | 0.482% |
| 61 | 12.451 | 1622 | 1626 | 1634 | rVB2 | 4046706  | 7287730  | 14.21%  | 0.497% |
| 62 | 12.598 | 1646 | 1651 | 1660 | rVB2 | 5546049  | 10729365 | 20.92%  | 0.732% |
| 63 | 12.810 | 1678 | 1687 | 1692 | rBV4 | 2292899  | 6120334  | 11.93%  | 0.417% |
| 64 | 12.904 | 1695 | 1703 | 1712 | rVB  | 11809981 | 24002047 | 46.79%  | 1.637% |
| 65 | 13.104 | 1734 | 1737 | 1747 | rVB6 | 3110315  | 6992668  | 13.63%  | 0.477% |
| 66 | 13.210 | 1747 | 1755 | 1757 | rBV4 | 5585198  | 11083789 | 21.61%  | 0.756% |
| 67 | 13.369 | 1778 | 1782 | 1786 | rVV  | 6088670  | 10825853 | 21.11%  | 0.738% |
| 68 | 13.422 | 1786 | 1791 | 1796 | rVV3 | 6642372  | 13286186 | 25.90%  | 0.906% |
| 69 | 13.469 | 1796 | 1799 | 1805 | rVB2 | 4465886  | 6843295  | 13.34%  | 0.467% |
| 70 | 13.569 | 1805 | 1816 | 1819 | rBV3 | 6529791  | 16540704 | 32.25%  | 1.128% |
| 71 | 13.704 | 1832 | 1839 | 1842 | rBV4 | 6055681  | 12102092 | 23.59%  | 0.825% |
| 72 | 13.998 | 1883 | 1889 | 1893 | rBV  | 13861136 | 27907999 | 54.41%  | 1.903% |
| 73 | 14.039 | 1893 | 1896 | 1904 | rVB5 | 3570753  | 7211033  | 14.06%  | 0.492% |
| 74 | 14.163 | 1911 | 1917 | 1921 | rBV4 | 6553606  | 10358707 | 20.20%  | 0.706% |
| 75 | 14.216 | 1921 | 1926 | 1930 | rVB  | 11198617 | 17650808 | 34.41%  | 1.204% |
| 76 | 14.292 | 1930 | 1939 | 1943 | rBV4 | 4084580  | 12442835 | 24.26%  | 0.848% |

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Integration Parameters: LSCINT.P

Integrator: RTE

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |          |          |        |        |
|-----|--------|------|------|------|------|----------|----------|--------|--------|
| 77  | 14.545 | 1978 | 1982 | 1987 | rVB2 | 5216648  | 7487764  | 14.60% | 0.511% |
| 78  | 14.610 | 1989 | 1993 | 1997 | rVV2 | 5783114  | 8492743  | 16.56% | 0.579% |
| 79  | 14.710 | 2002 | 2010 | 2014 | rVV2 | 13689875 | 35113045 | 68.46% | 2.394% |
| 80  | 14.839 | 2026 | 2032 | 2044 | rBV  | 12138978 | 25657227 | 50.02% | 1.749% |
| 81  | 14.957 | 2044 | 2052 | 2057 | rVB2 | 11445118 | 21449017 | 41.82% | 1.463% |
| 82  | 15.192 | 2084 | 2092 | 2095 | rBV5 | 3879429  | 9721035  | 18.95% | 0.663% |
| 83  | 15.357 | 2112 | 2120 | 2130 | rBV3 | 6311003  | 16014940 | 31.22% | 1.092% |
| 84  | 15.710 | 2175 | 2180 | 2187 | rBV3 | 6718180  | 11674400 | 22.76% | 0.796% |
| 85  | 15.821 | 2193 | 2199 | 2201 | rBV2 | 11297831 | 20300317 | 39.58% | 1.384% |
| 86  | 16.498 | 2305 | 2314 | 2315 | rBV  | 15338718 | 35929039 | 70.05% | 2.450% |
| 87  | 16.615 | 2329 | 2334 | 2338 | rVV  | 9309311  | 12693287 | 24.75% | 0.865% |
| 88  | 16.663 | 2338 | 2342 | 2352 | rVB2 | 7421200  | 13727721 | 26.76% | 0.936% |
| 89  | 16.927 | 2383 | 2387 | 2392 | rVB2 | 4304331  | 6958884  | 13.57% | 0.474% |
| 90  | 17.351 | 2454 | 2459 | 2463 | rVB  | 9688526  | 13769437 | 26.84% | 0.939% |
| 91  | 17.521 | 2484 | 2488 | 2497 | rVB2 | 8566820  | 13365276 | 26.06% | 0.911% |
| 92  | 17.745 | 2522 | 2526 | 2531 | rVB2 | 3946843  | 6157814  | 12.01% | 0.420% |
| 93  | 18.039 | 2571 | 2576 | 2580 | rBV  | 8741490  | 12068586 | 23.53% | 0.823% |
| 94  | 18.686 | 2680 | 2686 | 2688 | rBV  | 7796510  | 13145365 | 25.63% | 0.896% |
| 95  | 19.268 | 2781 | 2785 | 2790 | rVB  | 5818669  | 7021560  | 13.69% | 0.479% |
| 96  | 19.809 | 2872 | 2877 | 2884 | rVB2 | 6283800  | 7894567  | 15.39% | 0.538% |
| 97  | 20.780 | 3038 | 3042 | 3052 | rVB  | 9991877  | 12965821 | 25.28% | 0.884% |
| 98  | 21.256 | 3119 | 3123 | 3131 | rVB  | 4315678  | 5938709  | 11.58% | 0.405% |
| 99  | 21.668 | 3188 | 3193 | 3204 | rVB  | 3848237  | 6781181  | 13.22% | 0.462% |
| 100 | 21.780 | 3206 | 3212 | 3218 | rVB3 | 4974440  | 8484780  | 16.54% | 0.579% |

Sum of corrected areas: 1466587272

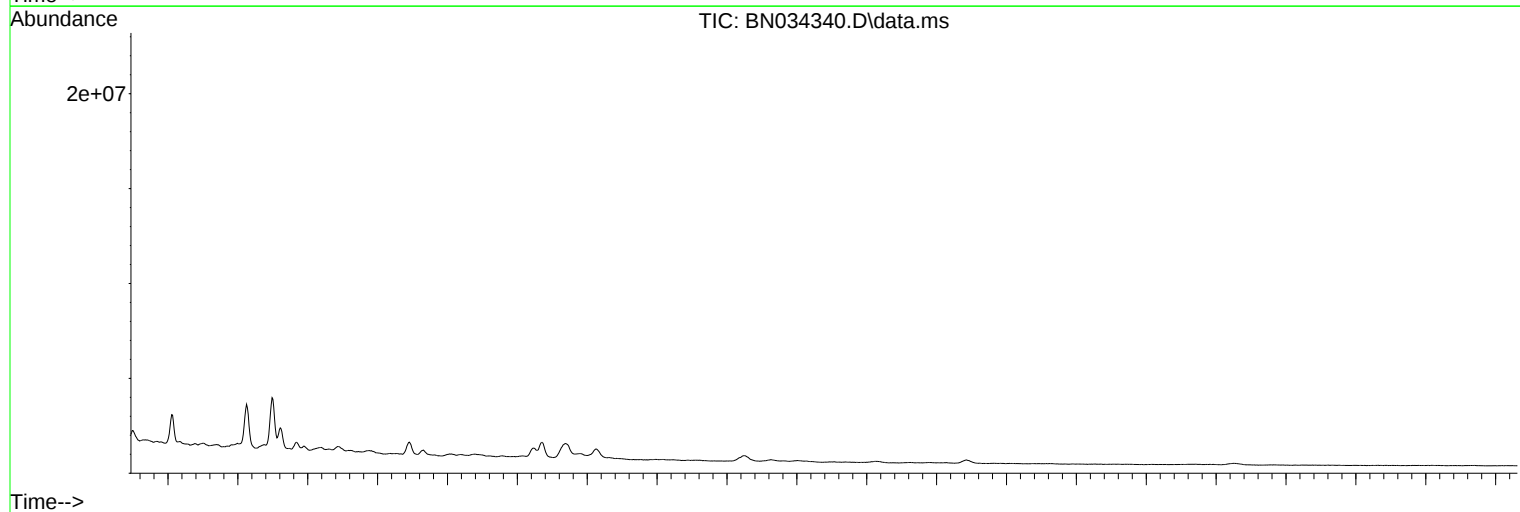
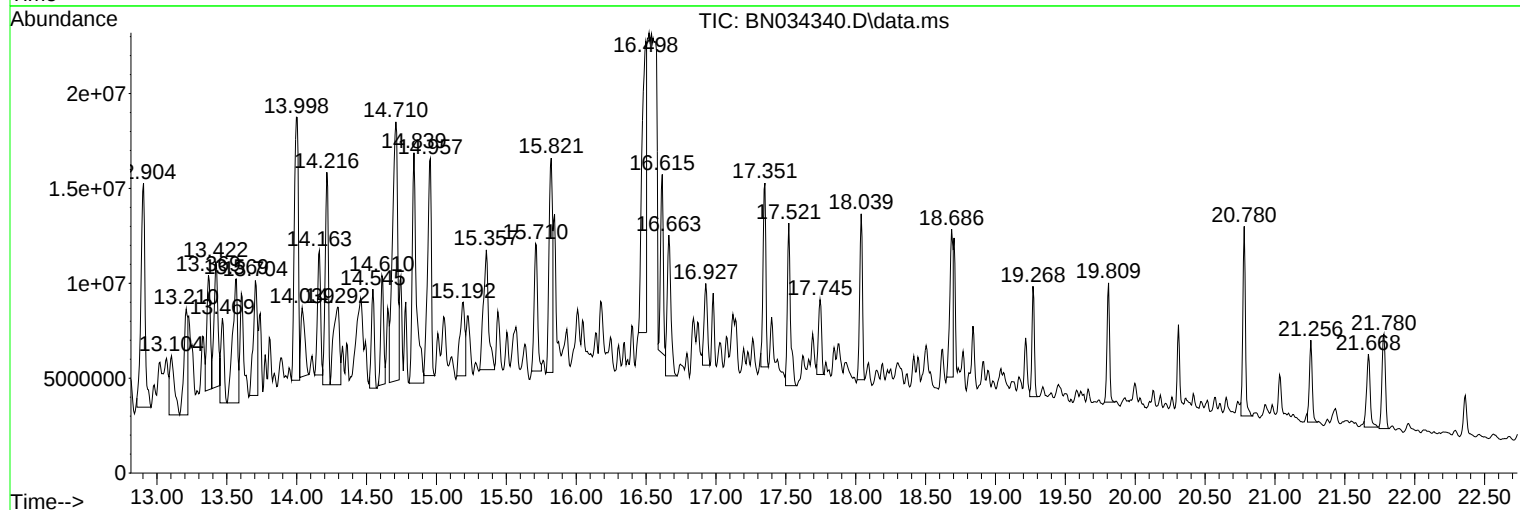
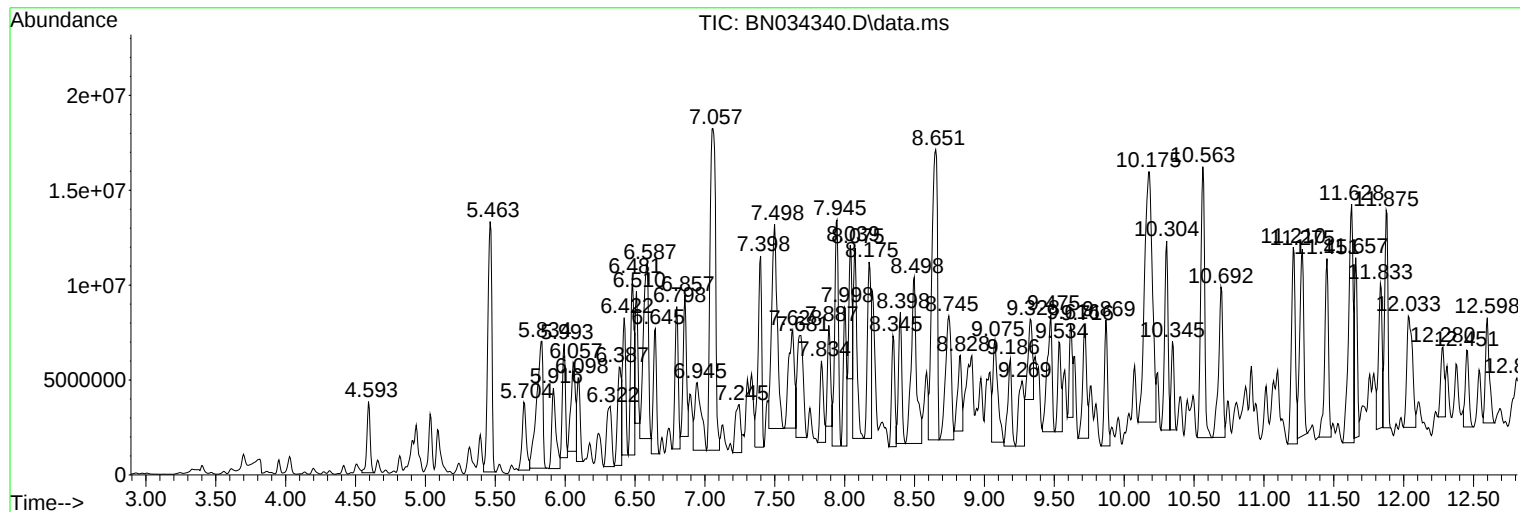
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P



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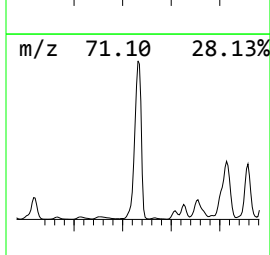
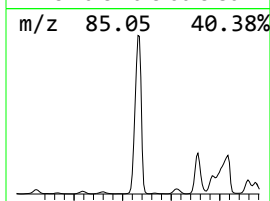
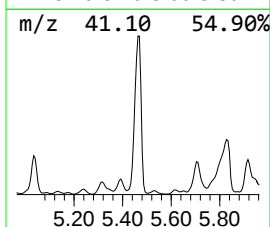
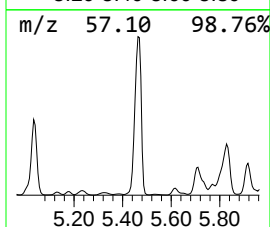
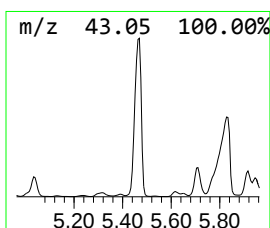
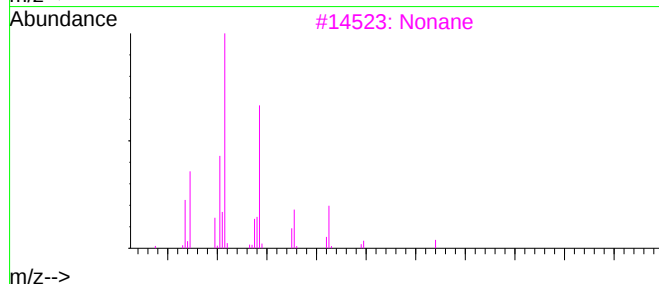
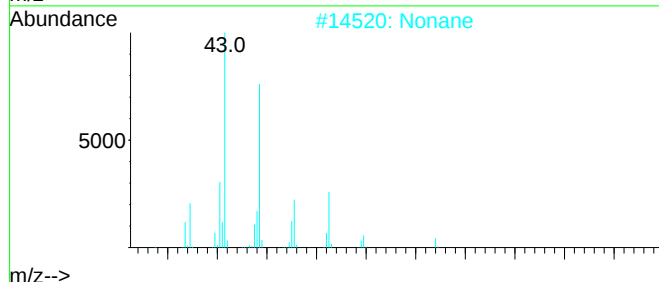
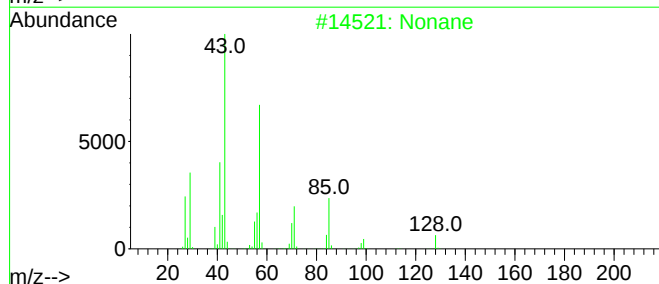
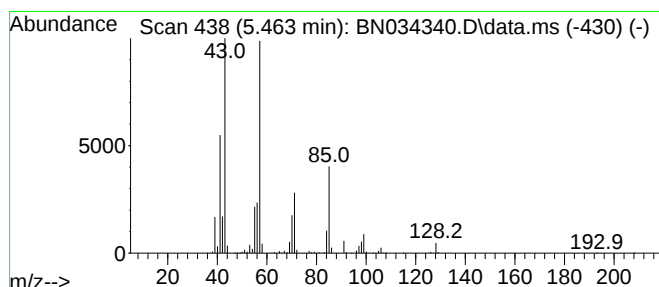
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Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 23

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 5.463 | 29.63 ng/ul | 26836400 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|-----------|--------------|-----|---------|-------------|------|
| 1         | Nonane       | 128 | C9H20   | 000111-84-2 | 94   |
| 2         | Nonane       | 128 | C9H20   | 000111-84-2 | 93   |
| 3         | Nonane       | 128 | C9H20   | 000111-84-2 | 91   |
| 4         | Nonane       | 128 | C9H20   | 000111-84-2 | 91   |
| 5         | Octane       | 114 | C8H18   | 000111-65-9 | 72   |



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TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

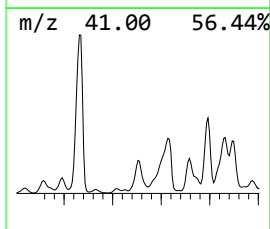
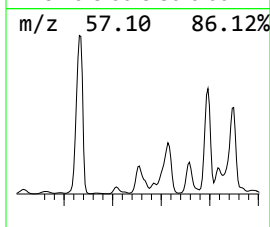
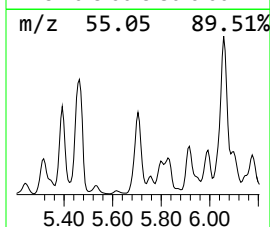
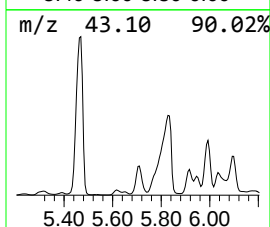
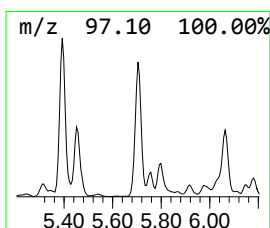
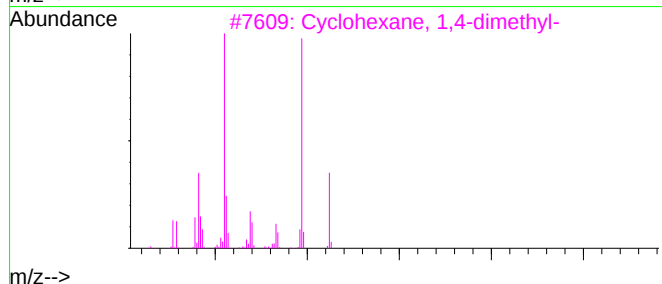
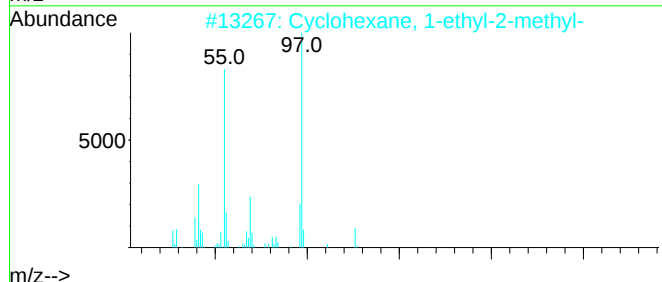
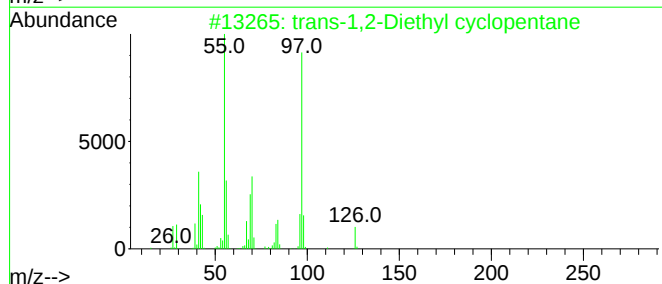
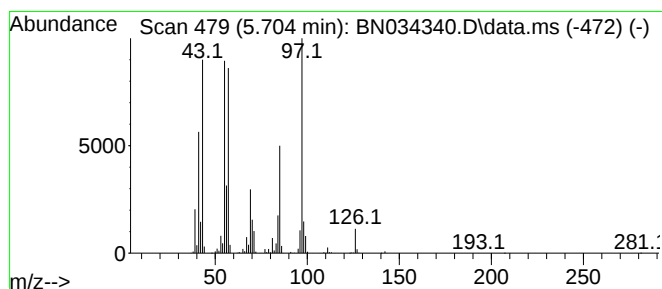
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Peak Number 3 (DEL) Alkane: Cyclic5.704 Concentration Rank 58

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 5.704 | 8.22 ng/ul | 7448920 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|-----------|-------------------------------------|-----|---------|-------------|------|
| 1         | trans-1,2-Diethyl cyclopentane      | 126 | C9H18   | 000932-40-1 | 58   |
| 2         | Cyclohexane, 1-ethyl-2-methyl-      | 126 | C9H18   | 003728-54-9 | 55   |
| 3         | Cyclohexane, 1,4-dimethyl-          | 112 | C8H16   | 000589-90-2 | 50   |
| 4         | Cyclohexane, 1-ethyl-2-methyl-, ... | 126 | C9H18   | 004923-78-8 | 49   |
| 5         | Cyclohexane, 1-ethyl-4-methyl-, ... | 126 | C9H18   | 004926-78-7 | 46   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

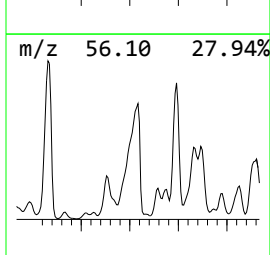
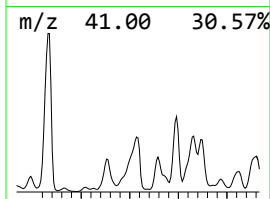
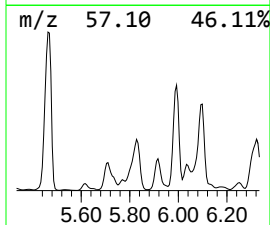
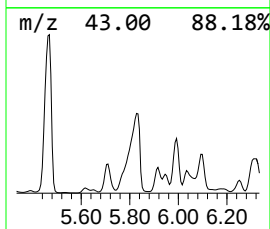
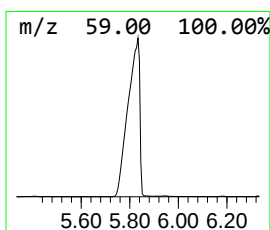
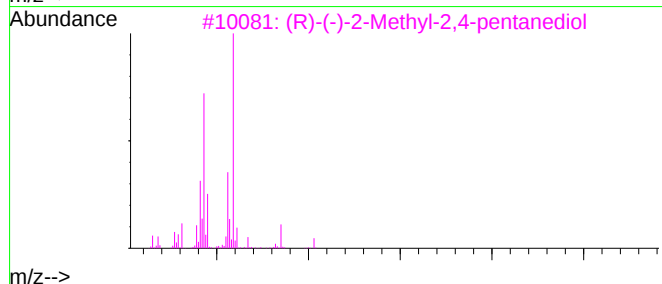
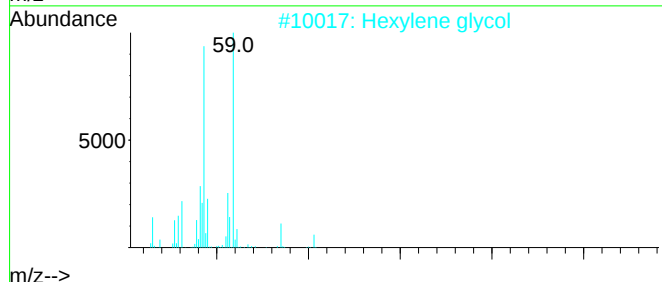
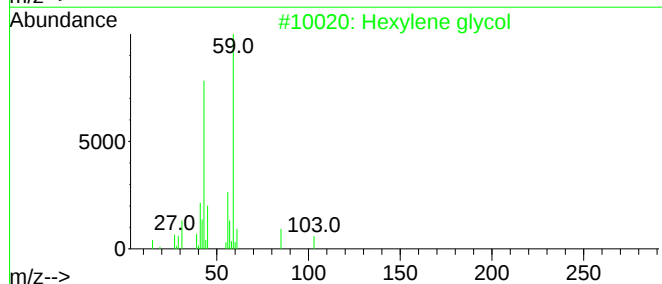
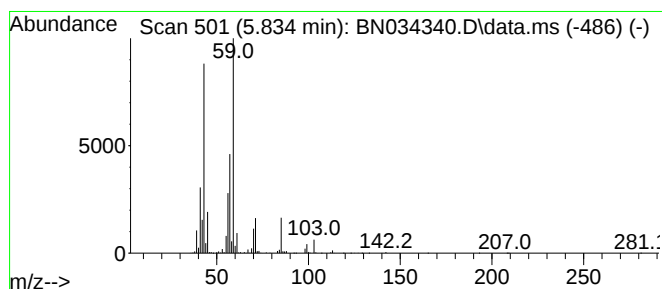
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Hexylene glycol Concentration Rank 28

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 5.834 | 23.17 ng/ul | 20984500 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Hexylene glycol                     | 118 | C6H14O2 | 000107-41-5  | 80   |
| 2    |    |   | Hexylene glycol                     | 118 | C6H14O2 | 000107-41-5  | 72   |
| 3    |    |   | (R)-(-)-2-Methyl-2,4-pentanediol    | 118 | C6H14O2 | 099210-90-9  | 72   |
| 4    |    |   | Hexylene glycol                     | 118 | C6H14O2 | 000107-41-5  | 72   |
| 5    |    |   | Carbonic acid, 2-methoxyethyl ne... | 190 | C9H18O4 | 1000331-42-7 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

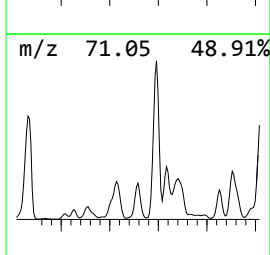
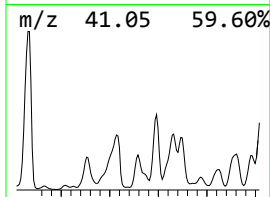
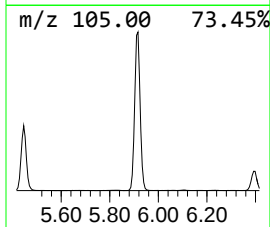
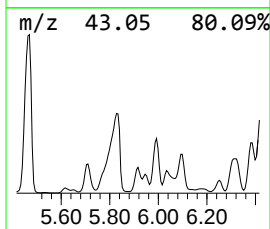
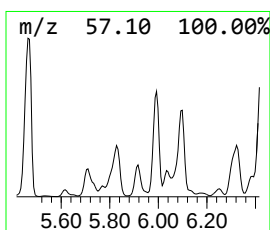
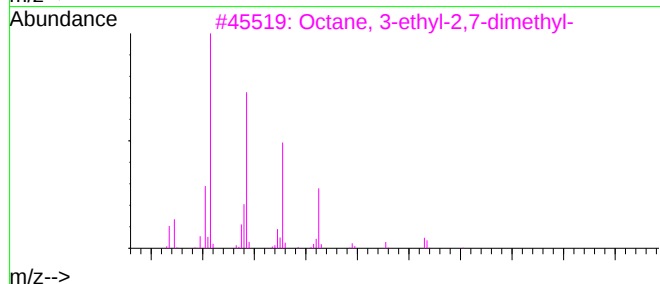
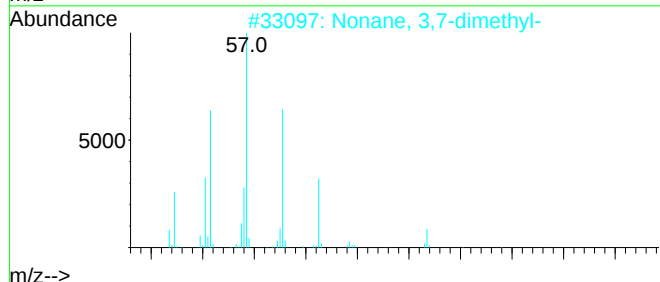
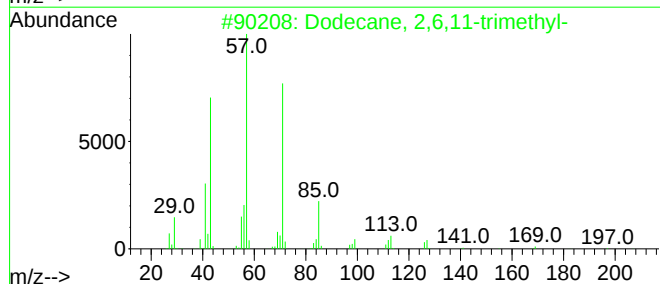
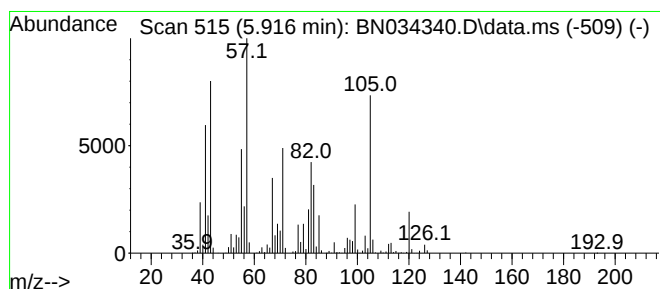
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 53

| R.T.      | EstConc                       | Area    | Relative to ISTD       | R.T.        |      |
|-----------|-------------------------------|---------|------------------------|-------------|------|
| 5.916     | 9.88 ng/ul                    | 8950850 | 1,4-Dichlorobenzene-d4 | 7.392       |      |
| Hit# of 5 | Tentative ID                  | MW      | MolForm                | CAS#        | Qual |
| 1         | Dodecane, 2,6,11-trimethyl-   | 212     | C15H32                 | 031295-56-4 | 27   |
| 2         | Nonane, 3,7-dimethyl-         | 156     | C11H24                 | 017302-32-8 | 27   |
| 3         | Octane, 3-ethyl-2,7-dimethyl- | 170     | C12H26                 | 062183-55-5 | 22   |
| 4         | Decane, 2-methyl-             | 156     | C11H24                 | 006975-98-0 | 22   |
| 5         | Dodecane, 2,6,11-trimethyl-   | 212     | C15H32                 | 031295-56-4 | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

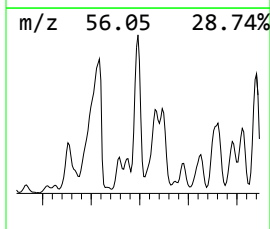
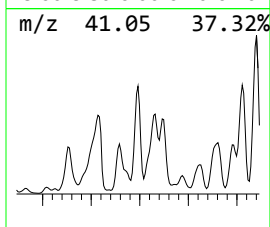
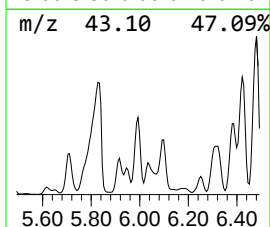
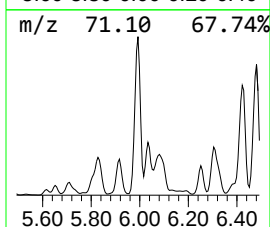
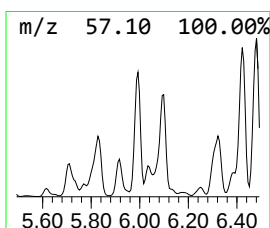
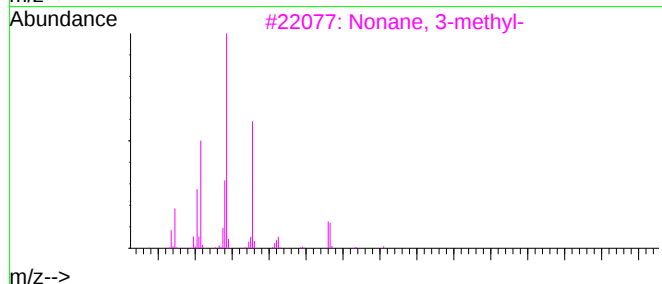
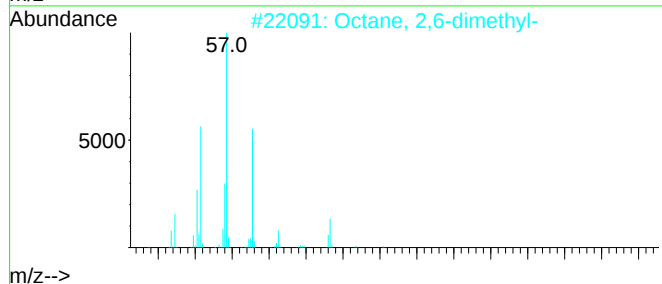
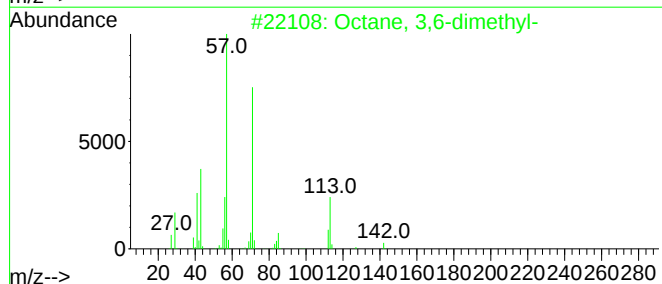
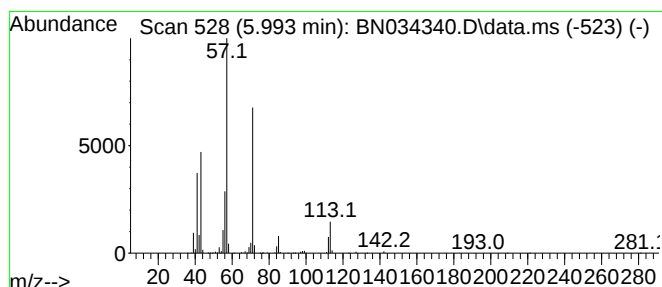
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Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 52

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 5.993 | 10.15 ng/ul | 9189080 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# of 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|-----------|-----------------------|-----|---------|-------------|------|
| 1         | Octane, 3,6-dimethyl- | 142 | C10H22  | 015869-94-0 | 91   |
| 2         | Octane, 2,6-dimethyl- | 142 | C10H22  | 002051-30-1 | 91   |
| 3         | Nonane, 3-methyl-     | 142 | C10H22  | 005911-04-6 | 90   |
| 4         | Octane, 3,6-dimethyl- | 142 | C10H22  | 015869-94-0 | 90   |
| 5         | Octane, 2,6-dimethyl- | 142 | C10H22  | 002051-30-1 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

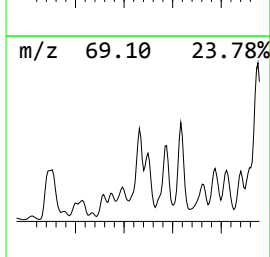
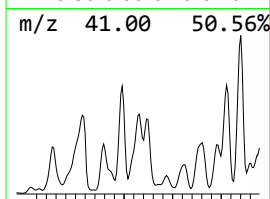
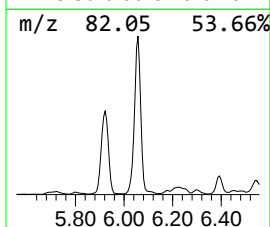
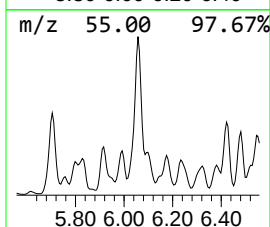
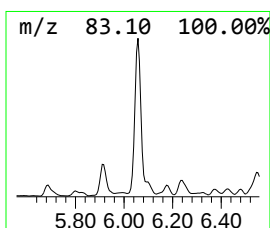
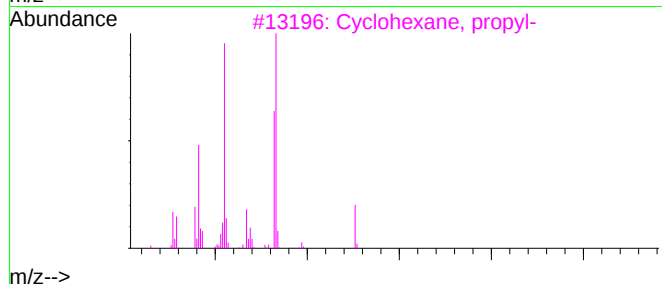
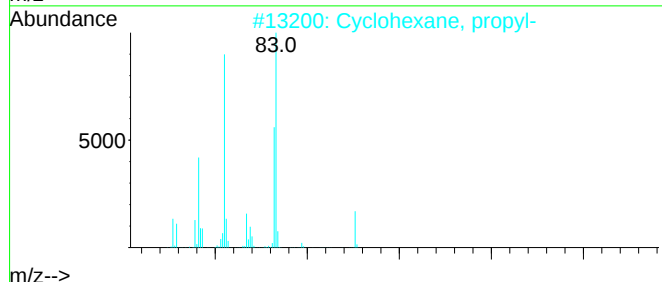
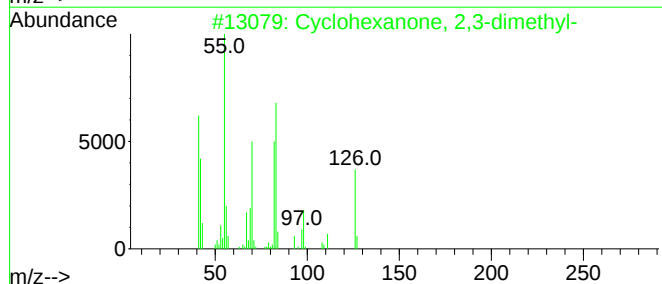
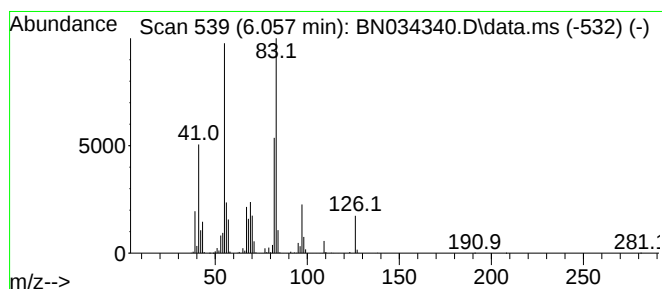
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Peak Number 7 Cyclohexanone, 2,3-dimethyl- Concentration Rank 46

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 6.057 | 12.06 ng/ul | 10925600 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# of 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|-----------|------------------------------|-----|---------|-------------|------|
| 1         | Cyclohexanone, 2,3-dimethyl- | 126 | C8H14O  | 013395-76-1 | 87   |
| 2         | Cyclohexane, propyl-         | 126 | C9H18   | 001678-92-8 | 74   |
| 3         | Cyclohexane, propyl-         | 126 | C9H18   | 001678-92-8 | 72   |
| 4         | Cyclohexane, propyl-         | 126 | C9H18   | 001678-92-8 | 64   |
| 5         | Cyclohexane, propyl-         | 126 | C9H18   | 001678-92-8 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

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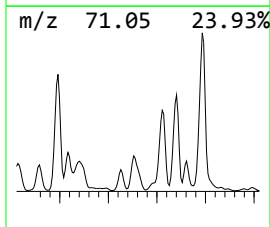
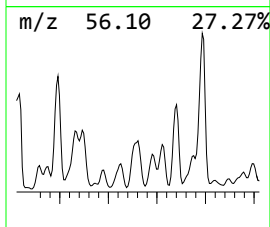
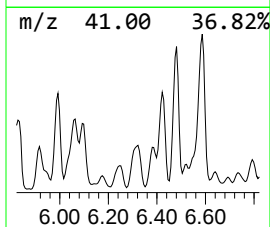
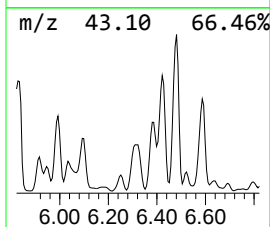
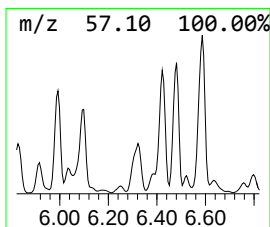
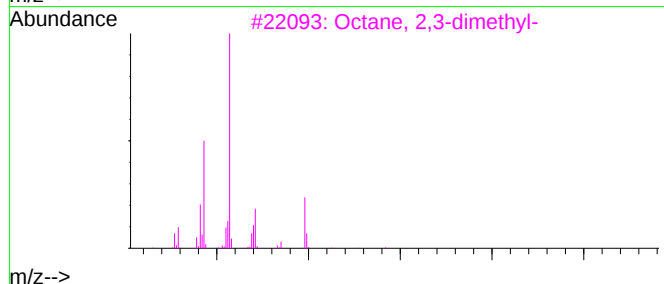
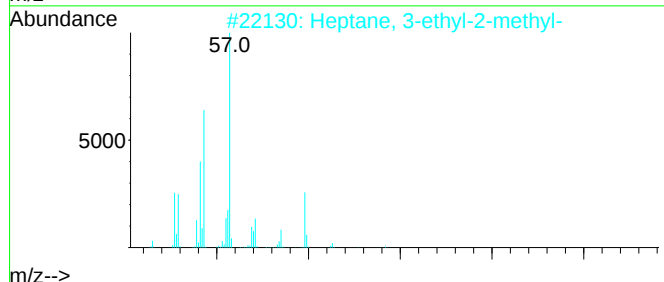
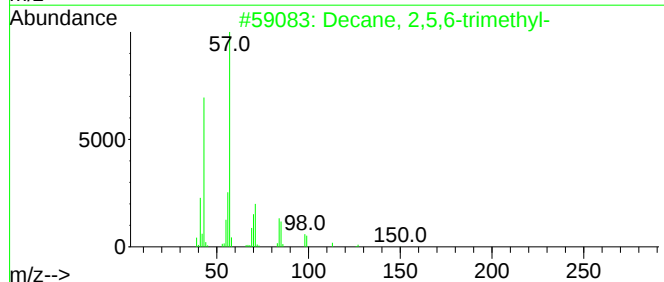
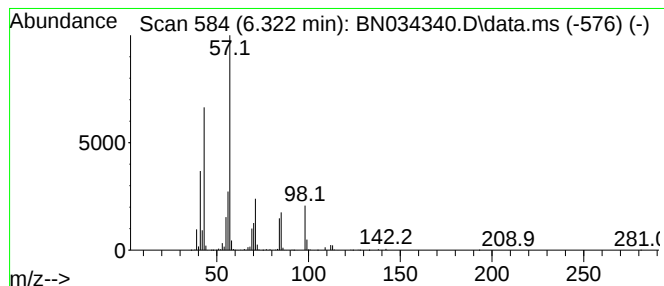
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Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 56

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 6.322 | 8.79 ng/ul | 7962340 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# of 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|-----------|-------------------------------|-----|---------|-------------|------|
| 1         | Decane, 2,5,6-trimethyl-      | 184 | C13H28  | 062108-23-0 | 80   |
| 2         | Heptane, 3-ethyl-2-methyl-    | 142 | C10H22  | 014676-29-0 | 64   |
| 3         | Octane, 2,3-dimethyl-         | 142 | C10H22  | 007146-60-3 | 64   |
| 4         | Hexane, 3-ethyl-2,5-dimethyl- | 142 | C10H22  | 052897-04-8 | 52   |
| 5         | Heptane, 3-ethyl-2-methyl-    | 142 | C10H22  | 014676-29-0 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

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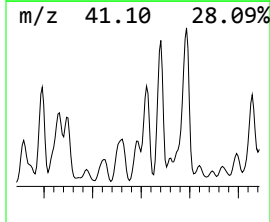
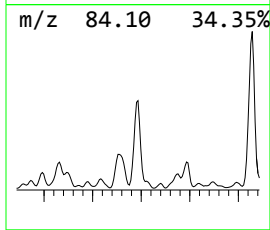
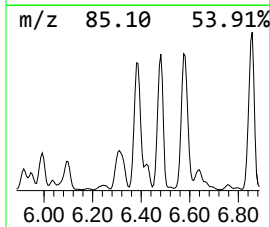
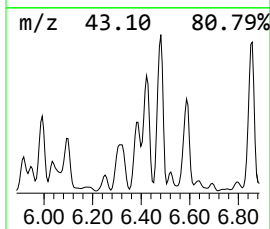
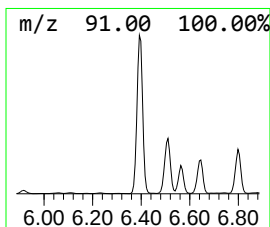
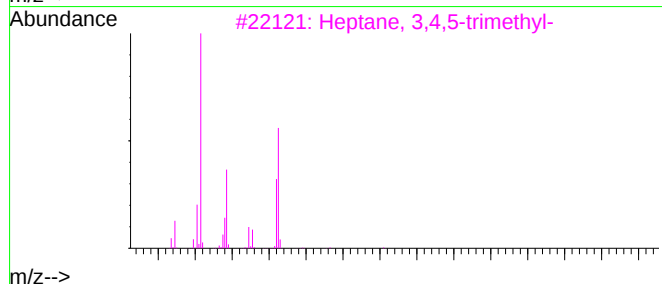
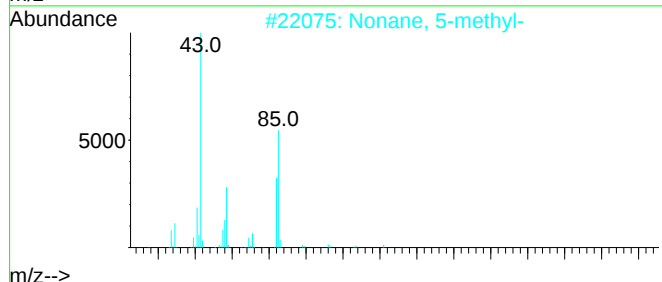
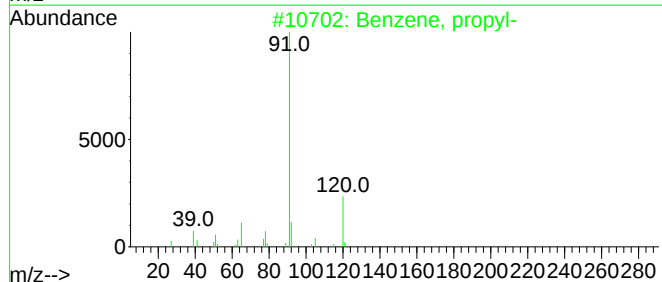
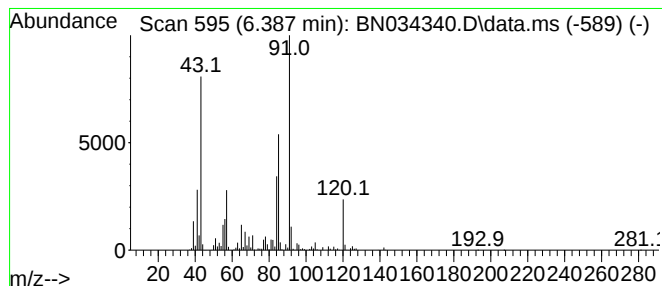
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Benzene, propyl- Concentration Rank 48

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 6.387 | 11.66 ng/ul | 10559800 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# of 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|-----------|----------------------------|-----|---------|-------------|------|
| 1         | Benzene, propyl-           | 120 | C9H12   | 000103-65-1 | 70   |
| 2         | Nonane, 5-methyl-          | 142 | C10H22  | 015869-85-9 | 60   |
| 3         | Heptane, 3,4,5-trimethyl-  | 142 | C10H22  | 020278-89-1 | 47   |
| 4         | Nonane, 5-methyl-          | 142 | C10H22  | 015869-85-9 | 47   |
| 5         | Pentane, 3-ethyl-3-methyl- | 114 | C8H18   | 001067-08-9 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

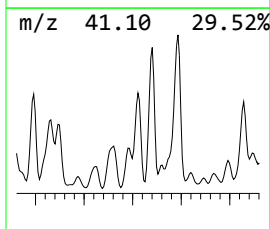
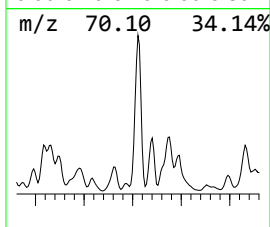
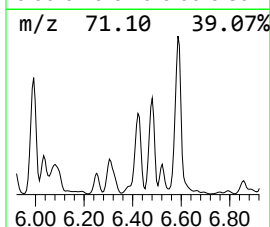
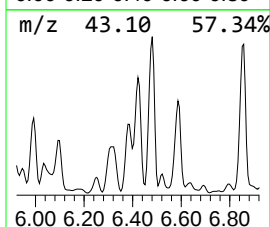
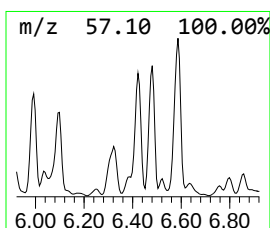
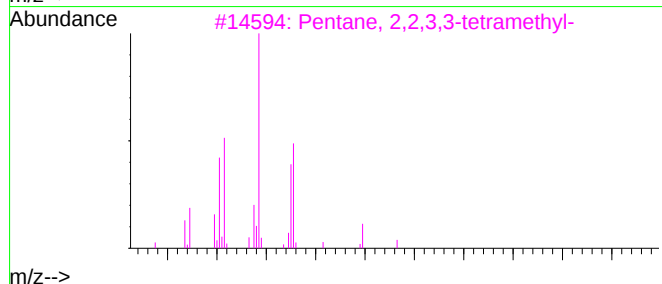
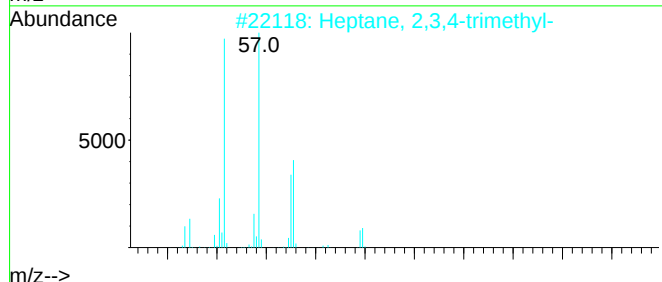
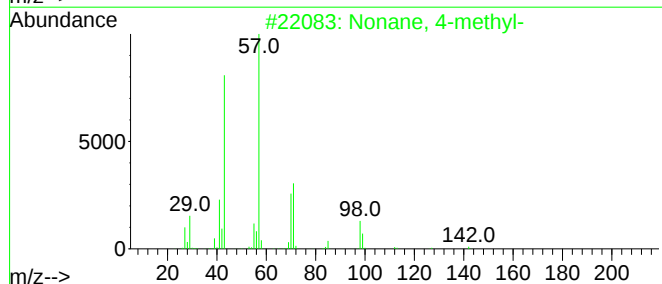
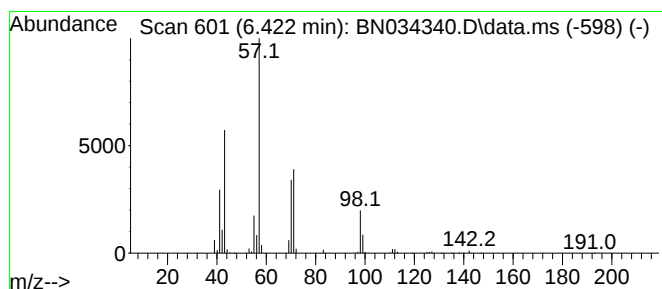
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Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 45

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 6.422 | 12.08 ng/ul | 10941100 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# of 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|-----------|-------------------------------|-----|---------|-------------|------|
| 1         | Nonane, 4-methyl-             | 142 | C10H22  | 017301-94-9 | 91   |
| 2         | Heptane, 2,3,4-trimethyl-     | 142 | C10H22  | 052896-95-4 | 64   |
| 3         | Pentane, 2,2,3,3-tetramethyl- | 128 | C9H20   | 007154-79-2 | 59   |
| 4         | Nonane, 4-methyl-             | 142 | C10H22  | 017301-94-9 | 59   |
| 5         | Pentane, 2,2,3,3-tetramethyl- | 128 | C9H20   | 007154-79-2 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

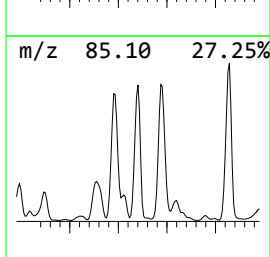
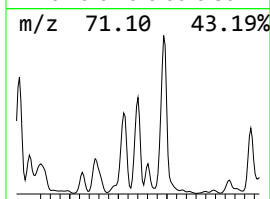
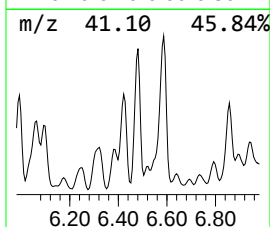
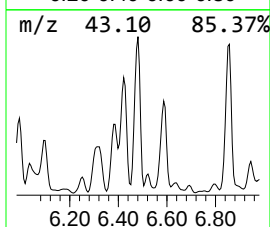
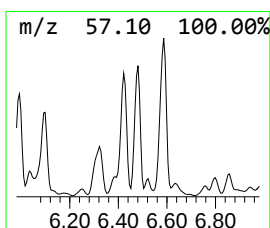
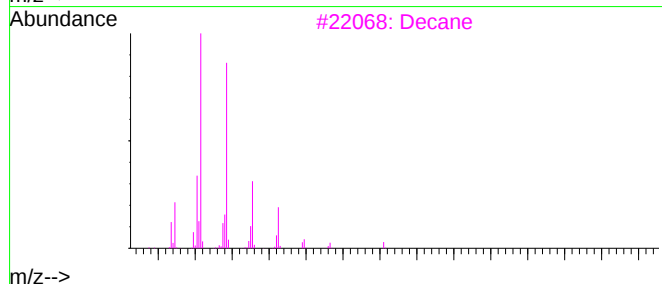
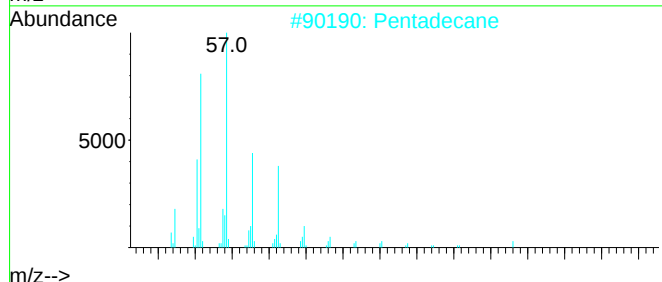
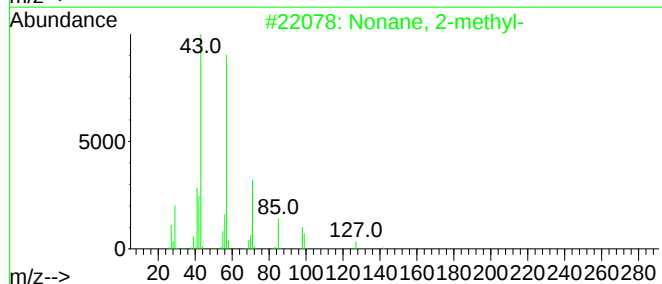
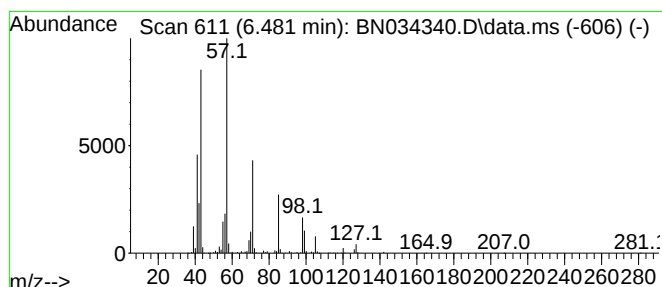
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Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 36

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 6.481 | 18.92 ng/ul | 17136400 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# of 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|-----------|----------------------|-----|---------|-------------|------|
| 1         | Nonane, 2-methyl-    | 142 | C10H22  | 000871-83-0 | 81   |
| 2         | Pentadecane          | 212 | C15H32  | 000629-62-9 | 64   |
| 3         | Decane               | 142 | C10H22  | 000124-18-5 | 64   |
| 4         | Tridecane, 6-methyl- | 198 | C14H30  | 013287-21-3 | 59   |
| 5         | Undecane, 5-methyl-  | 170 | C12H26  | 001632-70-8 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

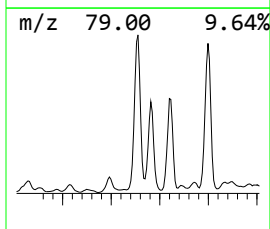
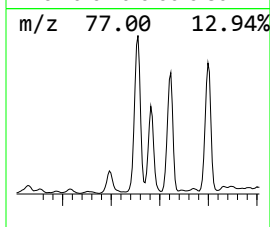
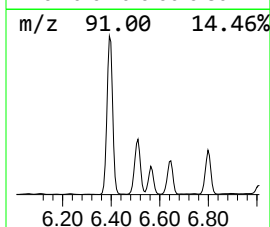
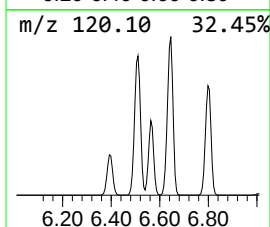
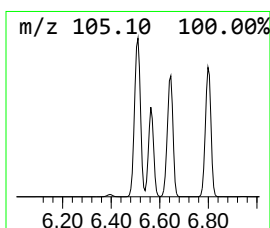
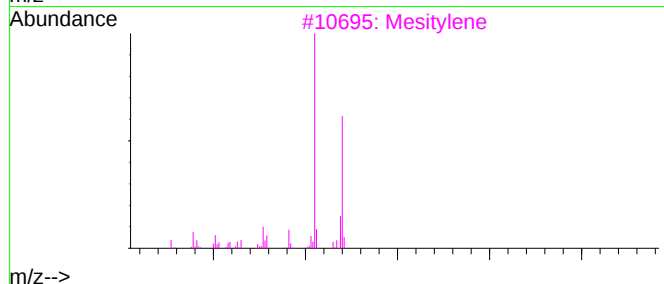
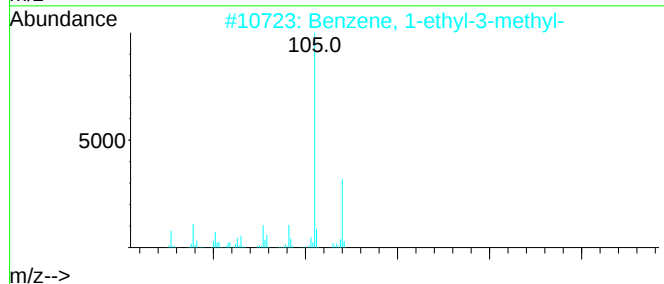
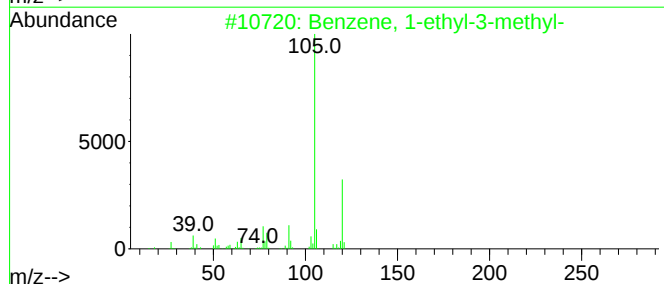
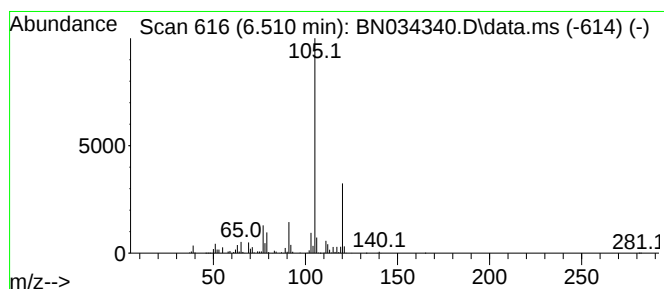
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 Benzene, 1-ethyl-3-methyl- Concentration Rank 55

| R.T.      | EstConc                    | Area    | Relative to ISTD       |         |             | R.T.  |
|-----------|----------------------------|---------|------------------------|---------|-------------|-------|
| 6.510     | 9.10 ng/ul                 | 8241200 | 1,4-Dichlorobenzene-d4 |         |             | 7.392 |
| Hit# of 5 | Tentative ID               |         | MW                     | MolForm | CAS#        | Qual  |
| 1         | Benzene, 1-ethyl-3-methyl- |         | 120                    | C9H12   | 000620-14-4 | 94    |
| 2         | Benzene, 1-ethyl-3-methyl- |         | 120                    | C9H12   | 000620-14-4 | 93    |
| 3         | Mesitylene                 |         | 120                    | C9H12   | 000108-67-8 | 90    |
| 4         | Benzene, 1-ethyl-4-methyl- |         | 120                    | C9H12   | 000622-96-8 | 90    |
| 5         | Benzene, 1-ethyl-2-methyl- |         | 120                    | C9H12   | 000611-14-3 | 87    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

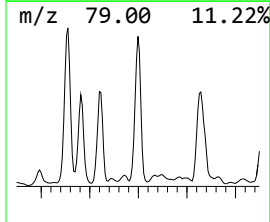
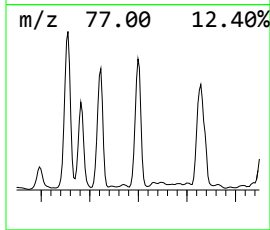
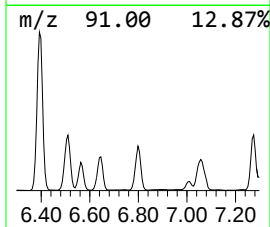
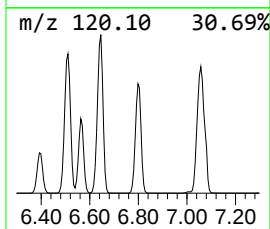
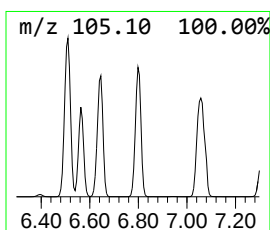
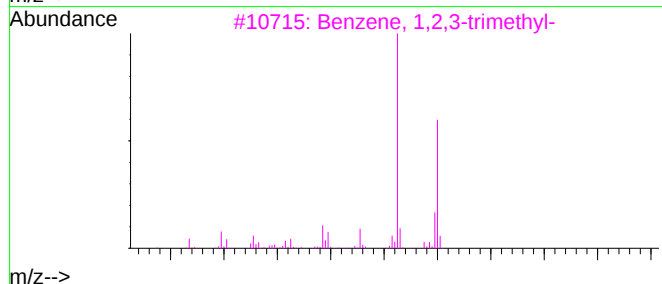
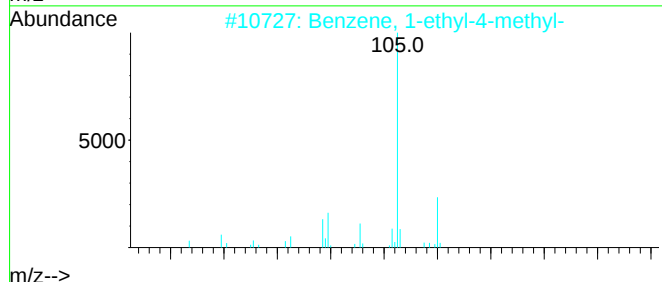
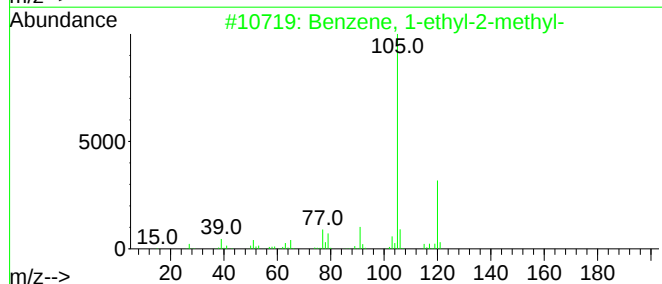
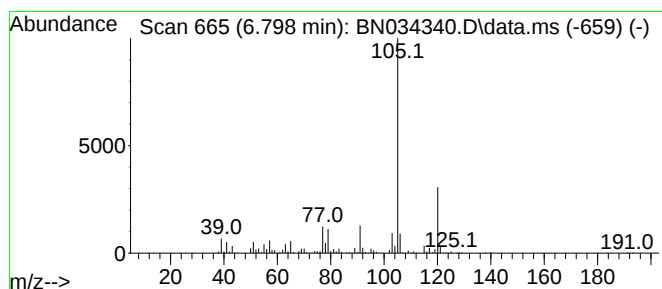
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TIC Integration Parameters: LSCINT.P

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Peak Number 14 Benzene, 1-ethyl-2-methyl- Concentration Rank 42

| R.T.      | EstConc                    | Area     | Relative to ISTD       |         |             | R.T.  |
|-----------|----------------------------|----------|------------------------|---------|-------------|-------|
| 6.798     | 13.46 ng/ul                | 12195100 | 1,4-Dichlorobenzene-d4 |         |             | 7.392 |
| Hit# of 5 | Tentative ID               |          | MW                     | MolForm | CAS#        | Qual  |
| 1         | Benzene, 1-ethyl-2-methyl- |          | 120                    | C9H12   | 000611-14-3 | 94    |
| 2         | Benzene, 1-ethyl-4-methyl- |          | 120                    | C9H12   | 000622-96-8 | 94    |
| 3         | Benzene, 1,2,3-trimethyl-  |          | 120                    | C9H12   | 000526-73-8 | 91    |
| 4         | Mesitylene                 |          | 120                    | C9H12   | 000108-67-8 | 91    |
| 5         | Mesitylene                 |          | 120                    | C9H12   | 000108-67-8 | 91    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

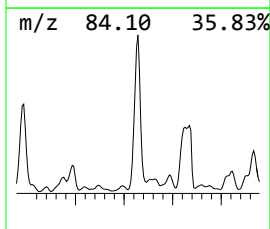
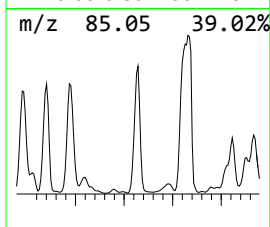
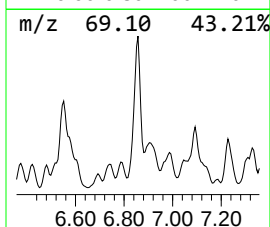
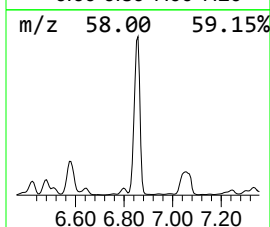
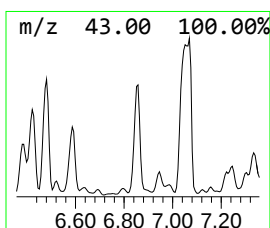
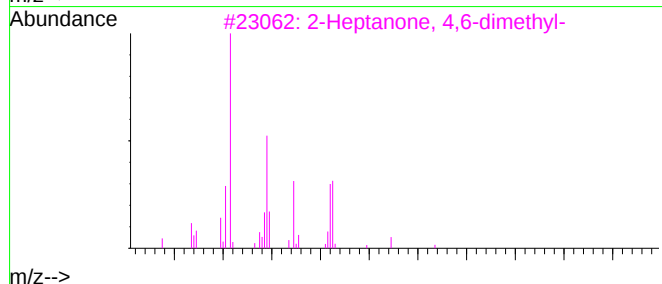
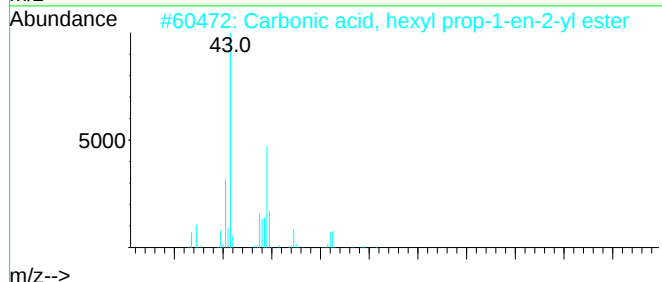
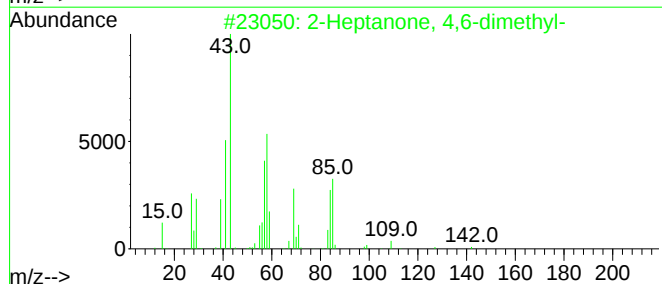
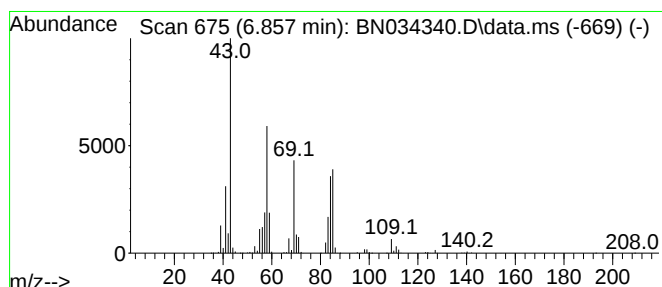
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Peak Number 15 2-Heptanone, 4,6-dimethyl- Concentration Rank 41

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 6.857 | 13.55 ng/ul | 12274100 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# of 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|-----------|-------------------------------------|-----|----------|--------------|------|
| 1         | 2-Heptanone, 4,6-dimethyl-          | 142 | C9H18O   | 019549-80-5  | 83   |
| 2         | Carbonic acid, hexyl prop-1-en-2... | 186 | C10H18O3 | 1000382-54-2 | 74   |
| 3         | 2-Heptanone, 4,6-dimethyl-          | 142 | C9H18O   | 019549-80-5  | 72   |
| 4         | 2-Pentanone, 5,5'-oxybis-           | 186 | C10H18O3 | 093677-66-8  | 59   |
| 5         | Pentane, 3-ethyl-2,4-dimethyl-      | 128 | C9H20    | 001068-87-7  | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
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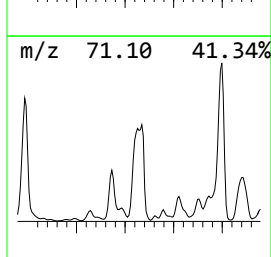
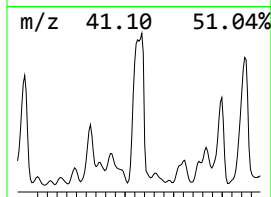
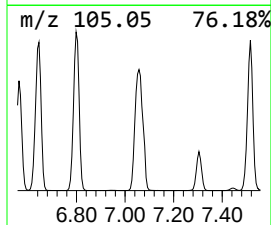
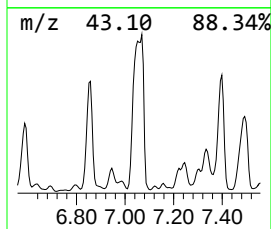
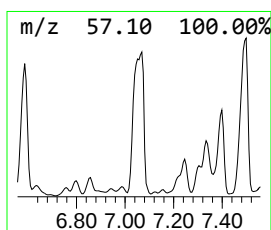
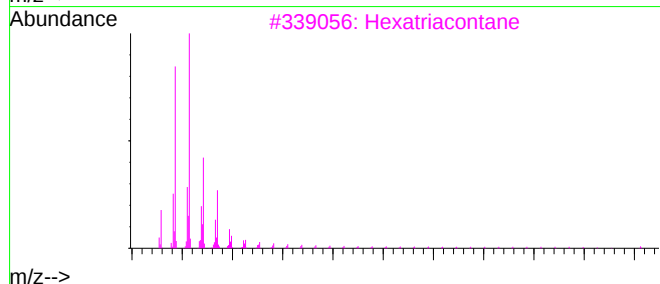
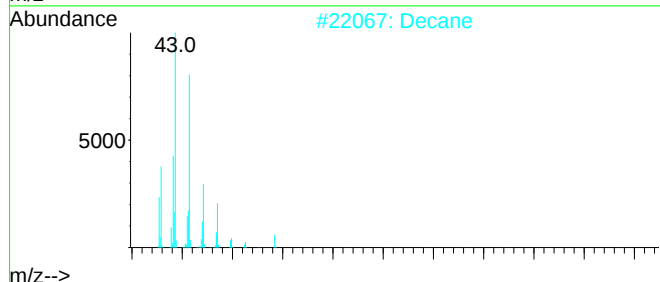
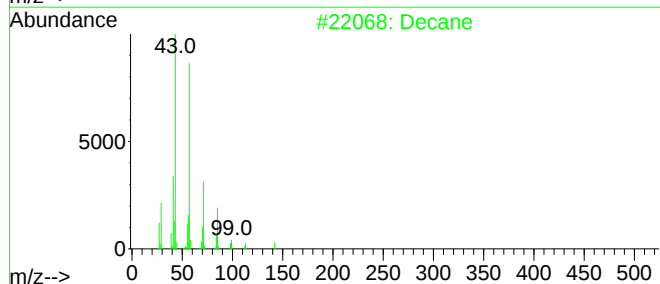
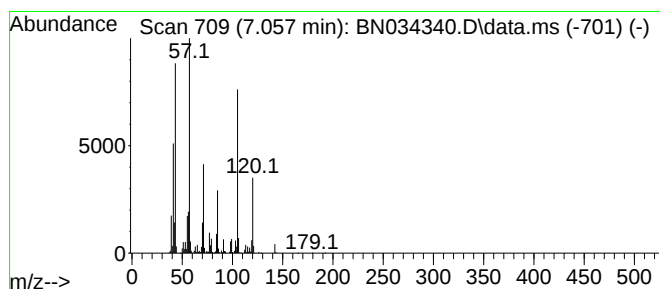
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TIC Integration Parameters: LSCINT.P

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Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 5

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 7.057 | 51.75 ng/ul | 46871800 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# of 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|-----------|---------------------------|-----|---------|-------------|------|
| 1         | Decane                    | 142 | C10H22  | 000124-18-5 | 83   |
| 2         | Decane                    | 142 | C10H22  | 000124-18-5 | 52   |
| 3         | Hexatriacontane           | 507 | C36H74  | 000630-06-8 | 50   |
| 4         | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 44   |
| 5         | Nonane                    | 128 | C9H20   | 000111-84-2 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

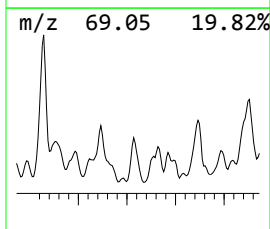
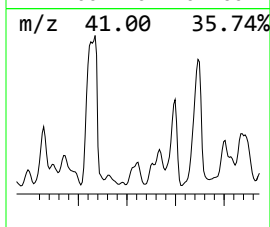
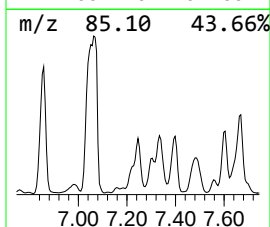
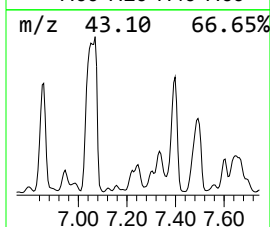
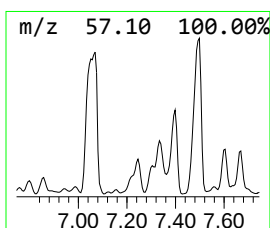
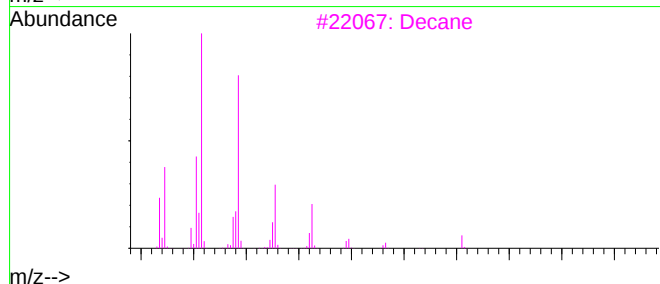
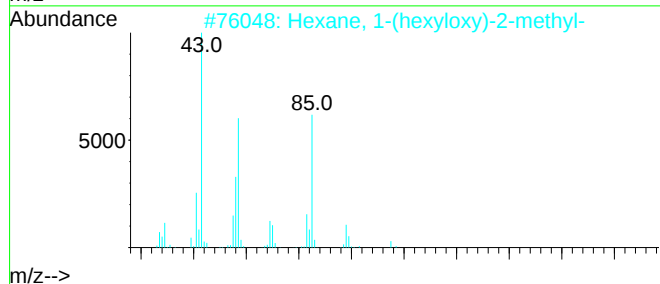
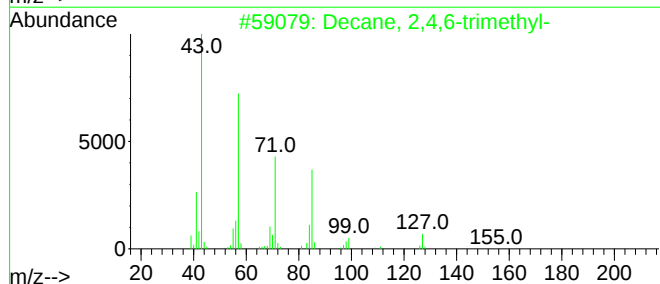
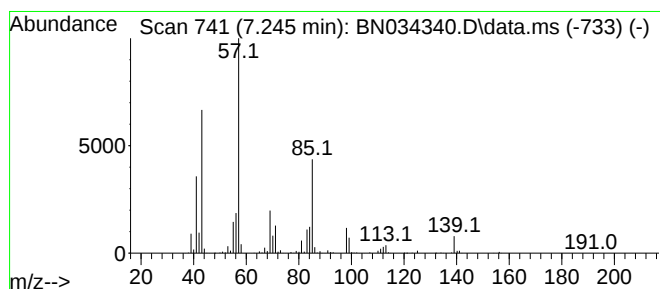
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Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 61

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 7.245 | 7.38 ng/ul | 6684190 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 2,4,6-trimethyl-       | 184 | C13H28  | 062108-27-4 | 53   |
| 2    |    |   | Hexane, 1-(hexyloxy)-2-methyl- | 200 | C13H28O | 074421-17-3 | 50   |
| 3    |    |   | Decane                         | 142 | C10H22  | 000124-18-5 | 47   |
| 4    |    |   | Hexane, 1,1'-oxybis-           | 186 | C12H26O | 000112-58-3 | 47   |
| 5    |    |   | Heptane, 4,4-dimethyl-         | 128 | C9H20   | 001068-19-5 | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

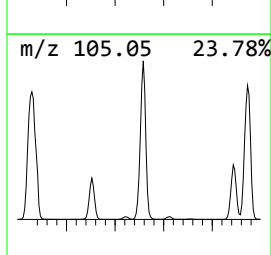
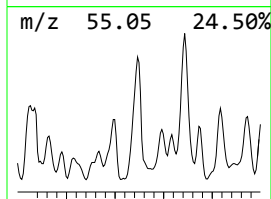
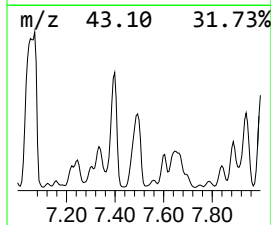
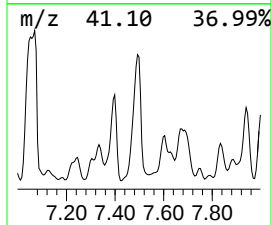
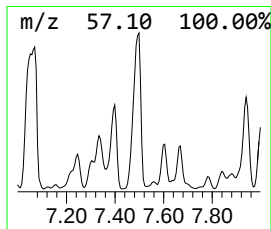
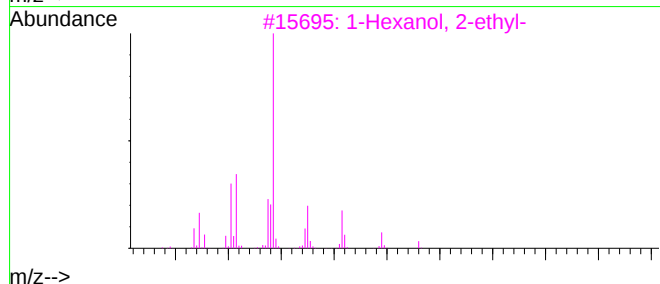
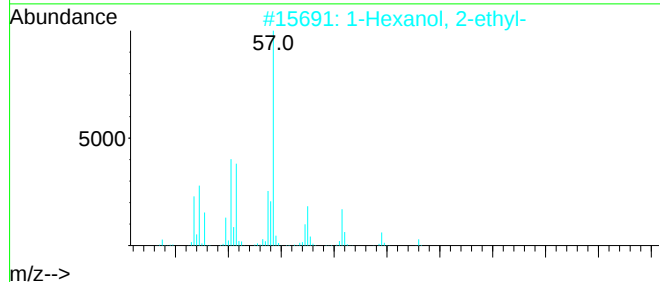
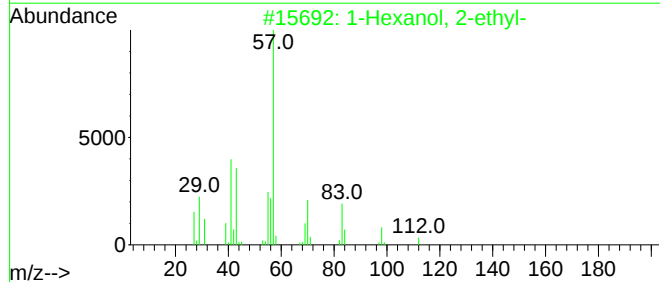
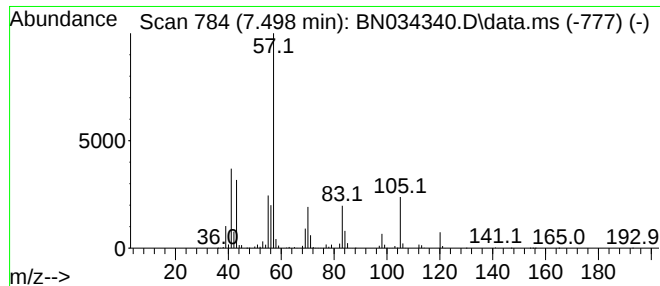
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ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 1-Hexanol, 2-ethyl- Concentration Rank 19

| R.T.      | EstConc                       | Area     | Relative to ISTD       |          | R.T.         |      |
|-----------|-------------------------------|----------|------------------------|----------|--------------|------|
| 7.498     | 32.62 ng/ul                   | 29546800 | 1,4-Dichlorobenzene-d4 |          | 7.392        |      |
| Hit# of 5 | Tentative ID                  |          | MW                     | MolForm  | CAS#         | Qual |
| 1         | 1-Hexanol, 2-ethyl-           |          | 130                    | C8H18O   | 000104-76-7  | 64   |
| 2         | 1-Hexanol, 2-ethyl-           |          | 130                    | C8H18O   | 000104-76-7  | 59   |
| 3         | 1-Hexanol, 2-ethyl-           |          | 130                    | C8H18O   | 000104-76-7  | 59   |
| 4         | Isobutyl octan-2-yl carbonate |          | 230                    | C13H26O3 | 1000372-74-7 | 59   |
| 5         | 1-Hexanol, 2-ethyl-           |          | 130                    | C8H18O   | 000104-76-7  | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

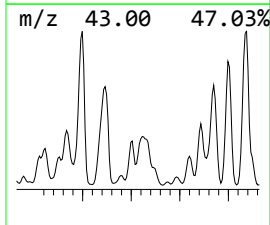
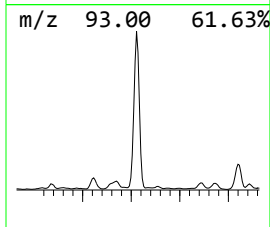
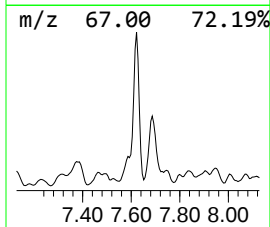
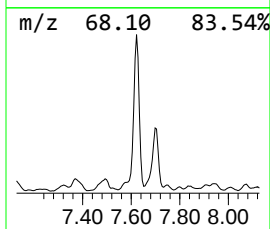
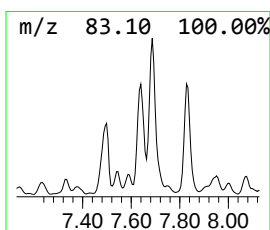
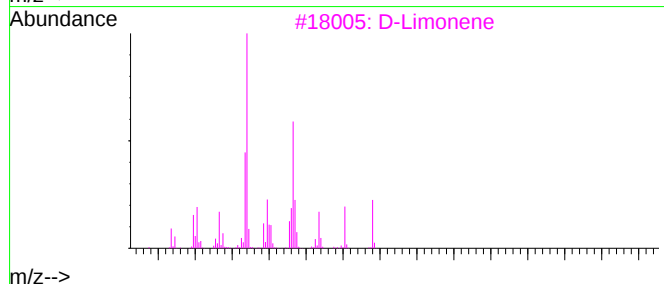
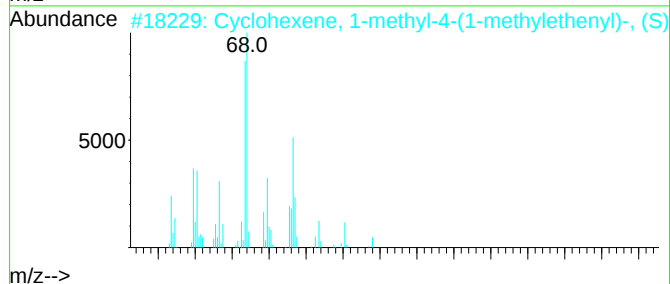
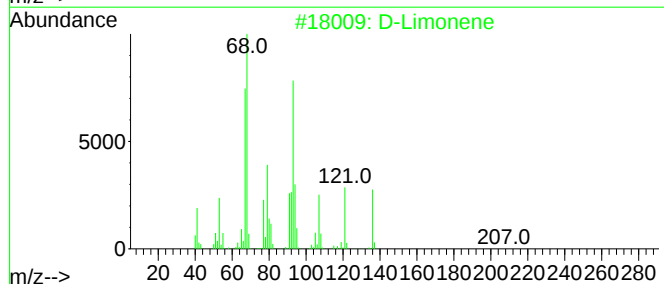
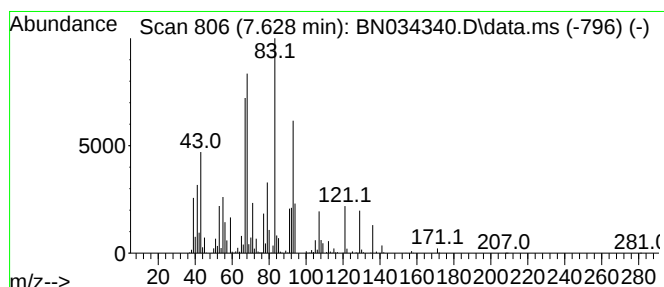
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Peak Number 19 D-Limonene Concentration Rank 40

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 7.628 | 16.85 ng/ul | 15263000 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | D-Limonene                          | 136 | C10H16  | 005989-27-5 | 96   |
| 2    |    |   | Cyclohexene, 1-methyl-4-(1-methy... | 136 | C10H16  | 005989-54-8 | 93   |
| 3    |    |   | D-Limonene                          | 136 | C10H16  | 005989-27-5 | 64   |
| 4    |    |   | Limonene                            | 136 | C10H16  | 000138-86-3 | 62   |
| 5    |    |   | D-Limonene                          | 136 | C10H16  | 005989-27-5 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L

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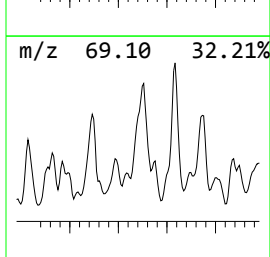
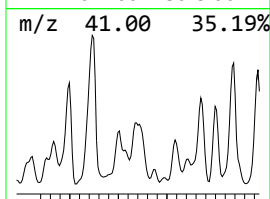
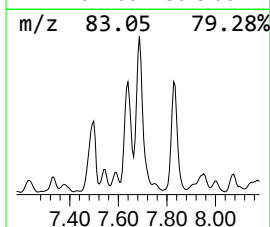
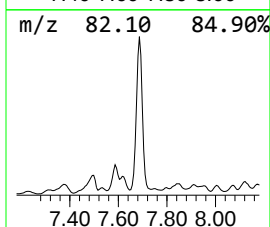
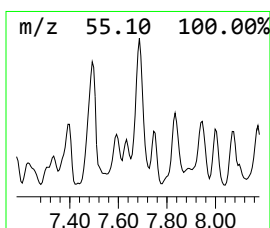
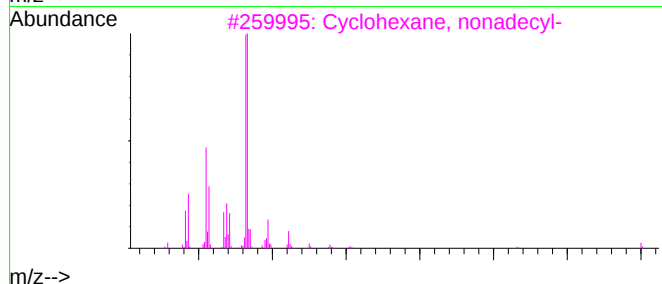
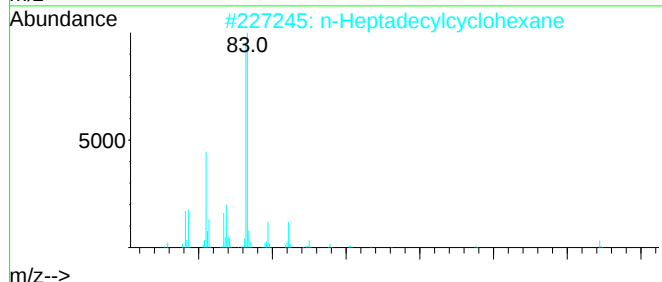
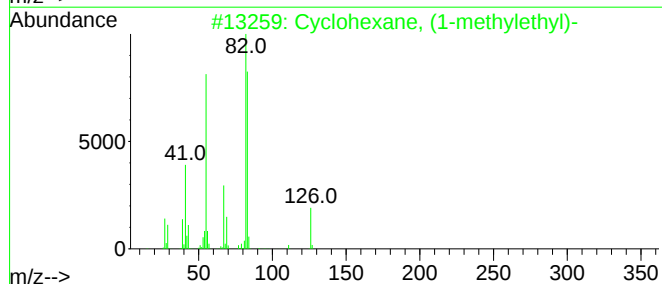
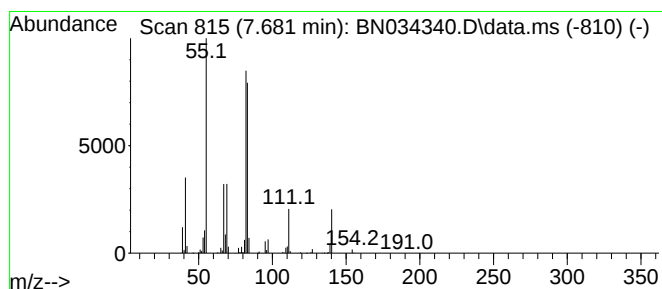
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Peak Number 20 (DEL) Alkane: Cyclic7.681 Concentration Rank 38

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 7.681 | 17.38 ng/ul | 15742900 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 64   |
| 2    |    |   | n-Heptadecylcyclohexane       | 322 | C23H46  | 019781-73-8 | 64   |
| 3    |    |   | Cyclohexane, nonadecyl-       | 350 | C25H50  | 022349-03-7 | 64   |
| 4    |    |   | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 59   |
| 5    |    |   | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 59   |



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Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

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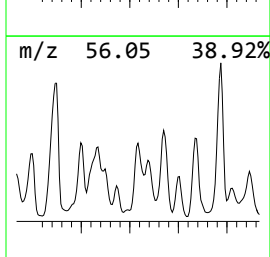
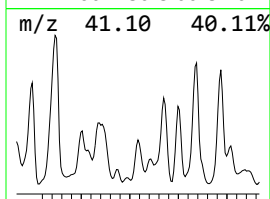
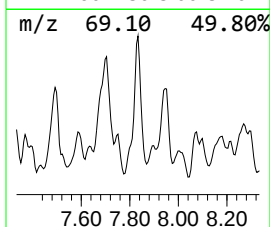
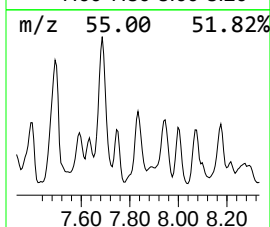
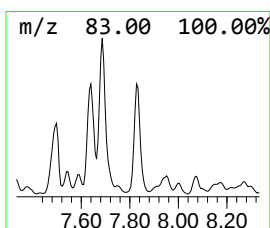
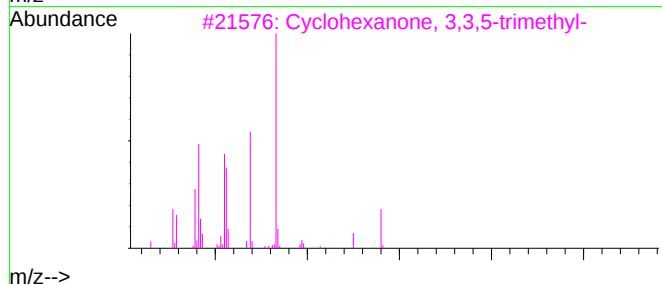
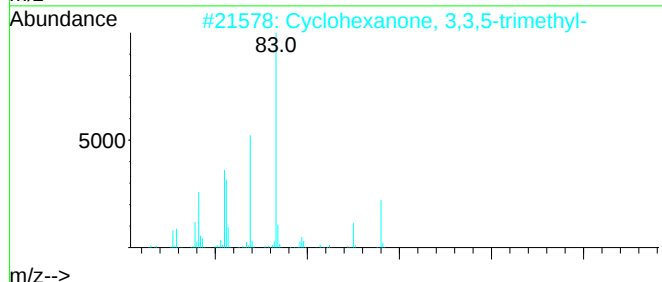
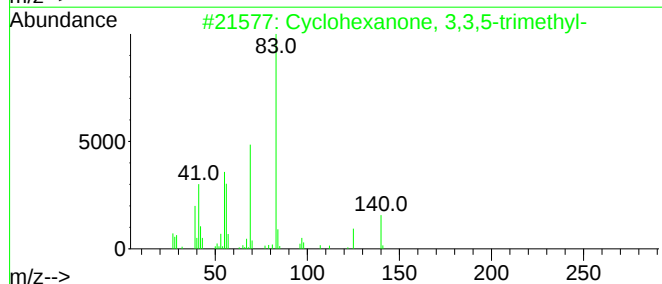
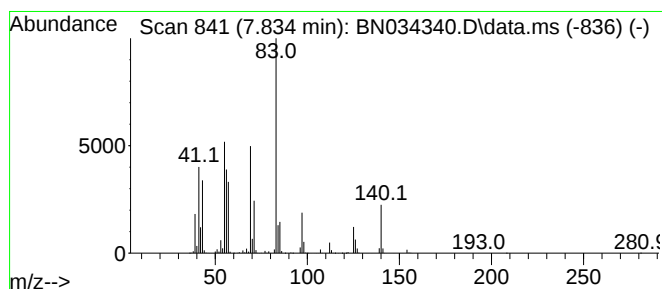
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Peak Number 21 Cyclohexanone, 3,3,5-trimet... Concentration Rank 57

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 7.834 | 8.26 ng/ul | 7480650 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# of 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|-----------|----------------------------------|-----|---------|-------------|------|
| 1         | Cyclohexanone, 3,3,5-trimethyl-  | 140 | C9H16O  | 000873-94-9 | 93   |
| 2         | Cyclohexanone, 3,3,5-trimethyl-  | 140 | C9H16O  | 000873-94-9 | 81   |
| 3         | Cyclohexanone, 3,3,5-trimethyl-  | 140 | C9H16O  | 000873-94-9 | 76   |
| 4         | Cyclohexanone, 3,3,5-trimethyl-  | 140 | C9H16O  | 000873-94-9 | 59   |
| 5         | Cyclopentanone, 2,4,4-trimethyl- | 126 | C8H14O  | 004694-12-6 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

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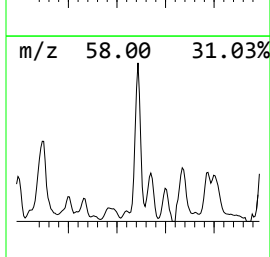
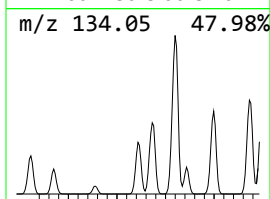
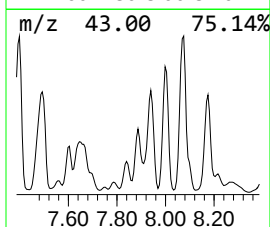
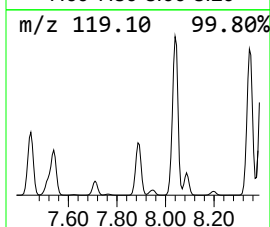
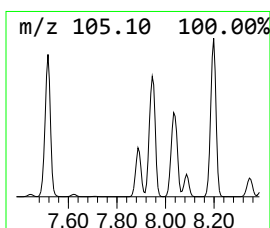
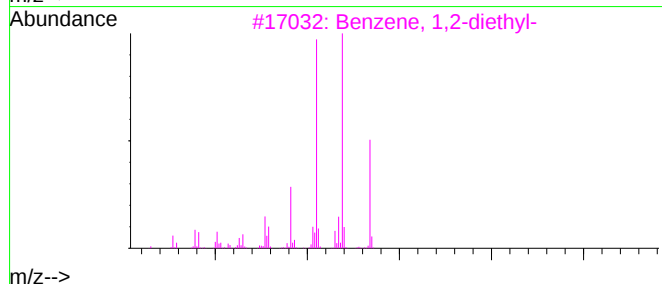
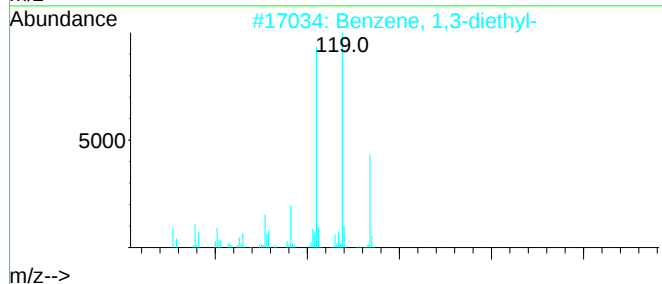
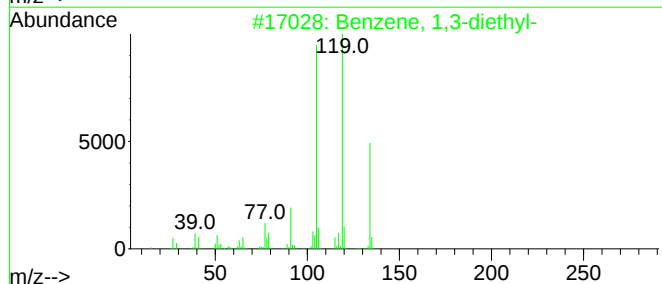
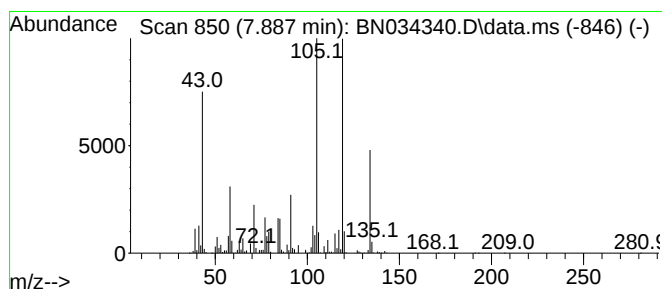
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Peak Number 22 Benzene, 1,3-diethyl- Concentration Rank 54

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 7.887 | 9.18 ng/ul | 8310120 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# of 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|-----------|-----------------------|-----|---------|-------------|------|
| 1         | Benzene, 1,3-diethyl- | 134 | C10H14  | 000141-93-5 | 96   |
| 2         | Benzene, 1,3-diethyl- | 134 | C10H14  | 000141-93-5 | 96   |
| 3         | Benzene, 1,2-diethyl- | 134 | C10H14  | 000135-01-3 | 96   |
| 4         | Benzene, 1,4-diethyl- | 134 | C10H14  | 000105-05-5 | 94   |
| 5         | Benzene, 1,4-diethyl- | 134 | C10H14  | 000105-05-5 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
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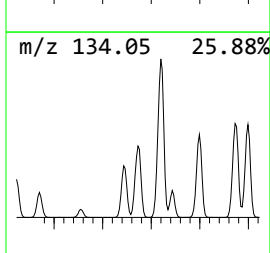
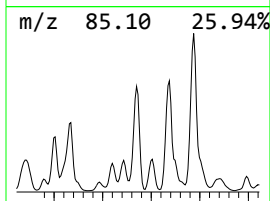
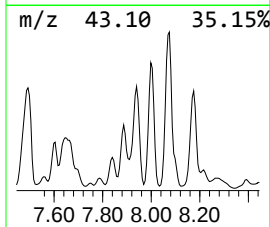
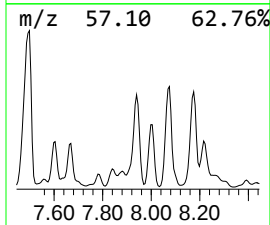
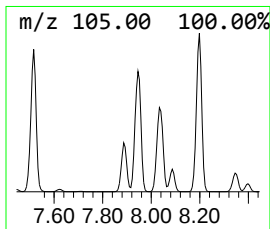
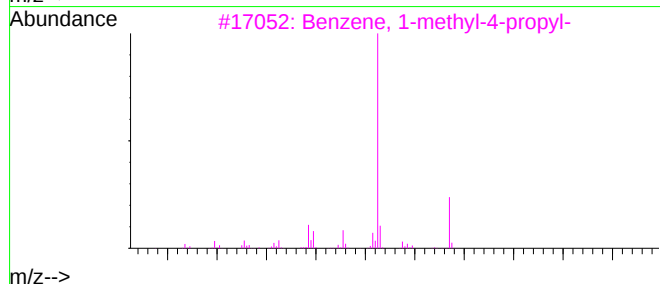
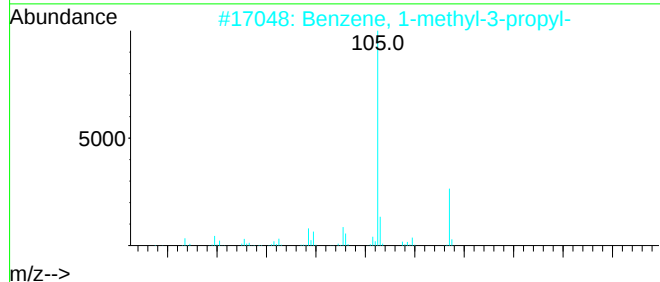
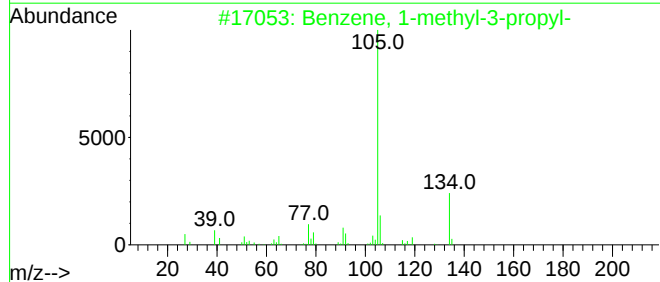
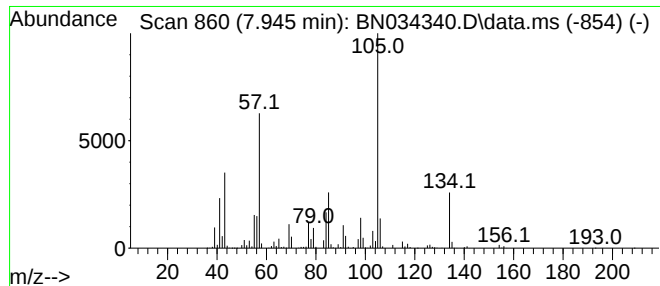
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TIC Integration Parameters: LSCINT.P

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Peak Number 23 Benzene, 1-methyl-3-propyl- Concentration Rank 26

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 7.945 | 24.35 ng/ul | 22055200 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-methyl-3-propyl- | 134 | C10H14  | 001074-43-7 | 55   |
| 2    |    |   | Benzene, 1-methyl-3-propyl- | 134 | C10H14  | 001074-43-7 | 55   |
| 3    |    |   | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 50   |
| 4    |    |   | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 50   |
| 5    |    |   | Benzene, (1-methylpropyl)-  | 134 | C10H14  | 000135-98-8 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

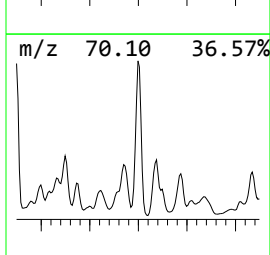
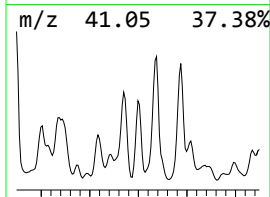
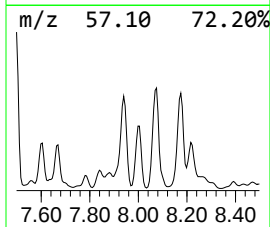
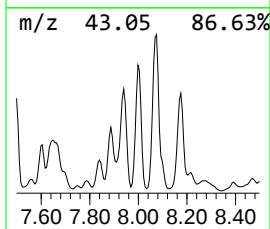
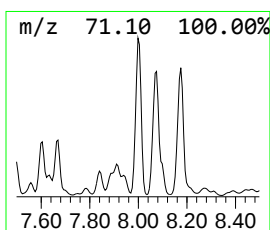
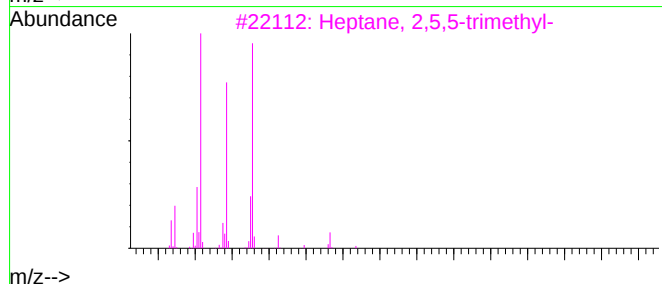
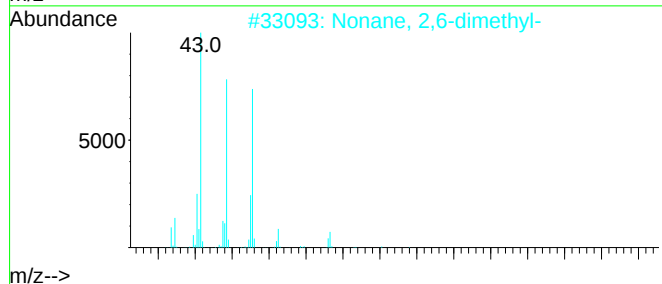
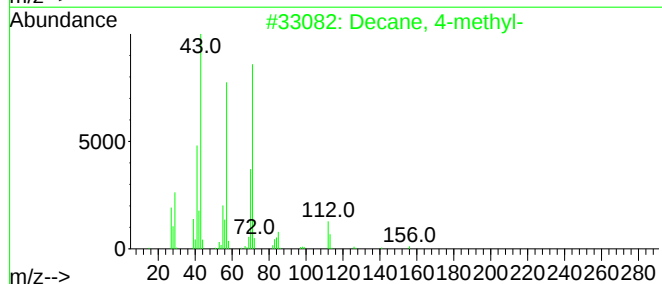
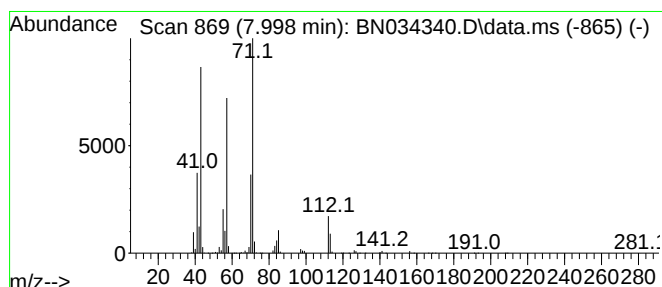
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 43

| R.T.      | EstConc                   | Area     | Relative to ISTD       | R.T.        |      |
|-----------|---------------------------|----------|------------------------|-------------|------|
| 7.998     | 12.65 ng/ul               | 11457600 | 1,4-Dichlorobenzene-d4 | 7.392       |      |
| Hit# of 5 | Tentative ID              | MW       | MolForm                | CAS#        | Qual |
| 1         | Decane, 4-methyl-         | 156      | C11H24                 | 002847-72-5 | 91   |
| 2         | Nonane, 2,6-dimethyl-     | 156      | C11H24                 | 017302-28-2 | 80   |
| 3         | Heptane, 2,5,5-trimethyl- | 142      | C10H22                 | 001189-99-7 | 78   |
| 4         | Decane, 4-methyl-         | 156      | C11H24                 | 002847-72-5 | 76   |
| 5         | Heptane, 3,3,5-trimethyl- | 142      | C10H22                 | 007154-80-5 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

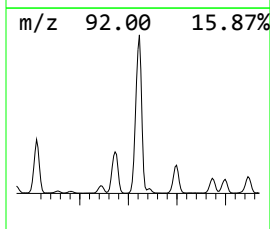
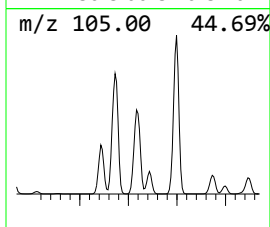
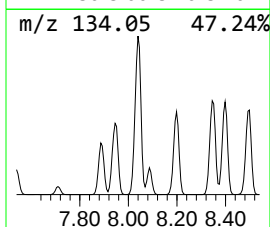
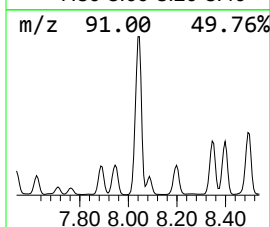
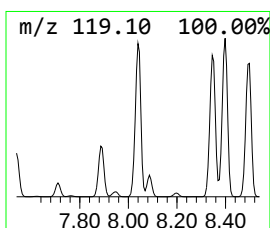
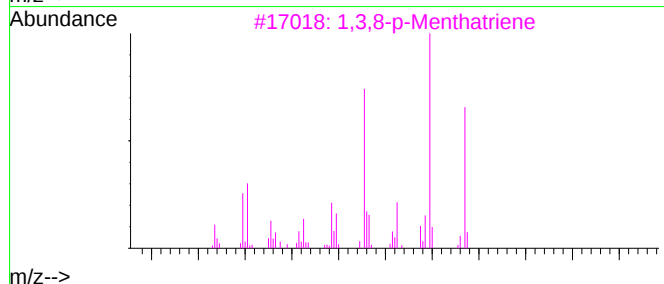
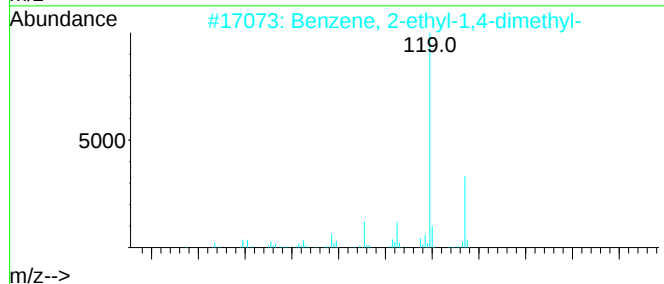
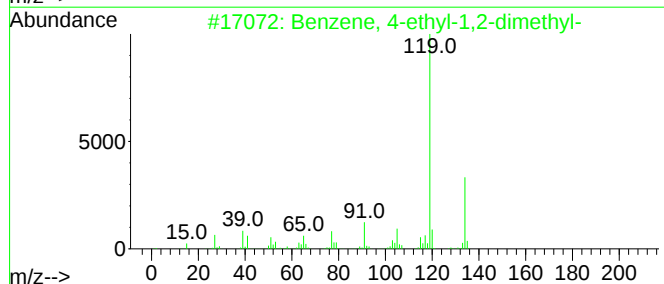
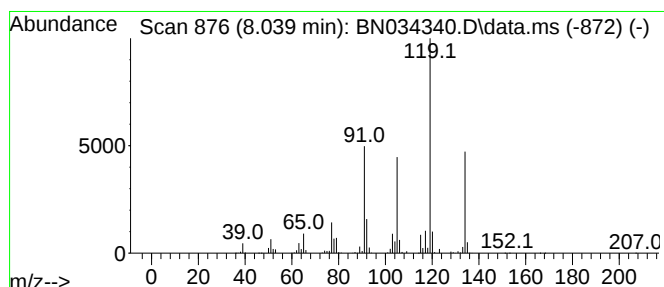
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Peak Number 25 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 47

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 8.039 | 11.69 ng/ul | 10585600 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# of 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|-----------|--------------------------------|-----|---------|-------------|------|
| 1         | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 95   |
| 2         | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 94   |
| 3         | 1,3,8-p-Menthatriene           | 134 | C10H14  | 018368-95-1 | 94   |
| 4         | Benzene, 1,4-diethyl-          | 134 | C10H14  | 000105-05-5 | 91   |
| 5         | Benzene, 1,2-diethyl-          | 134 | C10H14  | 000135-01-3 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

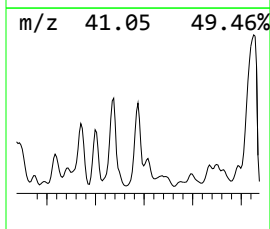
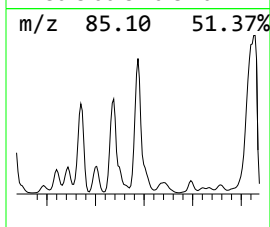
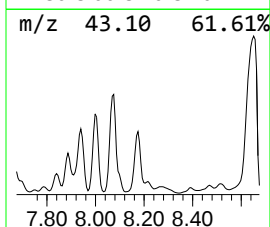
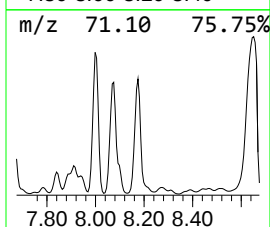
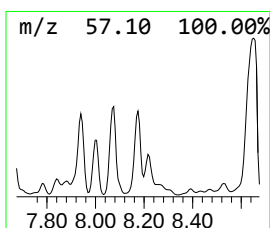
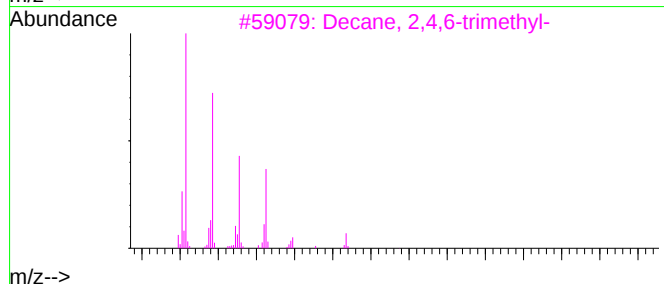
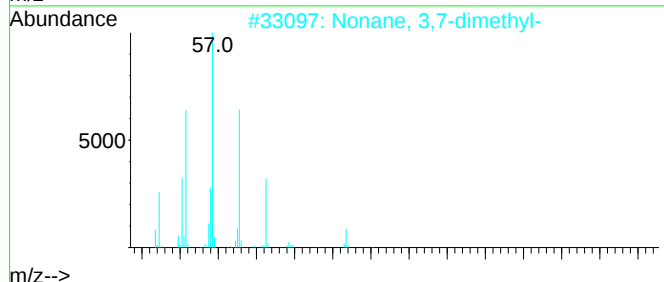
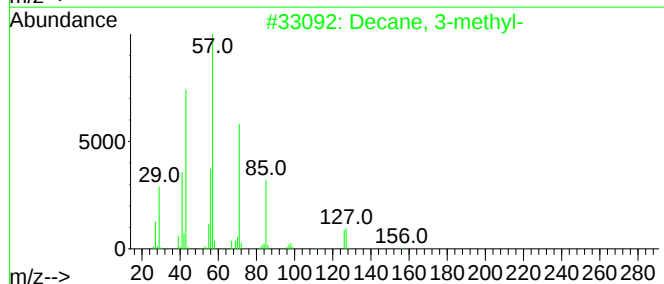
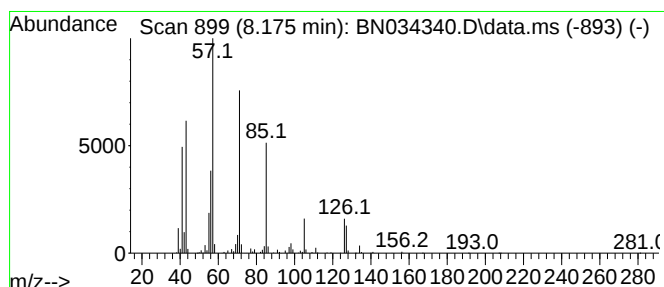
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Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 33

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 8.175 | 19.92 ng/ul | 18037700 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# of 5 | Tentative ID                      | MW  | MolForm   | CAS#         | Qual |
|-----------|-----------------------------------|-----|-----------|--------------|------|
| 1         | Decane, 3-methyl-                 | 156 | C11H24    | 013151-34-3  | 90   |
| 2         | Nonane, 3,7-dimethyl-             | 156 | C11H24    | 017302-32-8  | 64   |
| 3         | Decane, 2,4,6-trimethyl-          | 184 | C13H28    | 062108-27-4  | 59   |
| 4         | Sulfurous acid, butyl nonyl ester | 264 | C13H28O3S | 1000309-17-6 | 59   |
| 5         | Nonane, 1-iodo-                   | 254 | C9H19I    | 004282-42-2  | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

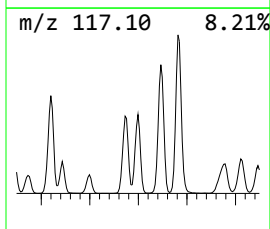
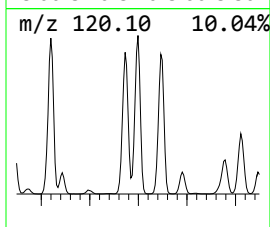
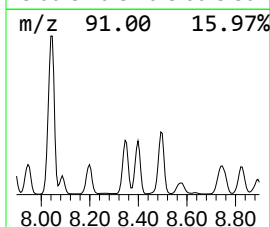
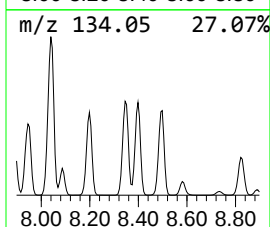
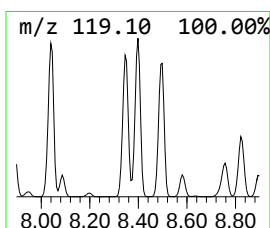
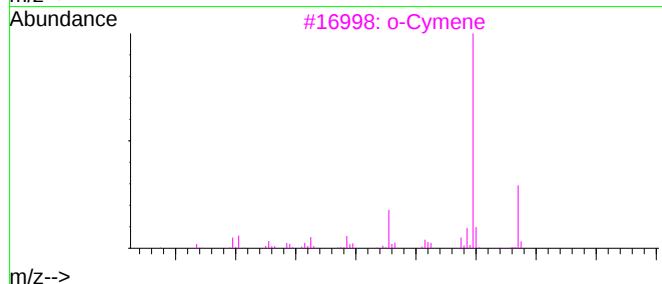
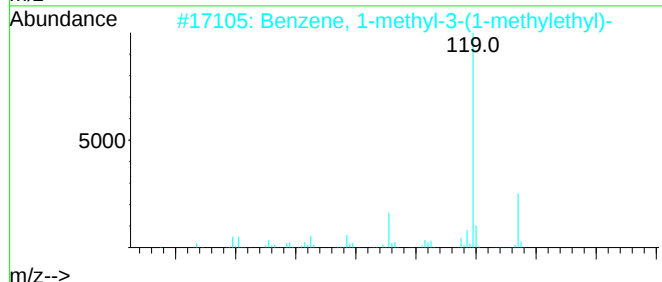
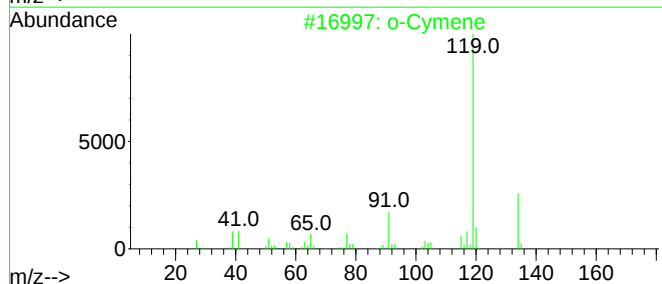
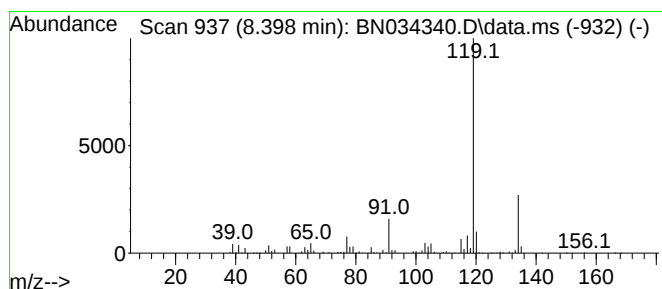
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Peak Number 28 o-Cymene Concentration Rank 44

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 8.398 | 12.29 ng/ul | 11129000 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|-----------|-------------------------------------|-----|---------|-------------|------|
| 1         | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 97   |
| 2         | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 95   |
| 3         | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 95   |
| 4         | Benzene, 1-ethyl-2,4-dimethyl-      | 134 | C10H14  | 000874-41-9 | 94   |
| 5         | p-Cymene                            | 134 | C10H14  | 000099-87-6 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

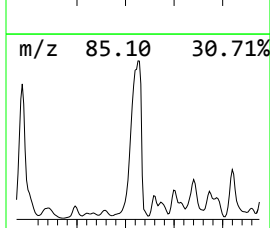
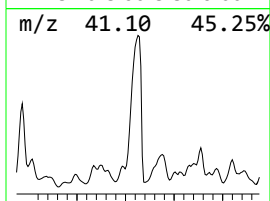
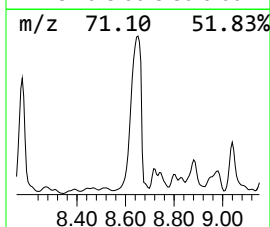
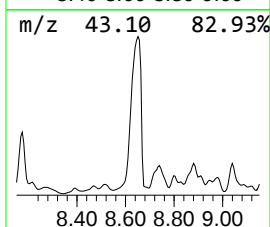
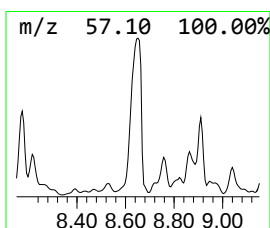
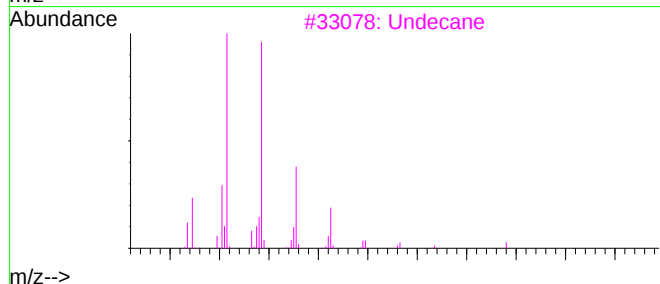
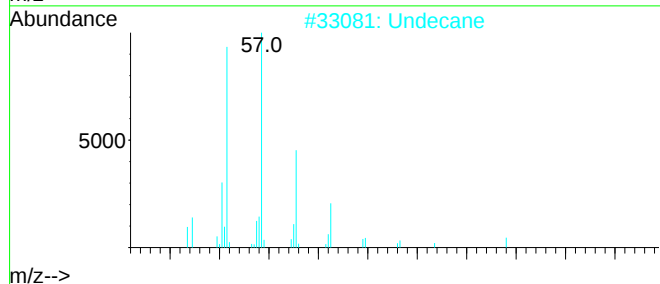
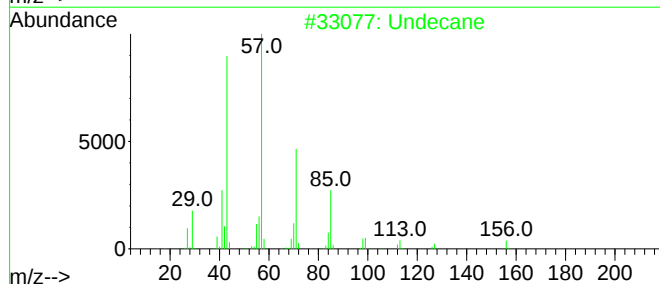
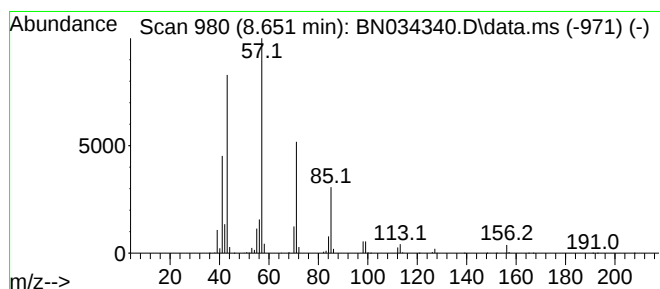
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Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 8.651 | 46.52 ng/ul | 42131700 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|-----------|--------------|-----|---------|-------------|------|
| 1         | Undecane     | 156 | C11H24  | 001120-21-4 | 97   |
| 2         | Undecane     | 156 | C11H24  | 001120-21-4 | 93   |
| 3         | Undecane     | 156 | C11H24  | 001120-21-4 | 91   |
| 4         | Tetradecane  | 198 | C14H30  | 000629-59-4 | 90   |
| 5         | Hexadecane   | 226 | C16H34  | 000544-76-3 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

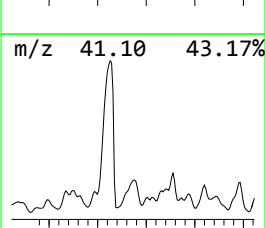
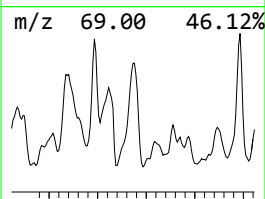
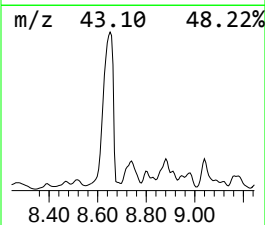
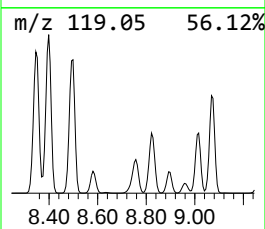
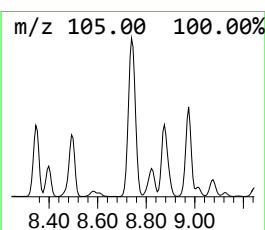
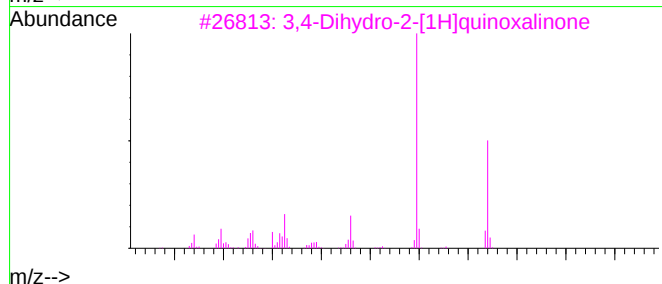
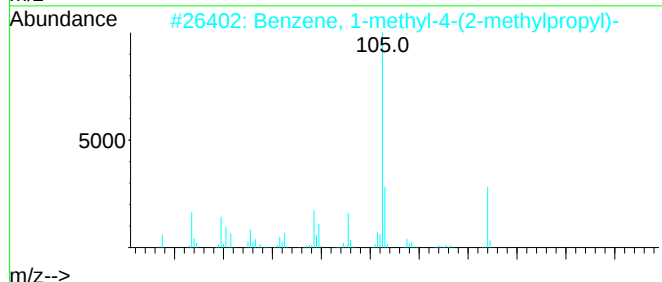
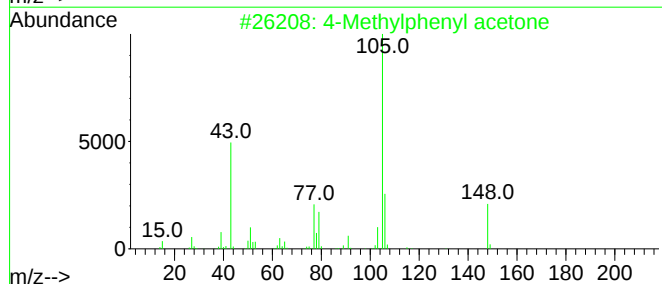
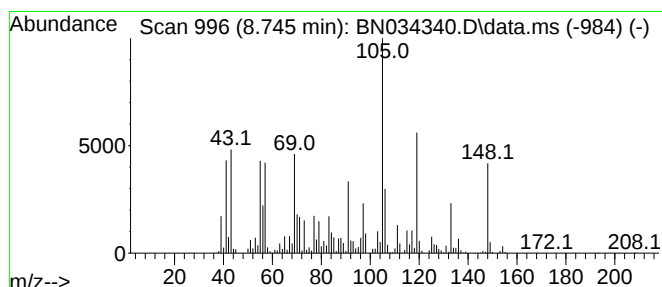
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Peak Number 30 unknown-01 Concentration Rank 30

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 8.745 | 20.90 ng/ul | 18927000 | 1,4-Dichlorobenzene-d4 | 7.392 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 4-Methylphenyl acetone              | 148 | C10H12O | 002096-86-8 | 46   |
| 2    |    |   | Benzene, 1-methyl-4-(2-methylpro... | 148 | C11H16  | 005161-04-6 | 41   |
| 3    |    |   | 3,4-Dihydro-2-[1H]quinoxalinone     | 148 | C8H8N2O | 059564-59-9 | 38   |
| 4    |    |   | 2-Methylindan-2-ol                  | 148 | C10H12O | 033223-84-6 | 38   |
| 5    |    |   | Benzene, (1,1-dimethylpropyl)-      | 148 | C11H16  | 002049-95-8 | 30   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

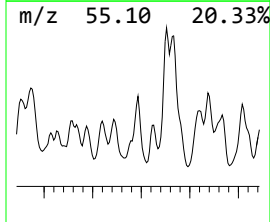
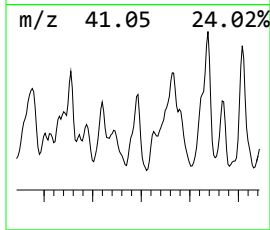
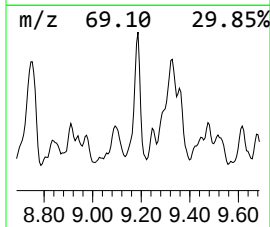
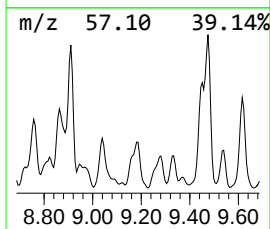
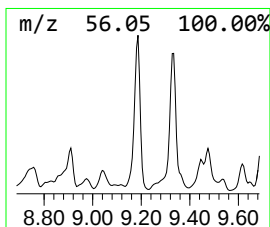
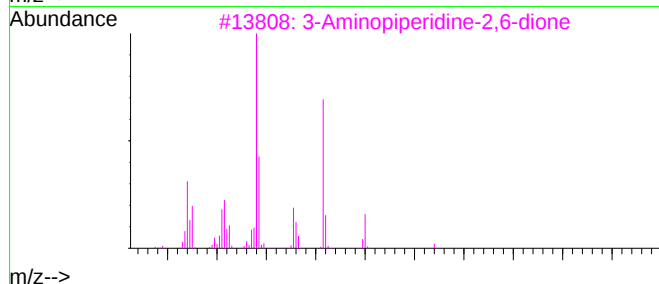
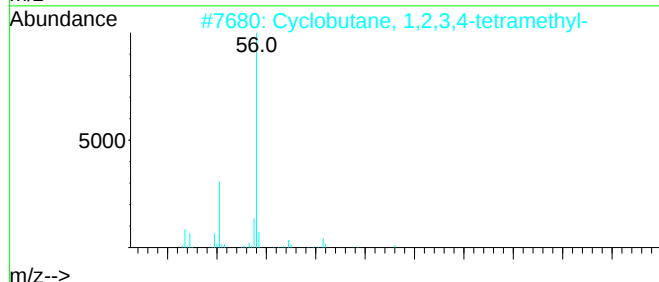
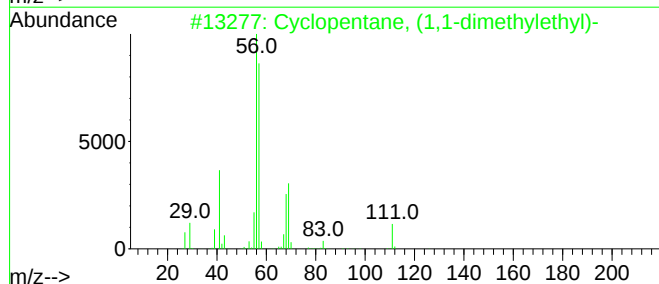
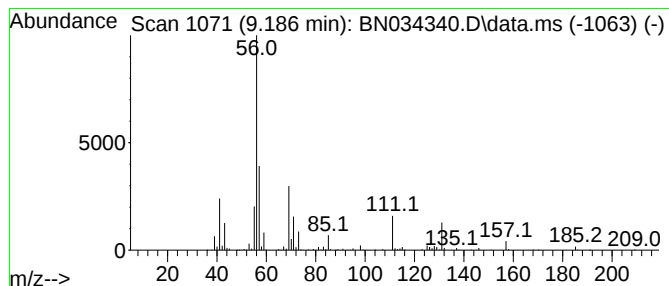
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

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Peak Number 33 (DEL) Alkane: Cyclic9.186 Concentration Rank 72

| R.T.      | EstConc                            | Area    | Relative to ISTD | R.T.        |      |
|-----------|------------------------------------|---------|------------------|-------------|------|
| 9.186     | 3.52 ng/ul                         | 9020800 | Naphthalene-d8   | 10.151      |      |
| Hit# of 5 | Tentative ID                       | MW      | MolForm          | CAS#        | Qual |
| 1         | Cyclopentane, (1,1-dimethylethyl)- | 126     | C9H18            | 003875-52-3 | 38   |
| 2         | Cyclobutane, 1,2,3,4-tetramethyl-  | 112     | C8H16            | 069531-57-3 | 38   |
| 3         | 3-Aminopiperidine-2,6-dione        | 128     | C5H8N2O2         | 002353-44-8 | 36   |
| 4         | 2,4-Dipropyl-5-ethyl-1,3-dioxane   | 200     | C12H24O2         | 006413-83-8 | 36   |
| 5         | 1-Decene, 2-methyl-                | 154     | C11H22           | 013151-27-4 | 28   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

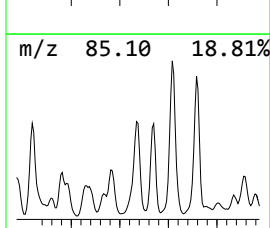
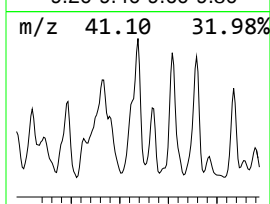
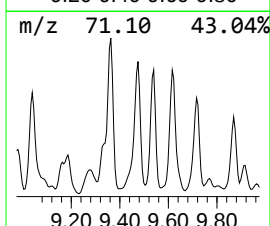
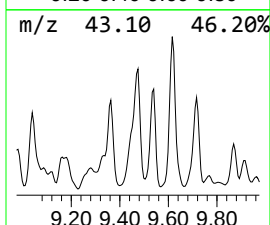
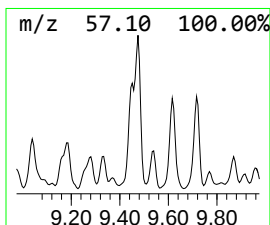
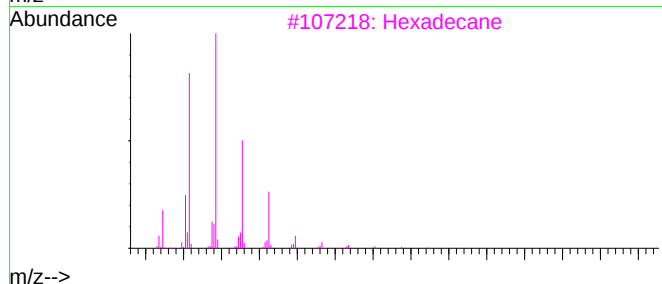
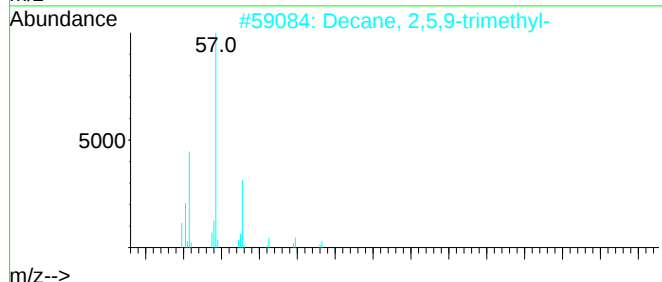
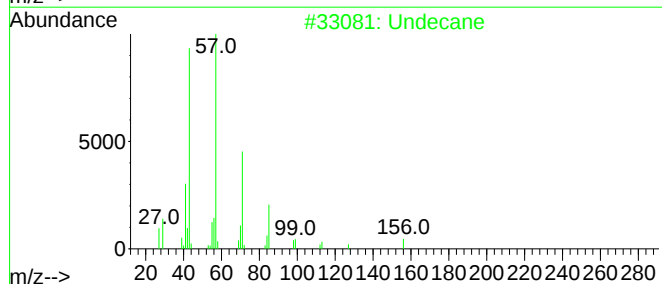
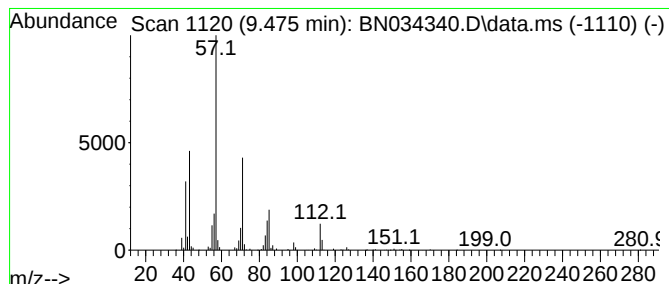
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Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 66

| R.T.  | EstConc    | Area     | Relative to ISTD | R.T.   |
|-------|------------|----------|------------------|--------|
| 9.475 | 5.79 ng/ul | 14837400 | Naphthalene-d8   | 10.151 |

| Hit# of 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|-----------|--------------------------|-----|---------|-------------|------|
| 1         | Undecane                 | 156 | C11H24  | 001120-21-4 | 72   |
| 2         | Decane, 2,5,9-trimethyl- | 184 | C13H28  | 062108-22-9 | 64   |
| 3         | Hexadecane               | 226 | C16H34  | 000544-76-3 | 64   |
| 4         | Pentadecane              | 212 | C15H32  | 000629-62-9 | 53   |
| 5         | Decane, 3,6-dimethyl-    | 170 | C12H26  | 017312-53-7 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

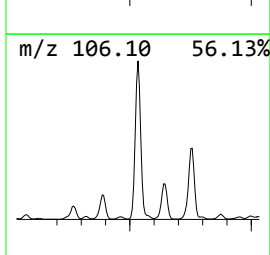
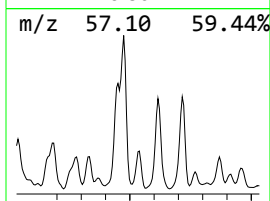
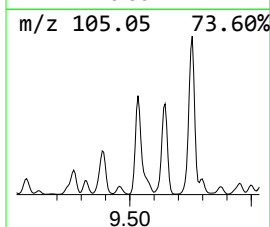
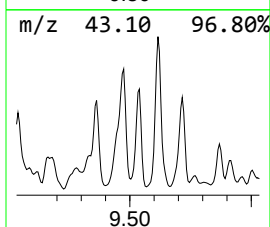
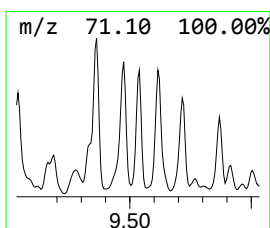
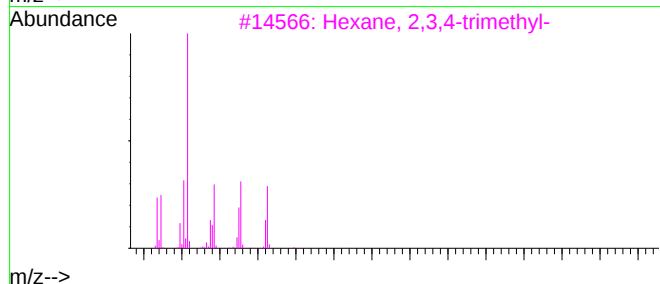
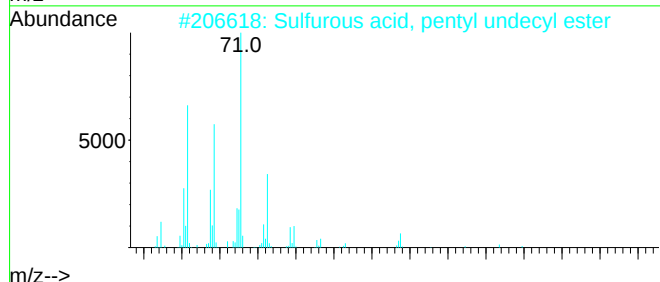
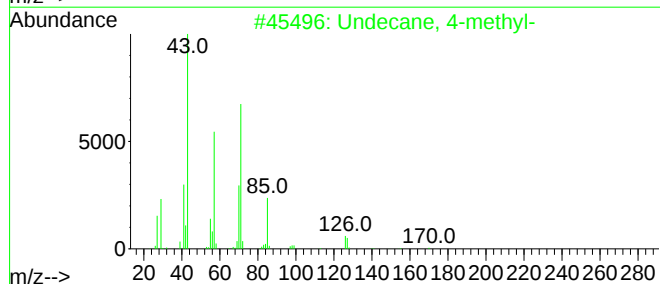
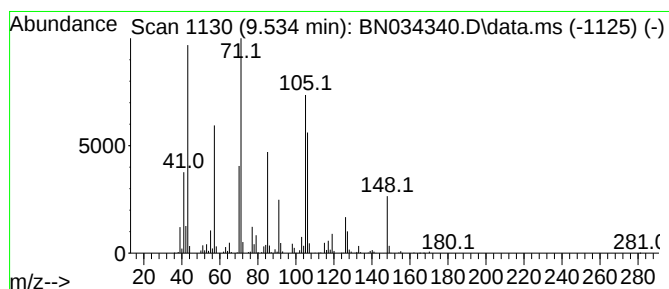
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 73

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|--------------|------|
| 9.534     | 3.27 ng/ul                          | 8392610 | Naphthalene-d8   | 10.151       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#         | Qual |
| 1         | Undecane, 4-methyl-                 | 170     | C12H26           | 002980-69-0  | 58   |
| 2         | Sulfurous acid, pentyl undecyl e... | 306     | C16H34O3S        | 1000309-14-4 | 38   |
| 3         | Hexane, 2,3,4-trimethyl-            | 128     | C9H20            | 000921-47-1  | 35   |
| 4         | 1,3-Dimethylcyclopentanol           | 114     | C7H14O           | 019550-46-0  | 35   |
| 5         | Succinic acid, heptyl tetrahydro... | 300     | C16H28O5         | 1000329-86-5 | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

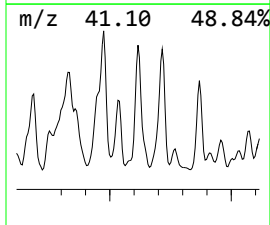
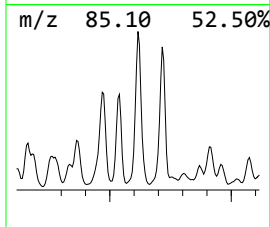
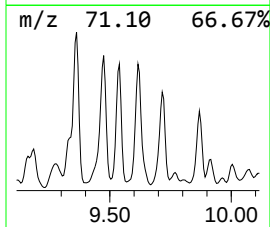
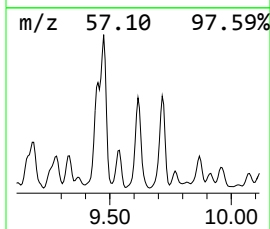
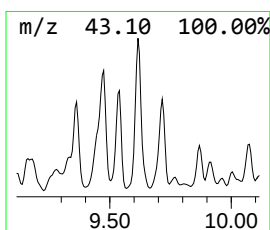
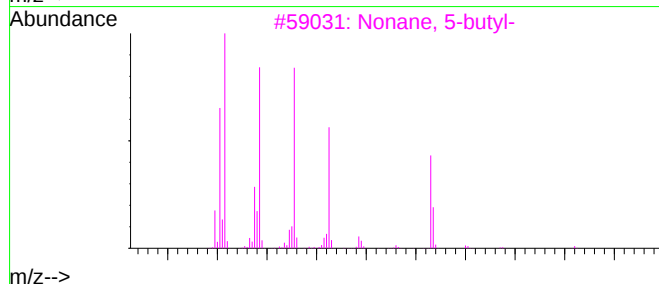
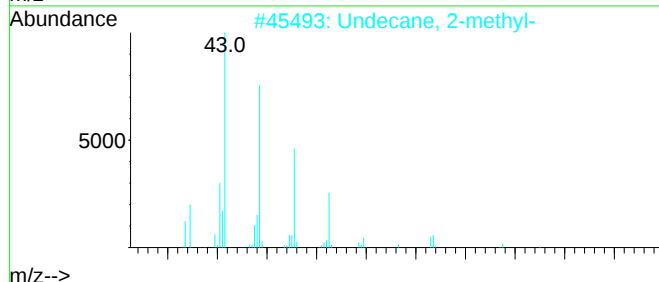
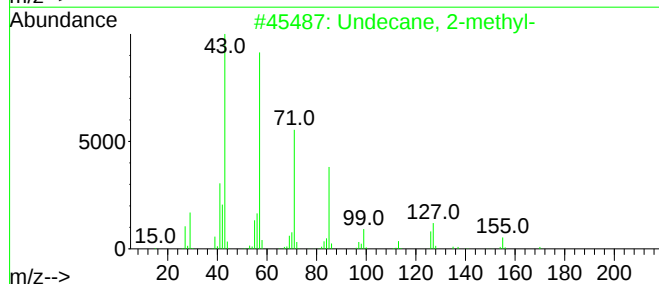
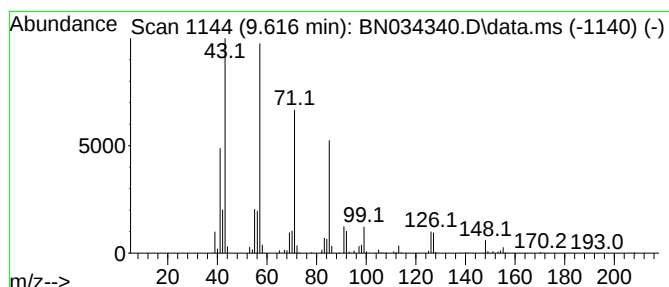
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

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Peak Number 37 (DEL) Alkane: Straight-Chai... Concentration Rank 75

| R.T.      | EstConc             | Area    | Relative to ISTD | R.T.        |      |
|-----------|---------------------|---------|------------------|-------------|------|
| 9.616     | 2.94 ng/ul          | 7532900 | Naphthalene-d8   | 10.151      |      |
| Hit# of 5 | Tentative ID        | MW      | MolForm          | CAS#        | Qual |
| 1         | Undecane, 2-methyl- | 170     | C12H26           | 007045-71-8 | 72   |
| 2         | Undecane, 2-methyl- | 170     | C12H26           | 007045-71-8 | 68   |
| 3         | Nonane, 5-butyl-    | 184     | C13H28           | 017312-63-9 | 64   |
| 4         | Heptadecane         | 240     | C17H36           | 000629-78-7 | 59   |
| 5         | Decane, 2-methyl-   | 156     | C11H24           | 006975-98-0 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

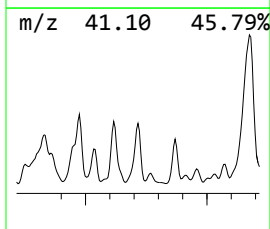
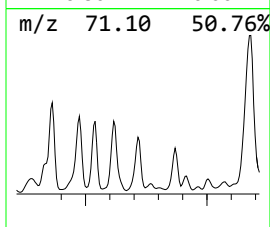
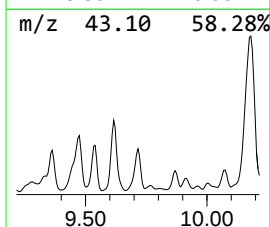
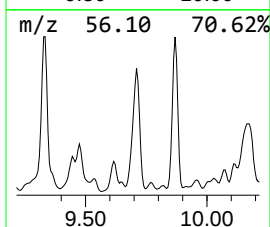
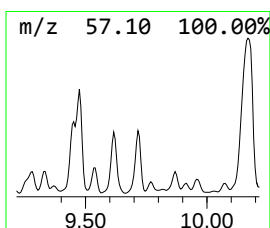
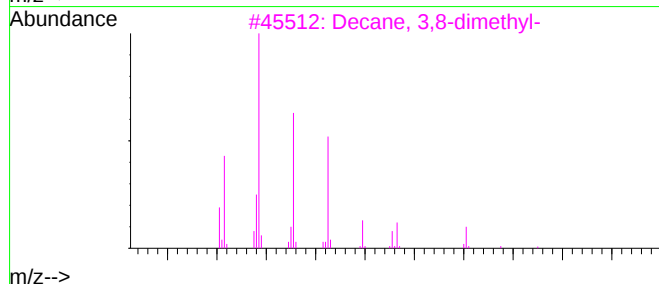
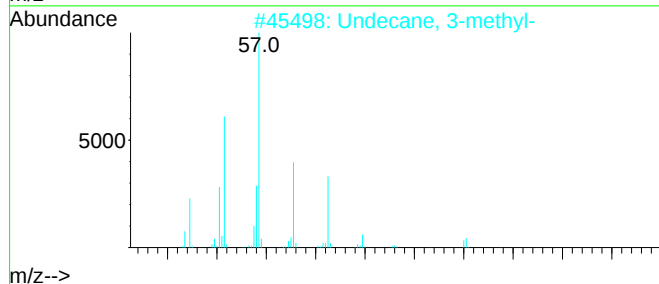
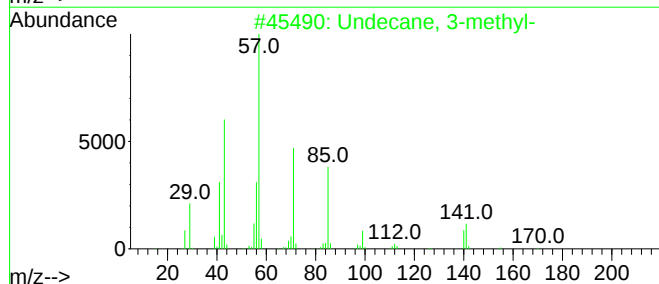
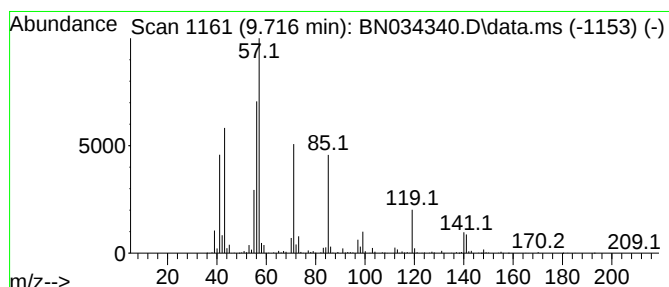
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Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 68

| R.T.  | EstConc    | Area     | Relative to ISTD | R.T.   |
|-------|------------|----------|------------------|--------|
| 9.716 | 4.75 ng/ul | 12185600 | Naphthalene-d8   | 10.151 |

| Hit# of 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|-----------|--------------------------|-----|---------|-------------|------|
| 1         | Undecane, 3-methyl-      | 170 | C12H26  | 001002-43-3 | 76   |
| 2         | Undecane, 3-methyl-      | 170 | C12H26  | 001002-43-3 | 59   |
| 3         | Decane, 3,8-dimethyl-    | 170 | C12H26  | 017312-55-9 | 49   |
| 4         | Pentane, 2,2-dimethyl-   | 100 | C7H16   | 000590-35-2 | 38   |
| 5         | Butane, 2,2,3-trimethyl- | 100 | C7H16   | 000464-06-2 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L

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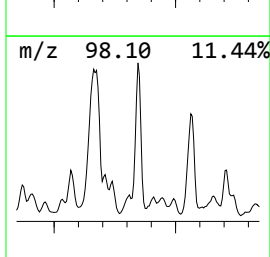
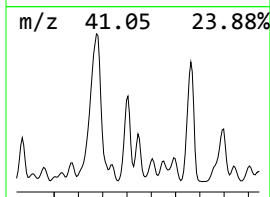
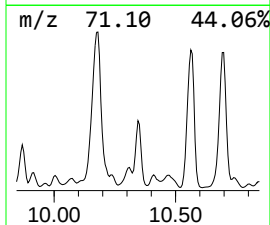
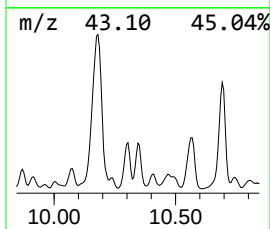
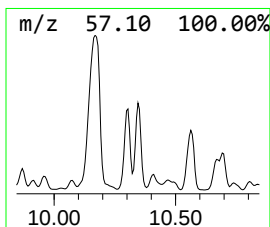
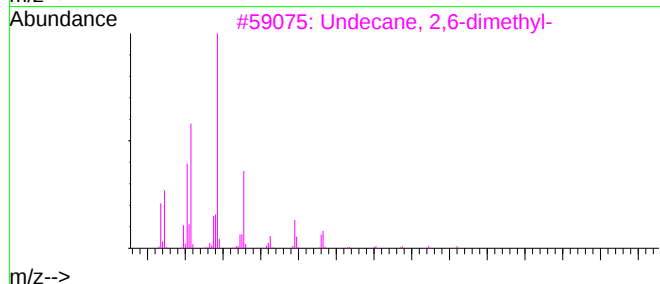
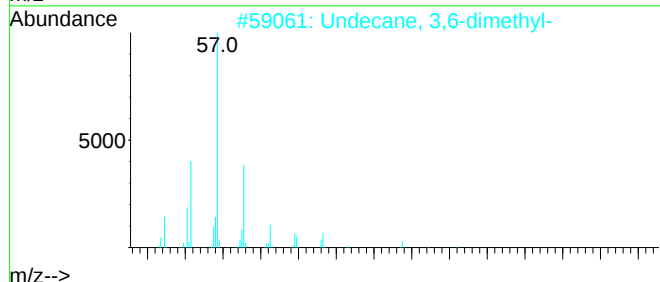
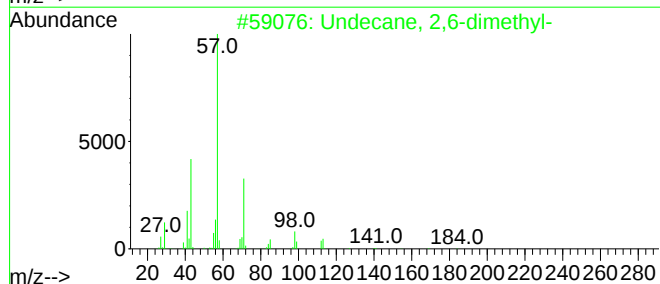
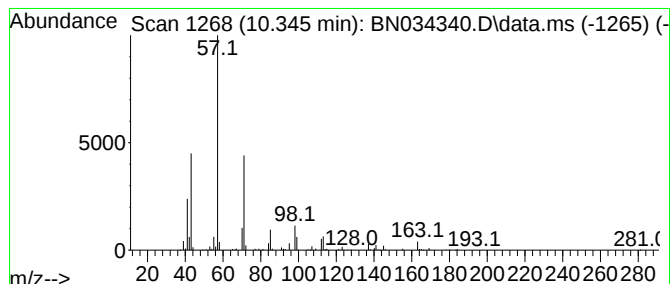
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Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 76

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 10.345 | 2.51 ng/ul | 6447510 | Naphthalene-d8   | 10.151 |

| Hit# of 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|-----------|-----------------------------|-----|---------|-------------|------|
| 1         | Undecane, 2,6-dimethyl-     | 184 | C13H28  | 017301-23-4 | 91   |
| 2         | Undecane, 3,6-dimethyl-     | 184 | C13H28  | 017301-28-9 | 91   |
| 3         | Undecane, 2,6-dimethyl-     | 184 | C13H28  | 017301-23-4 | 78   |
| 4         | Dodecane, 2,7,10-trimethyl- | 212 | C15H32  | 074645-98-0 | 64   |
| 5         | Octane, 2,6-dimethyl-       | 142 | C10H22  | 002051-30-1 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

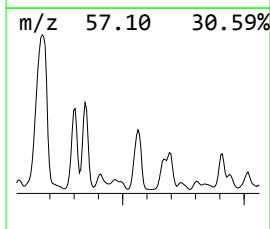
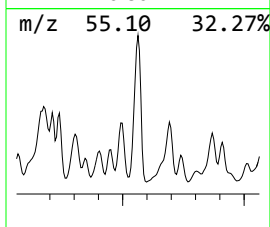
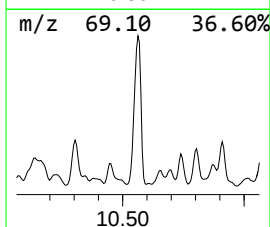
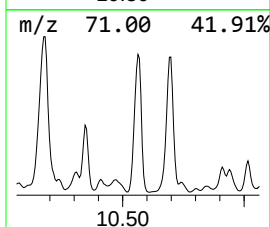
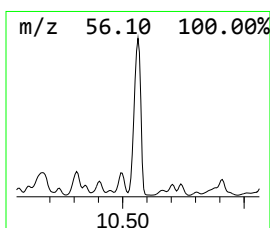
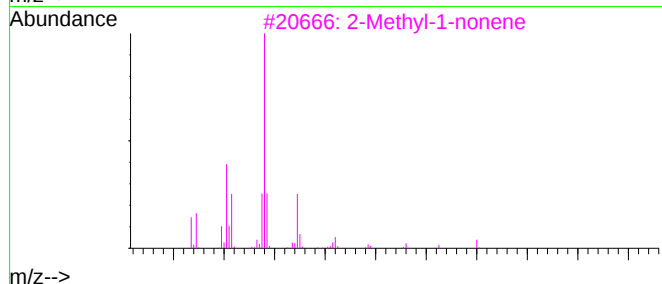
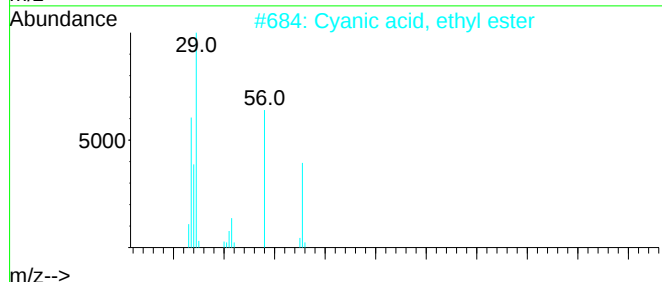
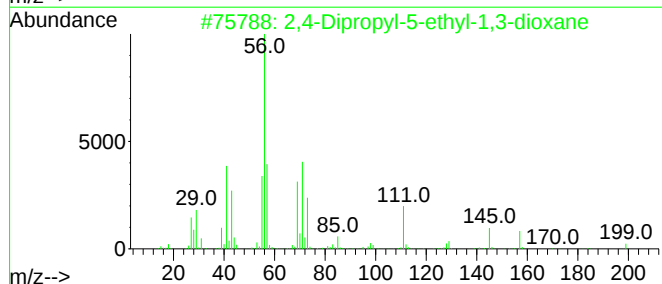
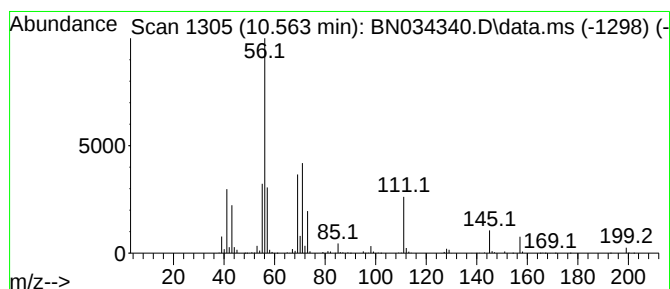
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TIC Integration Parameters: LSCINT.P

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Peak Number 41 unknown-02 Concentration Rank 50

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 10.563 | 10.99 ng/ul | 28183000 | Naphthalene-d8   | 10.151 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2,4-Dipropyl-5-ethyl-1,3-dioxane    | 200 | C12H24O2 | 006413-83-8 | 47   |
| 2    |    |   | Cyanoic acid, ethyl ester           | 71  | C3H5NO   | 000627-48-5 | 40   |
| 3    |    |   | 2-Methyl-1-nonene                   | 140 | C10H20   | 002980-71-4 | 27   |
| 4    |    |   | Cyclohexylamine                     | 99  | C6H13N   | 000108-91-8 | 27   |
| 5    |    |   | Oxirane, 2,3-bis(1-methylethyl)-... | 128 | C8H16O   | 054644-32-5 | 25   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

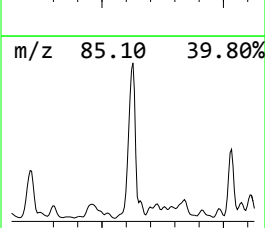
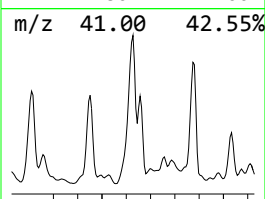
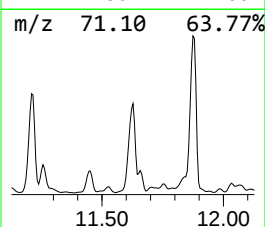
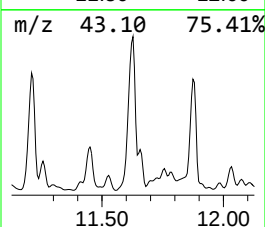
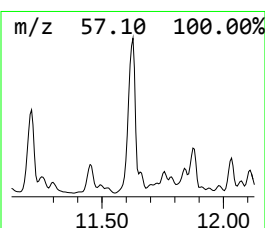
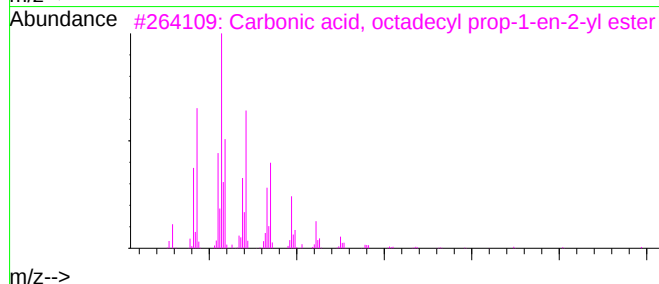
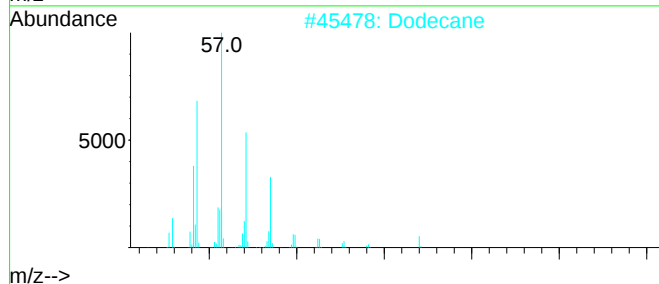
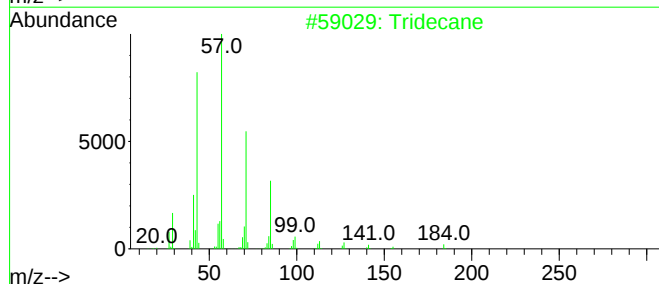
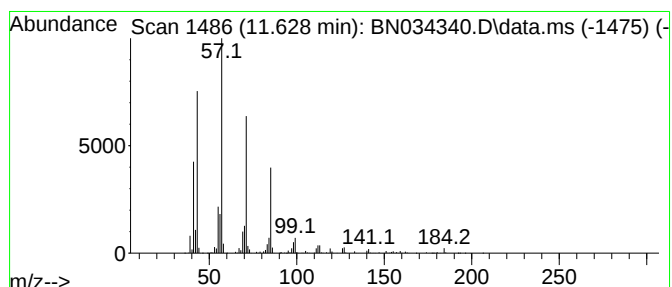
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Peak Number 46 (DEL) Alkane: Straight-Chai... Concentration Rank 49

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 11.628 | 11.62 ng/ul | 29795500 | Naphthalene-d8   | 10.151 |

| Hit# of 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|-----------|-------------------------------------|-----|----------|--------------|------|
| 1         | Tridecane                           | 184 | C13H28   | 000629-50-5  | 96   |
| 2         | Dodecane                            | 170 | C12H26   | 000112-40-3  | 90   |
| 3         | Carbonic acid, octadecyl prop-1-... | 354 | C22H42O3 | 1000383-11-5 | 90   |
| 4         | Tetradecane                         | 198 | C14H30   | 000629-59-4  | 90   |
| 5         | Tridecane                           | 184 | C13H28   | 000629-50-5  | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L

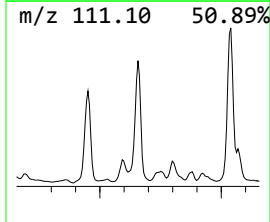
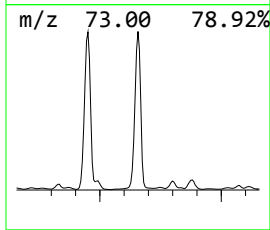
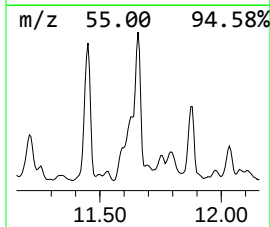
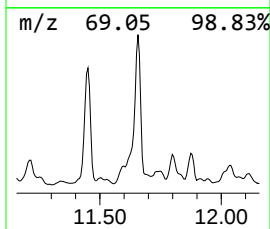
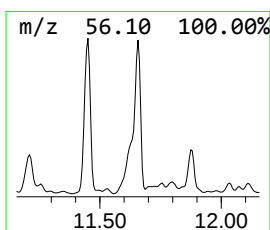
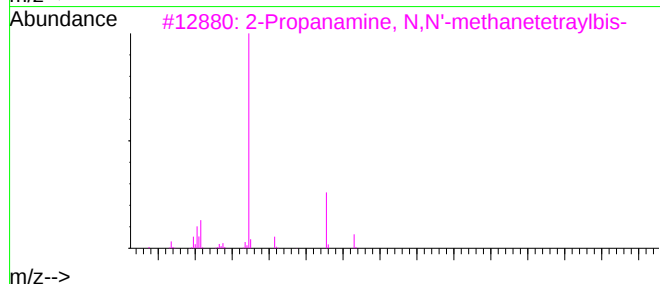
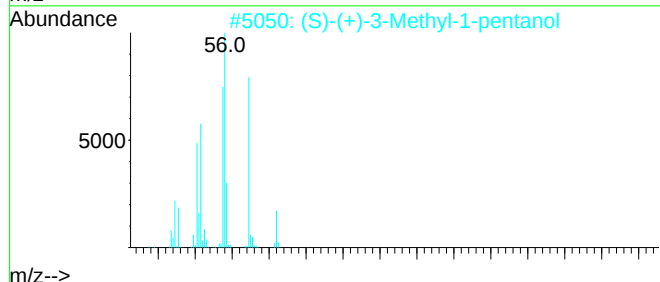
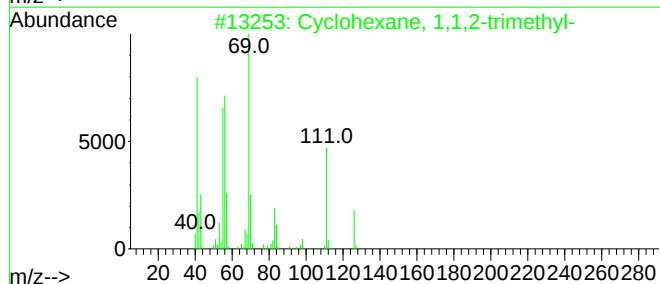
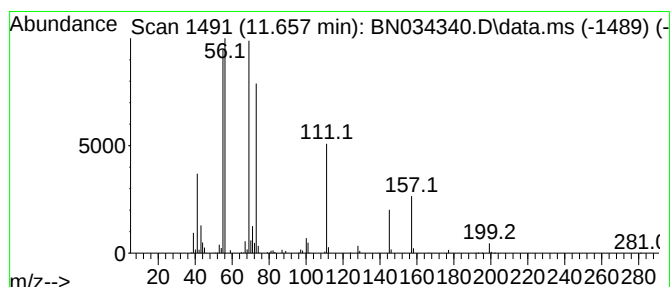
TIC Integration Parameters: LSCINT.P

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Peak Number 47 (DEL) Alkane: Cyclic11.657 Concentration Rank 69

| R.T.   | EstConc    | Area     | Relative to ISTD | R.T.   |
|--------|------------|----------|------------------|--------|
| 11.657 | 4.65 ng/ul | 11914700 | Naphthalene-d8   | 10.151 |

| Hit# of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|-----------|-------------------------------------|-----|---------|-------------|------|
| 1         | Cyclohexane, 1,1,2-trimethyl-       | 126 | C9H18   | 007094-26-0 | 38   |
| 2         | (S)-(+)-3-Methyl-1-pentanol         | 102 | C6H14O  | 042072-39-9 | 28   |
| 3         | 2-Propanamine, N,N'-methanetetra... | 126 | C7H14N2 | 000693-13-0 | 27   |
| 4         | Undecanol-4                         | 172 | C11H24O | 004272-06-4 | 17   |
| 5         | Neopentyl glycol                    | 104 | C5H12O2 | 000126-30-7 | 16   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

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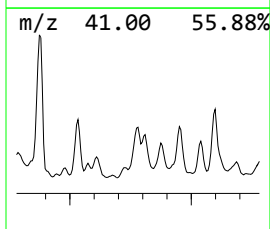
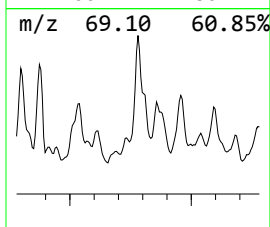
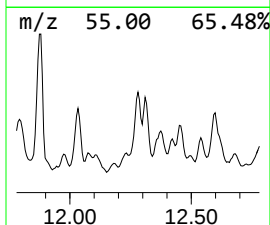
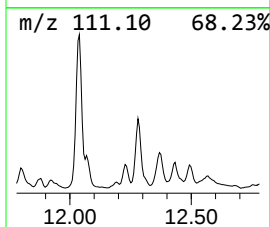
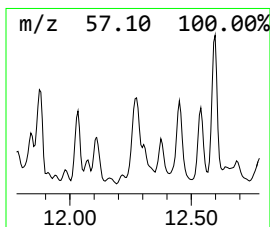
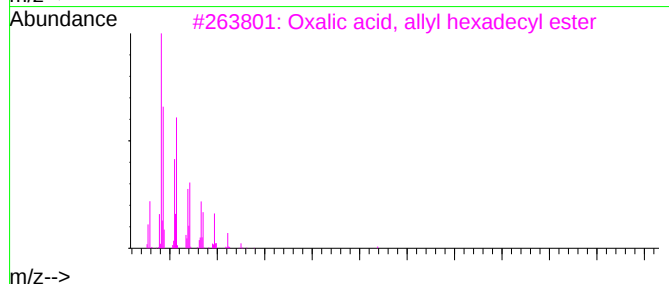
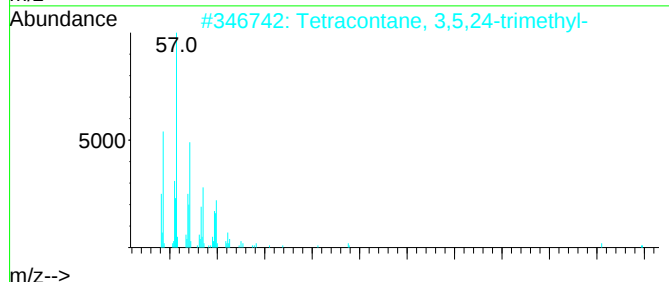
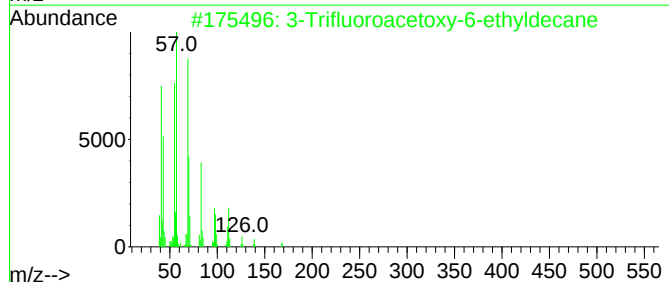
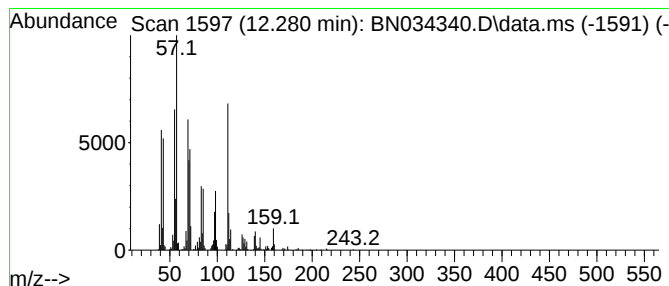
TIC Integration Parameters: LSCINT.P

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Peak Number 48 unknown-03

Concentration Rank 34

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|--------------|------|
| 12.280    | 19.60 ng/ul                         | 7066660 | Acenaphthene-d10 | 14.051       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#         | Qual |
| 1         | 3-Trifluoroacetoxy-6-ethyldecane    | 282     | C14H25F3O2       | 116436-59-0  | 41   |
| 2         | Tetracontane, 3,5,24-trimethyl-     | 605     | C43H88           | 055162-61-3  | 38   |
| 3         | Oxalic acid, allyl hexadecyl ester  | 354     | C21H38O4         | 1000309-24-4 | 35   |
| 4         | Cyclodecane                         | 140     | C10H20           | 000293-96-9  | 30   |
| 5         | 1H-Pyrazole, 3-ethyl-4,5-dihydro... | 126     | C7H14N2          | 075011-91-5  | 25   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

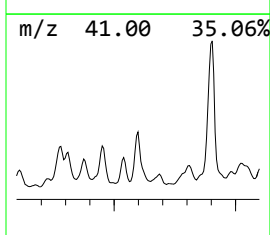
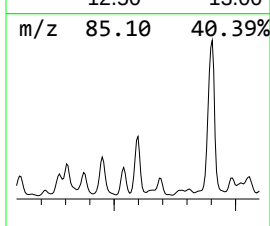
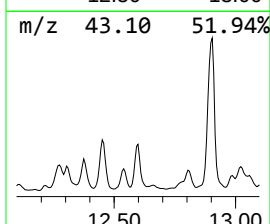
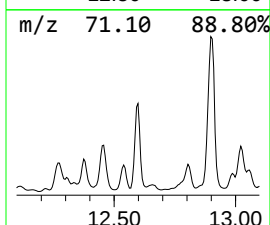
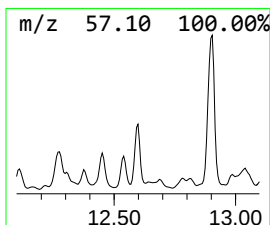
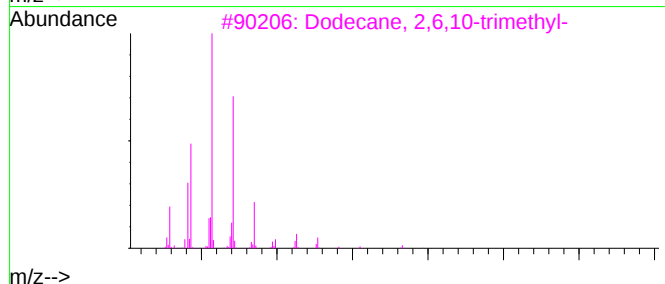
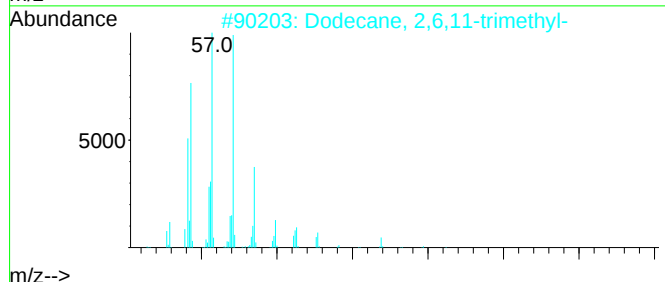
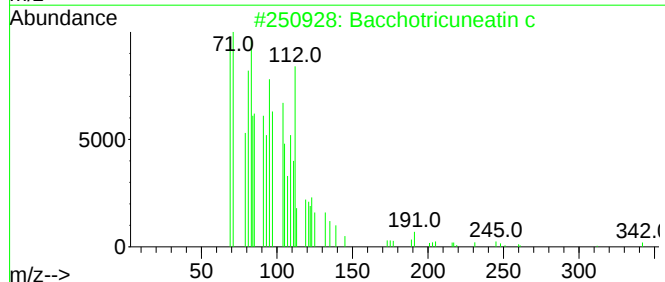
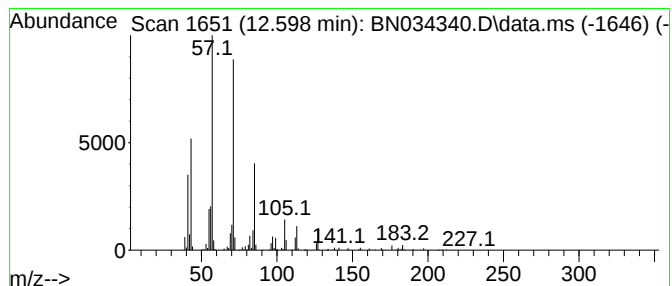
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

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Peak Number 49 Bacchotricuneatin c Concentration Rank 22

| R.T.      | EstConc                     | Area     | Relative to ISTD |          | R.T.        |      |
|-----------|-----------------------------|----------|------------------|----------|-------------|------|
| 12.598    | 29.76 ng/ul                 | 10729400 | Acenaphthene-d10 |          | 14.051      |      |
| Hit# of 5 | Tentative ID                |          | MW               | MolForm  | CAS#        | Qual |
| 1         | Bacchotricuneatin c         |          | 342              | C20H22O5 | 066563-30-2 | 72   |
| 2         | Dodecane, 2,6,11-trimethyl- |          | 212              | C15H32   | 031295-56-4 | 72   |
| 3         | Dodecane, 2,6,10-trimethyl- |          | 212              | C15H32   | 003891-98-3 | 72   |
| 4         | Decane, 5-propyl-           |          | 184              | C13H28   | 017312-62-8 | 64   |
| 5         | Dodecane, 2,7,10-trimethyl- |          | 212              | C15H32   | 074645-98-0 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

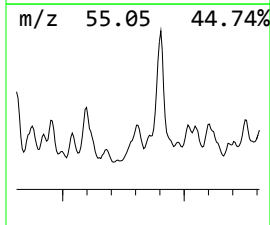
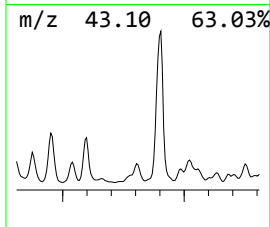
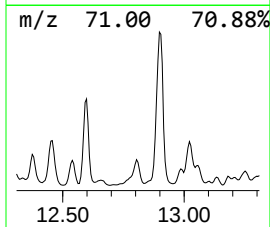
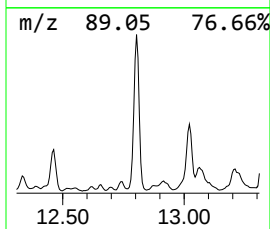
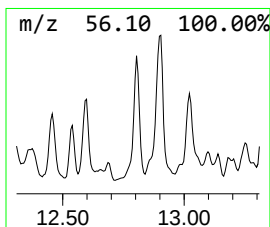
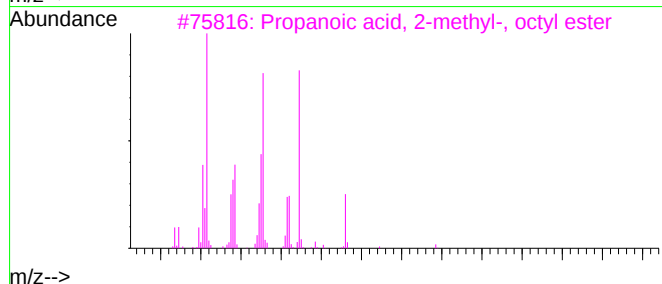
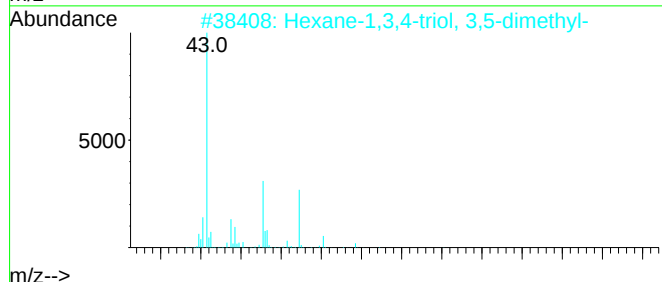
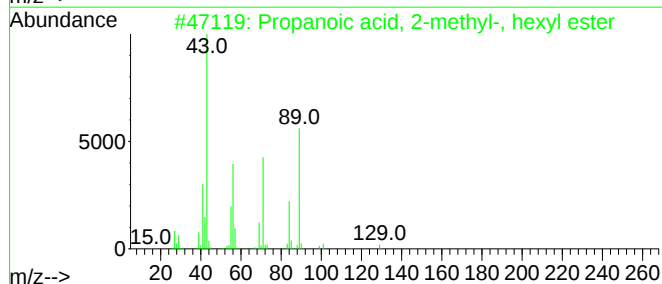
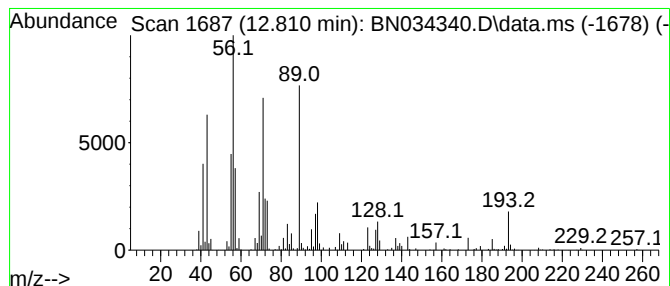
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

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Peak Number 50 Propanoic acid, 2-methyl-, ... Concentration Rank 39

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|--------------|------|
| 12.810    | 16.97 ng/ul                         | 6120330 | Acenaphthene-d10 | 14.051       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#         | Qual |
| 1         | Propanoic acid, 2-methyl-, hexyl... | 172     | C10H20O2         | 002349-07-7  | 59   |
| 2         | Hexane-1,3,4-triol, 3,5-dimethyl-   | 162     | C8H18O3          | 1000268-02-6 | 59   |
| 3         | Propanoic acid, 2-methyl-, octyl... | 200     | C12H24O2         | 000109-15-9  | 53   |
| 4         | Propanoic acid, 2-methyl-, hepty... | 186     | C11H22O2         | 002349-13-5  | 45   |
| 5         | Butyric acid, dodecyl ester         | 256     | C16H32O2         | 003724-61-6  | 42   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

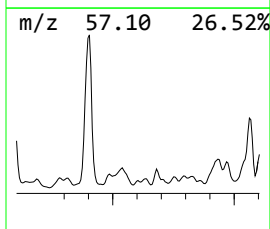
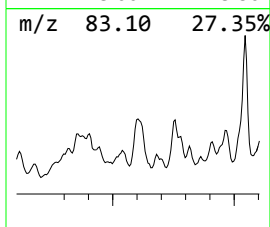
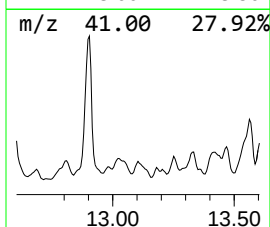
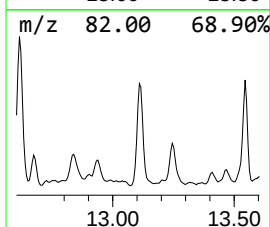
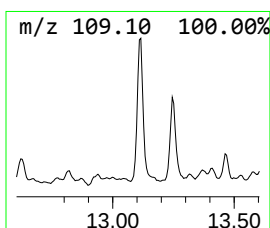
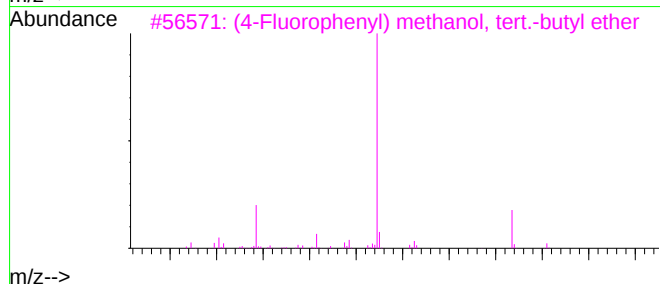
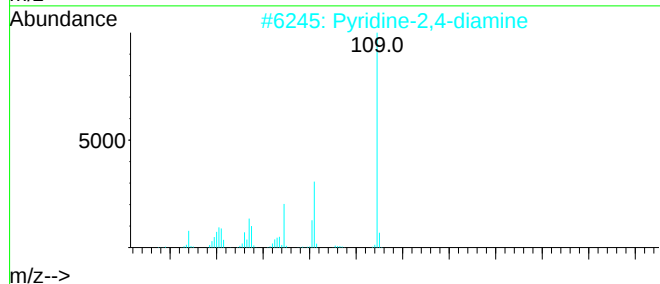
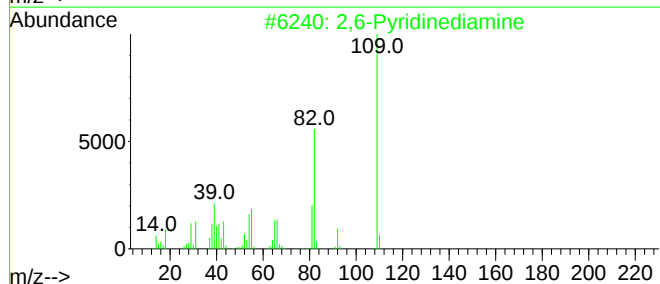
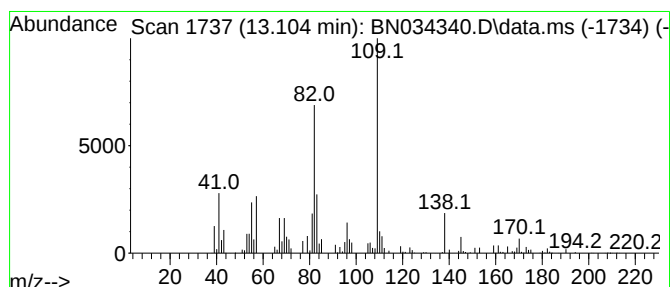
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 51 2,6-Pyridinediamine Concentration Rank 35

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|--------------|------|
| 13.104    | 19.39 ng/ul                         | 6992670 | Acenaphthene-d10 | 14.051       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#         | Qual |
| 1         | 2,6-Pyridinediamine                 | 109     | C5H7N3           | 000141-86-6  | 50   |
| 2         | Pyridine-2,4-diamine                | 109     | C5H7N3           | 000461-88-1  | 50   |
| 3         | (4-Fluorophenyl) methanol, tert...  | 182     | C11H15FO         | 1000374-64-1 | 43   |
| 4         | 2-Cyclopenten-1-one, 3,4,4-trime... | 124     | C8H12O           | 030434-65-2  | 43   |
| 5         | Phenol, 3-amino-                    | 109     | C6H7NO           | 000591-27-5  | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

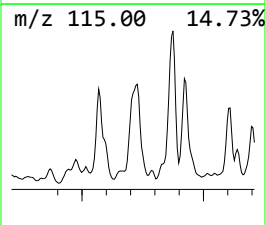
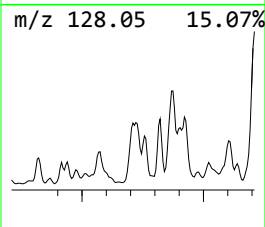
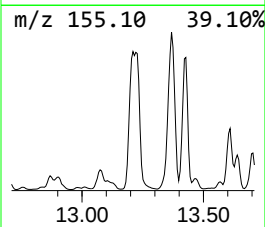
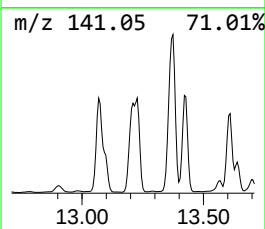
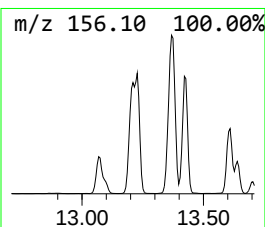
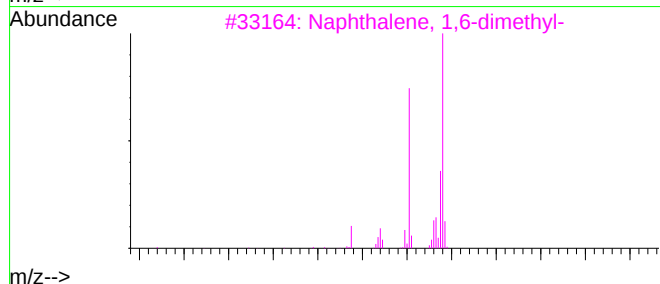
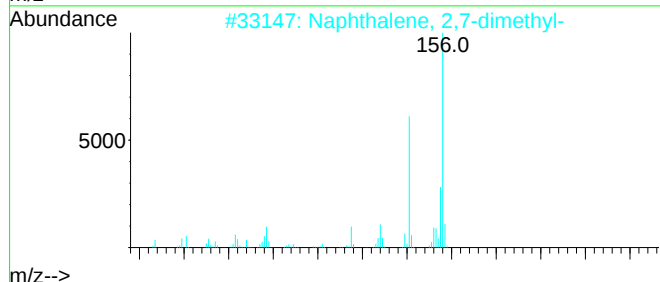
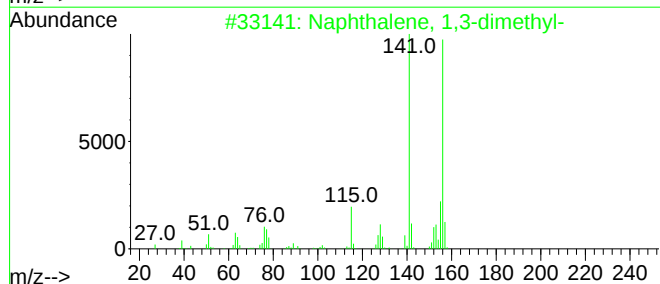
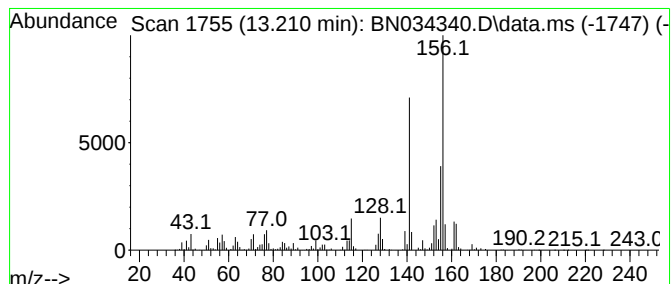
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TIC Integration Parameters: LSCINT.P

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Peak Number 52 Naphthalene, 1,3-dimethyl- Concentration Rank 20

| R.T.      | EstConc                    | Area     | Relative to ISTD | R.T.        |      |
|-----------|----------------------------|----------|------------------|-------------|------|
| 13.210    | 30.74 ng/ul                | 11083800 | Acenaphthene-d10 | 14.051      |      |
| Hit# of 5 | Tentative ID               | MW       | MolForm          | CAS#        | Qual |
| 1         | Naphthalene, 1,3-dimethyl- | 156      | C12H12           | 000575-41-7 | 96   |
| 2         | Naphthalene, 2,7-dimethyl- | 156      | C12H12           | 000582-16-1 | 96   |
| 3         | Naphthalene, 1,6-dimethyl- | 156      | C12H12           | 000575-43-9 | 96   |
| 4         | Naphthalene, 2,3-dimethyl- | 156      | C12H12           | 000581-40-8 | 96   |
| 5         | Naphthalene, 2,6-dimethyl- | 156      | C12H12           | 000581-42-0 | 95   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

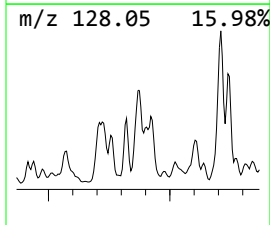
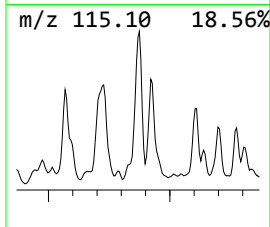
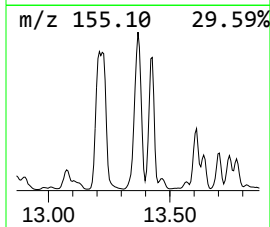
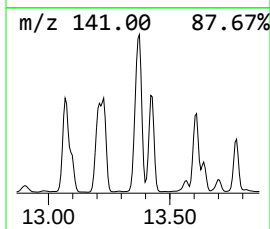
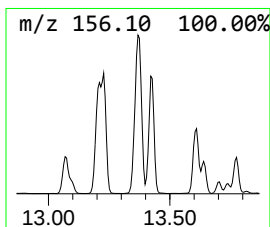
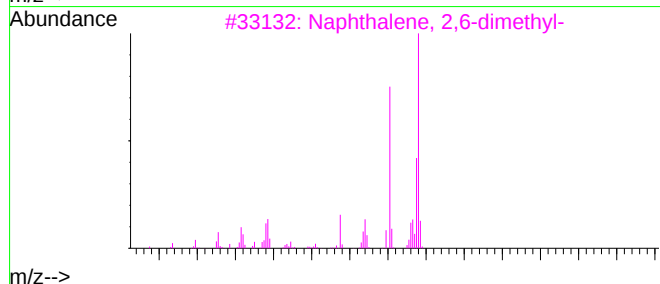
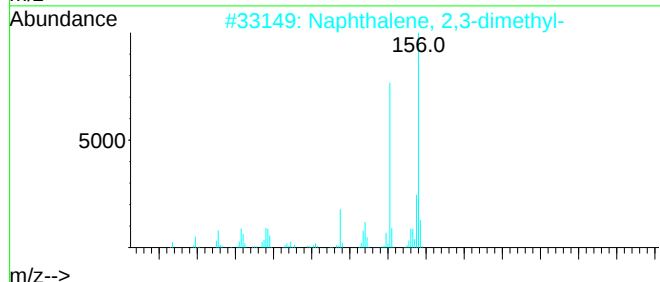
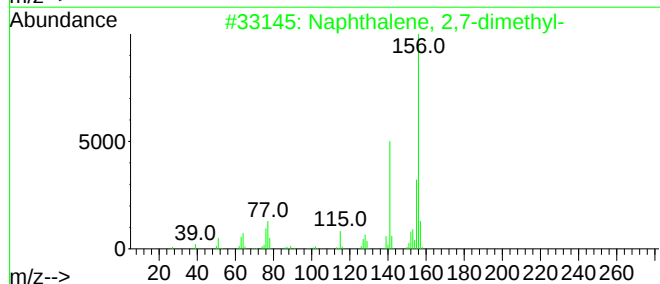
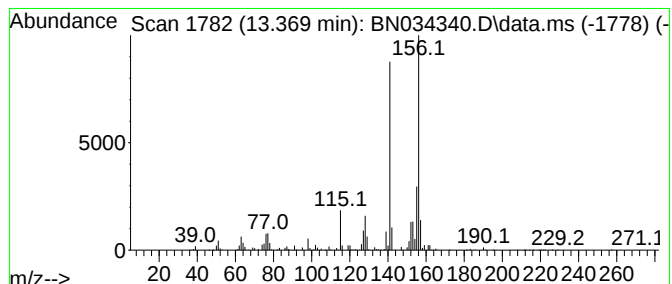
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

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Peak Number 53 Naphthalene, 2,7-dimethyl- Concentration Rank 21

| R.T.      | EstConc                    | Area     | Relative to ISTD | R.T.        |      |
|-----------|----------------------------|----------|------------------|-------------|------|
| 13.369    | 30.03 ng/ul                | 10825900 | Acenaphthene-d10 | 14.051      |      |
| Hit# of 5 | Tentative ID               | MW       | MolForm          | CAS#        | Qual |
| 1         | Naphthalene, 2,7-dimethyl- | 156      | C12H12           | 000582-16-1 | 97   |
| 2         | Naphthalene, 2,3-dimethyl- | 156      | C12H12           | 000581-40-8 | 97   |
| 3         | Naphthalene, 2,6-dimethyl- | 156      | C12H12           | 000581-42-0 | 97   |
| 4         | Naphthalene, 1,4-dimethyl- | 156      | C12H12           | 000571-58-4 | 97   |
| 5         | Naphthalene, 1,7-dimethyl- | 156      | C12H12           | 000575-37-1 | 97   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

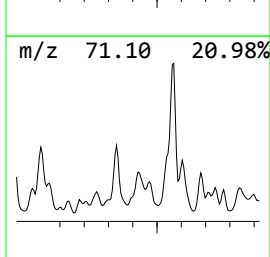
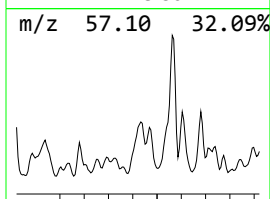
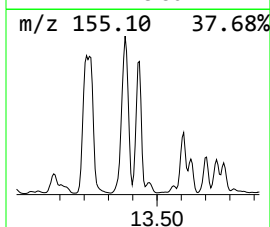
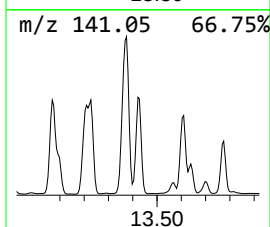
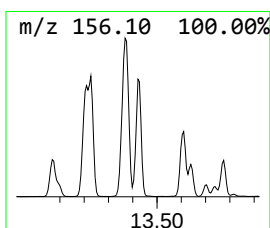
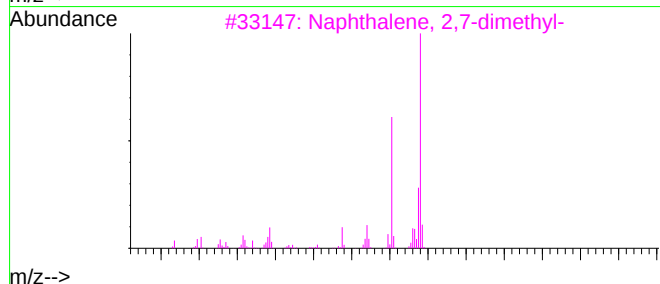
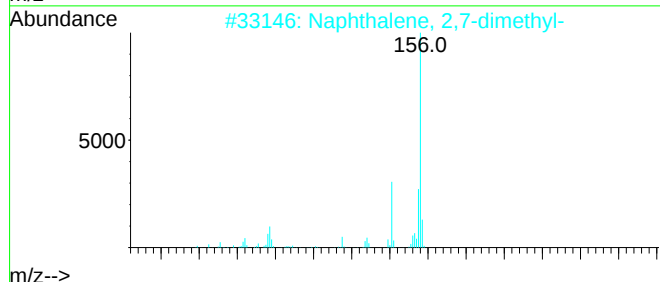
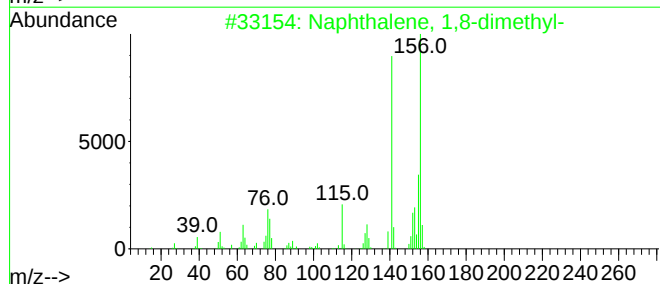
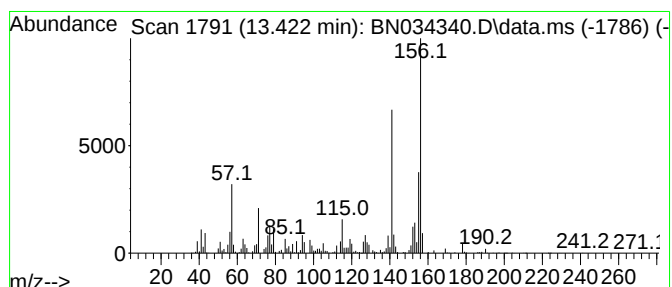
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

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Peak Number 54 Naphthalene, 1,8-dimethyl- Concentration Rank 15

| R.T.      | EstConc                    | Area     | Relative to ISTD | R.T.        |      |
|-----------|----------------------------|----------|------------------|-------------|------|
| 13.422    | 36.85 ng/ul                | 13286200 | Acenaphthene-d10 | 14.051      |      |
| Hit# of 5 | Tentative ID               | MW       | MolForm          | CAS#        | Qual |
| 1         | Naphthalene, 1,8-dimethyl- | 156      | C12H12           | 000569-41-5 | 93   |
| 2         | Naphthalene, 2,7-dimethyl- | 156      | C12H12           | 000582-16-1 | 90   |
| 3         | Naphthalene, 2,7-dimethyl- | 156      | C12H12           | 000582-16-1 | 90   |
| 4         | Naphthalene, 1,6-dimethyl- | 156      | C12H12           | 000575-43-9 | 89   |
| 5         | Naphthalene, 1,4-dimethyl- | 156      | C12H12           | 000571-58-4 | 89   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

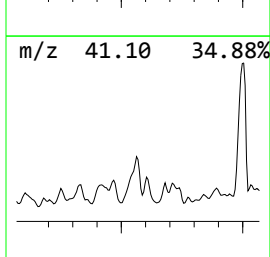
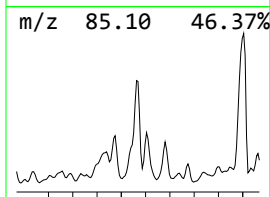
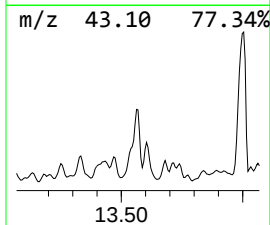
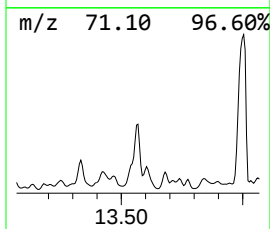
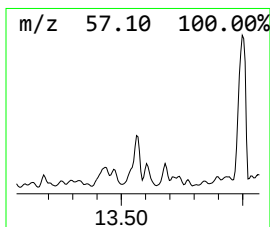
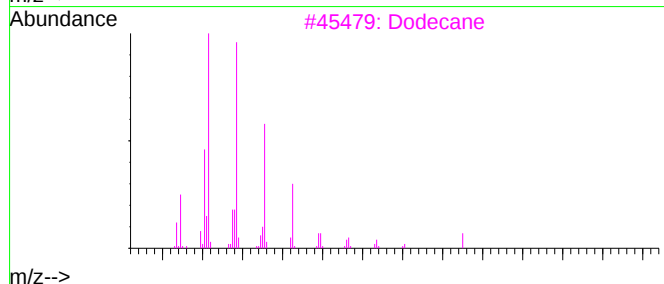
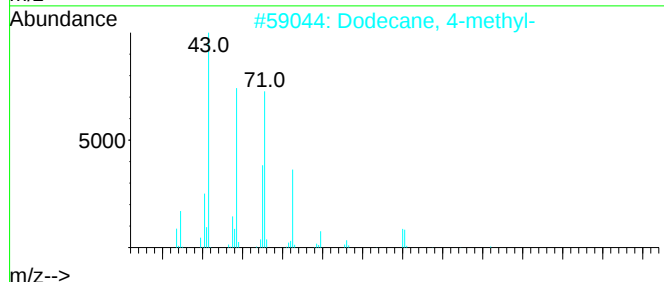
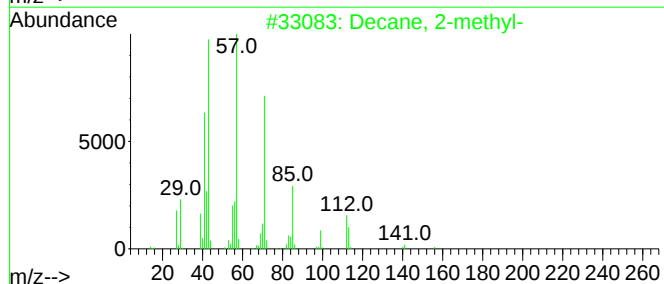
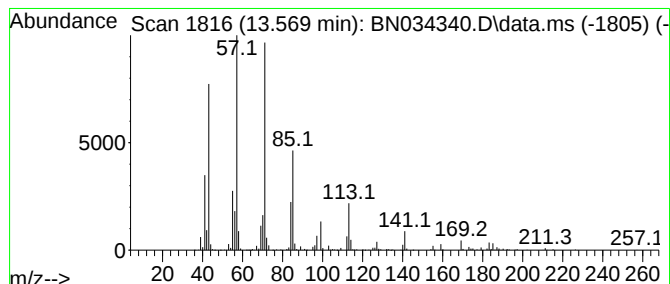
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 55 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.      | EstConc                            | Area     | Relative to ISTD | R.T.         |      |
|-----------|------------------------------------|----------|------------------|--------------|------|
| 13.569    | 45.88 ng/ul                        | 16540700 | Acenaphthene-d10 | 14.051       |      |
| Hit# of 5 | Tentative ID                       | MW       | MolForm          | CAS#         | Qual |
| 1         | Decane, 2-methyl-                  | 156      | C11H24           | 006975-98-0  | 72   |
| 2         | Dodecane, 4-methyl-                | 184      | C13H28           | 006117-97-1  | 68   |
| 3         | Dodecane                           | 170      | C12H26           | 000112-40-3  | 68   |
| 4         | Decane, 5-propyl-                  | 184      | C13H28           | 017312-62-8  | 68   |
| 5         | Carbonic acid, eicosyl vinyl ester | 368      | C23H44O3         | 1000382-54-3 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

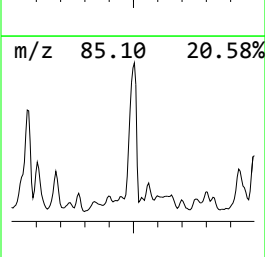
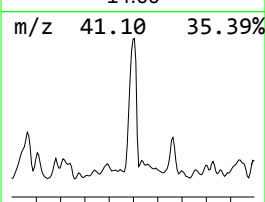
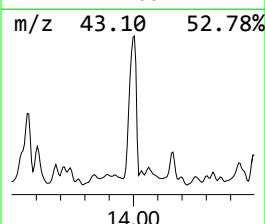
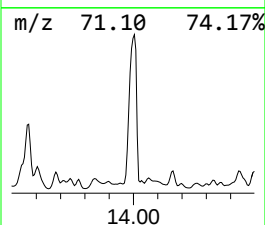
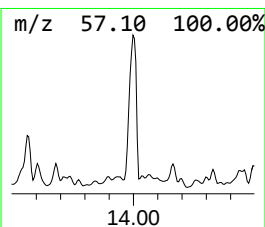
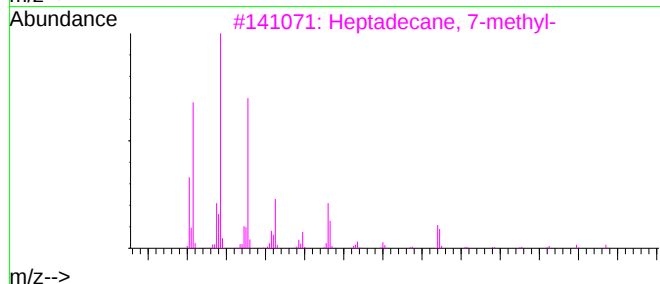
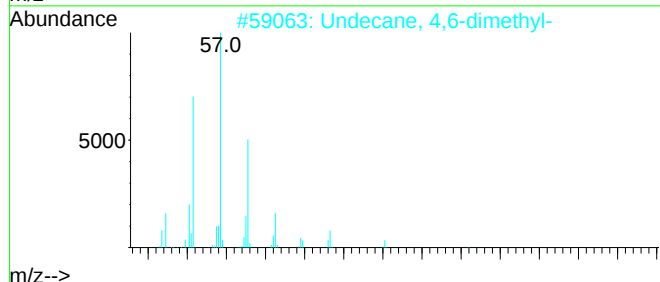
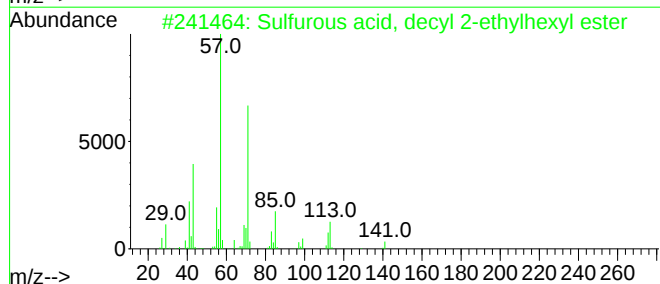
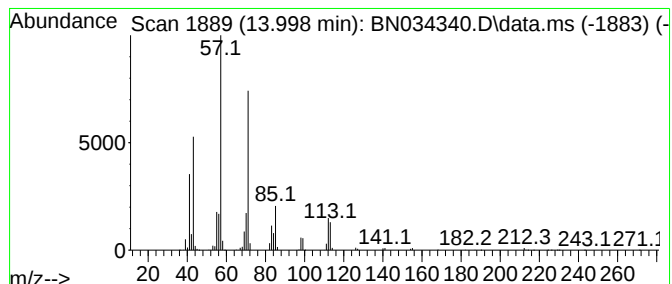
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 56 Sulfurous acid, decyl 2-eth... Concentration Rank 1

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 13.998 | 77.40 ng/ul | 27908000 | Acenaphthene-d10 | 14.051 |

| Hit# of 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|-----------|-------------------------------------|-----|-----------|--------------|------|
| 1         | Sulfurous acid, decyl 2-ethylhex... | 334 | C18H38O3S | 1000309-19-3 | 90   |
| 2         | Undecane, 4,6-dimethyl-             | 184 | C13H28    | 017312-82-2  | 86   |
| 3         | Heptadecane, 7-methyl-              | 254 | C18H38    | 020959-33-5  | 86   |
| 4         | Dodecane, 2,7,10-trimethyl-         | 212 | C15H32    | 074645-98-0  | 80   |
| 5         | Dodecane, 2,6,11-trimethyl-         | 212 | C15H32    | 031295-56-4  | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

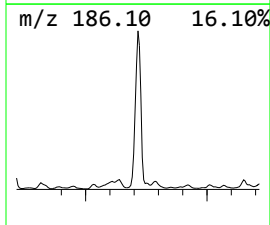
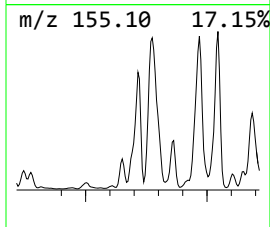
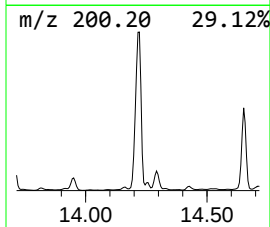
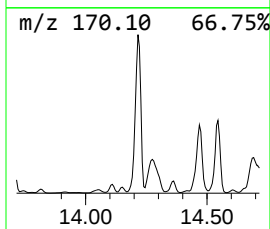
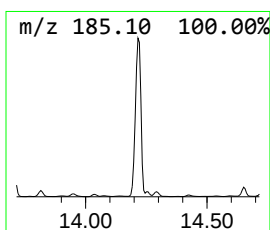
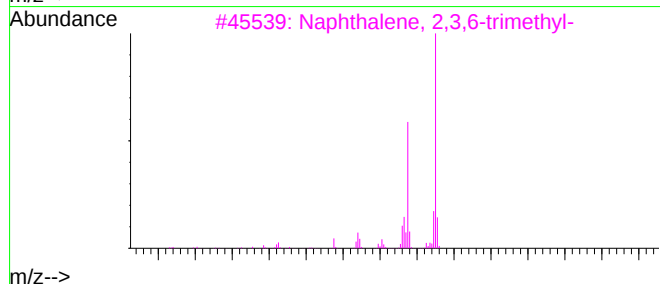
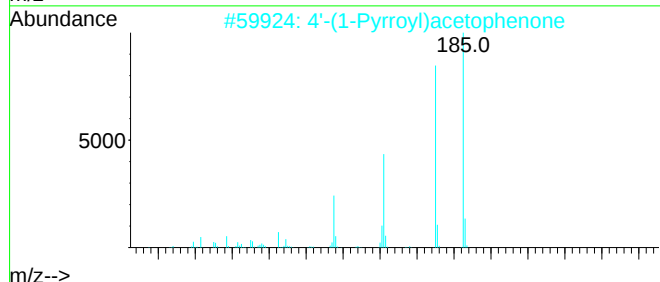
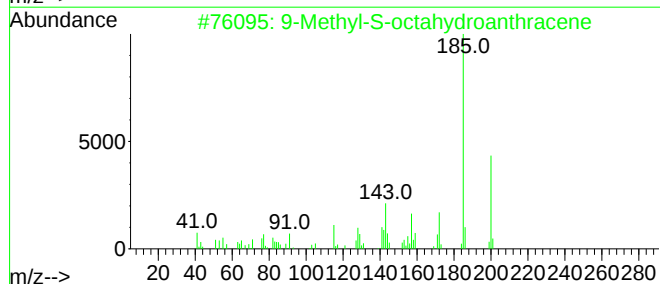
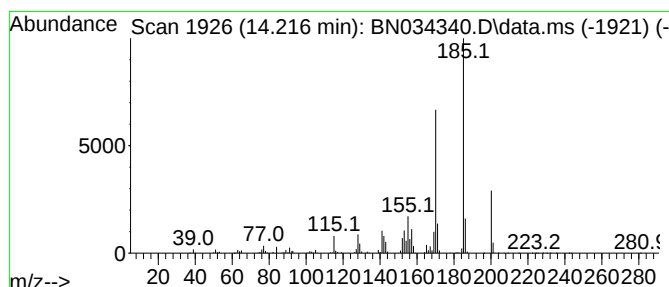
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Peak Number 57 9-Methyl-S-octahydroanthracene Concentration Rank 6

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 14.216 | 48.95 ng/ul | 17650800 | Acenaphthene-d10 | 14.051 |

| Hit# of 5 | Tentative ID                       | MW  | MolForm  | CAS#         | Qual |
|-----------|------------------------------------|-----|----------|--------------|------|
| 1         | 9-Methyl-S-octahydroanthracene     | 200 | C15H20   | 004948-51-0  | 60   |
| 2         | 4'-(1-Pyrrolyl)acetophenone        | 185 | C12H11NO | 022106-37-2  | 52   |
| 3         | Naphthalene, 2,3,6-trimethyl-      | 170 | C13H14   | 000829-26-5  | 51   |
| 4         | 6-tert-Butylquinoline              | 185 | C13H15N  | 068141-13-9  | 50   |
| 5         | 1H-Pyrazole, 5-(t-butyl)-3-phenyl- | 200 | C13H16N2 | 1000261-34-8 | 46   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

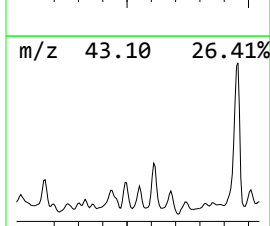
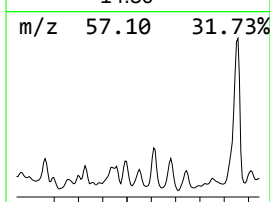
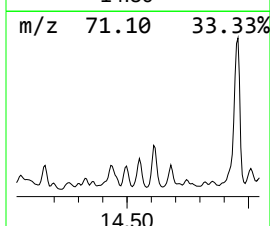
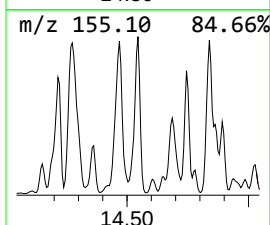
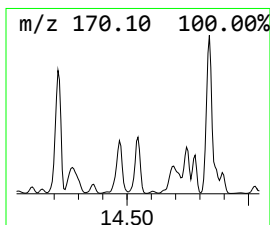
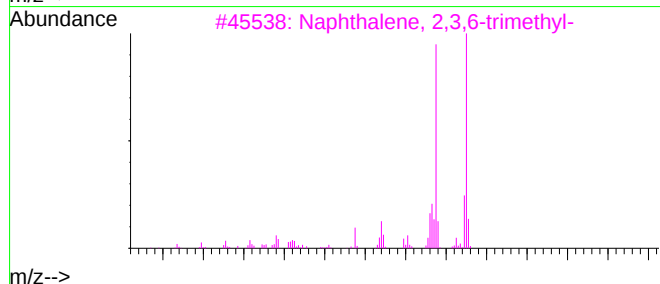
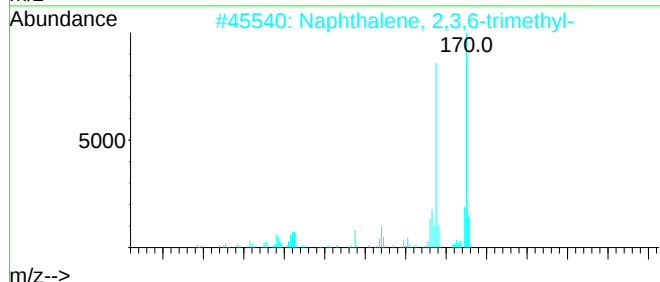
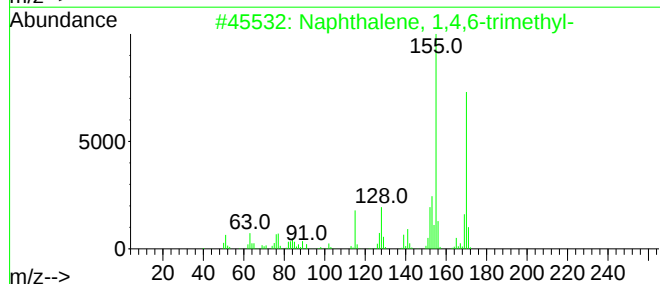
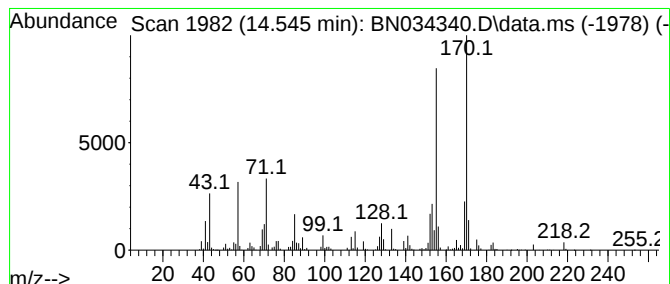
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Peak Number 58 Naphthalene, 1,4,6-trimethyl- Concentration Rank 31

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 14.545 | 20.77 ng/ul | 7487760 | Acenaphthene-d10 | 14.051 |

| Hit# of 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|-----------|-------------------------------|-----|---------|-------------|------|
| 1         | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 97   |
| 2         | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 96   |
| 3         | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 96   |
| 4         | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 96   |
| 5         | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

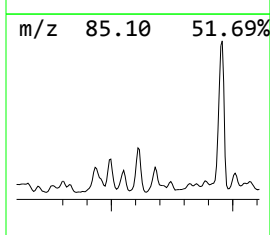
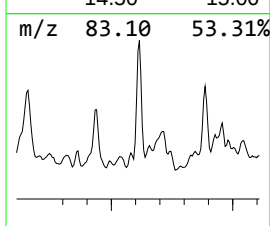
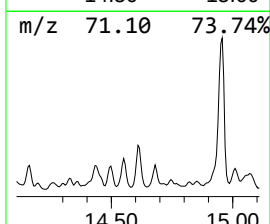
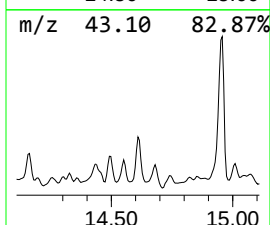
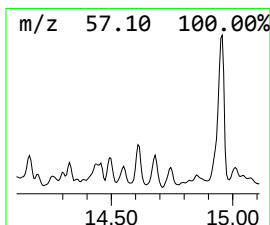
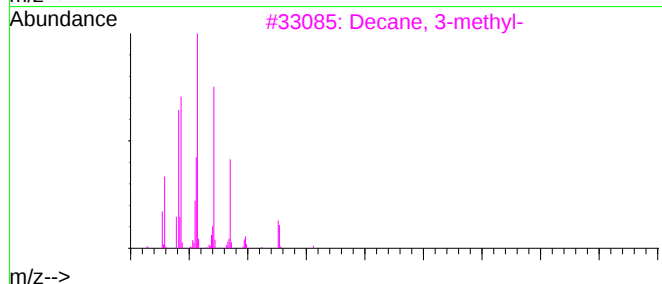
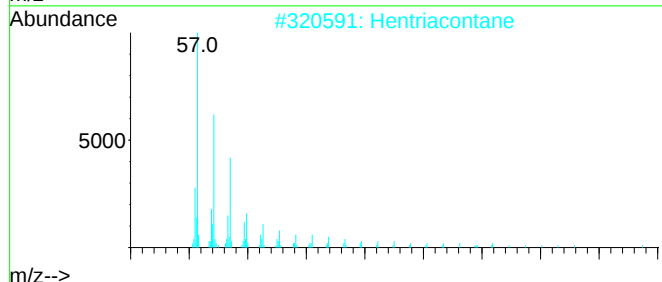
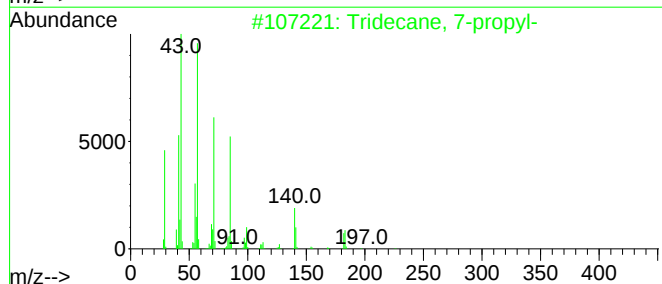
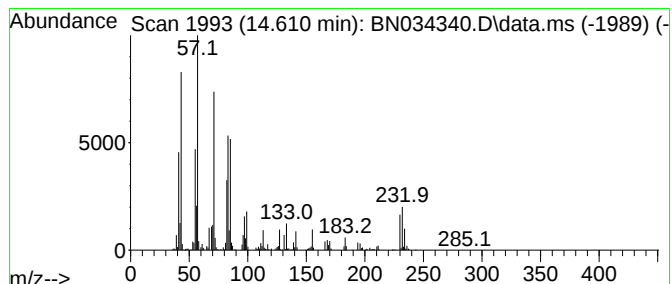
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 27

| R.T.      | EstConc              | Area    | Relative to ISTD | R.T.        |      |
|-----------|----------------------|---------|------------------|-------------|------|
| 14.610    | 23.55 ng/ul          | 8492740 | Acenaphthene-d10 | 14.051      |      |
| Hit# of 5 | Tentative ID         | MW      | MolForm          | CAS#        | Qual |
| 1         | Tridecane, 7-propyl- | 226     | C16H34           | 055045-09-5 | 64   |
| 2         | Hentriacontane       | 437     | C31H64           | 000630-04-6 | 53   |
| 3         | Decane, 3-methyl-    | 156     | C11H24           | 013151-34-3 | 50   |
| 4         | Heneicosane          | 296     | C21H44           | 000629-94-7 | 49   |
| 5         | Octacosane           | 394     | C28H58           | 000630-02-4 | 49   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

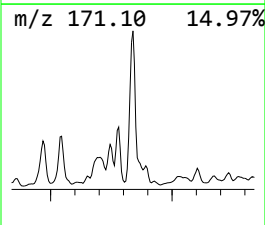
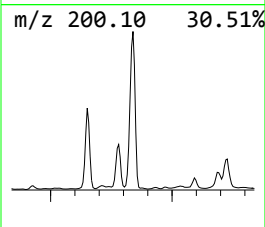
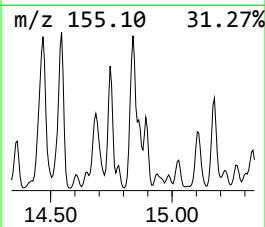
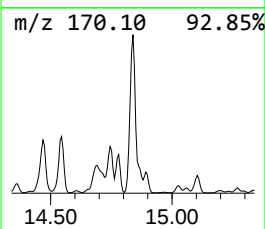
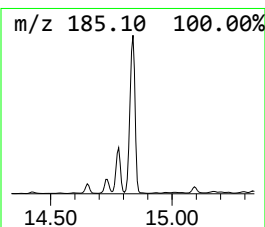
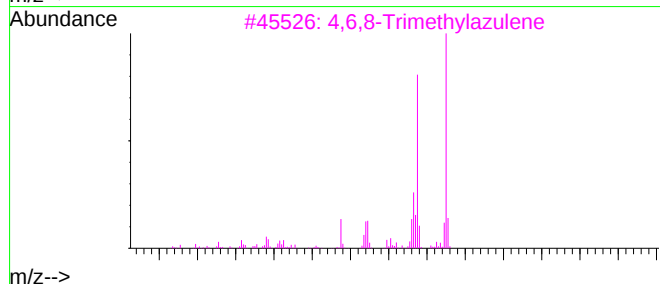
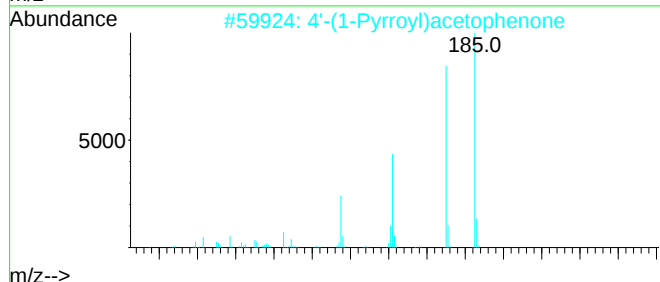
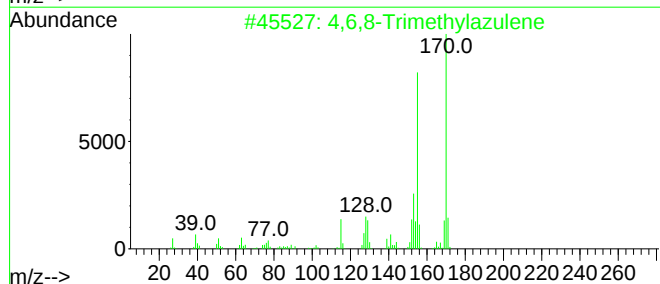
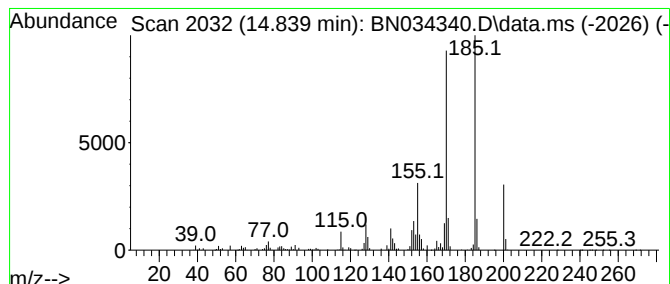
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TIC Integration Parameters: LSCINT.P

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Peak Number 60 4,6,8-Trimethylazulene Concentration Rank 2

| R.T.      | EstConc                       | Area     | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------|----------|------------------|-------------|------|
| 14.839    | 71.16 ng/ul                   | 25657200 | Acenaphthene-d10 | 14.051      |      |
| Hit# of 5 | Tentative ID                  | MW       | MolForm          | CAS#        | Qual |
| 1         | 4,6,8-Trimethylazulene        | 170      | C13H14           | 000941-81-1 | 70   |
| 2         | 4'-(1-Pyrroyl)acetophenone    | 185      | C12H11NO         | 022106-37-2 | 58   |
| 3         | 4,6,8-Trimethylazulene        | 170      | C13H14           | 000941-81-1 | 55   |
| 4         | Naphthalene, 1,4,6-trimethyl- | 170      | C13H14           | 002131-42-2 | 55   |
| 5         | Naphthalene, 2,3,6-trimethyl- | 170      | C13H14           | 000829-26-5 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

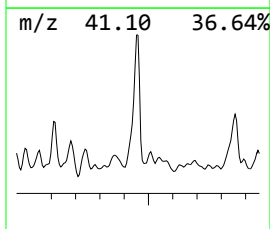
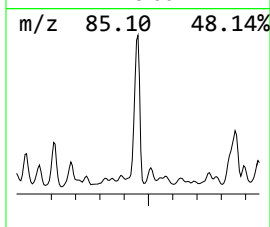
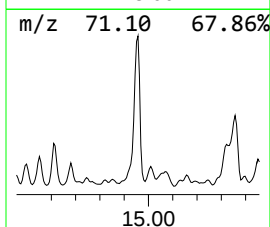
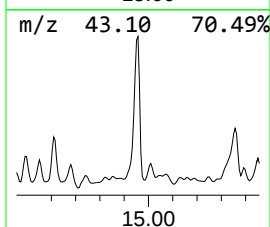
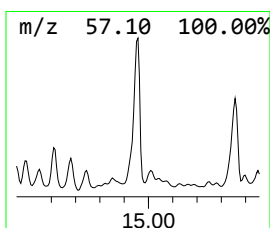
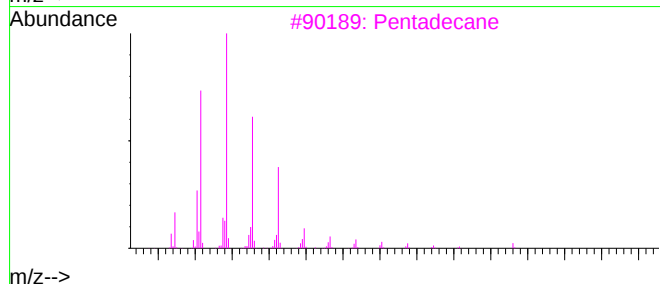
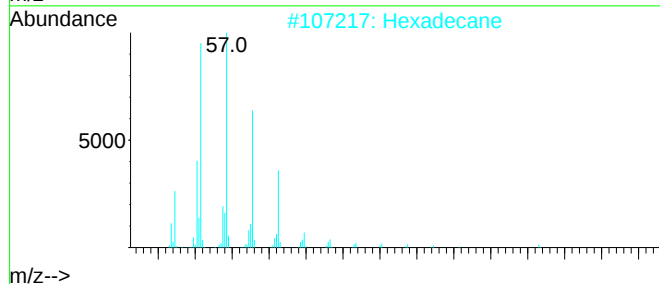
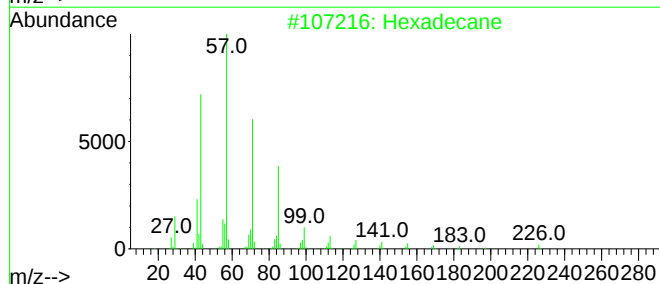
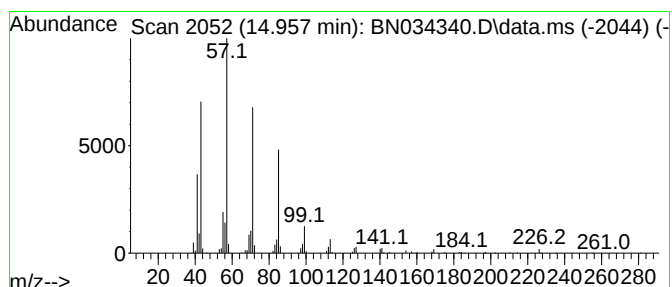
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 61 (DEL) Alkane: Straight-Chai... Concentration Rank 3

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 14.957 | 59.49 ng/ul | 21449000 | Acenaphthene-d10 | 14.051 |

| Hit# of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|-----------|--------------|-----|---------|-------------|------|
| 1         | Hexadecane   | 226 | C16H34  | 000544-76-3 | 93   |
| 2         | Hexadecane   | 226 | C16H34  | 000544-76-3 | 91   |
| 3         | Pentadecane  | 212 | C15H32  | 000629-62-9 | 90   |
| 4         | Tetradecane  | 198 | C14H30  | 000629-59-4 | 90   |
| 5         | Heptadecane  | 240 | C17H36  | 000629-78-7 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

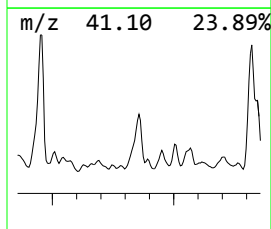
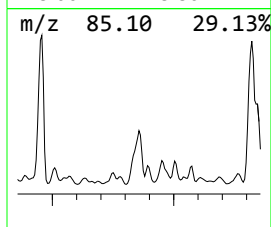
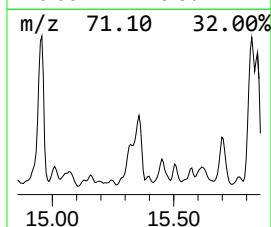
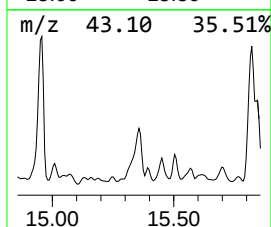
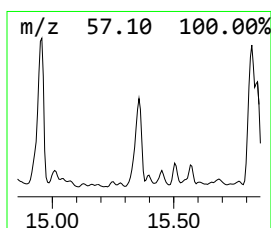
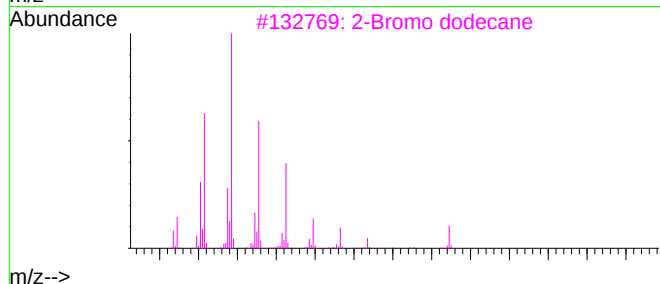
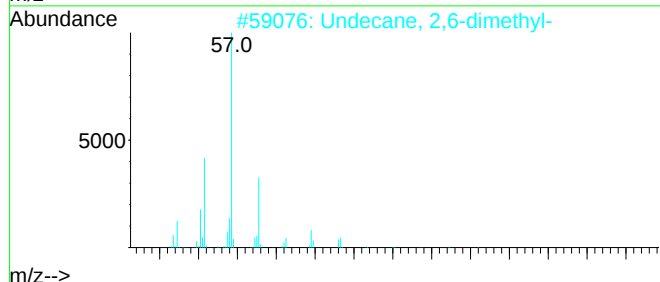
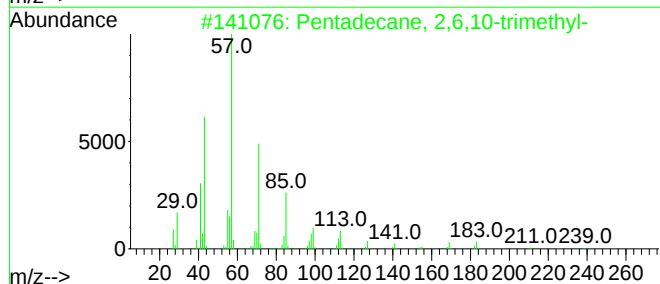
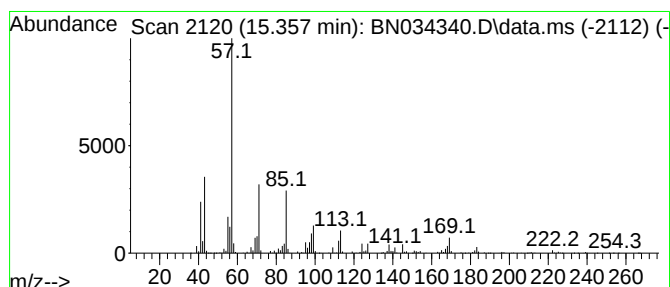
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 62 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 15.357 | 44.42 ng/ul | 16014900 | Acenaphthene-d10 | 14.051 |

| Hit# of 5 | Tentative ID                       | MW  | MolForm  | CAS#        | Qual |
|-----------|------------------------------------|-----|----------|-------------|------|
| 1         | Pentadecane, 2,6,10-trimethyl-     | 254 | C18H38   | 003892-00-0 | 83   |
| 2         | Undecane, 2,6-dimethyl-            | 184 | C13H28   | 017301-23-4 | 64   |
| 3         | 2-Bromo dodecane                   | 248 | C12H25Br | 013187-99-0 | 58   |
| 4         | Undecane, 3,5-dimethyl-            | 184 | C13H28   | 017312-81-1 | 52   |
| 5         | Nonane, 2,2,4,4,6,8,8-heptamethyl- | 226 | C16H34   | 004390-04-9 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

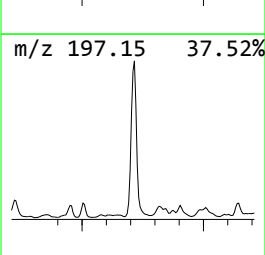
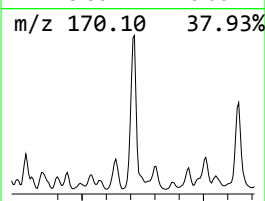
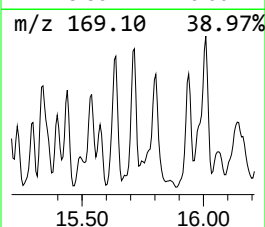
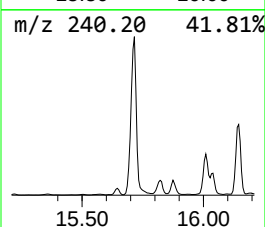
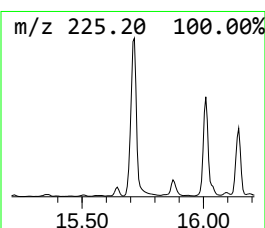
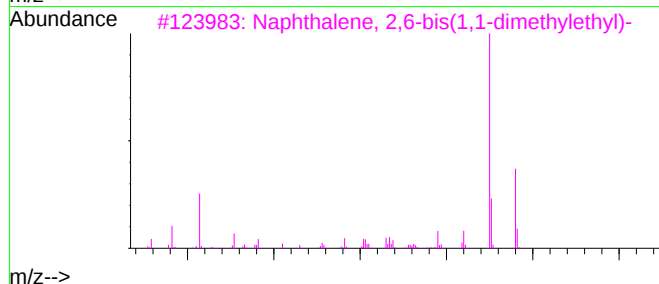
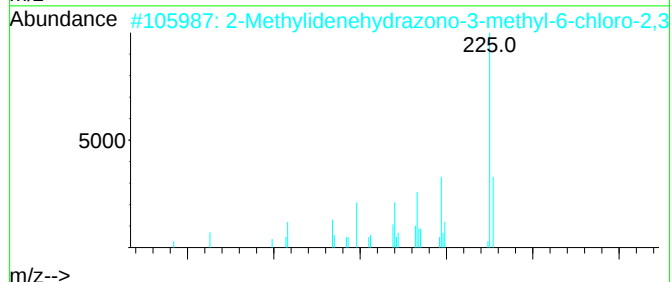
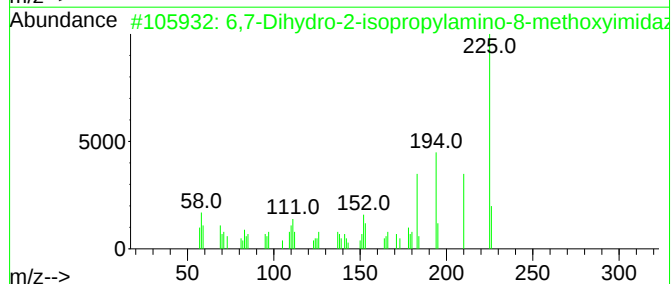
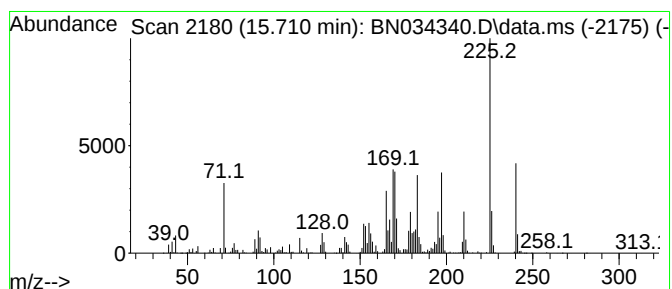
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 63 unknown-04 Concentration Rank 18

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 15.710 | 33.55 ng/ul | 11674400 | Phenanthrene-d10 | 16.833 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|--------------|--------------|------|
| 1    |    |   | 6,7-Dihydro-2-isopropylamino-8-m... | 225 | C9H15N5O2    | 1000101-98-2 | 46   |
| 2    |    |   | 2-Methylidenehydrazono-3-methyl-... | 225 | C9H8ClN3S    | 1000147-91-5 | 43   |
| 3    |    |   | Naphthalene, 2,6-bis(1,1-dimethy... | 240 | C18H24       | 003905-64-4  | 38   |
| 4    |    |   | Indolo[2,3-a]quinolizine, 1,2,3,... | 226 | C15H18N2     | 004802-79-3  | 38   |
| 5    |    |   | 4-Fluoro-1,3-benzothiazol-2-ylam... | 240 | C10H13FN2SSi | 1000478-75-9 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

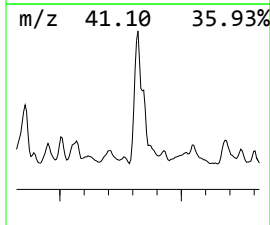
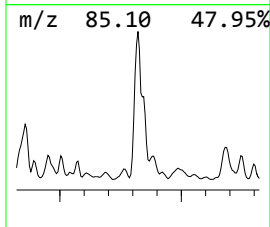
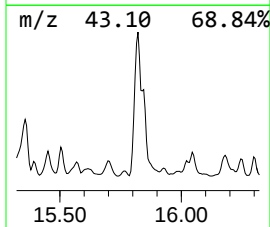
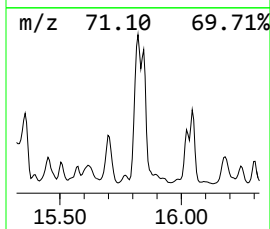
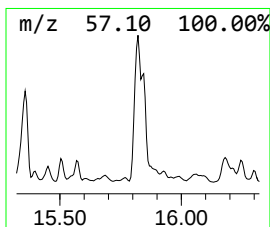
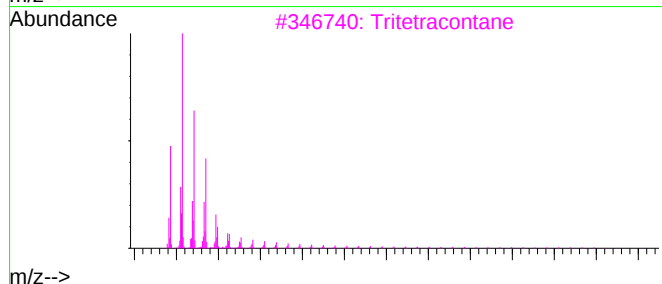
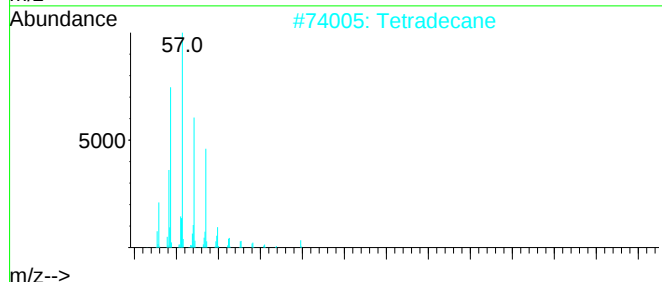
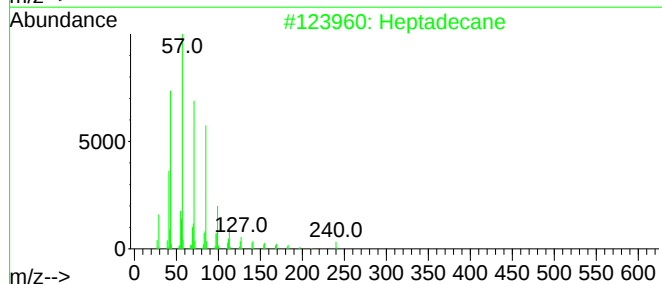
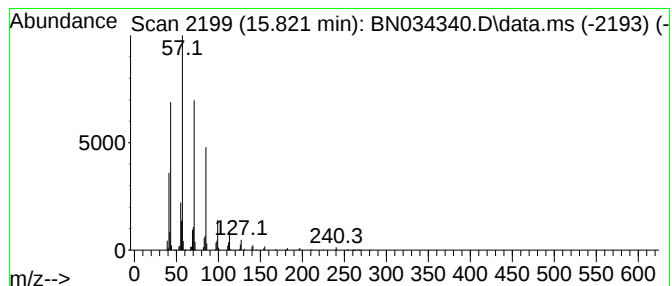
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Peak Number 64 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 15.821 | 58.34 ng/ul | 20300300 | Phenanthrene-d10 | 16.833 |

| Hit# of 5 | Tentative ID    | MW  | MolForm | CAS#        | Qual |
|-----------|-----------------|-----|---------|-------------|------|
| 1         | Heptadecane     | 240 | C17H36  | 000629-78-7 | 91   |
| 2         | Tetradecane     | 198 | C14H30  | 000629-59-4 | 91   |
| 3         | Tritetracontane | 605 | C43H88  | 007098-21-7 | 90   |
| 4         | Hexadecane      | 226 | C16H34  | 000544-76-3 | 90   |
| 5         | Hexadecane      | 226 | C16H34  | 000544-76-3 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

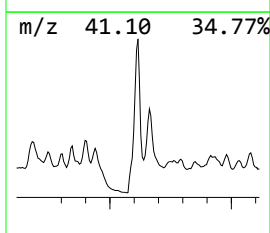
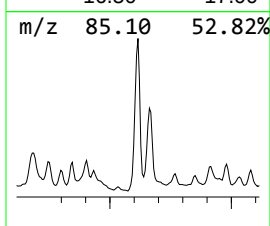
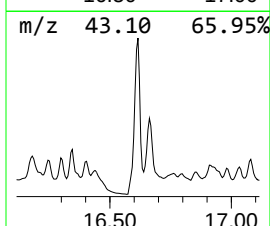
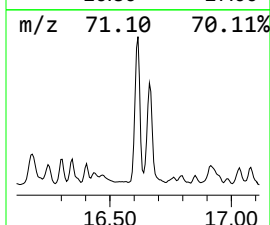
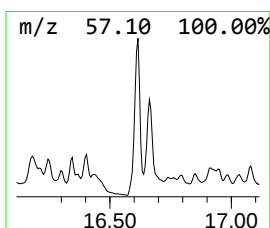
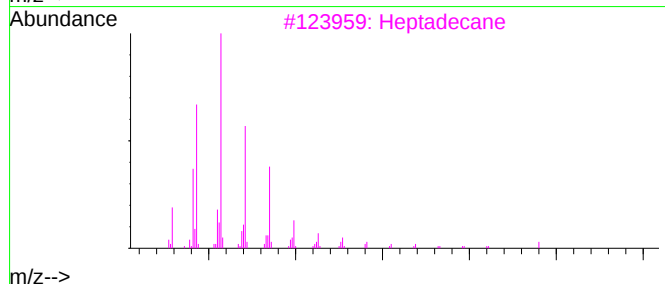
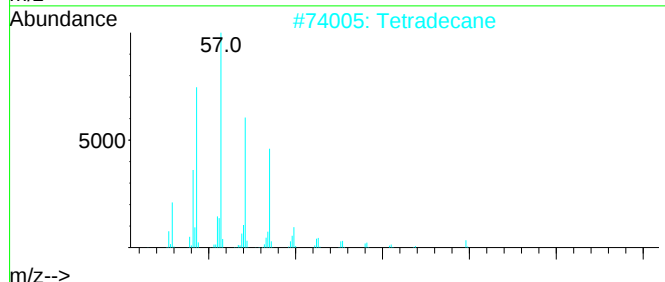
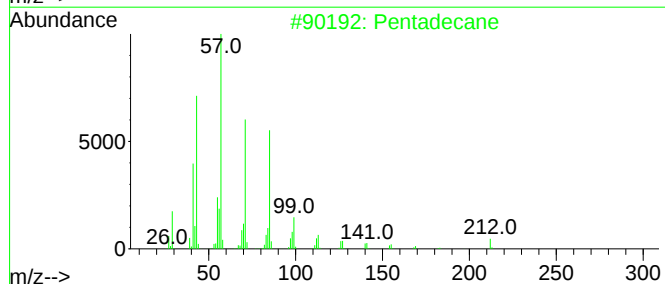
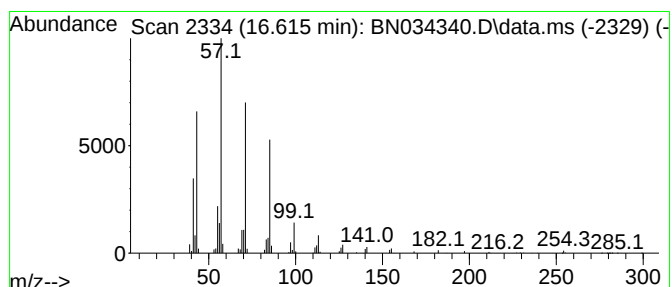
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 65 (DEL) Alkane: Straight-Chai... Concentration Rank 16

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 16.616 | 36.48 ng/ul | 12693300 | Phenanthrene-d10 | 16.833 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentadecane                         | 212 | C15H32  | 000629-62-9 | 90   |
| 2    |    |   | Tetradecane                         | 198 | C14H30  | 000629-59-4 | 87   |
| 3    |    |   | Heptadecane                         | 240 | C17H36  | 000629-78-7 | 86   |
| 4    |    |   | Dodecane, 2-methyl-6-propyl-        | 226 | C16H34  | 055045-08-4 | 78   |
| 5    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

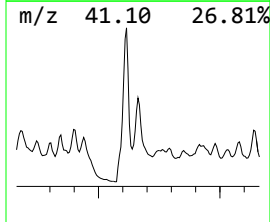
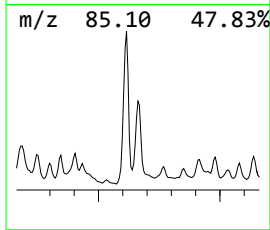
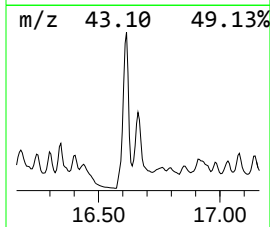
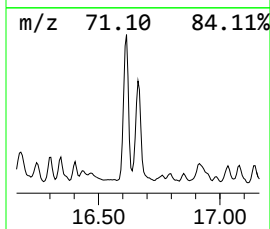
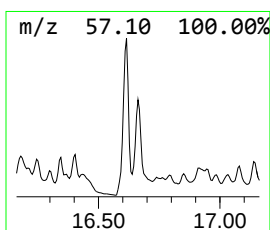
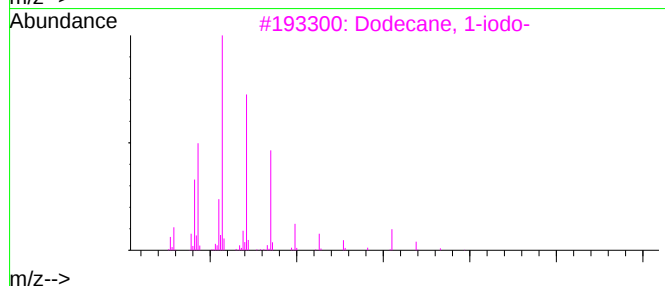
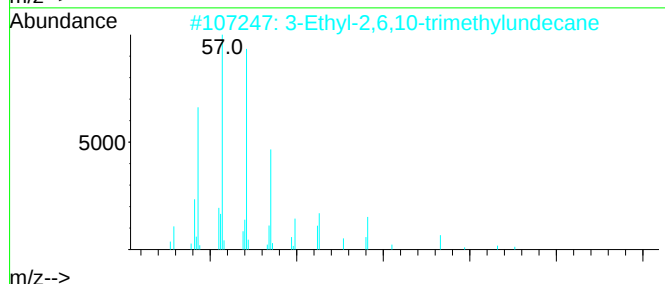
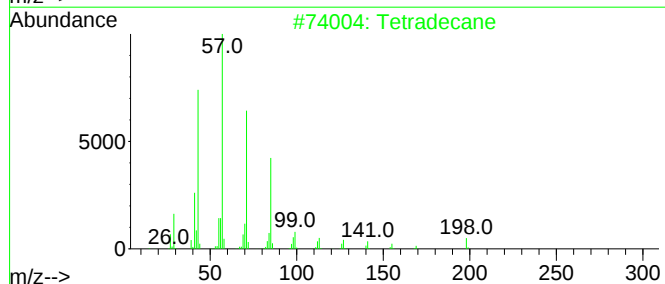
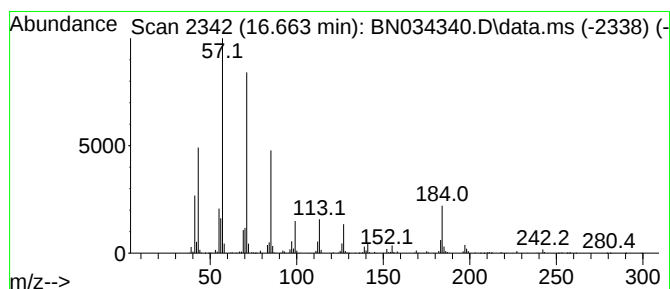
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TIC Integration Parameters: LSCINT.P

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Peak Number 66 (DEL) Alkane: Straight-Chai... Concentration Rank 12

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 16.663 | 39.45 ng/ul | 13727700 | Phenanthrene-d10 | 16.833 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm  | CAS#         | Qual |
|------|----|---|----------------------------------|-----|----------|--------------|------|
| 1    |    |   | Tetradecane                      | 198 | C14H30   | 000629-59-4  | 81   |
| 2    |    |   | 3-Ethyl-2,6,10-trimethylundecane | 226 | C16H34   | 1000432-25-9 | 72   |
| 3    |    |   | Dodecane, 1-iodo-                | 296 | C12H25I  | 004292-19-7  | 72   |
| 4    |    |   | 2-Bromotetradecane               | 276 | C14H29Br | 074036-95-6  | 72   |
| 5    |    |   | Decane, 3-methyl-                | 156 | C11H24   | 013151-34-3  | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

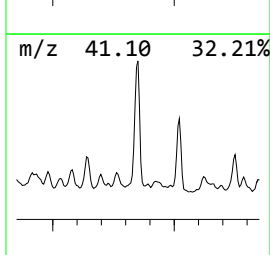
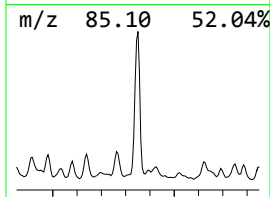
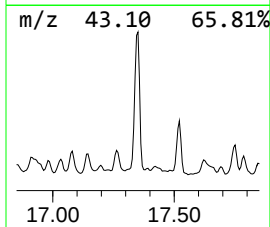
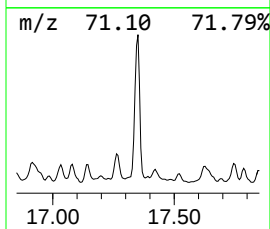
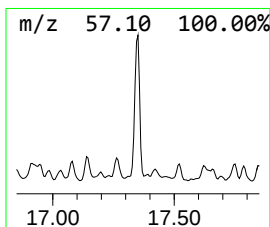
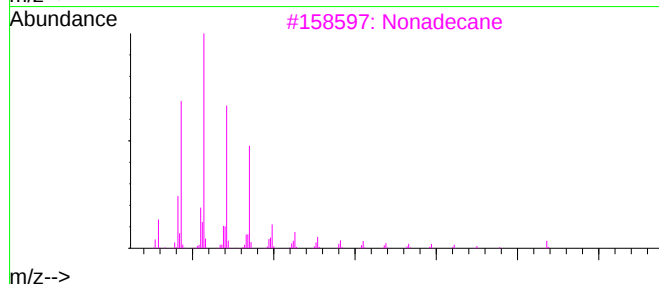
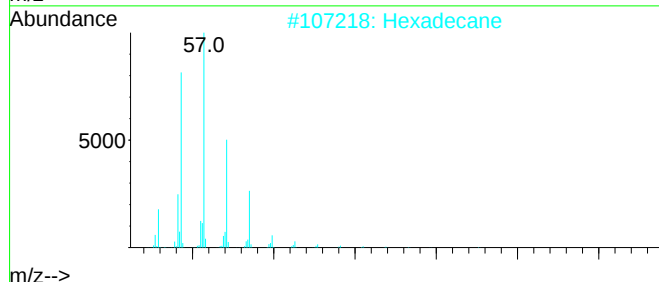
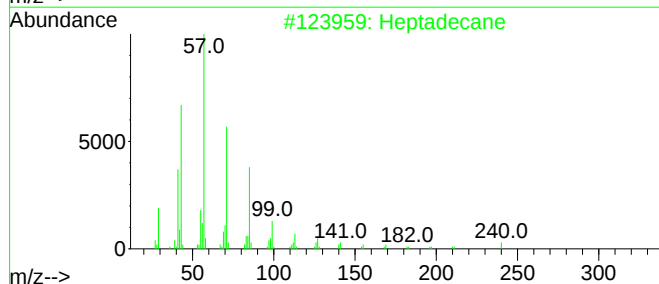
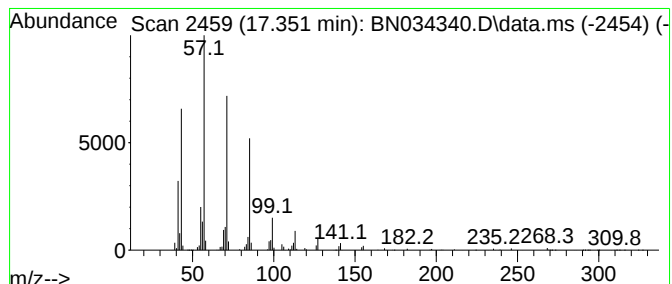
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 67 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 17.351 | 39.57 ng/ul | 13769400 | Phenanthrene-d10 | 16.833 |

| Hit# of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|-----------|--------------|-----|---------|-------------|------|
| 1         | Heptadecane  | 240 | C17H36  | 000629-78-7 | 91   |
| 2         | Hexadecane   | 226 | C16H34  | 000544-76-3 | 90   |
| 3         | Nonadecane   | 268 | C19H40  | 000629-92-5 | 90   |
| 4         | Tetradecane  | 198 | C14H30  | 000629-59-4 | 87   |
| 5         | Hexadecane   | 226 | C16H34  | 000544-76-3 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

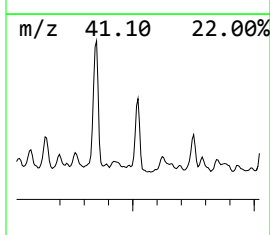
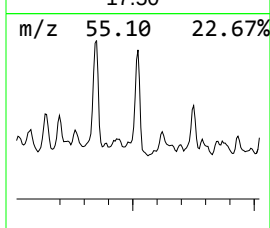
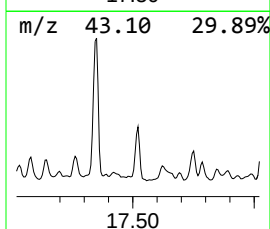
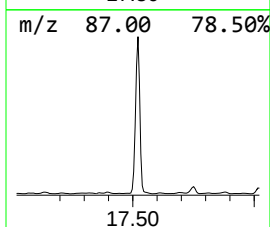
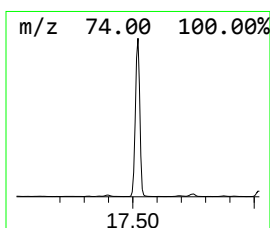
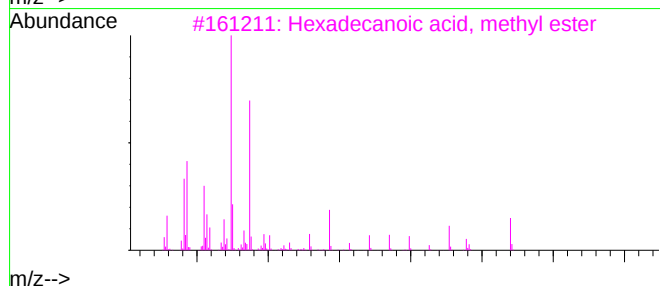
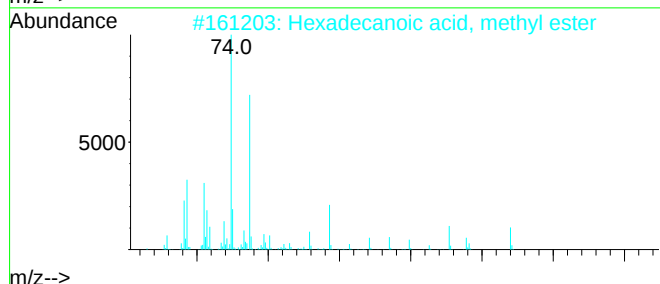
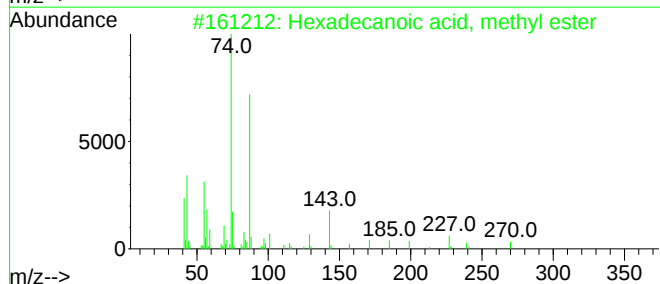
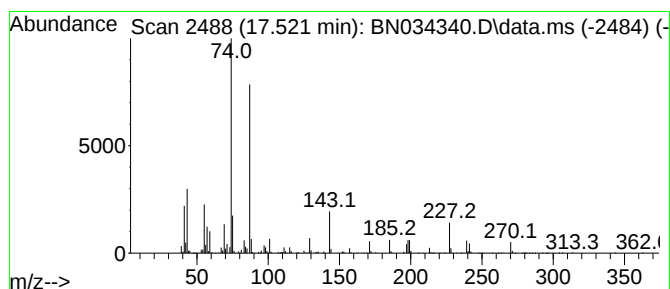
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Peak Number 68 Hexadecanoic acid, methyl e... Concentration Rank 13

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 17.521 | 38.41 ng/ul | 13365300 | Phenanthrene-d10 | 16.833 |

| Hit# of 5 | Tentative ID                    | MW  | MolForm  | CAS#        | Qual |
|-----------|---------------------------------|-----|----------|-------------|------|
| 1         | Hexadecanoic acid, methyl ester | 270 | C17H34O2 | 000112-39-0 | 98   |
| 2         | Hexadecanoic acid, methyl ester | 270 | C17H34O2 | 000112-39-0 | 95   |
| 3         | Hexadecanoic acid, methyl ester | 270 | C17H34O2 | 000112-39-0 | 93   |
| 4         | Tridecanoic acid, methyl ester  | 228 | C14H28O2 | 001731-88-0 | 90   |
| 5         | Hexadecanoic acid, methyl ester | 270 | C17H34O2 | 000112-39-0 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

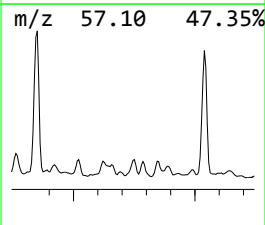
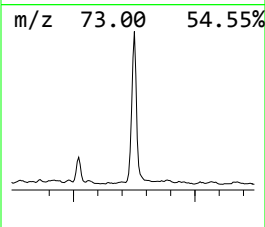
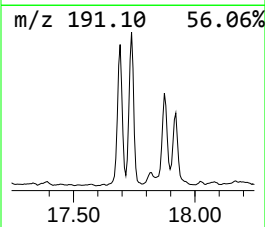
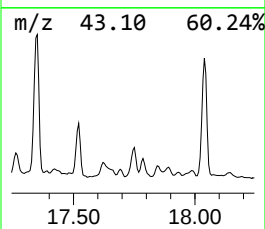
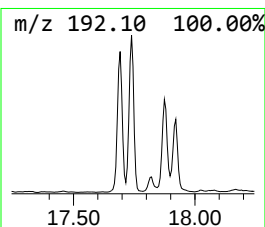
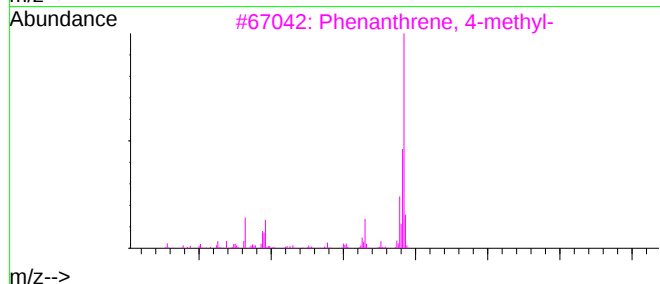
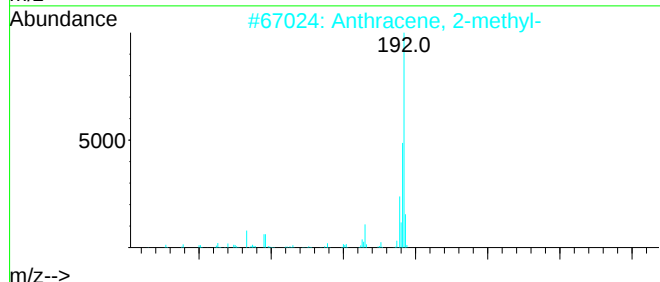
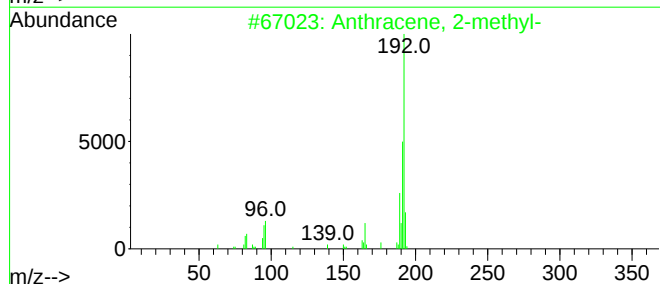
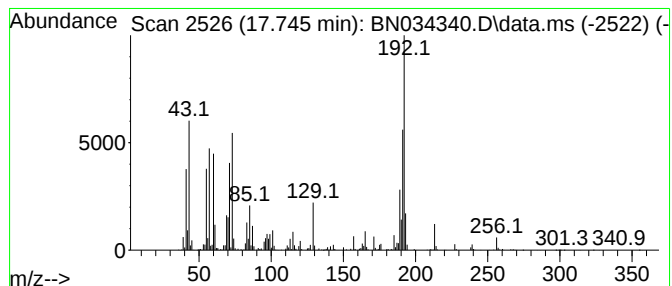
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 69 Anthracene, 2-methyl- Concentration Rank 37

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 17.745    | 17.70 ng/ul                         | 6157810 | Phenanthrene-d10 | 16.833      |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | Anthracene, 2-methyl-               | 192     | C15H12           | 000613-12-7 | 90   |
| 2         | Anthracene, 2-methyl-               | 192     | C15H12           | 000613-12-7 | 84   |
| 3         | Phenanthrene, 4-methyl-             | 192     | C15H12           | 000832-64-4 | 84   |
| 4         | Phenanthrene, 1-methyl-             | 192     | C15H12           | 000832-69-9 | 84   |
| 5         | 1H-Cyclopropa[1]phenanthrene,1a,... | 192     | C15H12           | 000949-41-7 | 84   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

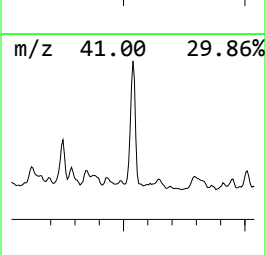
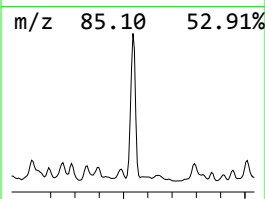
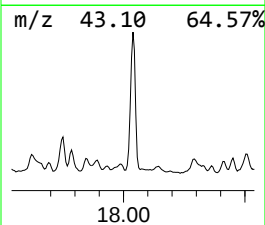
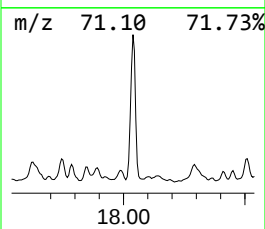
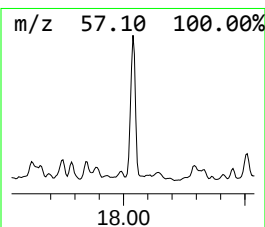
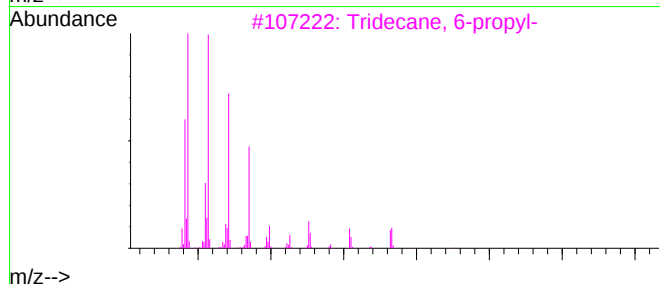
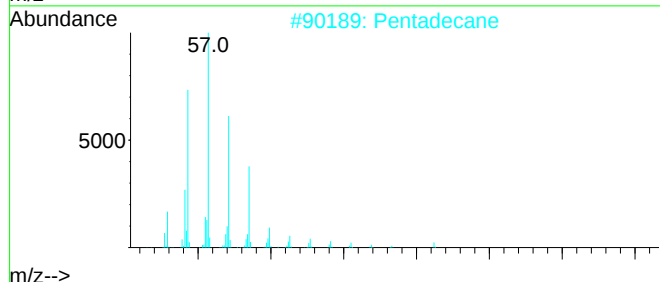
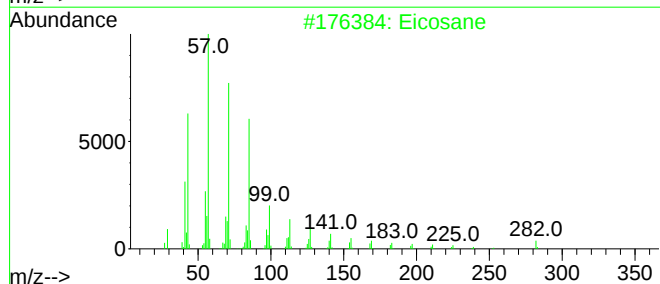
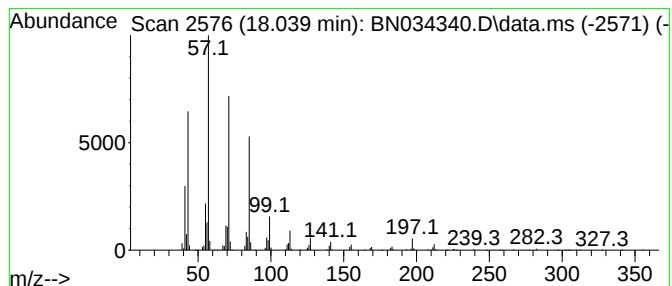
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 70 (DEL) Alkane: Straight-Chai... Concentration Rank 17

| R.T.      | EstConc              | Area     | Relative to ISTD | R.T.        |      |
|-----------|----------------------|----------|------------------|-------------|------|
| 18.039    | 34.69 ng/ul          | 12068600 | Phenanthrene-d10 | 16.833      |      |
| Hit# of 5 | Tentative ID         | MW       | MolForm          | CAS#        | Qual |
| 1         | Eicosane             | 282      | C20H42           | 000112-95-8 | 97   |
| 2         | Pentadecane          | 212      | C15H32           | 000629-62-9 | 96   |
| 3         | Tridecane, 6-propyl- | 226      | C16H34           | 055045-10-8 | 91   |
| 4         | Heneicosane          | 296      | C21H44           | 000629-94-7 | 91   |
| 5         | Nonadecane           | 268      | C19H40           | 000629-92-5 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L

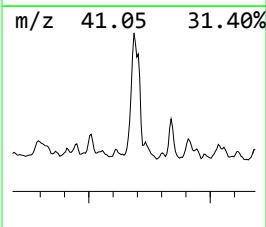
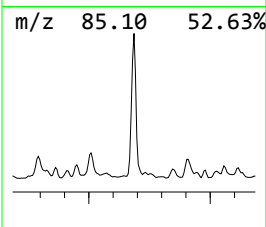
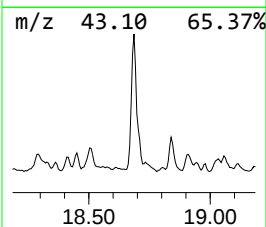
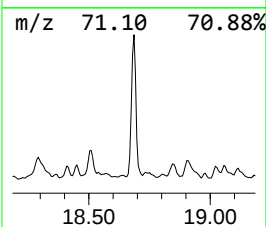
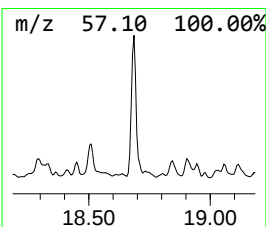
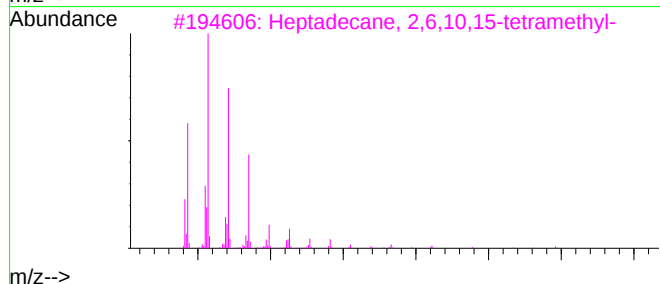
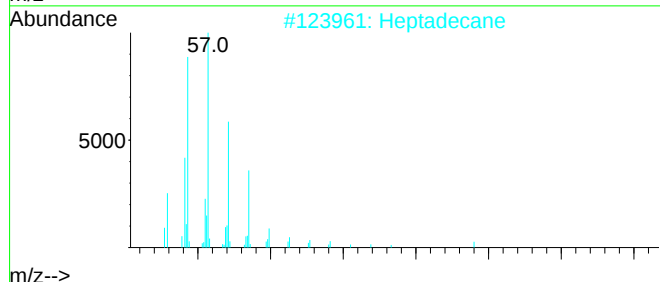
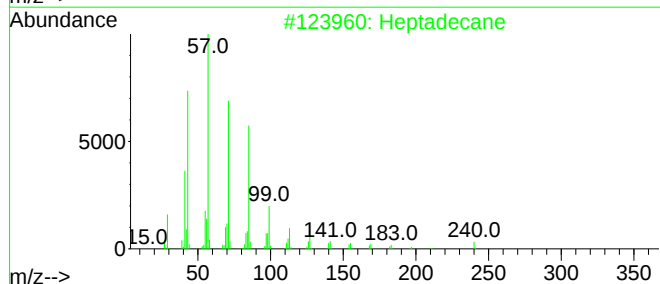
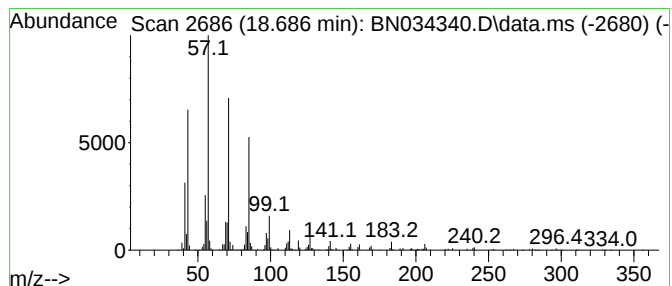
TIC Integration Parameters: LSCINT.P

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Peak Number 71 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 18.686 | 37.78 ng/ul | 13145400 | Phenanthrene-d10 | 16.833 |

| Hit# of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|-----------|-------------------------------------|-----|---------|-------------|------|
| 1         | Heptadecane                         | 240 | C17H36  | 000629-78-7 | 97   |
| 2         | Heptadecane                         | 240 | C17H36  | 000629-78-7 | 96   |
| 3         | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 94   |
| 4         | Tetracosane                         | 338 | C24H50  | 000646-31-1 | 91   |
| 5         | Octadecane                          | 254 | C18H38  | 000593-45-3 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

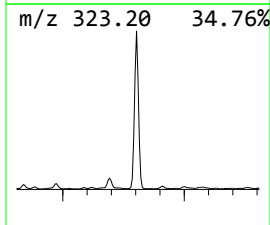
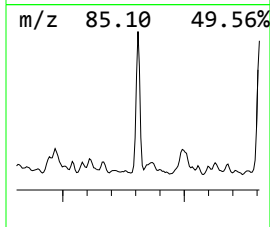
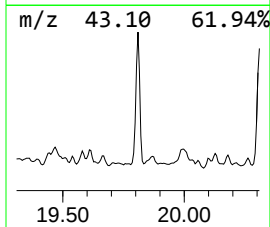
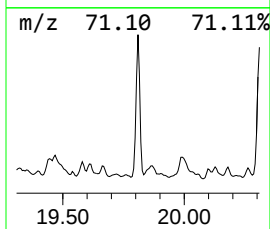
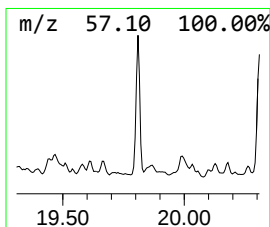
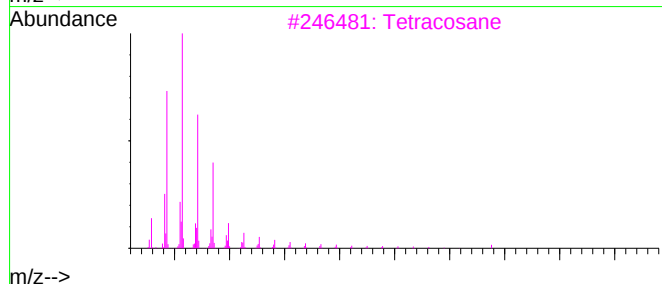
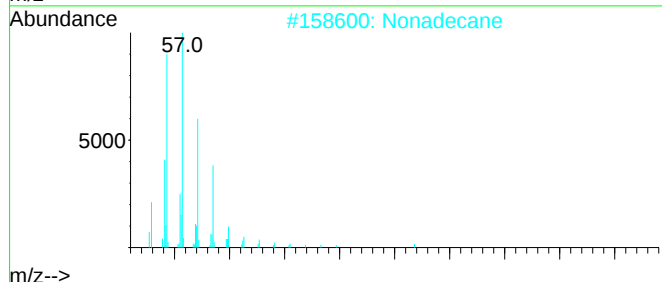
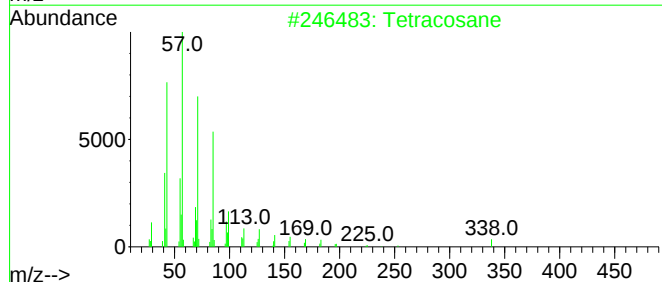
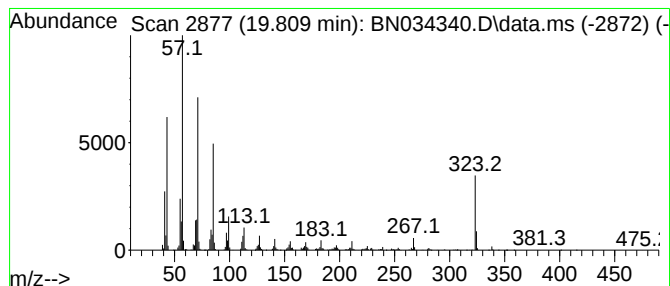
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Peak Number 72 (DEL) Alkane: Straight-Chai... Concentration Rank 25

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 19.809 | 26.59 ng/ul | 7894570 | Chrysene-d12     | 21.033 |

| Hit# of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|-----------|--------------|-----|---------|-------------|------|
| 1         | Tetracosane  | 338 | C24H50  | 000646-31-1 | 98   |
| 2         | Nonadecane   | 268 | C19H40  | 000629-92-5 | 96   |
| 3         | Tetracosane  | 338 | C24H50  | 000646-31-1 | 96   |
| 4         | Tetracosane  | 338 | C24H50  | 000646-31-1 | 93   |
| 5         | Tricosane    | 324 | C23H48  | 000638-67-5 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

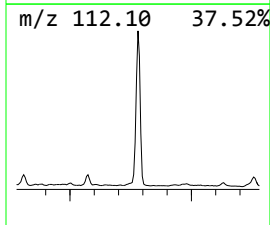
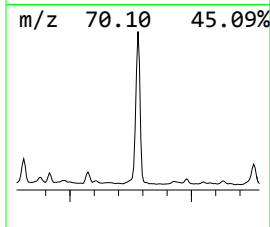
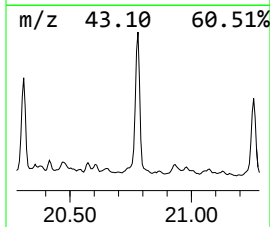
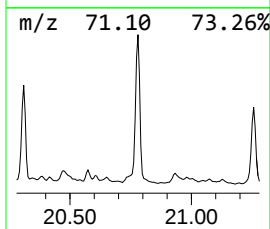
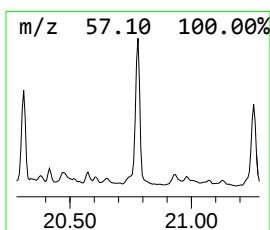
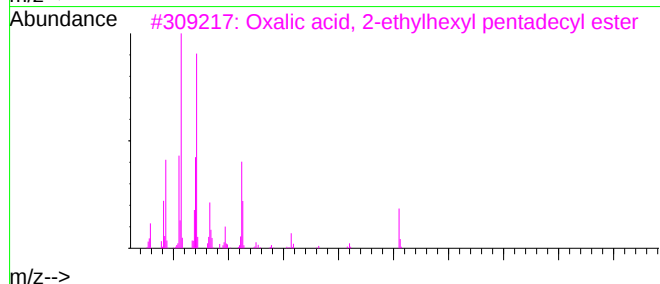
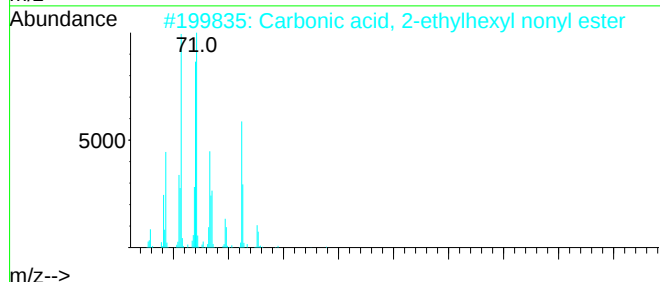
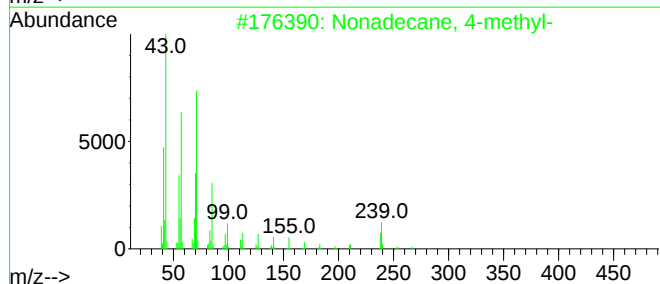
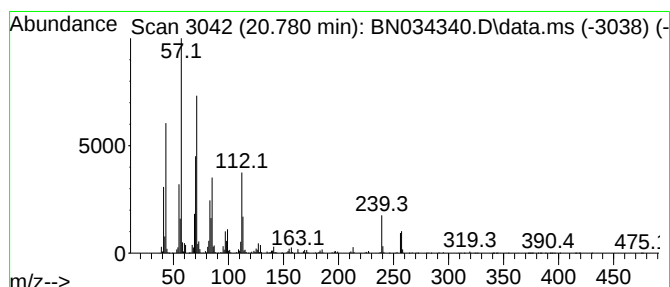
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

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Peak Number 73 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.      | EstConc                               | Area     | Relative to ISTD | R.T.         |      |
|-----------|---------------------------------------|----------|------------------|--------------|------|
| 20.780    | 43.67 ng/ul                           | 12965800 | Chrysene-d12     | 21.033       |      |
| Hit# of 5 | Tentative ID                          | MW       | MolForm          | CAS#         | Qual |
| 1         | Nonadecane, 4-methyl-                 | 282      | C20H42           | 025117-27-5  | 81   |
| 2         | Carbonic acid, 2-ethylhexyl nonyl...  | 300      | C18H36O3         | 1000383-13-6 | 68   |
| 3         | Oxalic acid, 2-ethylhexyl pentadec... | 412      | C25H48O4         | 1000309-39-8 | 64   |
| 4         | Nonane, 4-methyl-5-propyl-            | 184      | C13H28           | 062185-55-1  | 62   |
| 5         | Carbonic acid, 2-ethylhexyl octyl...  | 286      | C17H34O3         | 1000383-13-5 | 62   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

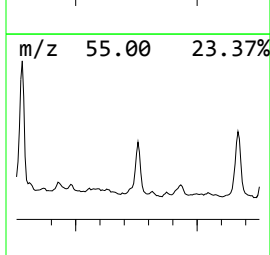
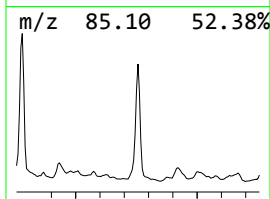
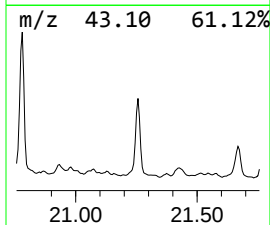
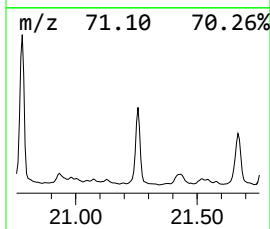
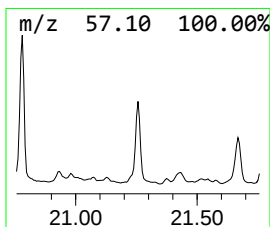
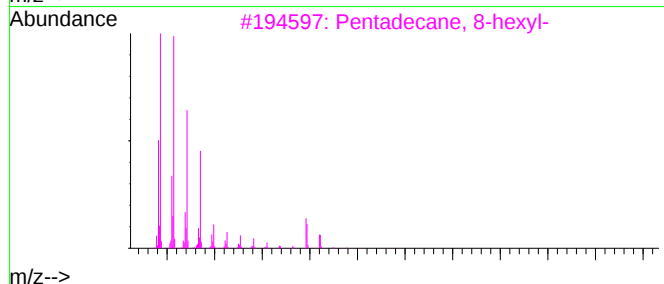
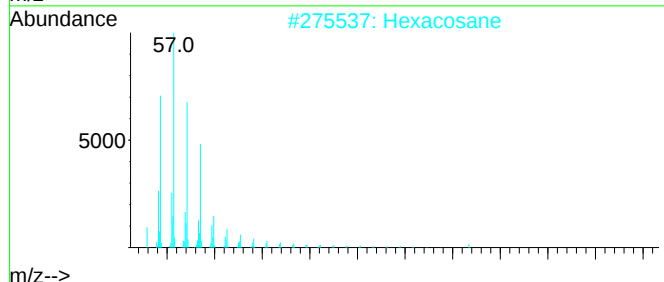
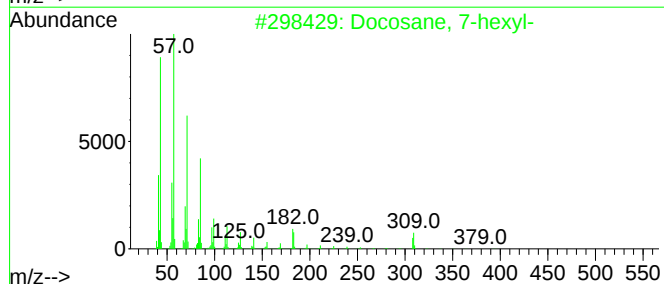
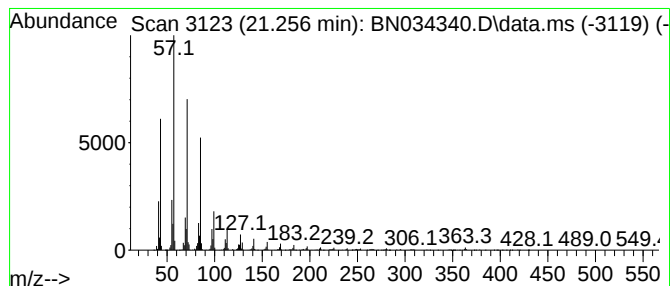
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Peak Number 74 (DEL) Alkane: Straight-Chai... Concentration Rank 32

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 21.256 | 20.00 ng/ul | 5938710 | Chrysene-d12     | 21.033 |

| Hit# of 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|-----------|-----------------------|-----|---------|-------------|------|
| 1         | Docosane, 7-hexyl-    | 394 | C28H58  | 055373-86-9 | 93   |
| 2         | Hexacosane            | 366 | C26H54  | 000630-01-3 | 93   |
| 3         | Pentadecane, 8-hexyl- | 296 | C21H44  | 013475-75-7 | 93   |
| 4         | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 91   |
| 5         | Heneicosane           | 296 | C21H44  | 000629-94-7 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

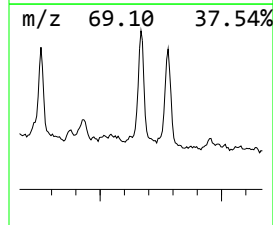
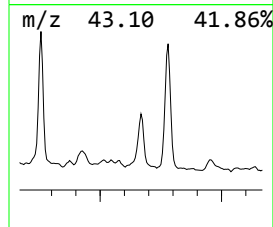
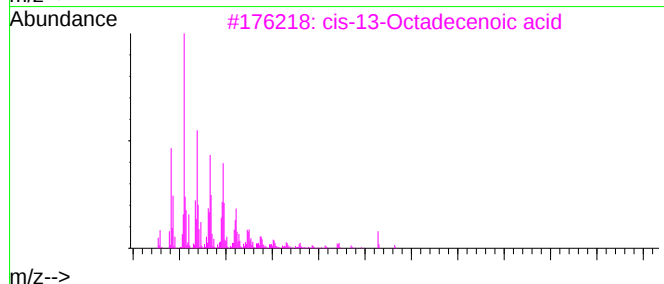
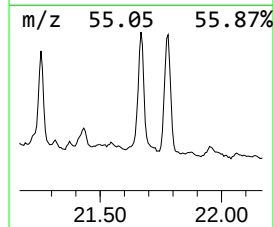
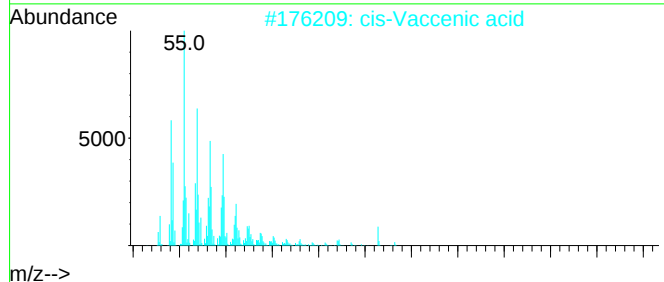
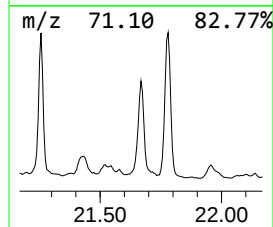
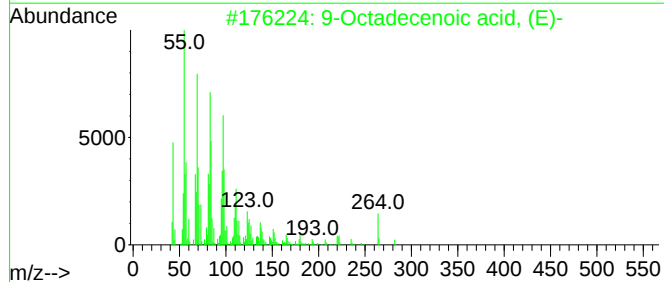
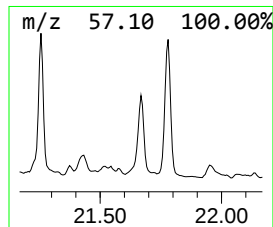
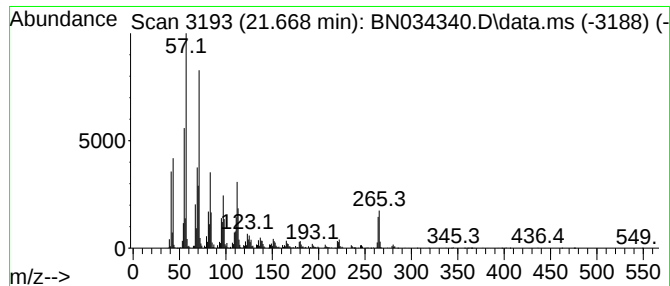
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 75 9-Octadecenoic acid, (E)- Concentration Rank 29

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 21.668 | 22.84 ng/ul | 6781180 | Chrysene-d12     | 21.033 |

| Hit# of 5 | Tentative ID               | MW  | MolForm  | CAS#        | Qual |
|-----------|----------------------------|-----|----------|-------------|------|
| 1         | 9-Octadecenoic acid, (E)-  | 282 | C18H34O2 | 000112-79-8 | 91   |
| 2         | cis-Vaccenic acid          | 282 | C18H34O2 | 000506-17-2 | 66   |
| 3         | cis-13-Octadecenoic acid   | 282 | C18H34O2 | 013126-39-1 | 62   |
| 4         | trans-13-Octadecenoic acid | 282 | C18H34O2 | 000693-71-0 | 53   |
| 5         | 9-Eicosene, (E)-           | 280 | C20H40   | 074685-29-3 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
Data File : BN034340.D  
Acq On : 02 Oct 2024 16:57  
Operator : RC/JU  
Sample : P4016-18  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

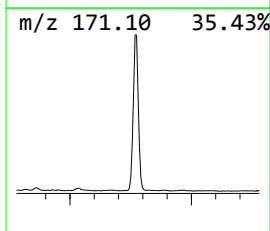
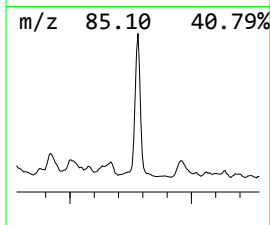
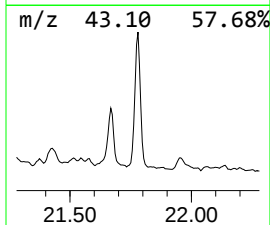
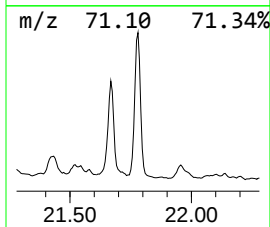
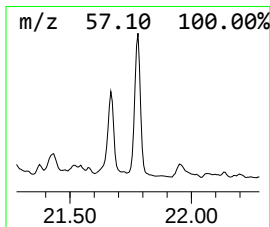
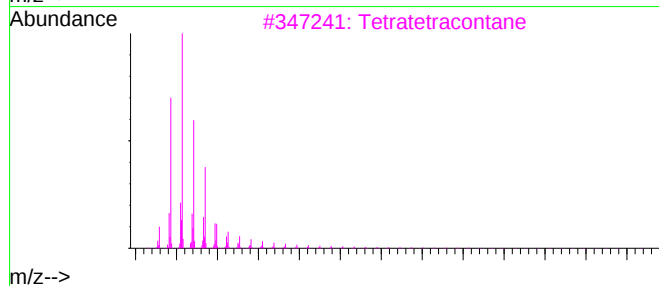
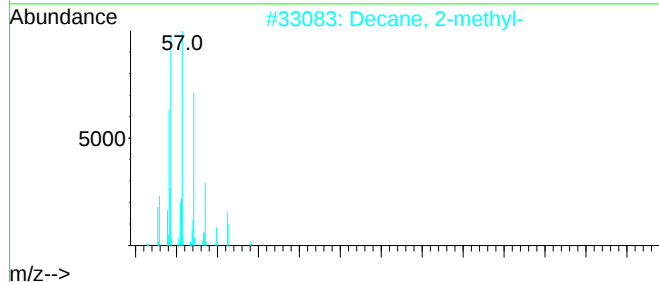
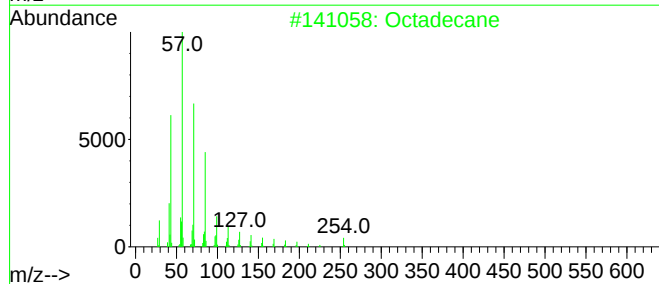
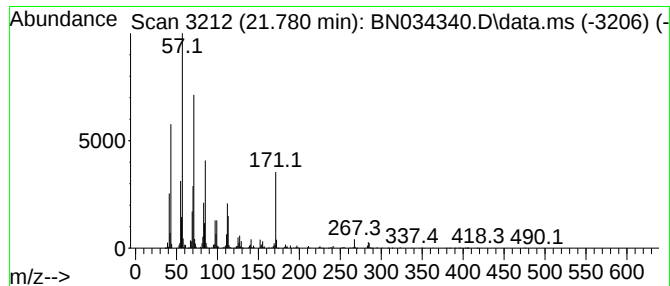
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 76 (DEL) Alkane: Straight-Chai... Concentration Rank 24

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 21.780 | 28.57 ng/ul | 8484780 | Chrysene-d12     | 21.033 |

| Hit# of 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|-----------|-------------------------------------|-----|----------|--------------|------|
| 1         | Octadecane                          | 254 | C18H38   | 000593-45-3  | 78   |
| 2         | Decane, 2-methyl-                   | 156 | C11H24   | 006975-98-0  | 62   |
| 3         | Tetratetracontane                   | 619 | C44H90   | 007098-22-8  | 58   |
| 4         | Oxalic acid, 2-ethylhexyl tetrad... | 398 | C24H46O4 | 1000309-39-7 | 58   |
| 5         | Heptadecane, 9-octyl-               | 352 | C25H52   | 007225-64-1  | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
 Data File : BN034340.D  
 Acq On : 02 Oct 2024 16:57  
 Operator : RC/JU  
 Sample : P4016-18  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 DDC23

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
 Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |          |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|----------|------|
|                    |        |         |       |          | #                     | RT     | Resp     | Conc |
| (DEL) Alkane: S... | 5.463  | 29.6    | ng/u1 | 26836400 | 1                     | 7.392  | 18114600 | 20.0 |
| (DEL) Alkane: C... | 5.704  | 8.2     | ng/u1 | 7448920  | 1                     | 7.392  | 18114600 | 20.0 |
| Hexylene glycol    | 5.834  | 23.2    | ng/u1 | 20984500 | 1                     | 7.392  | 18114600 | 20.0 |
| (DEL) Alkane: S... | 5.916  | 9.9     | ng/u1 | 8950850  | 1                     | 7.392  | 18114600 | 20.0 |
| (DEL) Alkane: S... | 5.993  | 10.2    | ng/u1 | 9189080  | 1                     | 7.392  | 18114600 | 20.0 |
| Cyclohexanone, ... | 6.057  | 12.1    | ng/u1 | 10925600 | 1                     | 7.392  | 18114600 | 20.0 |
| (DEL) Alkane: S... | 6.322  | 8.8     | ng/u1 | 7962340  | 1                     | 7.392  | 18114600 | 20.0 |
| Benzene, propyl-   | 6.387  | 11.7    | ng/u1 | 10559800 | 1                     | 7.392  | 18114600 | 20.0 |
| (DEL) Alkane: S... | 6.422  | 12.1    | ng/u1 | 10941100 | 1                     | 7.392  | 18114600 | 20.0 |
| (DEL) Alkane: S... | 6.481  | 18.9    | ng/u1 | 17136400 | 1                     | 7.392  | 18114600 | 20.0 |
| Benzene, 1-ethy... | 6.510  | 9.1     | ng/u1 | 8241200  | 1                     | 7.392  | 18114600 | 20.0 |
| Benzene, 1-ethy... | 6.798  | 13.5    | ng/u1 | 12195100 | 1                     | 7.392  | 18114600 | 20.0 |
| 2-Heptanone, 4,... | 6.857  | 13.6    | ng/u1 | 12274100 | 1                     | 7.392  | 18114600 | 20.0 |
| (DEL) Alkane: S... | 7.057  | 51.8    | ng/u1 | 46871800 | 1                     | 7.392  | 18114600 | 20.0 |
| (DEL) Alkane: S... | 7.245  | 7.4     | ng/u1 | 6684190  | 1                     | 7.392  | 18114600 | 20.0 |
| 1-Hexanol, 2-et... | 7.498  | 32.6    | ng/u1 | 29546800 | 1                     | 7.392  | 18114600 | 20.0 |
| D-Limonene         | 7.628  | 16.9    | ng/u1 | 15263000 | 1                     | 7.392  | 18114600 | 20.0 |
| (DEL) Alkane: C... | 7.681  | 17.4    | ng/u1 | 15742900 | 1                     | 7.392  | 18114600 | 20.0 |
| Cyclohexanone, ... | 7.834  | 8.3     | ng/u1 | 7480650  | 1                     | 7.392  | 18114600 | 20.0 |
| Benzene, 1,3-di... | 7.887  | 9.2     | ng/u1 | 8310120  | 1                     | 7.392  | 18114600 | 20.0 |
| Benzene, 1-meth... | 7.945  | 24.4    | ng/u1 | 22055200 | 1                     | 7.392  | 18114600 | 20.0 |
| (DEL) Alkane: S... | 7.998  | 12.7    | ng/u1 | 11457600 | 1                     | 7.392  | 18114600 | 20.0 |
| Benzene, 4-ethy... | 8.039  | 11.7    | ng/u1 | 10585600 | 1                     | 7.392  | 18114600 | 20.0 |
| (DEL) Alkane: S... | 8.175  | 19.9    | ng/u1 | 18037700 | 1                     | 7.392  | 18114600 | 20.0 |
| o-Cymene           | 8.398  | 12.3    | ng/u1 | 11129000 | 1                     | 7.392  | 18114600 | 20.0 |
| (DEL) Alkane: S... | 8.651  | 46.5    | ng/u1 | 42131700 | 1                     | 7.392  | 18114600 | 20.0 |
| unknown-01         | 8.745  | 20.9    | ng/u1 | 18927000 | 1                     | 7.392  | 18114600 | 20.0 |
| (DEL) Alkane: C... | 9.186  | 3.5     | ng/u1 | 9020800  | 2                     | 10.151 | 51293200 | 20.0 |
| (DEL) Alkane: S... | 9.475  | 5.8     | ng/u1 | 14837400 | 2                     | 10.151 | 51293200 | 20.0 |
| (DEL) Alkane: S... | 9.534  | 3.3     | ng/u1 | 8392610  | 2                     | 10.151 | 51293200 | 20.0 |
| (DEL) Alkane: S... | 9.616  | 2.9     | ng/u1 | 7532900  | 2                     | 10.151 | 51293200 | 20.0 |
| (DEL) Alkane: S... | 9.716  | 4.8     | ng/u1 | 12185600 | 2                     | 10.151 | 51293200 | 20.0 |
| (DEL) Alkane: S... | 10.345 | 2.5     | ng/u1 | 6447510  | 2                     | 10.151 | 51293200 | 20.0 |
| unknown-02         | 10.563 | 11.0    | ng/u1 | 28183000 | 2                     | 10.151 | 51293200 | 20.0 |
| (DEL) Alkane: S... | 11.628 | 11.6    | ng/u1 | 29795500 | 2                     | 10.151 | 51293200 | 20.0 |
| (DEL) Alkane: C... | 11.657 | 4.7     | ng/u1 | 11914700 | 2                     | 10.151 | 51293200 | 20.0 |
| unknown-03         | 12.280 | 19.6    | ng/u1 | 7066660  | 3                     | 14.051 | 7211030  | 20.0 |
| Bacchotricuneat... | 12.598 | 29.8    | ng/u1 | 10729400 | 3                     | 14.051 | 7211030  | 20.0 |
| Propanoic acid,... | 12.810 | 17.0    | ng/u1 | 6120330  | 3                     | 14.051 | 7211030  | 20.0 |
| 2,6-Pyridinedia... | 13.104 | 19.4    | ng/u1 | 6992670  | 3                     | 14.051 | 7211030  | 20.0 |
| Naphthalene, 1,... | 13.210 | 30.7    | ng/u1 | 11083800 | 3                     | 14.051 | 7211030  | 20.0 |
| Naphthalene, 2,... | 13.369 | 30.0    | ng/u1 | 10825900 | 3                     | 14.051 | 7211030  | 20.0 |
| Naphthalene, 1,... | 13.422 | 36.9    | ng/u1 | 13286200 | 3                     | 14.051 | 7211030  | 20.0 |
| (DEL) Alkane: S... | 13.569 | 45.9    | ng/u1 | 16540700 | 3                     | 14.051 | 7211030  | 20.0 |
| Sulfurous acid,... | 13.998 | 77.4    | ng/u1 | 27908000 | 3                     | 14.051 | 7211030  | 20.0 |
| 9-Methyl-S-octa... | 14.216 | 49.0    | ng/u1 | 17650800 | 3                     | 14.051 | 7211030  | 20.0 |
| Naphthalene, 1,... | 14.545 | 20.8    | ng/u1 | 7487760  | 3                     | 14.051 | 7211030  | 20.0 |
| (DEL) Alkane: S... | 14.610 | 23.6    | ng/u1 | 8492740  | 3                     | 14.051 | 7211030  | 20.0 |
| 4,6,8-Trimethyl... | 14.839 | 71.2    | ng/u1 | 25657200 | 3                     | 14.051 | 7211030  | 20.0 |
| (DEL) Alkane: S... | 14.957 | 59.5    | ng/u1 | 21449000 | 3                     | 14.051 | 7211030  | 20.0 |
| (DEL) Alkane: S... | 15.357 | 44.4    | ng/u1 | 16014900 | 3                     | 14.051 | 7211030  | 20.0 |
| unknown-04         | 15.710 | 33.5    | ng/u1 | 11674400 | 4                     | 16.833 | 6958880  | 20.0 |

|                    |        |      |       |          |   |        |         |      |
|--------------------|--------|------|-------|----------|---|--------|---------|------|
| (DEL) Alkane: S... | 15.821 | 58.3 | ng/ul | 20300300 | 4 | 16.833 | 6958880 | 20.0 |
| (DEL) Alkane: S... | 16.616 | 36.5 | ng/ul | 12693300 | 4 | 16.833 | 6958880 | 20.0 |
| (DEL) Alkane: S... | 16.663 | 39.5 | ng/ul | 13727700 | 4 | 16.833 | 6958880 | 20.0 |
| (DEL) Alkane: S... | 17.351 | 39.6 | ng/ul | 13769400 | 4 | 16.833 | 6958880 | 20.0 |
| Hexadecanoic ac... | 17.521 | 38.4 | ng/ul | 13365300 | 4 | 16.833 | 6958880 | 20.0 |
| Anthracene, 2-m... | 17.745 | 17.7 | ng/ul | 6157810  | 4 | 16.833 | 6958880 | 20.0 |
| (DEL) Alkane: S... | 18.039 | 34.7 | ng/ul | 12068600 | 4 | 16.833 | 6958880 | 20.0 |
| (DEL) Alkane: S... | 18.686 | 37.8 | ng/ul | 13145400 | 4 | 16.833 | 6958880 | 20.0 |
| (DEL) Alkane: S... | 19.809 | 26.6 | ng/ul | 7894570  | 5 | 21.033 | 5938710 | 20.0 |
| (DEL) Alkane: S... | 20.780 | 43.7 | ng/ul | 12965800 | 5 | 21.033 | 5938710 | 20.0 |
| (DEL) Alkane: S... | 21.256 | 20.0 | ng/ul | 5938710  | 5 | 21.033 | 5938710 | 20.0 |
| 9-Octadecenoic ... | 21.668 | 22.8 | ng/ul | 6781180  | 5 | 21.033 | 5938710 | 20.0 |
| (DEL) Alkane: S... | 21.780 | 28.6 | ng/ul | 8484780  | 5 | 21.033 | 5938710 | 20.0 |

Instrument :

BNAN

ClientSampleId :

DDC23