

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
 Data File : BN033874.D  
 Acq On : 11 Sep 2024 23:28  
 Operator : JU/RC  
 Sample : P3878-02DL 20X  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 DDC46DL

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

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: BN033874.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         | 5.563       | 431           | 438         | 446          | rVB2     | 304899         | 469729        | 2.39%           | 0.586%        |
| 2         | 6.093       | 523           | 528         | 533          | rBV      | 130513         | 186709        | 0.95%           | 0.233%        |
| 3         | 6.163       | 533           | 540         | 543          | rVV4     | 85322          | 196035        | 1.00%           | 0.244%        |
| 4         | 6.493       | 590           | 596         | 597          | rBV      | 126128         | 195686        | 1.00%           | 0.244%        |
| 5         | 6.516       | 597           | 600         | 605          | rVV      | 200434         | 305845        | 1.56%           | 0.381%        |
| 6         | 6.575       | 605           | 610         | 612          | rVV      | 254775         | 353556        | 1.80%           | 0.441%        |
| 7         | 6.605       | 612           | 615         | 620          | rVV      | 340817         | 524610        | 2.67%           | 0.654%        |
| 8         | 6.681       | 620           | 628         | 633          | rVV4     | 367580         | 837510        | 4.26%           | 1.045%        |
| 9         | 6.740       | 633           | 638         | 644          | rVB      | 184738         | 271028        | 1.38%           | 0.338%        |
| 10        | 6.899       | 660           | 665         | 669          | rVV      | 235953         | 360343        | 1.83%           | 0.449%        |
| 11        | 6.946       | 669           | 673         | 679          | rVV      | 377497         | 598415        | 3.04%           | 0.746%        |
| 12        | 7.140       | 701           | 706         | 717          | rVV2     | 1569982        | 2717651       | 13.82%          | 3.389%        |
| 13        | 7.493       | 760           | 766         | 772          | rVV2     | 800292         | 1285838       | 6.54%           | 1.604%        |
| 14        | 7.575       | 776           | 780         | 783          | rVV2     | 181893         | 331775        | 1.69%           | 0.414%        |
| 15        | 7.610       | 783           | 786         | 796          | rVV3     | 199660         | 391935        | 1.99%           | 0.489%        |
| 16        | 7.775       | 809           | 814         | 823          | rVV6     | 128036         | 387163        | 1.97%           | 0.483%        |
| 17        | 7.922       | 834           | 839         | 845          | rVV      | 358484         | 598564        | 3.04%           | 0.747%        |
| 18        | 7.981       | 845           | 849         | 854          | rVV2     | 243069         | 420930        | 2.14%           | 0.525%        |
| 19        | 8.040       | 854           | 859         | 864          | rVB2     | 363227         | 602339        | 3.06%           | 0.751%        |
| 20        | 8.093       | 864           | 868         | 871          | rBV      | 213542         | 295505        | 1.50%           | 0.369%        |
| 21        | 8.134       | 871           | 875         | 878          | rVV2     | 340078         | 575647        | 2.93%           | 0.718%        |
| 22        | 8.163       | 878           | 880         | 892          | rVB3     | 354388         | 603110        | 3.07%           | 0.752%        |
| 23        | 8.269       | 892           | 898         | 900          | rBV      | 293751         | 459184        | 2.34%           | 0.573%        |
| 24        | 8.293       | 900           | 902         | 910          | rVB      | 224415         | 335687        | 1.71%           | 0.419%        |
| 25        | 8.440       | 923           | 927         | 931          | rBV      | 176318         | 262320        | 1.33%           | 0.327%        |
| 26        | 8.493       | 931           | 936         | 943          | rVB      | 239915         | 368257        | 1.87%           | 0.459%        |
| 27        | 8.593       | 944           | 953         | 963          | rBV4     | 234986         | 603255        | 3.07%           | 0.752%        |
| 28        | 8.728       | 971           | 976         | 982          | rVV      | 1294820        | 1849684       | 9.41%           | 2.307%        |
| 29        | 8.840       | 990           | 995         | 1001         | rVB5     | 161447         | 366903        | 1.87%           | 0.458%        |
| 30        | 8.922       | 1001          | 1009        | 1013         | rBV3     | 124425         | 266996        | 1.36%           | 0.333%        |
| 31        | 8.981       | 1013          | 1019        | 1026         | rVV5     | 73316          | 213949        | 1.09%           | 0.267%        |
| 32        | 9.163       | 1047          | 1050        | 1063         | rVB2     | 121440         | 299613        | 1.52%           | 0.374%        |
| 33        | 9.281       | 1063          | 1070        | 1076         | rVB2     | 94548          | 183347        | 0.93%           | 0.229%        |
| 34        | 9.363       | 1077          | 1084        | 1088         | rBV5     | 82012          | 212731        | 1.08%           | 0.265%        |
| 35        | 9.422       | 1088          | 1094        | 1098         | rBV3     | 99528          | 179858        | 0.91%           | 0.224%        |
| 36        | 9.569       | 1109          | 1119        | 1125         | rVB2     | 150495         | 354786        | 1.80%           | 0.442%        |

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 ClientSampleId :  
 DDC46DL

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

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

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 9.634  | 1125 | 1130 | 1133 | rBV3 | 118885  | 191088  | 0.97%  | 0.238% |
| 38 | 9.822  | 1156 | 1162 | 1166 | rBV2 | 116113  | 210969  | 1.07%  | 0.263% |
| 39 | 9.969  | 1181 | 1187 | 1192 | rBV2 | 123199  | 188242  | 0.96%  | 0.235% |
| 40 | 10.263 | 1226 | 1237 | 1253 | rVB3 | 1946098 | 4318738 | 21.97% | 5.386% |
| 41 | 10.393 | 1253 | 1259 | 1265 | rBV2 | 453624  | 758856  | 3.86%  | 0.946% |
| 42 | 10.657 | 1297 | 1304 | 1317 | rBV  | 356675  | 618224  | 3.14%  | 0.771% |
| 43 | 10.787 | 1317 | 1326 | 1333 | rBV  | 292009  | 560271  | 2.85%  | 0.699% |
| 44 | 11.193 | 1393 | 1395 | 1408 | rVB7 | 92379   | 195312  | 0.99%  | 0.244% |
| 45 | 11.304 | 1408 | 1414 | 1419 | rBV2 | 294731  | 489773  | 2.49%  | 0.611% |
| 46 | 11.545 | 1451 | 1455 | 1460 | rVB2 | 130397  | 202105  | 1.03%  | 0.252% |
| 47 | 11.710 | 1475 | 1483 | 1487 | rBV  | 325718  | 599823  | 3.05%  | 0.748% |
| 48 | 11.928 | 1510 | 1520 | 1524 | rVV  | 230696  | 416192  | 2.12%  | 0.519% |
| 49 | 11.969 | 1524 | 1527 | 1532 | rVB  | 171819  | 216654  | 1.10%  | 0.270% |
| 50 | 12.128 | 1548 | 1554 | 1564 | rVB3 | 381404  | 745395  | 3.79%  | 0.930% |
| 51 | 12.551 | 1619 | 1626 | 1634 | rVB4 | 100737  | 206654  | 1.05%  | 0.258% |
| 52 | 12.692 | 1645 | 1650 | 1658 | rVB3 | 120617  | 283412  | 1.44%  | 0.353% |
| 53 | 12.892 | 1675 | 1684 | 1692 | rBV5 | 83756   | 229321  | 1.17%  | 0.286% |
| 54 | 12.981 | 1692 | 1699 | 1705 | rBV  | 407262  | 692706  | 3.52%  | 0.864% |
| 55 | 13.192 | 1732 | 1735 | 1745 | rVB3 | 177616  | 307583  | 1.56%  | 0.384% |
| 56 | 13.298 | 1745 | 1753 | 1765 | rBV6 | 230110  | 779189  | 3.96%  | 0.972% |
| 57 | 13.416 | 1769 | 1773 | 1776 | rBV2 | 144983  | 211337  | 1.07%  | 0.264% |
| 58 | 13.457 | 1776 | 1780 | 1784 | rVV  | 259416  | 421034  | 2.14%  | 0.525% |
| 59 | 13.510 | 1784 | 1789 | 1794 | rVV4 | 215195  | 368849  | 1.88%  | 0.460% |
| 60 | 13.563 | 1794 | 1798 | 1803 | rVB  | 156110  | 230101  | 1.17%  | 0.287% |
| 61 | 13.651 | 1804 | 1813 | 1817 | rBV2 | 141377  | 316330  | 1.61%  | 0.395% |
| 62 | 13.692 | 1817 | 1820 | 1825 | rVV2 | 178105  | 231632  | 1.18%  | 0.289% |
| 63 | 13.828 | 1835 | 1843 | 1846 | rVV3 | 154701  | 368548  | 1.87%  | 0.460% |
| 64 | 14.069 | 1880 | 1884 | 1888 | rBV  | 463202  | 585369  | 2.98%  | 0.730% |
| 65 | 14.139 | 1888 | 1896 | 1902 | rVV  | 1459425 | 2211820 | 11.25% | 2.759% |
| 66 | 14.245 | 1909 | 1914 | 1917 | rVB4 | 141884  | 188048  | 0.96%  | 0.235% |
| 67 | 14.357 | 1926 | 1933 | 1939 | rBV4 | 155234  | 432394  | 2.20%  | 0.539% |
| 68 | 14.545 | 1957 | 1965 | 1968 | rBV3 | 124241  | 265530  | 1.35%  | 0.331% |
| 69 | 14.628 | 1973 | 1979 | 1983 | rVB2 | 160890  | 252800  | 1.29%  | 0.315% |
| 70 | 14.692 | 1985 | 1990 | 1994 | rBV2 | 188181  | 260889  | 1.33%  | 0.325% |
| 71 | 14.775 | 1994 | 2004 | 2009 | rVV  | 1401502 | 2199728 | 11.19% | 2.744% |
| 72 | 14.822 | 2009 | 2012 | 2017 | rVB2 | 179366  | 248798  | 1.27%  | 0.310% |
| 73 | 14.939 | 2029 | 2032 | 2039 | rVV7 | 101376  | 208418  | 1.06%  | 0.260% |
| 74 | 15.028 | 2043 | 2047 | 2052 | rVV  | 477256  | 599608  | 3.05%  | 0.748% |
| 75 | 15.139 | 2062 | 2066 | 2070 | rVB  | 155505  | 213220  | 1.08%  | 0.266% |
| 76 | 15.257 | 2080 | 2086 | 2089 | rBV5 | 114841  | 227776  | 1.16%  | 0.284% |

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Instrument :  
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 ClientSampleId :  
 DDC46DL

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

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

|     |        |      |      |      |      |          |          |         |         |
|-----|--------|------|------|------|------|----------|----------|---------|---------|
| 77  | 15.433 | 2109 | 2116 | 2122 | rBV4 | 168069   | 407713   | 2.07%   | 0.509%  |
| 78  | 15.898 | 2189 | 2195 | 2198 | rBV  | 526177   | 785461   | 4.00%   | 0.980%  |
| 79  | 16.245 | 2249 | 2254 | 2260 | rBV5 | 80585    | 182387   | 0.93%   | 0.227%  |
| 80  | 16.557 | 2299 | 2307 | 2318 | rBV  | 13537530 | 19660443 | 100.00% | 24.521% |
|     |        |      |      |      |      |          |          |         |         |
| 81  | 16.686 | 2325 | 2329 | 2335 | rBV2 | 500683   | 686958   | 3.49%   | 0.857%  |
| 82  | 16.739 | 2335 | 2338 | 2347 | rVB  | 194926   | 323486   | 1.65%   | 0.403%  |
| 83  | 16.892 | 2359 | 2364 | 2369 | rBV  | 1925160  | 2591337  | 13.18%  | 3.232%  |
| 84  | 16.992 | 2376 | 2381 | 2390 | rBV  | 205664   | 347953   | 1.77%   | 0.434%  |
| 85  | 17.427 | 2451 | 2455 | 2459 | rBV  | 494395   | 588639   | 2.99%   | 0.734%  |
|     |        |      |      |      |      |          |          |         |         |
| 86  | 17.598 | 2480 | 2484 | 2493 | rVB2 | 795099   | 1066178  | 5.42%   | 1.330%  |
| 87  | 17.816 | 2517 | 2521 | 2528 | rVB2 | 208682   | 295038   | 1.50%   | 0.368%  |
| 88  | 18.122 | 2568 | 2573 | 2577 | rBV  | 439297   | 519326   | 2.64%   | 0.648%  |
| 89  | 18.780 | 2676 | 2685 | 2690 | rBV2 | 643846   | 1236052  | 6.29%   | 1.542%  |
| 90  | 18.922 | 2705 | 2709 | 2714 | rVB  | 301565   | 368842   | 1.88%   | 0.460%  |
|     |        |      |      |      |      |          |          |         |         |
| 91  | 19.298 | 2769 | 2773 | 2778 | rBV2 | 200806   | 282049   | 1.43%   | 0.352%  |
| 92  | 19.351 | 2778 | 2782 | 2788 | rVB  | 289435   | 309525   | 1.57%   | 0.386%  |
| 93  | 19.892 | 2870 | 2874 | 2878 | rBV  | 231819   | 262605   | 1.34%   | 0.328%  |
| 94  | 20.392 | 2955 | 2959 | 2964 | rBV  | 192850   | 233702   | 1.19%   | 0.291%  |
| 95  | 20.863 | 3035 | 3039 | 3044 | rVB  | 436554   | 505155   | 2.57%   | 0.630%  |
|     |        |      |      |      |      |          |          |         |         |
| 96  | 21.127 | 3079 | 3084 | 3093 | rVB  | 2644358  | 3478568  | 17.69%  | 4.339%  |
| 97  | 21.351 | 3118 | 3122 | 3129 | rVB  | 168679   | 241285   | 1.23%   | 0.301%  |
| 98  | 21.880 | 3207 | 3212 | 3221 | rVB3 | 259982   | 460992   | 2.34%   | 0.575%  |
| 99  | 23.715 | 3519 | 3524 | 3532 | rVB  | 165627   | 363249   | 1.85%   | 0.453%  |
| 100 | 23.909 | 3548 | 3557 | 3565 | rBV  | 1850559  | 4262397  | 21.68%  | 5.316%  |

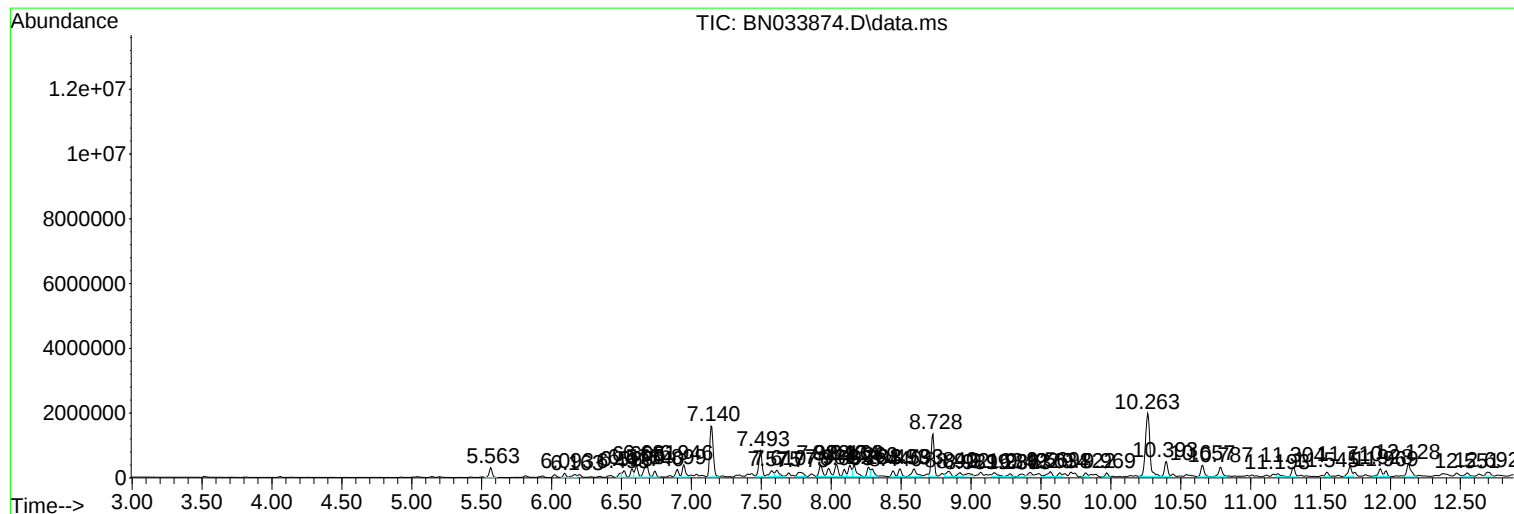
Sum of corrected areas: 80178569

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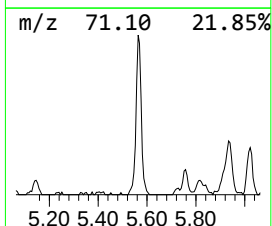
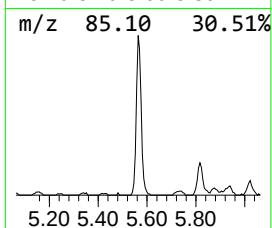
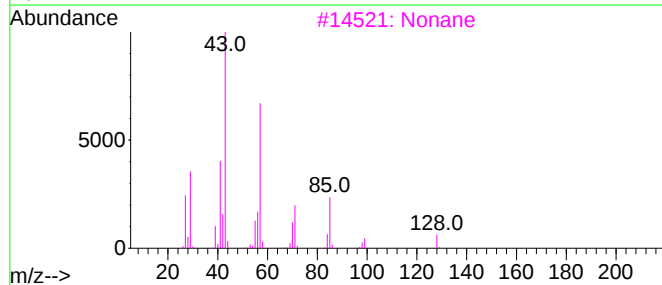
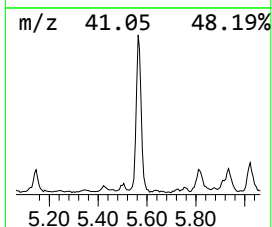
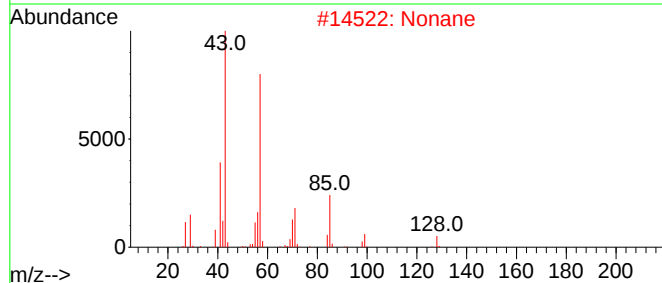
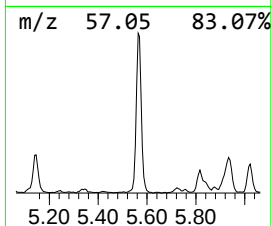
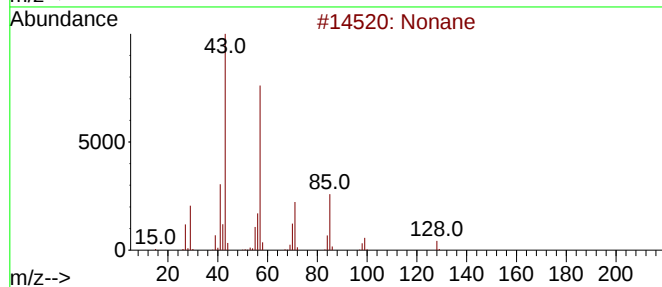
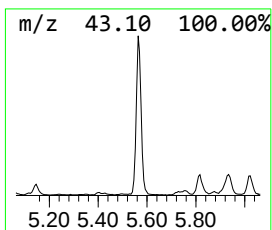
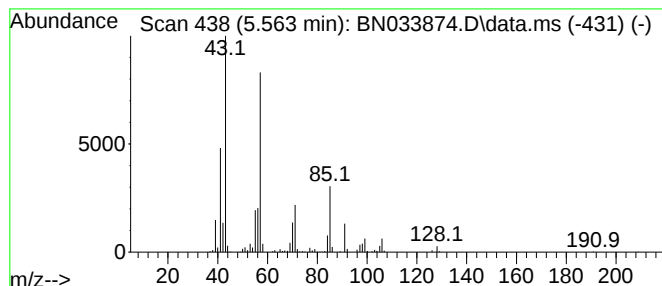
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 5.563 | 7.31 ng/ul | 469729 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of     | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|--------|---|--------------|-----|---------|-------------|------|
| 1    | Nonane |   |              | 128 | C9H20   | 000111-84-2 | 91   |
| 2    | Nonane |   |              | 128 | C9H20   | 000111-84-2 | 72   |
| 3    | Nonane |   |              | 128 | C9H20   | 000111-84-2 | 72   |
| 4    | Nonane |   |              | 128 | C9H20   | 000111-84-2 | 72   |
| 5    | Decane |   |              | 142 | C10H22  | 000124-18-5 | 53   |



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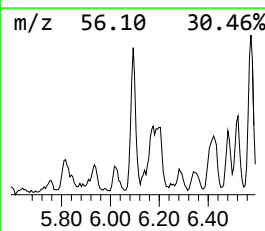
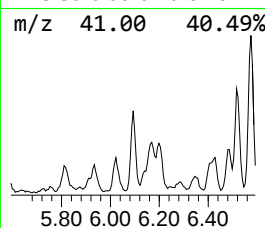
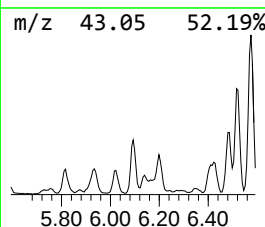
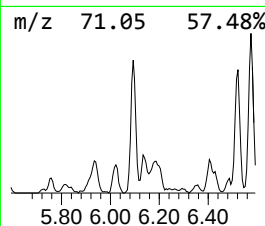
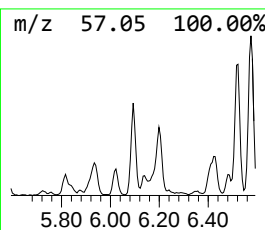
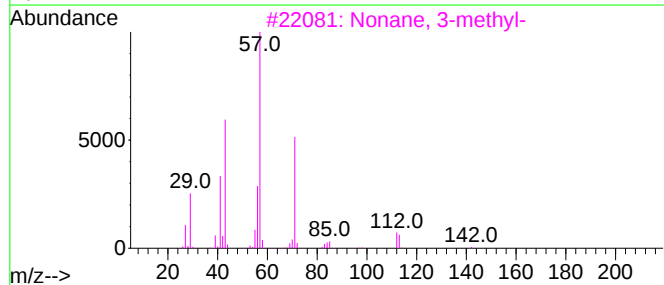
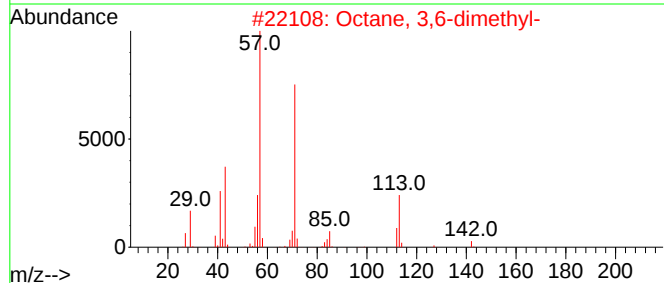
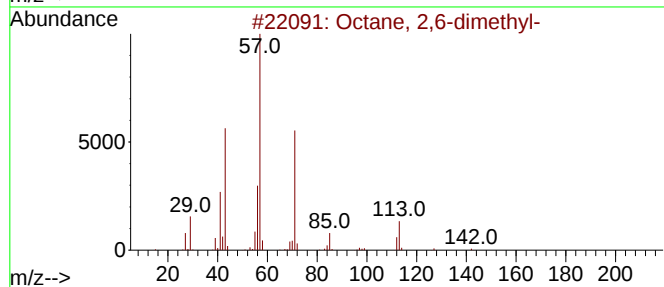
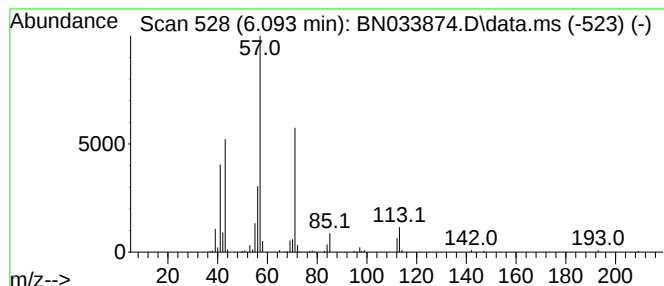
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 39

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 6.093 | 2.90 ng/ul | 186709 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Octane, 2,6-dimethyl- | 142 | C10H22  | 002051-30-1 | 93   |
| 2    |    |   | Octane, 3,6-dimethyl- | 142 | C10H22  | 015869-94-0 | 91   |
| 3    |    |   | Nonane, 3-methyl-     | 142 | C10H22  | 005911-04-6 | 90   |
| 4    |    |   | Octane, 3,6-dimethyl- | 142 | C10H22  | 015869-94-0 | 90   |
| 5    |    |   | Nonane, 3-methyl-     | 142 | C10H22  | 005911-04-6 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

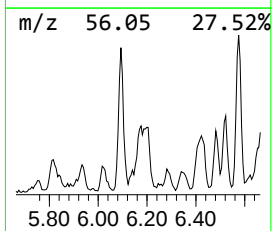
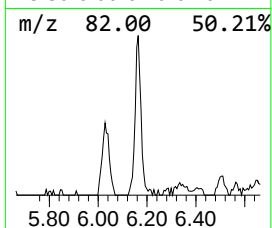
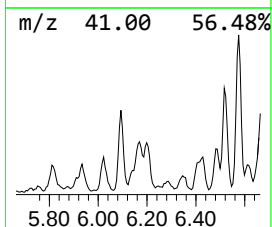
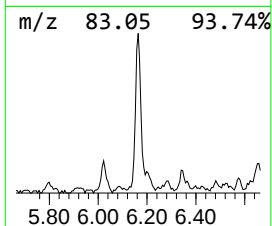
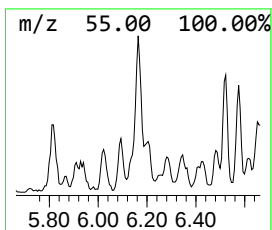
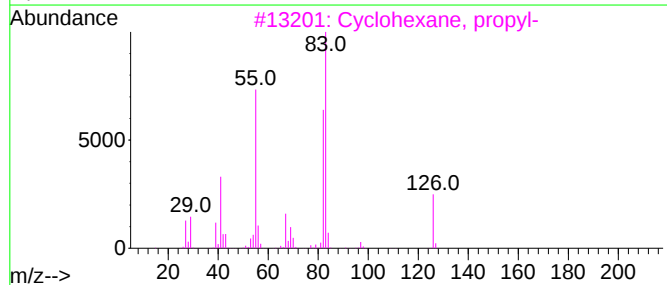
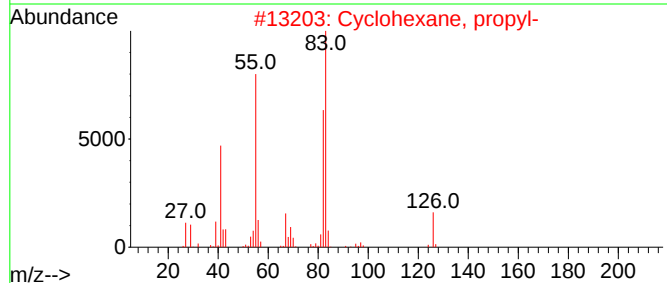
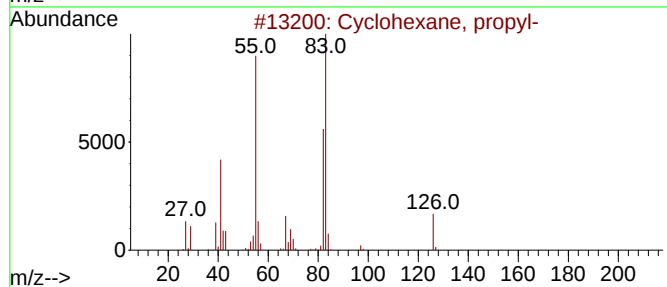
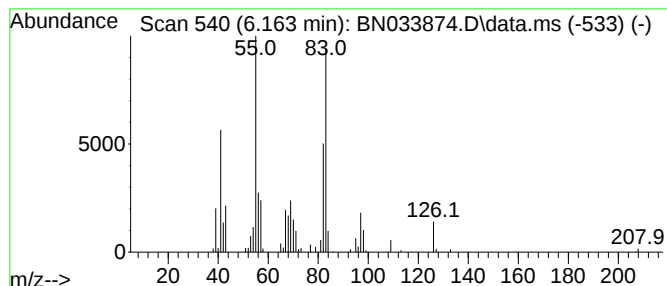
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 (DEL) Alkane: Cyclic6.163 Concentration Rank 36

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 6.163 | 3.05 ng/ul | 196035 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, propyl-           | 126 | C9H18   | 001678-92-8 | 80   |
| 2    |    |   | Cyclohexane, propyl-           | 126 | C9H18   | 001678-92-8 | 80   |
| 3    |    |   | Cyclohexane, propyl-           | 126 | C9H18   | 001678-92-8 | 80   |
| 4    |    |   | Cyclohexane, (3-methylpentyl)- | 168 | C12H24  | 061142-38-9 | 78   |
| 5    |    |   | Cyclohexane, propyl-           | 126 | C9H18   | 001678-92-8 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

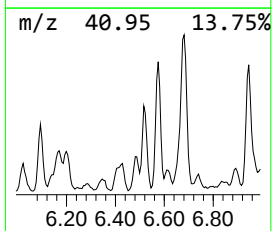
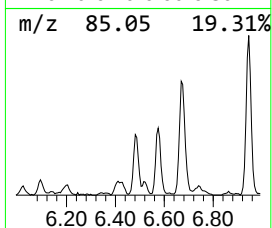
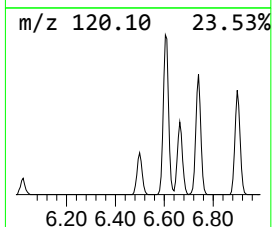
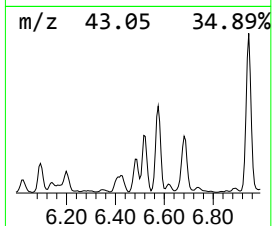
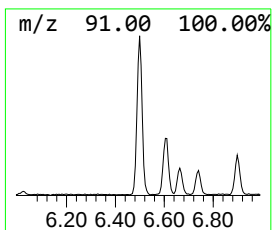
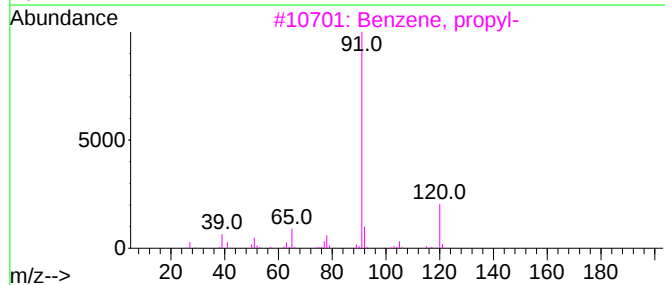
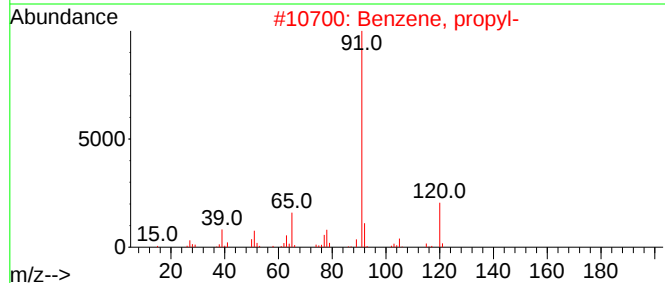
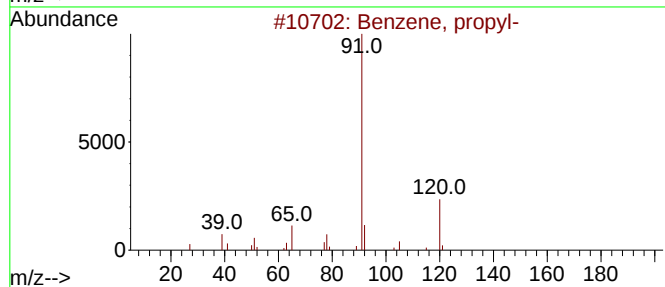
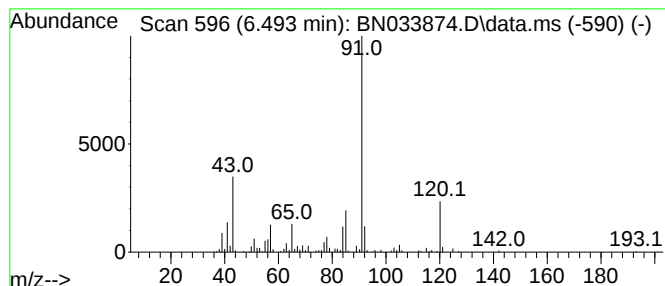
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Quant Title : SVOA CALIBRATION

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

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Peak Number 4 Benzene, propyl- Concentration Rank 37

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 6.493 | 3.04 ng/ul | 195686 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm   | CAS#        | Qual |
|------|----|---|------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | Benzene, propyl-                   | 120 | C9H12     | 000103-65-1 | 90   |
| 2    |    |   | Benzene, propyl-                   | 120 | C9H12     | 000103-65-1 | 90   |
| 3    |    |   | Benzene, propyl-                   | 120 | C9H12     | 000103-65-1 | 55   |
| 4    |    |   | Phenethylamine, N-benzyl-p-chloro- | 245 | C15H16ClN | 013622-43-0 | 50   |
| 5    |    |   | Benzene, propyl-                   | 120 | C9H12     | 000103-65-1 | 49   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

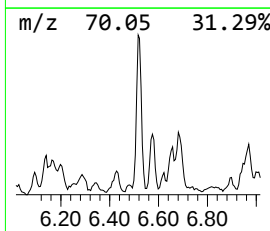
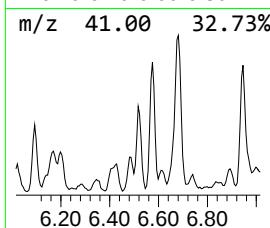
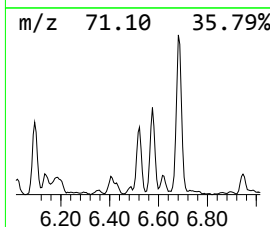
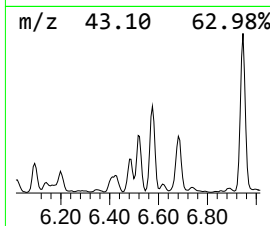
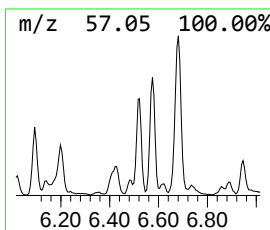
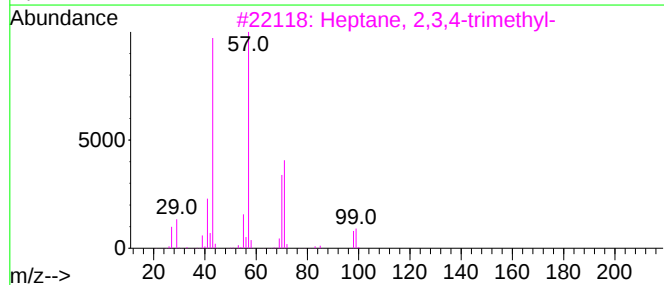
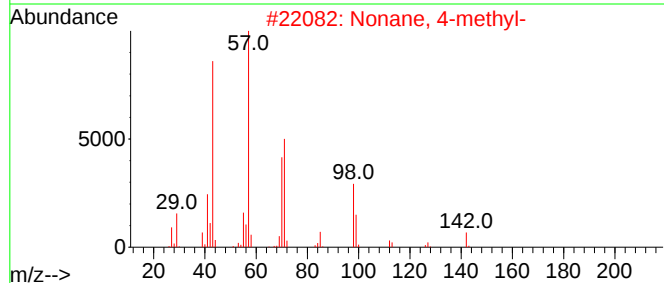
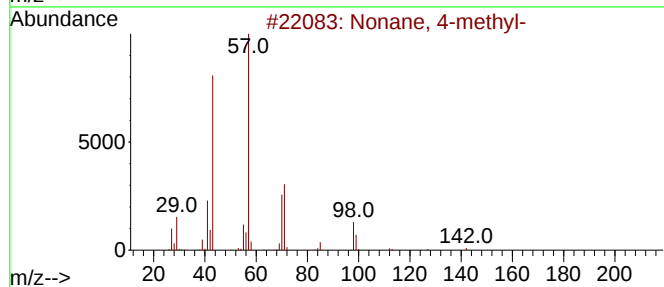
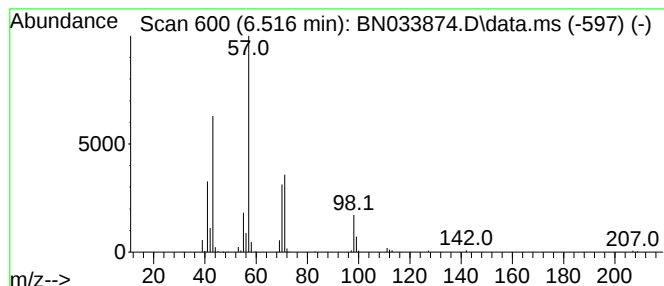
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 27

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 6.516 | 4.76 ng/ul | 305845 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Nonane, 4-methyl-             | 142 | C10H22  | 017301-94-9 | 91   |
| 2    |    |   | Nonane, 4-methyl-             | 142 | C10H22  | 017301-94-9 | 78   |
| 3    |    |   | Heptane, 2,3,4-trimethyl-     | 142 | C10H22  | 052896-95-4 | 72   |
| 4    |    |   | Pentane, 2,2,3,3-tetramethyl- | 128 | C9H20   | 007154-79-2 | 59   |
| 5    |    |   | Undecane, 4,5-dimethyl-       | 184 | C13H28  | 017312-79-7 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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Quant Title : SVOA CALIBRATION

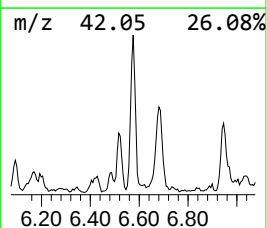
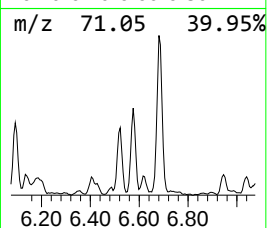
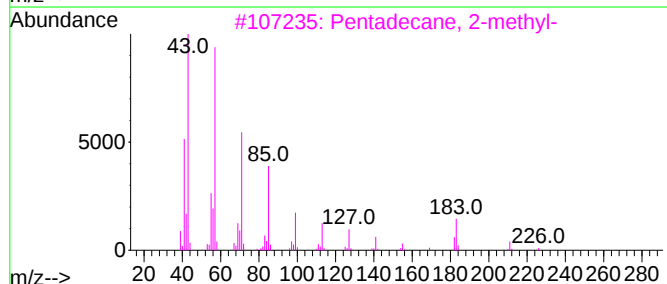
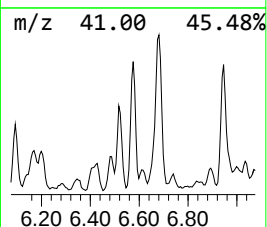
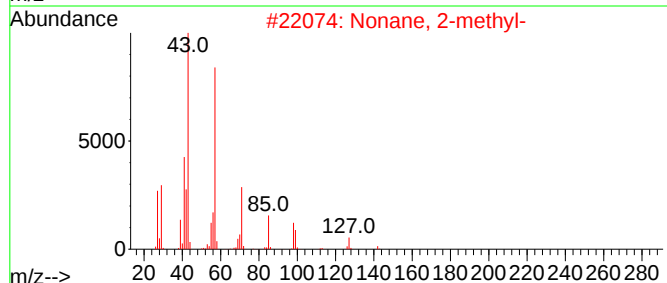
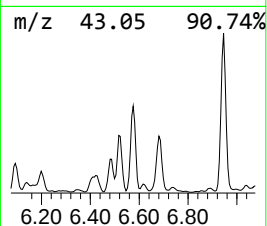
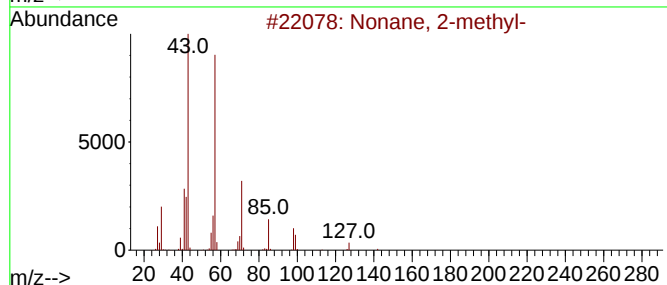
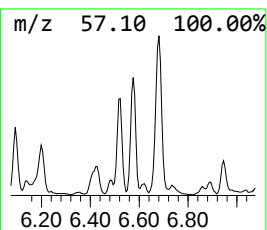
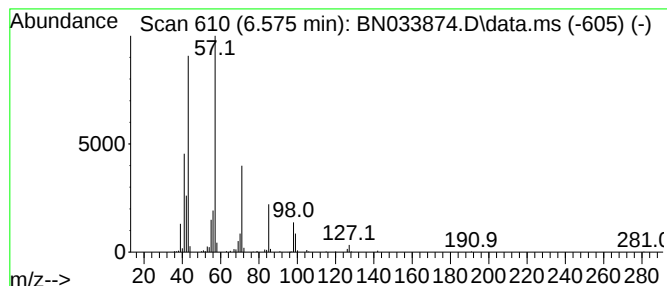
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 21

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 6.575 | 5.50 ng/ul | 353556 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Nonane, 2-methyl-      | 142 | C10H22  | 000871-83-0 | 94   |
| 2    |    |   | Nonane, 2-methyl-      | 142 | C10H22  | 000871-83-0 | 83   |
| 3    |    |   | Pentadecane, 2-methyl- | 226 | C16H34  | 001560-93-6 | 72   |
| 4    |    |   | Decane, 2,9-dimethyl-  | 170 | C12H26  | 001002-17-1 | 64   |
| 5    |    |   | Octane, 2-methyl-      | 128 | C9H20   | 003221-61-2 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

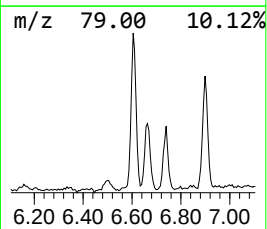
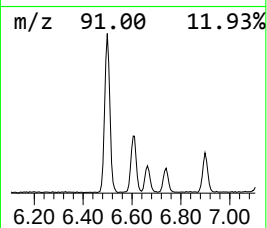
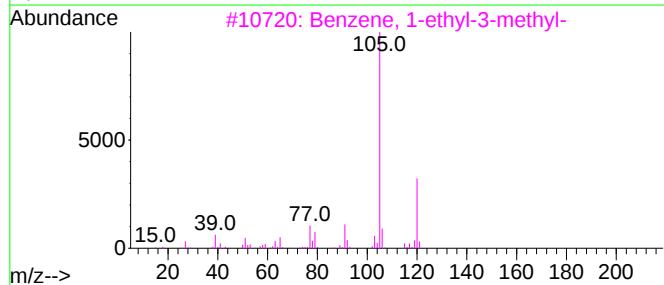
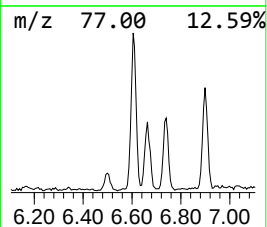
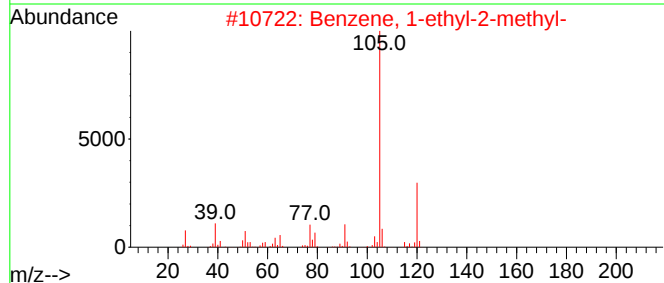
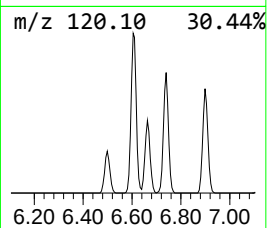
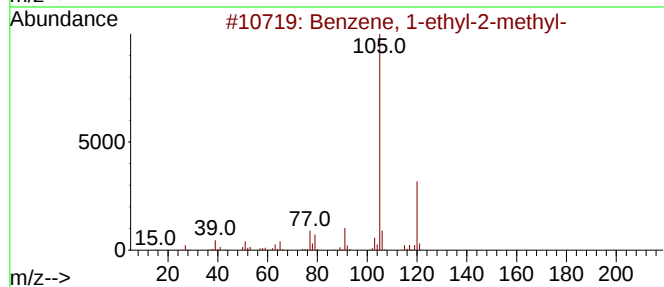
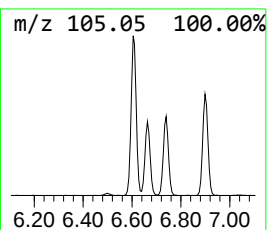
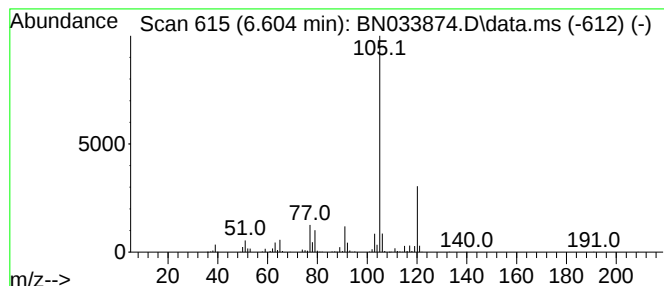
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 6.604 | 8.16 ng/ul | 524610 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 95   |
| 2    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 95   |
| 3    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 95   |
| 4    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 94   |
| 5    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

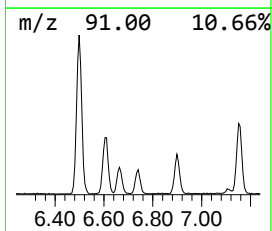
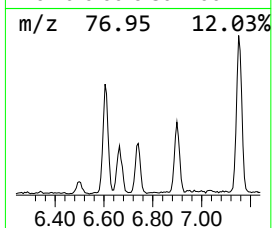
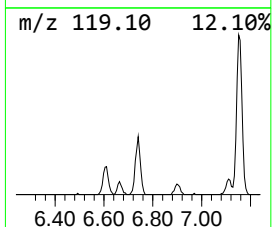
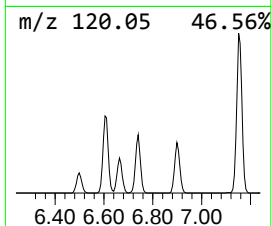
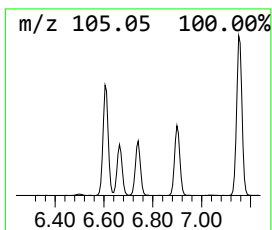
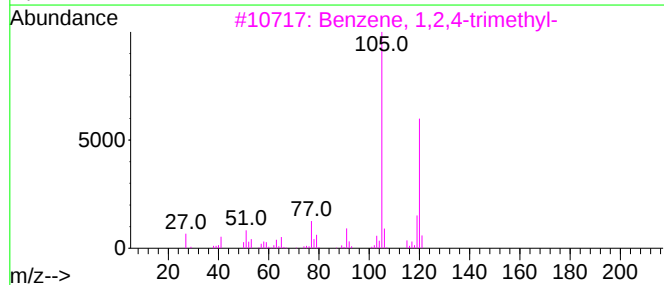
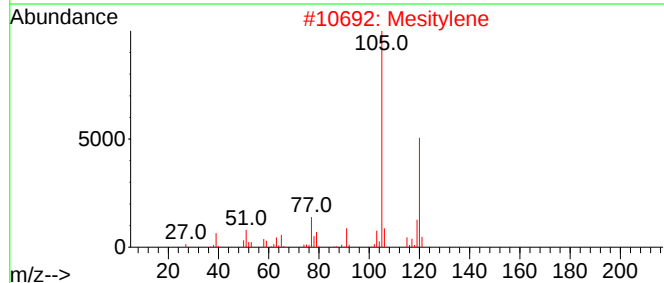
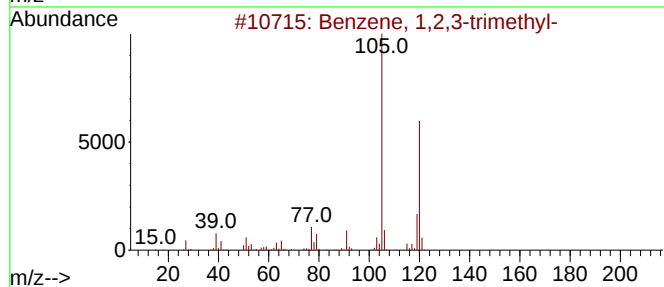
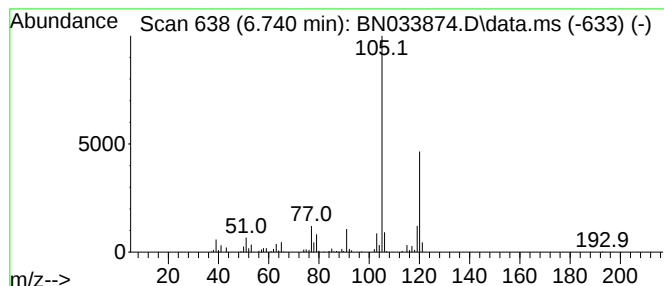
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Quant Title : SVOA CALIBRATION

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

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Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 30

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 6.740 | 4.22 ng/ul | 271028 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 97   |
| 2    |    |   | Mesitylene                | 120 | C9H12   | 000108-67-8 | 97   |
| 3    |    |   | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 95   |
| 4    |    |   | Mesitylene                | 120 | C9H12   | 000108-67-8 | 94   |
| 5    |    |   | Mesitylene                | 120 | C9H12   | 000108-67-8 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

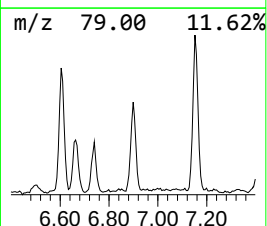
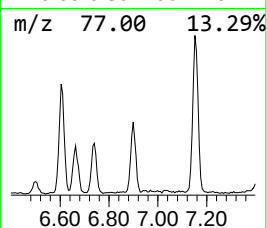
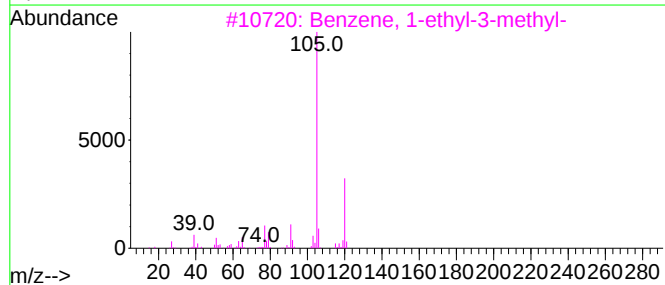
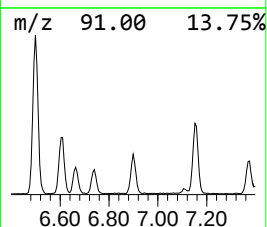
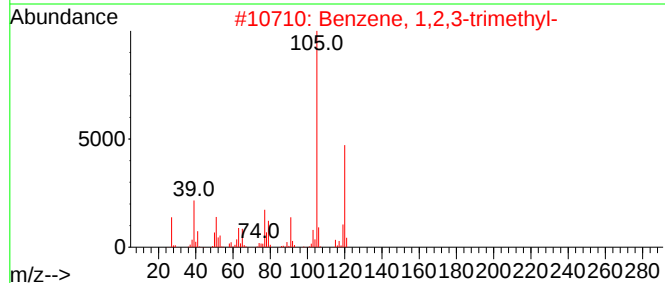
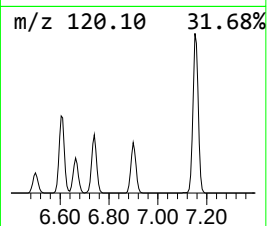
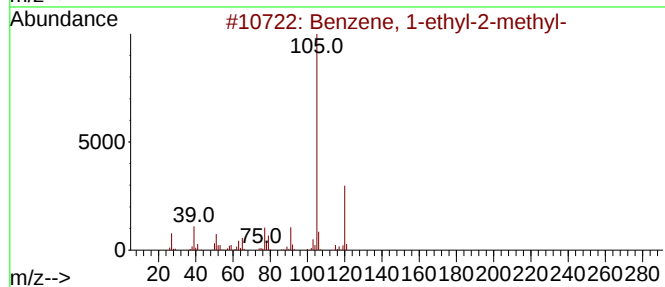
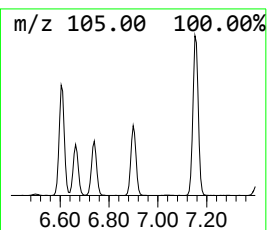
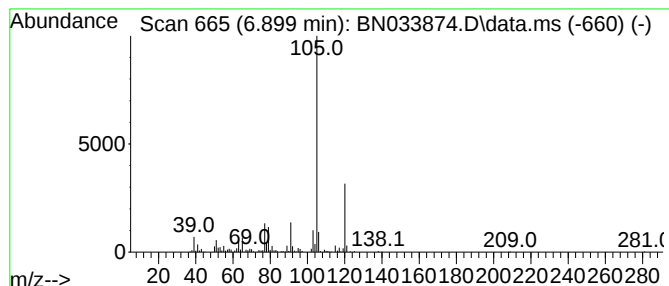
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 9 Benzene, 1-ethyl-3-methyl- Concentration Rank 20

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 6.899 | 5.60 ng/ul | 360343 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 94   |
| 2    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |
| 3    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |
| 4    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |
| 5    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

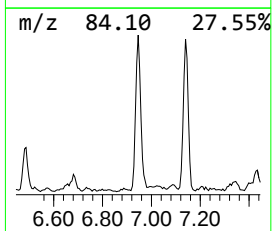
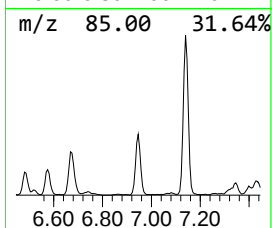
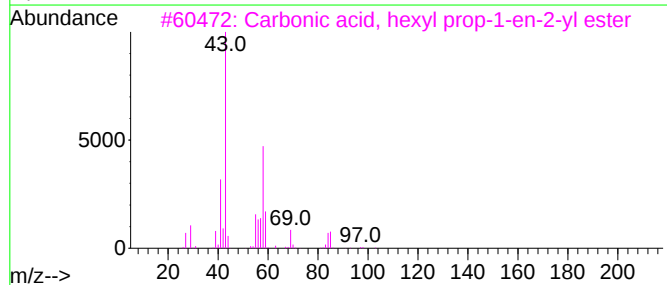
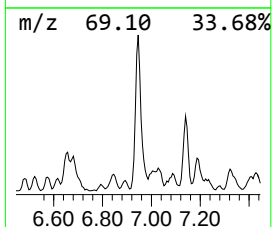
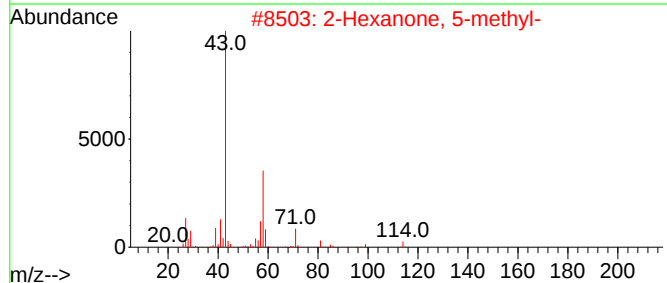
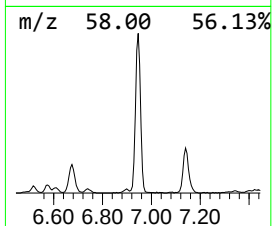
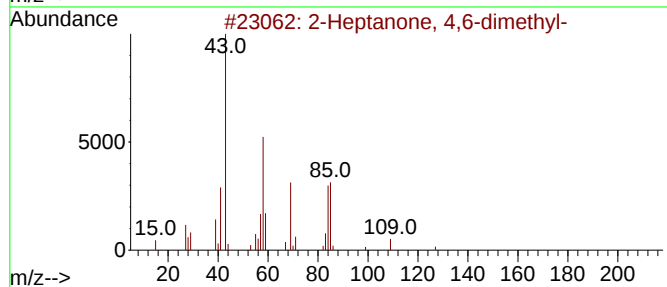
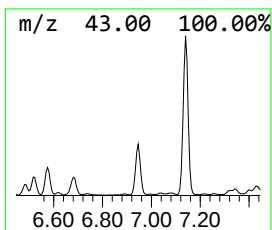
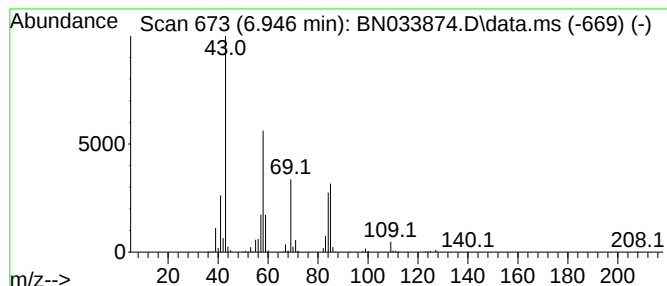
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Quant Title : SVOA CALIBRATION

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

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Peak Number 10 2-Heptanone, 4,6-dimethyl- Concentration Rank 6

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 6.946 | 9.31 ng/ul | 598415 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 2-Heptanone, 4,6-dimethyl-          | 142 | C9H18O   | 019549-80-5  | 90   |
| 2    |    |   | 2-Hexanone, 5-methyl-               | 114 | C7H14O   | 000110-12-3  | 50   |
| 3    |    |   | Carbonic acid, hexyl prop-1-en-2... | 186 | C10H18O3 | 1000382-54-2 | 50   |
| 4    |    |   | 2-Hexanone, 5-methyl-               | 114 | C7H14O   | 000110-12-3  | 50   |
| 5    |    |   | 2-Hexanone, 4-methyl-               | 114 | C7H14O   | 000105-42-0  | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

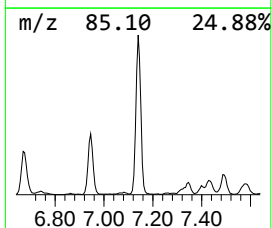
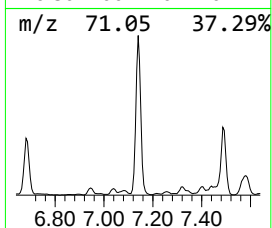
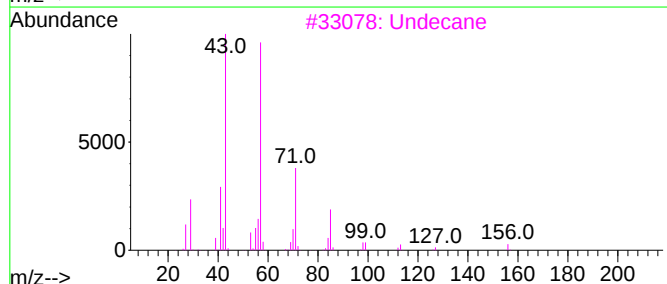
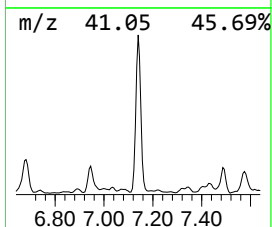
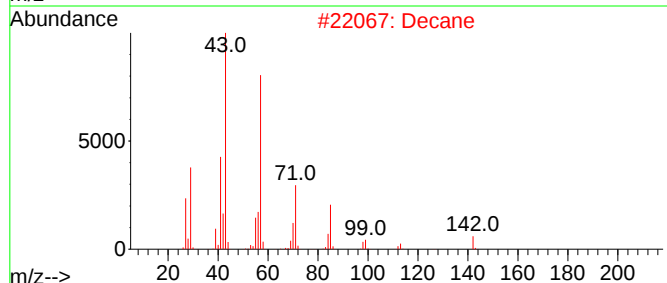
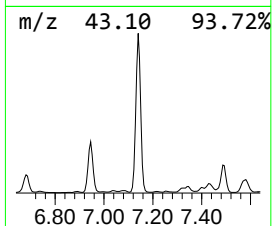
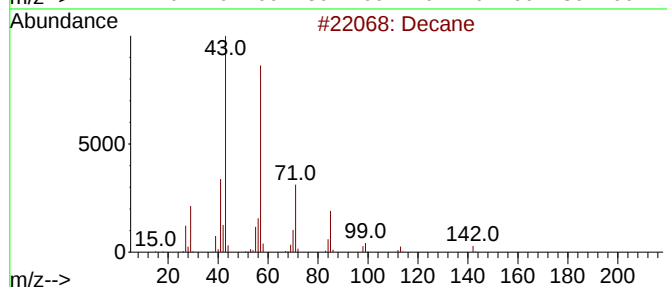
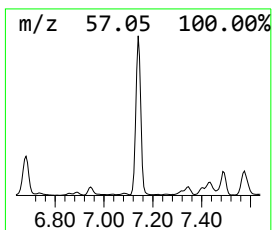
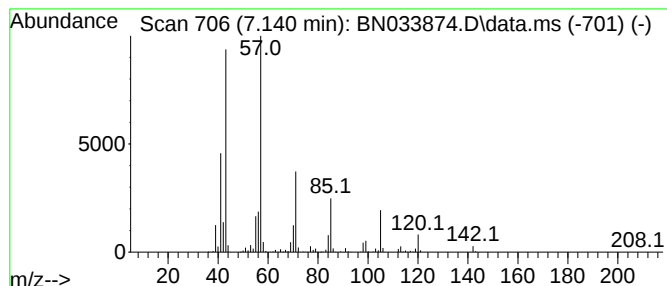
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 1

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 7.140 | 42.27 ng/ul | 2717650 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of       | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----------|---|--------------|-----|---------|-------------|------|
| 1    | Decane   |   |              | 142 | C10H22  | 000124-18-5 | 94   |
| 2    | Decane   |   |              | 142 | C10H22  | 000124-18-5 | 72   |
| 3    | Undecane |   |              | 156 | C11H24  | 001120-21-4 | 72   |
| 4    | Nonane   |   |              | 128 | C9H20   | 000111-84-2 | 72   |
| 5    | Nonane   |   |              | 128 | C9H20   | 000111-84-2 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

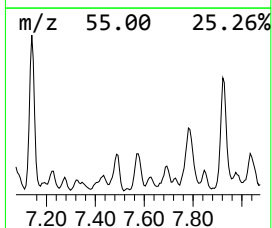
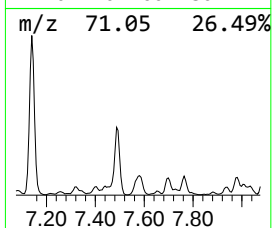
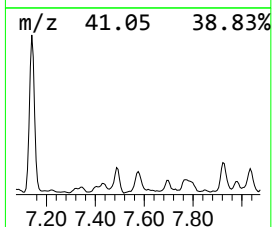
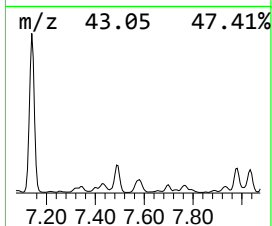
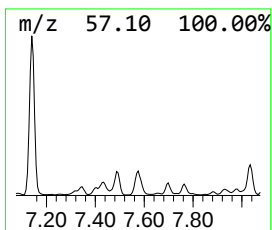
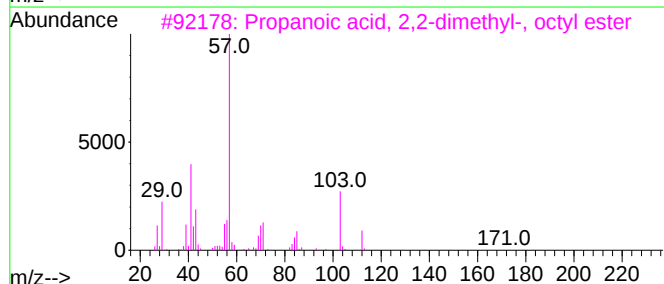
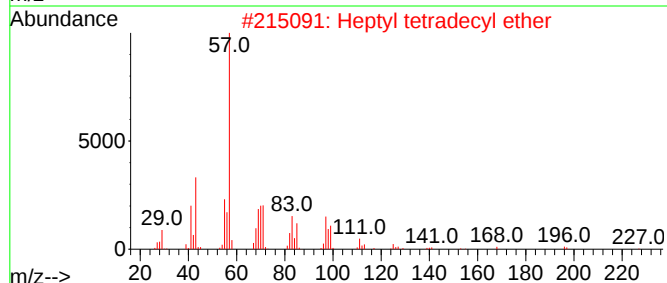
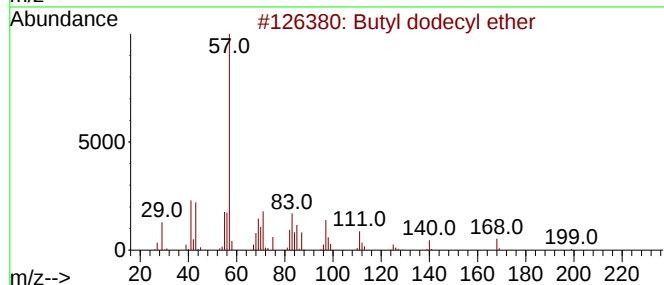
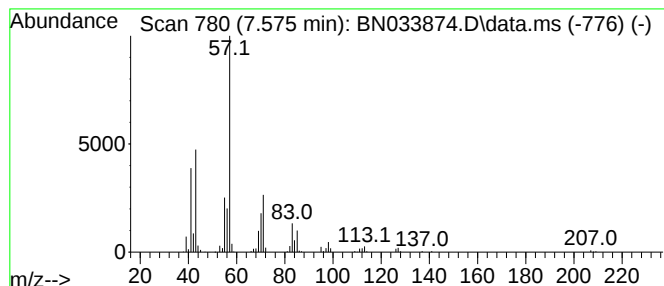
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 Butyl dodecyl ether Concentration Rank 26

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 7.575 | 5.16 ng/ul | 331775 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Butyl dodecyl ether                 | 242 | C16H34O  | 1000406-40-3 | 72   |
| 2    |    |   | Heptyl tetradecyl ether             | 312 | C21H44O  | 1000406-39-3 | 64   |
| 3    |    |   | Propanoic acid, 2,2-dimethyl-, o... | 214 | C13H26O2 | 027751-88-8  | 59   |
| 4    |    |   | 1-Octanol, 2-butyl-                 | 186 | C12H26O  | 003913-02-8  | 59   |
| 5    |    |   | 1-Hexanol, 2-ethyl-                 | 130 | C8H18O   | 000104-76-7  | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

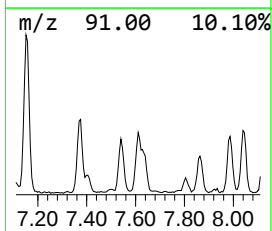
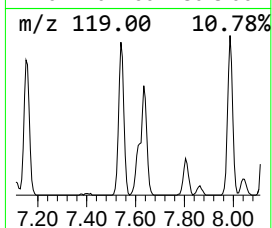
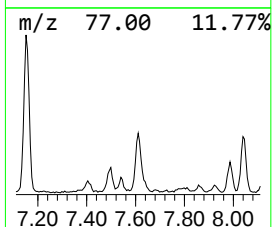
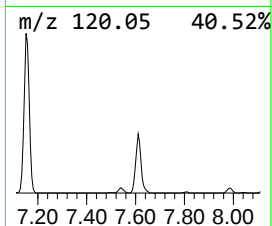
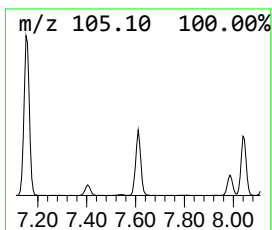
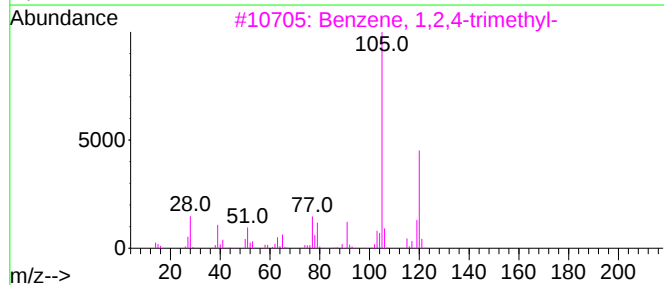
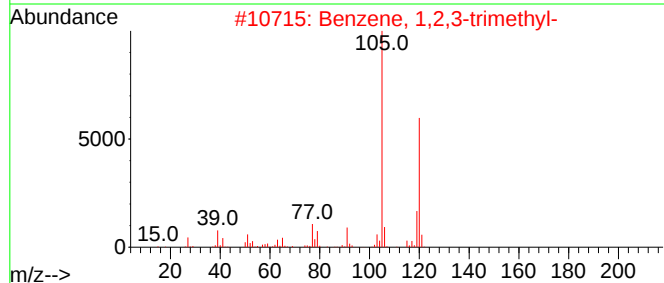
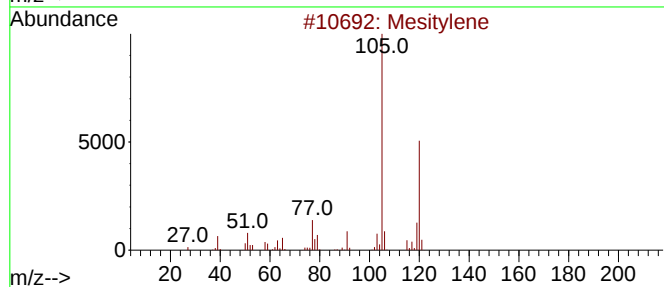
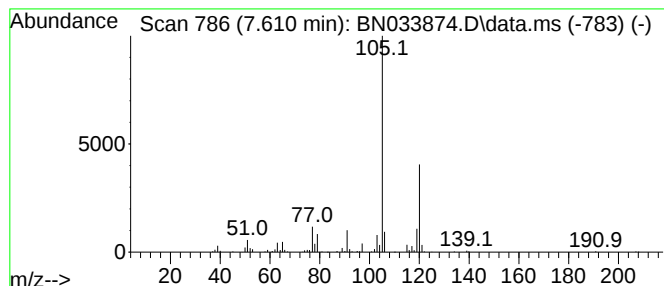
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Quant Title : SVOA CALIBRATION

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

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Peak Number 13 Mesitylene Concentration Rank 15

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 7.610 | 6.10 ng/ul | 391935 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Mesitylene                | 120 | C9H12   | 000108-67-8 | 95   |
| 2    |    |   | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 95   |
| 3    |    |   | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 94   |
| 4    |    |   | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 94   |
| 5    |    |   | Mesitylene                | 120 | C9H12   | 000108-67-8 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

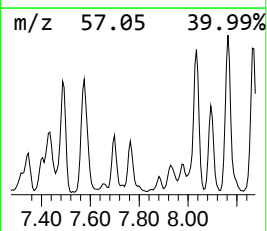
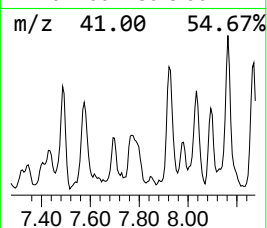
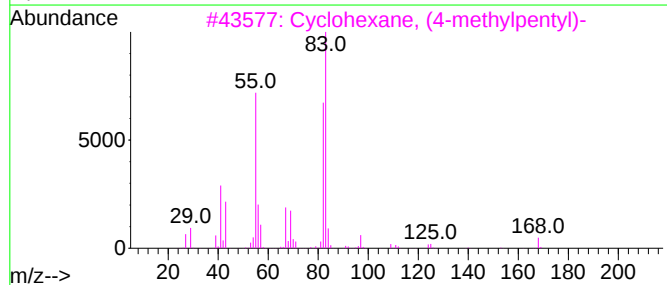
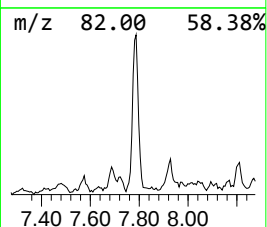
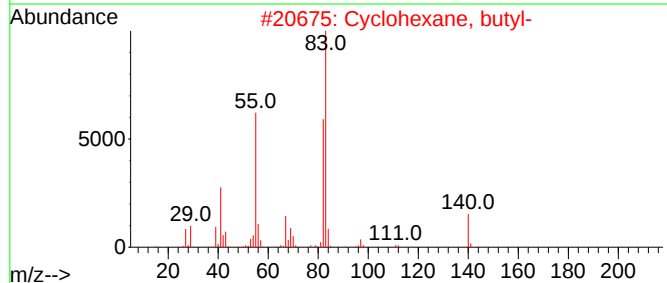
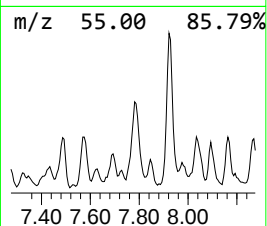
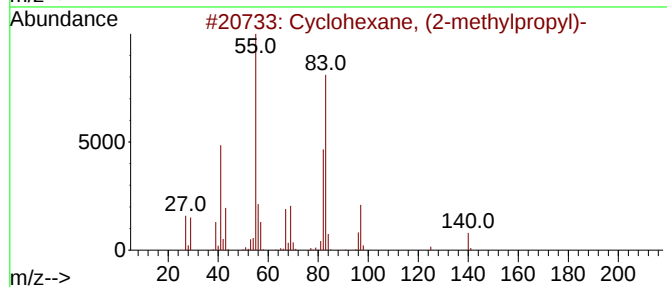
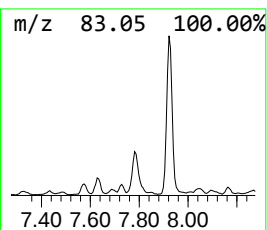
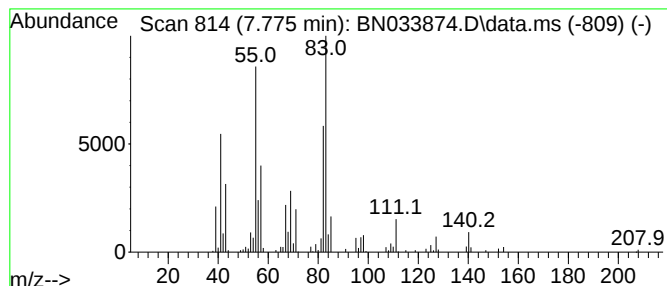
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Quant Title : SVOA CALIBRATION

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

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Peak Number 14 (DEL) Alkane: Cyclic7.775 Concentration Rank 17

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 7.775 | 6.02 ng/ul | 387163 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|------|--------------------------------|-----|---------|-------------|------|
| 1    |      | Cyclohexane, (2-methylpropyl)- | 140 | C10H20  | 001678-98-4 | 62   |
| 2    |      | Cyclohexane, butyl-            | 140 | C10H20  | 001678-93-9 | 59   |
| 3    |      | Cyclohexane, (4-methylpentyl)- | 168 | C12H24  | 061142-20-9 | 59   |
| 4    |      | Cyclohexane, pentyl-           | 154 | C11H22  | 004292-92-6 | 53   |
| 5    |      | Cyclohexane, pentyl-           | 154 | C11H22  | 004292-92-6 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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Quant Title : SVOA CALIBRATION

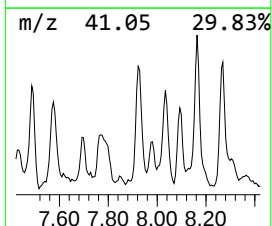
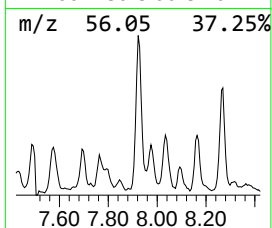
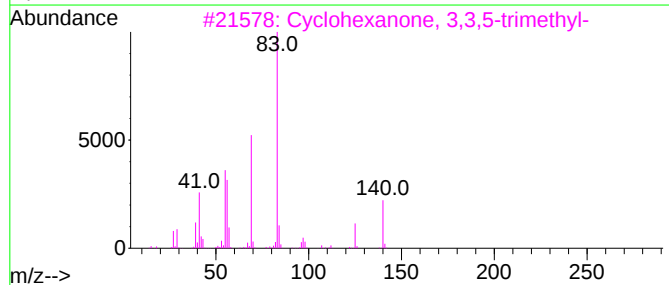
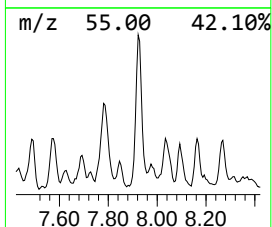
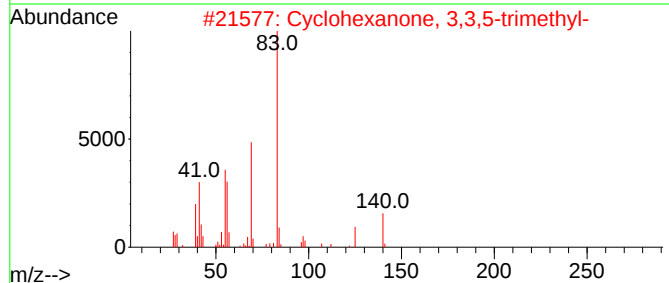
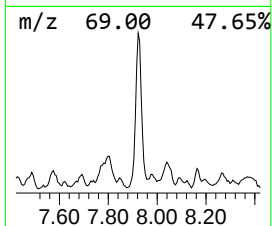
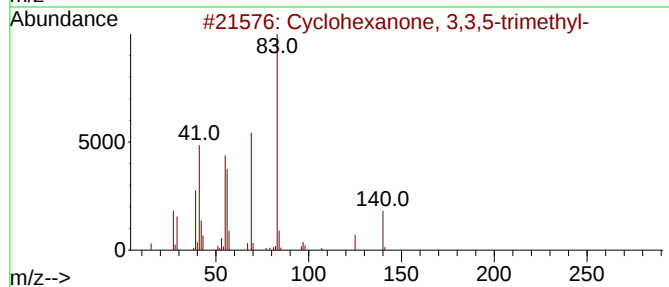
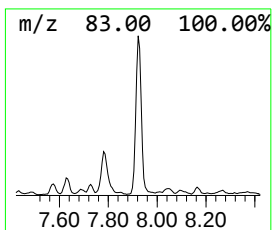
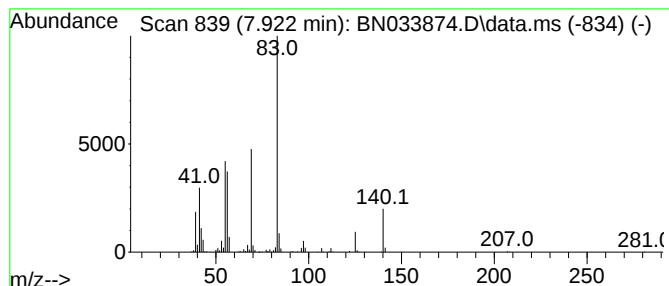
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TIC Integration Parameters: LSCINT.P

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Peak Number 15 Cyclohexanone, 3,3,5-trimet... Concentration Rank 5

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 7.922 | 9.31 ng/ul | 598564 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Cyclohexanone, 3,3,5-trimethyl-     | 140 | C9H16O  | 000873-94-9 | 95   |
| 2    |      | Cyclohexanone, 3,3,5-trimethyl-     | 140 | C9H16O  | 000873-94-9 | 94   |
| 3    |      | Cyclohexanone, 3,3,5-trimethyl-     | 140 | C9H16O  | 000873-94-9 | 93   |
| 4    |      | Cyclohexanone, 3,3,5-trimethyl-     | 140 | C9H16O  | 000873-94-9 | 91   |
| 5    |      | 1,3-Cyclohexanedione, 5,5-dimethyl- | 140 | C8H12O2 | 000126-81-8 | 56   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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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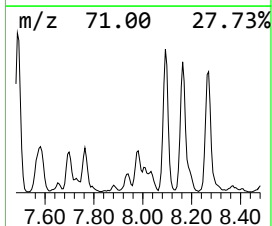
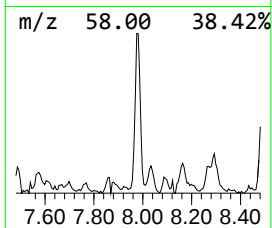
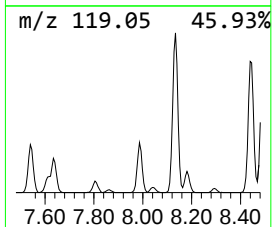
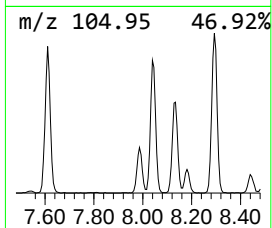
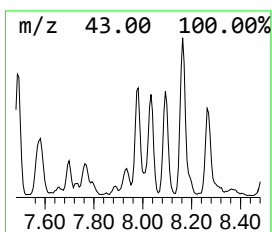
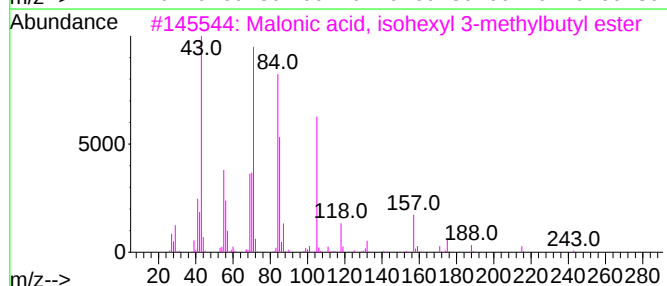
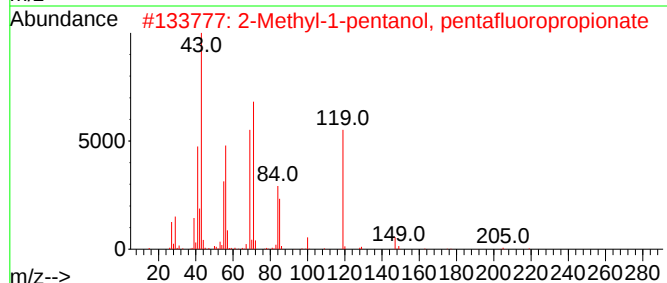
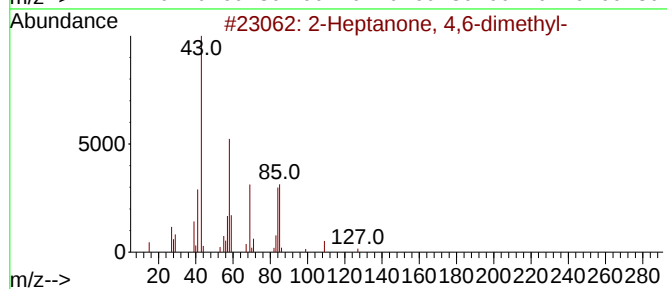
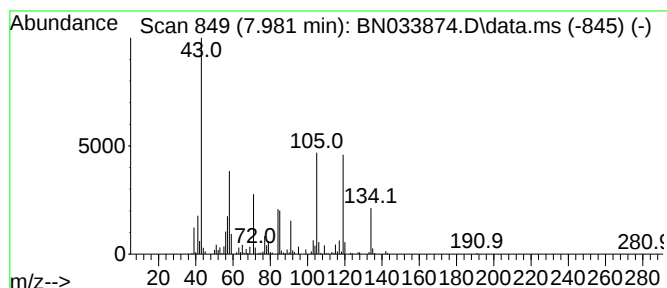
Peak Number 16 unknown-01

Concentration Rank 13

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 7.981 | 6.55 ng/ul | 420930 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of                                   | 5 | Tentative ID | MW         | MolForm | CAS#         | Qual |
|------|--------------------------------------|---|--------------|------------|---------|--------------|------|
| 1    | 2-Heptanone, 4,6-dimethyl-           |   | 142          | C9H18O     |         | 019549-80-5  | 38   |
| 2    | 2-Methyl-1-pentanol, pentafluoro...  |   | 248          | C9H13F5O2  |         | 1000352-35-9 | 16   |
| 3    | Malonic acid, isohexyl 3-methylb...  |   | 258          | C14H26O4   |         | 1010348-98-0 | 14   |
| 4    | Dichloroacetic acid, 4-methylpen...  |   | 212          | C8H14Cl2O2 |         | 1000282-43-7 | 14   |
| 5    | Thioglycolic acid, S-isopropyllox... |   | 220          | C9H16O4S   |         | 1000476-02-2 | 12   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

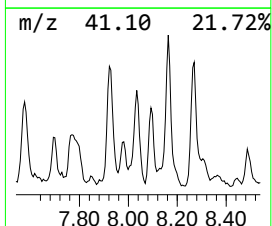
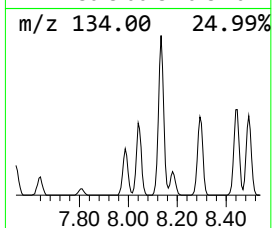
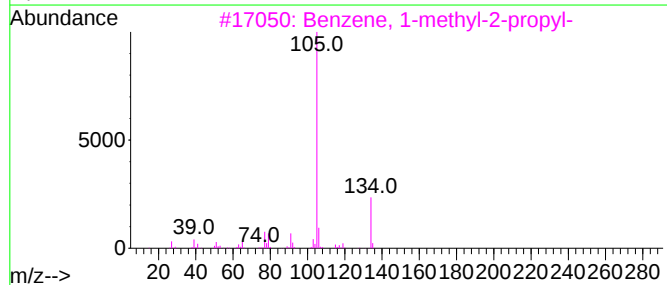
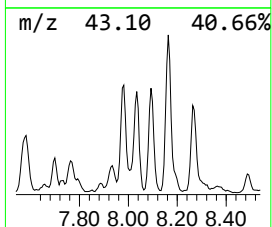
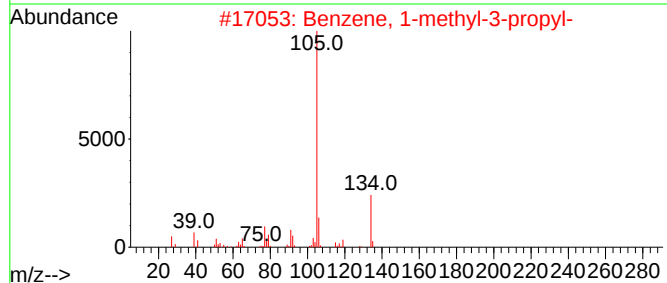
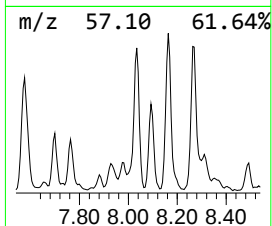
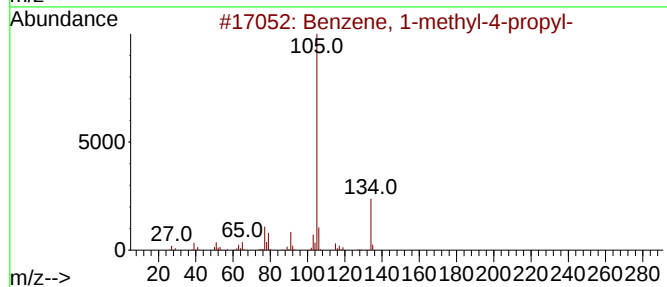
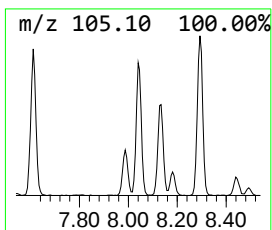
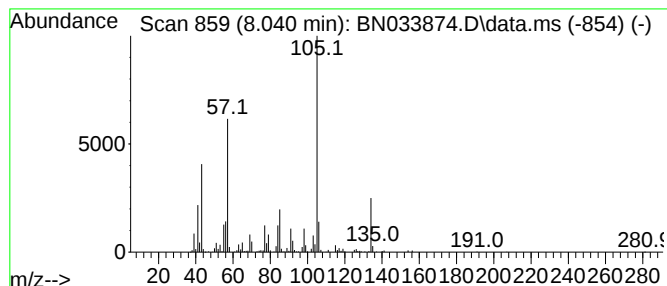
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 17 Benzene, 1-methyl-4-propyl- Concentration Rank 4

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.040 | 9.37 ng/ul | 602339 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 90   |
| 2    |    |   | Benzene, 1-methyl-3-propyl- | 134 | C10H14  | 001074-43-7 | 83   |
| 3    |    |   | Benzene, 1-methyl-2-propyl- | 134 | C10H14  | 001074-17-5 | 64   |
| 4    |    |   | Benzene, 1-methyl-3-propyl- | 134 | C10H14  | 001074-43-7 | 64   |
| 5    |    |   | Benzene, 1-methyl-3-propyl- | 134 | C10H14  | 001074-43-7 | 60   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

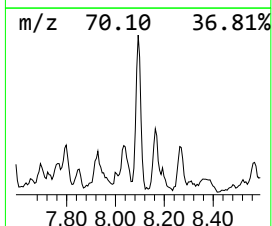
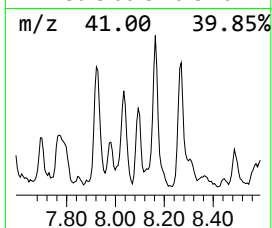
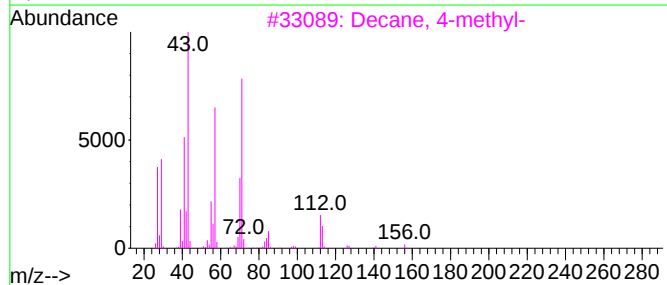
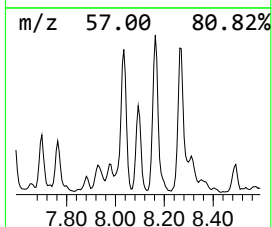
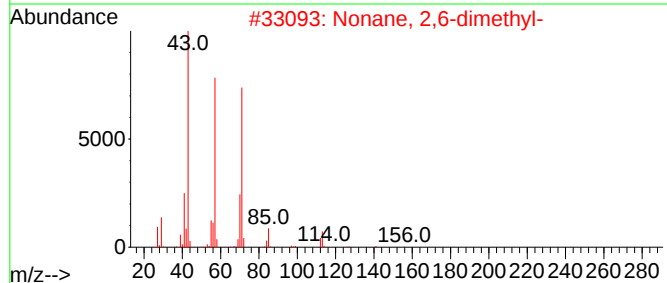
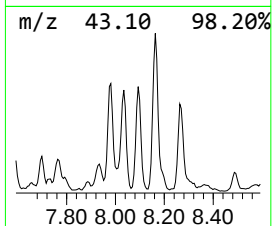
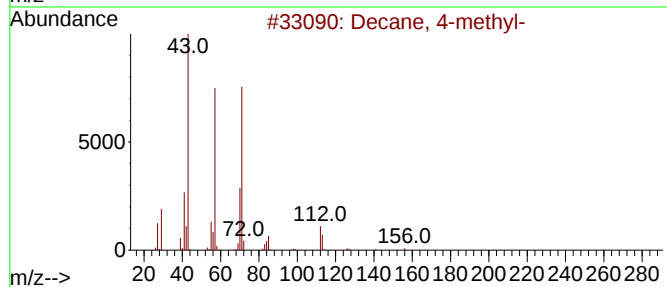
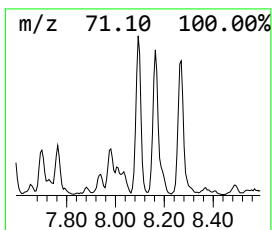
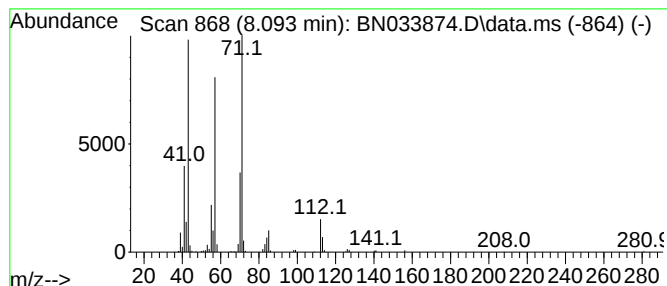
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TIC Integration Parameters: LSCINT.P

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Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 28

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.093 | 4.60 ng/ul | 295505 | 1,4-Dichlorobenzene-d4 | 7.499 |

| 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  | 90   |
| 3    |      | Decane, 4-methyl-                   | 156 | C11H24     | 002847-72-5  | 83   |
| 4    |      | 4-Methyl-3-heptanol, pentafluoro... | 276 | C11H17F5O2 | 1000365-51-7 | 78   |
| 5    |      | Heptane, 2,5,5-trimethyl-           | 142 | C10H22     | 001189-99-7  | 78   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

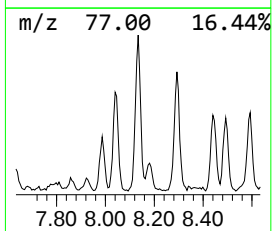
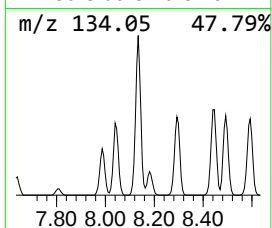
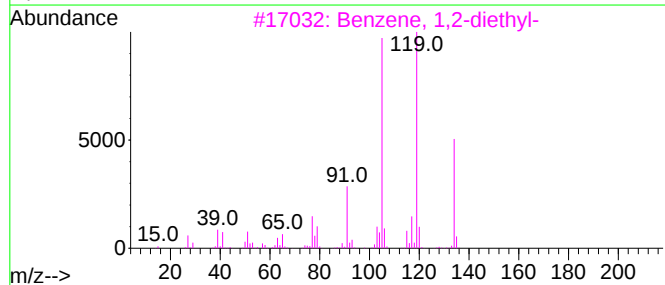
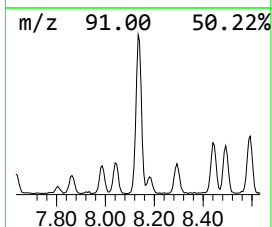
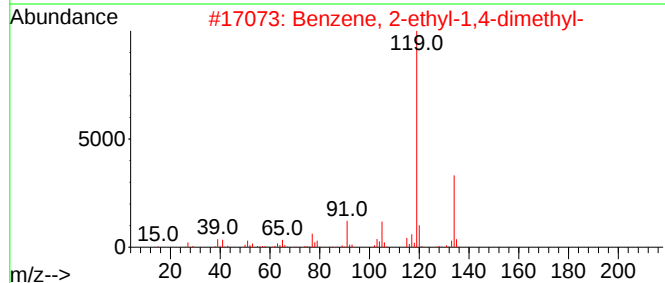
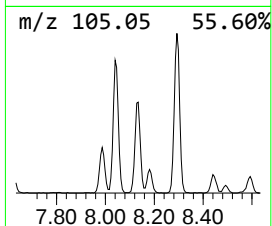
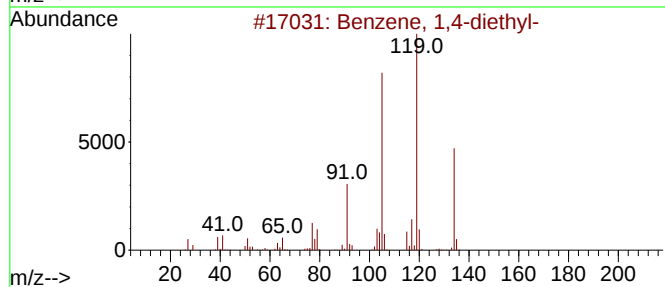
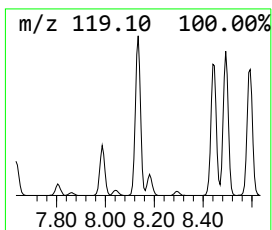
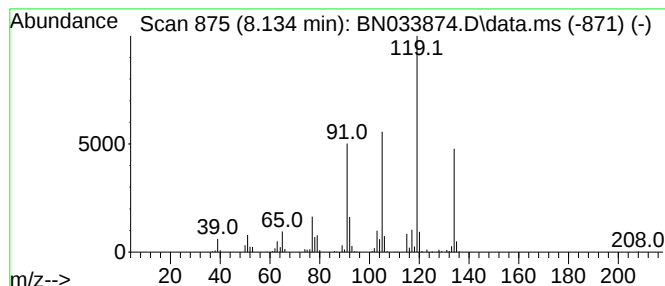
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Quant Title : SVOA CALIBRATION

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

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Peak Number 19 Benzene, 1,4-diethyl- Concentration Rank 7

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.134 | 8.95 ng/ul | 575647 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,4-diethyl-          | 134 | C10H14  | 000105-05-5 | 96   |
| 2    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 94   |
| 3    |    |   | Benzene, 1,2-diethyl-          | 134 | C10H14  | 000135-01-3 | 94   |
| 4    |    |   | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 91   |
| 5    |    |   | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

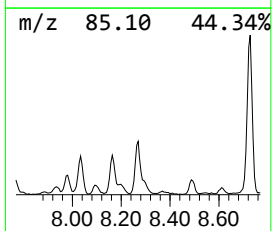
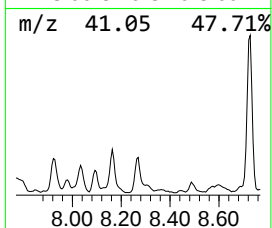
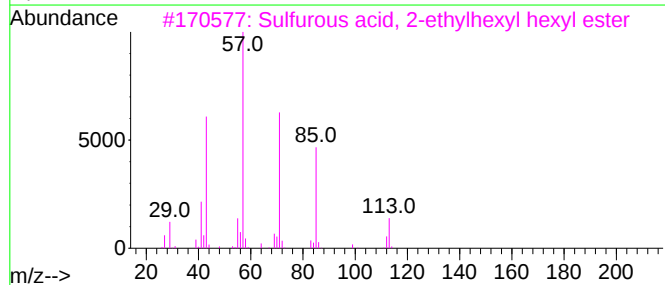
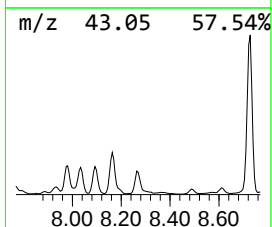
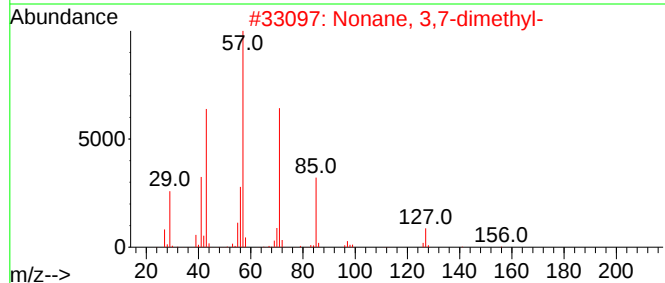
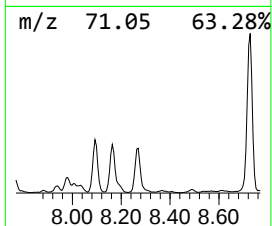
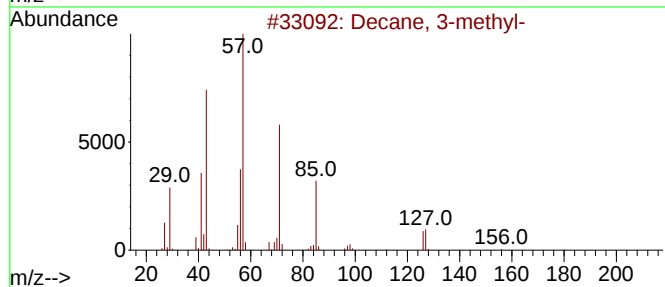
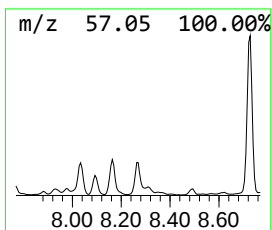
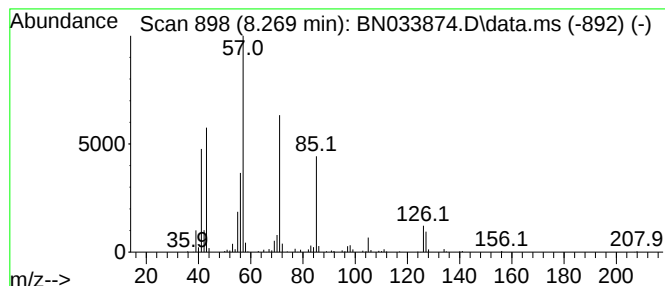
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.269 | 7.14 ng/ul | 459184 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Decane, 3-methyl-                   | 156 | C11H24    | 013151-34-3  | 93   |
| 2    |    |   | Nonane, 3,7-dimethyl-               | 156 | C11H24    | 017302-32-8  | 64   |
| 3    |    |   | Sulfurous acid, 2-ethylhexyl hex... | 278 | C14H30O3S | 1000309-20-2 | 53   |
| 4    |    |   | Dodecane, 1-iodo-                   | 296 | C12H25I   | 004292-19-7  | 53   |
| 5    |    |   | Decane, 2,3,8-trimethyl-            | 184 | C13H28    | 062238-14-6  | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

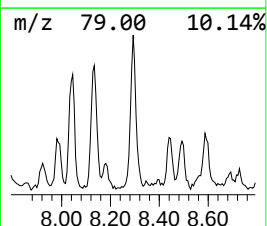
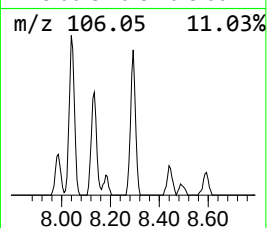
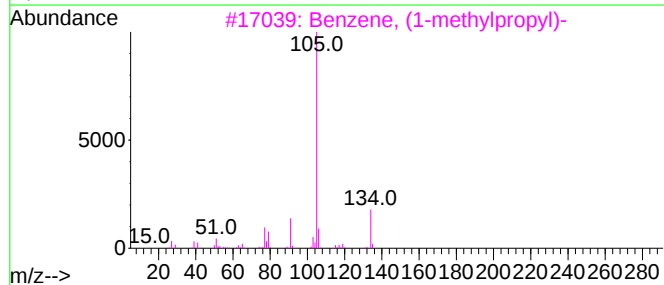
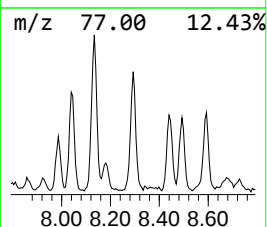
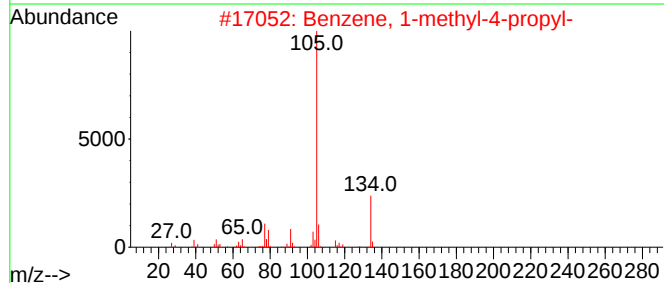
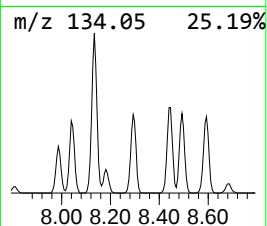
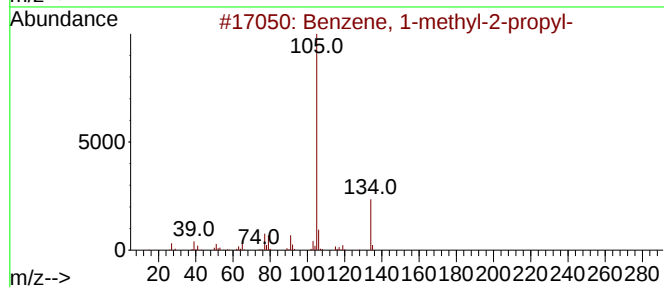
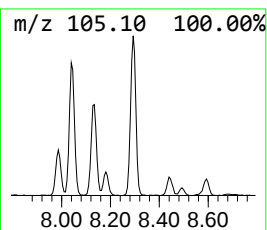
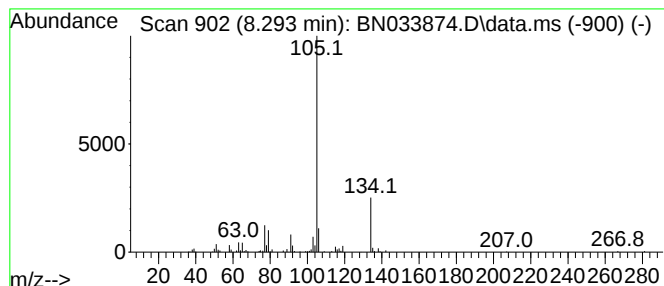
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Quant Title : SVOA CALIBRATION

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

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Peak Number 21 Benzene, 1-methyl-2-propyl- Concentration Rank 25

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.293 | 5.22 ng/ul | 335687 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-methyl-2-propyl- | 134 | C10H14  | 001074-17-5 | 94   |
| 2    |    |   | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 93   |
| 3    |    |   | Benzene, (1-methylpropyl)-  | 134 | C10H14  | 000135-98-8 | 91   |
| 4    |    |   | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 91   |
| 5    |    |   | Benzene, 1-methyl-3-propyl- | 134 | C10H14  | 001074-43-7 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

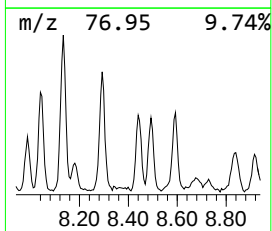
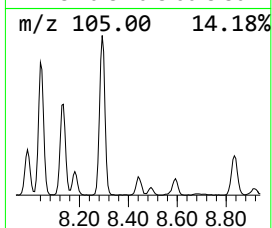
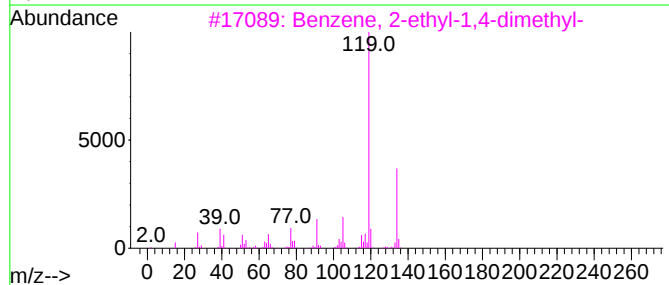
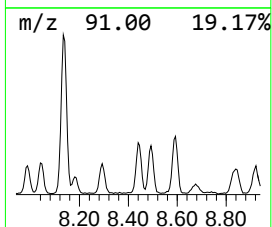
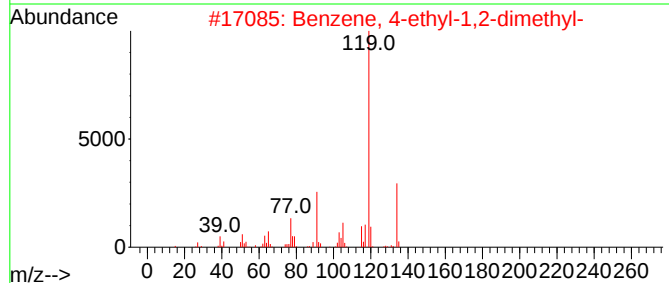
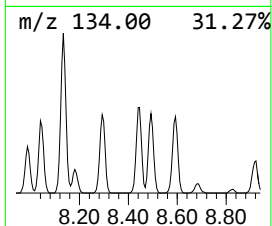
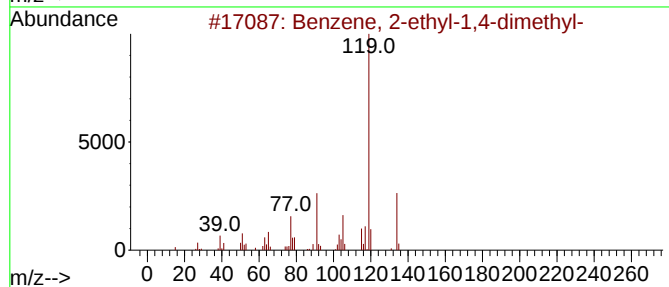
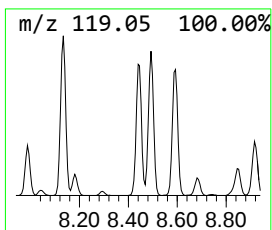
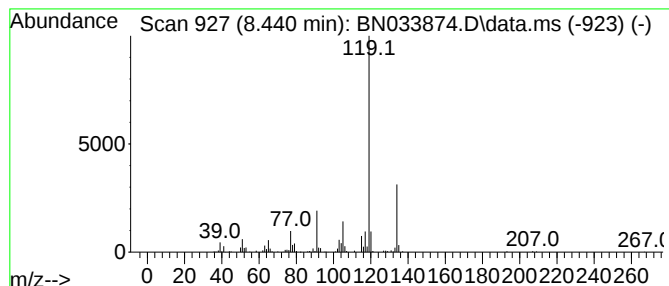
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TIC Integration Parameters: LSCINT.P

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Peak Number 22 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 31

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.440 | 4.08 ng/ul | 262320 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 96   |
| 2    |    |   | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 96   |
| 3    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 95   |
| 4    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 95   |
| 5    |    |   | p-Cymene                       | 134 | C10H14  | 000099-87-6 | 95   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

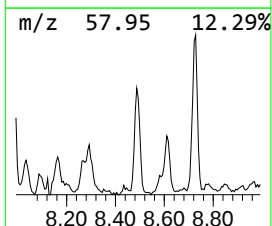
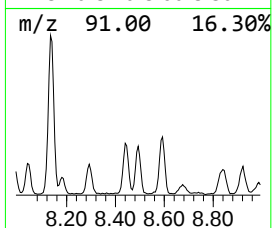
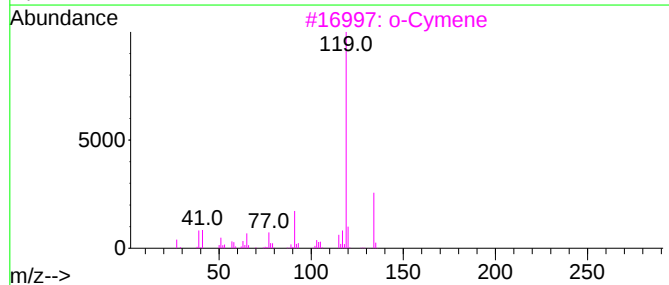
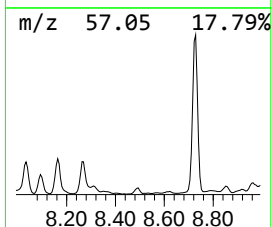
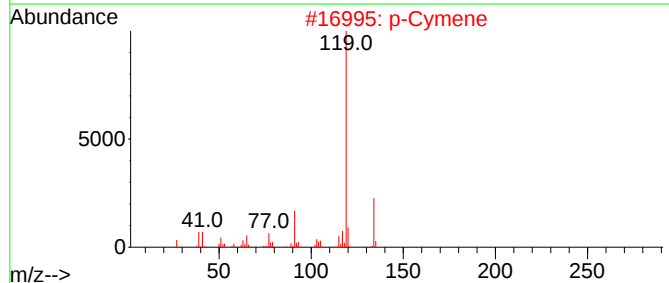
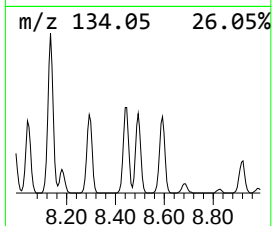
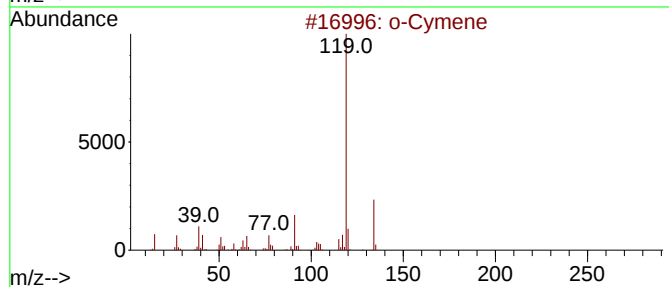
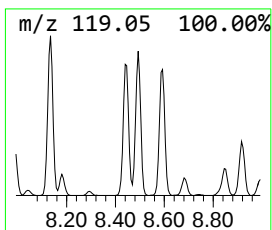
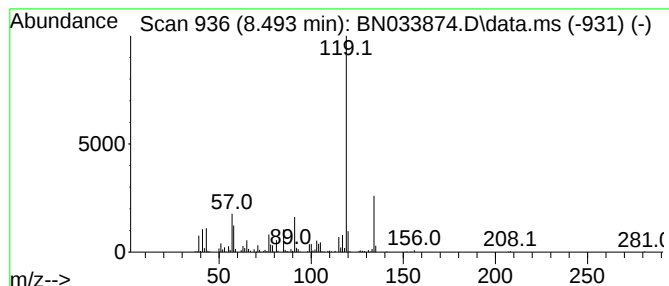
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Quant Title : SVOA CALIBRATION

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

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Peak Number 23 o-Cymene Concentration Rank 18

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.493 | 5.73 ng/ul | 368257 | 1,4-Dichlorobenzene-d4 | 7.499 |

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



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

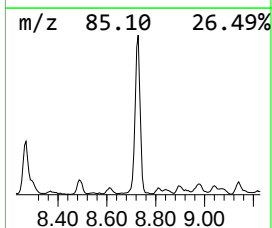
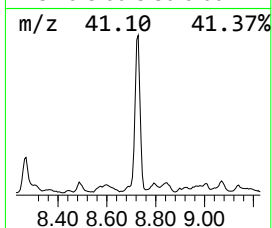
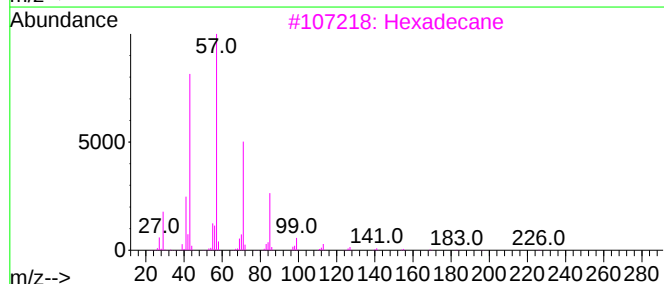
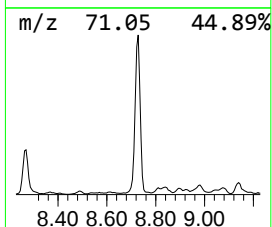
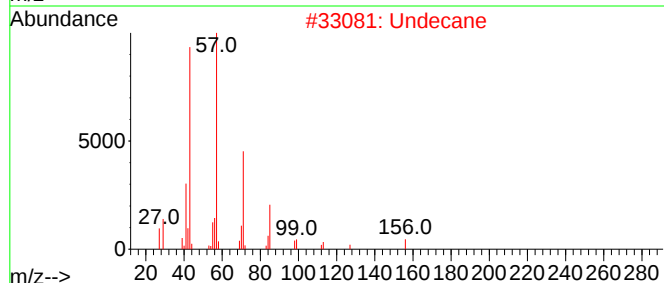
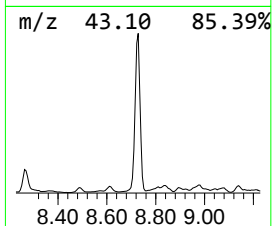
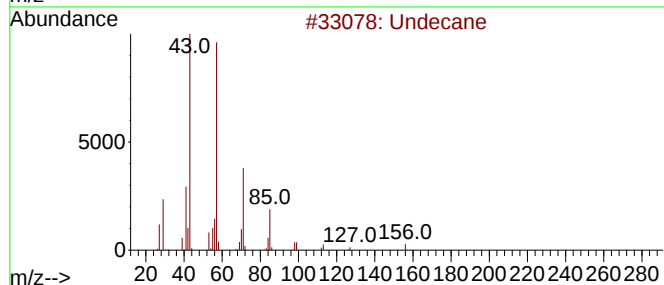
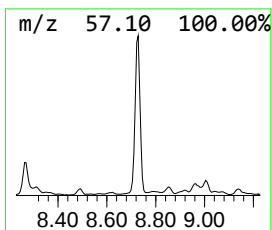
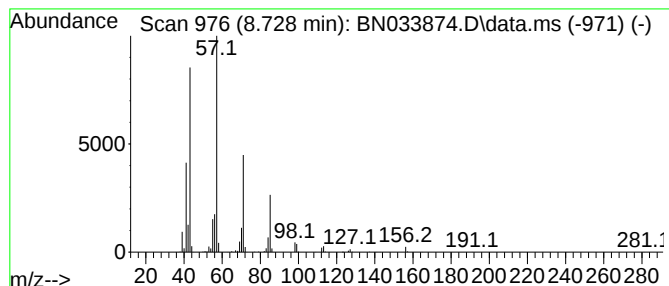
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 2

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 8.728 | 28.77 ng/ul | 1849680 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of         | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------------|---|--------------|-----|---------|-------------|------|
| 1    | Undecane   |   |              | 156 | C11H24  | 001120-21-4 | 93   |
| 2    | Undecane   |   |              | 156 | C11H24  | 001120-21-4 | 87   |
| 3    | Hexadecane |   |              | 226 | C16H34  | 000544-76-3 | 86   |
| 4    | Tridecane  |   |              | 184 | C13H28  | 000629-50-5 | 83   |
| 5    | Tridecane  |   |              | 184 | C13H28  | 000629-50-5 | 78   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

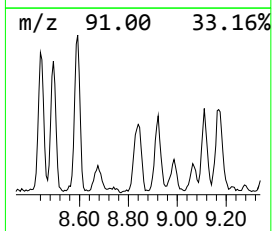
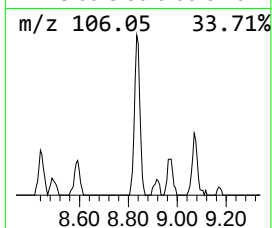
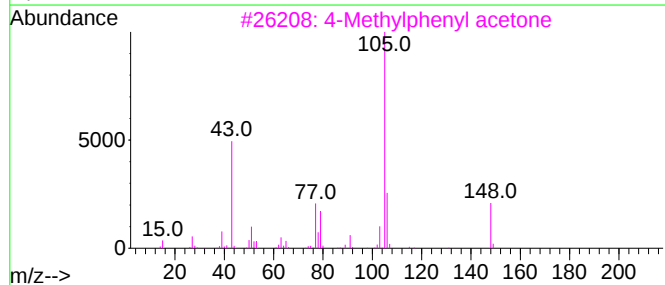
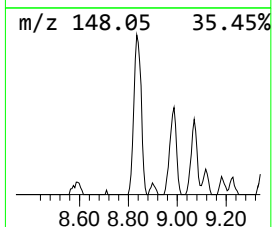
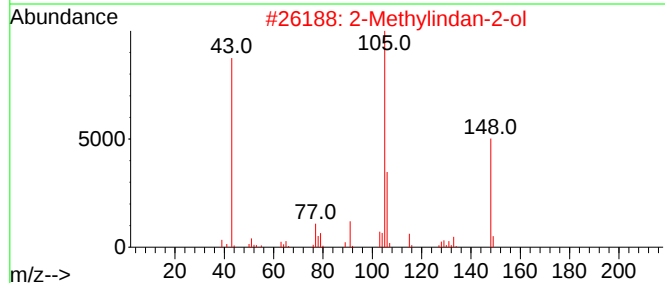
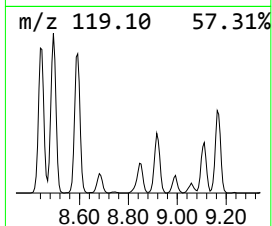
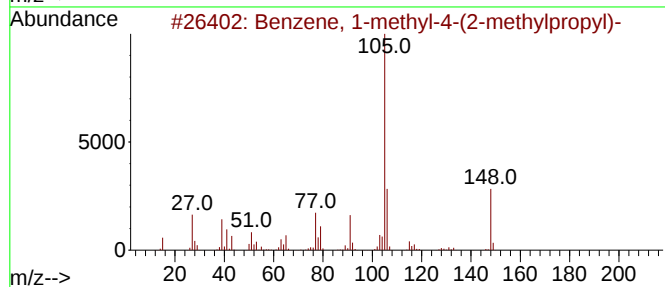
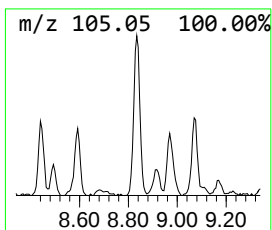
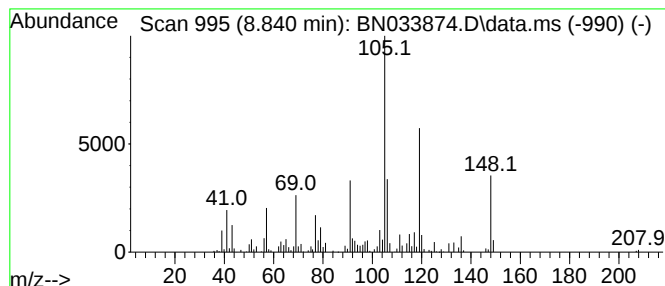
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Peak Number 25 Benzene, 1-methyl-4-(2-meth... Concentration Rank 19

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.840 | 5.71 ng/ul | 366903 | 1,4-Dichlorobenzene-d4 | 7.499 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-methyl-4-(2-methylpro... | 148 | C11H16  | 005161-04-6 | 55   |
| 2    |    |   | 2-Methylindan-2-ol                  | 148 | C10H12O | 033223-84-6 | 55   |
| 3    |    |   | 4-Methylphenyl acetone              | 148 | C10H12O | 002096-86-8 | 46   |
| 4    |    |   | Benzeneethanol, 2-methyl-           | 136 | C9H12O  | 019819-98-8 | 45   |
| 5    |    |   | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

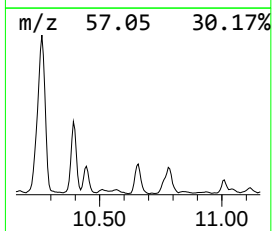
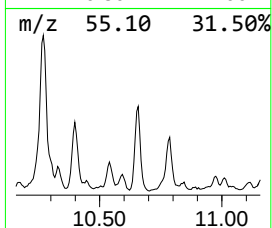
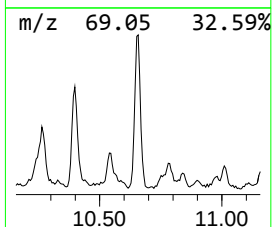
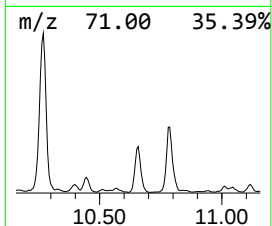
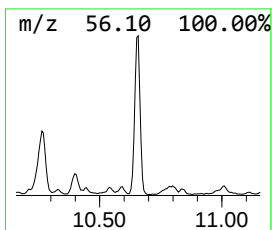
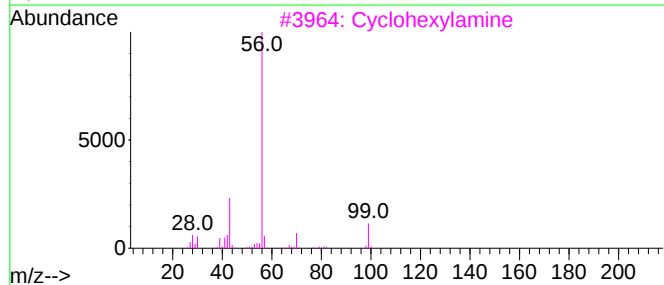
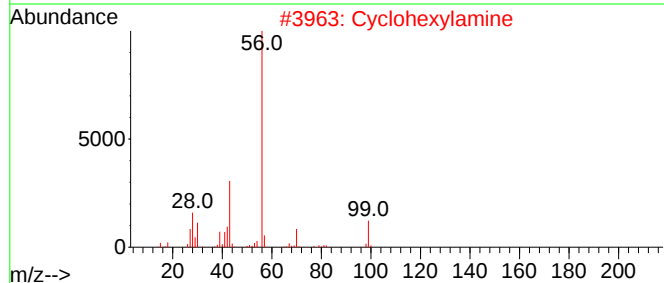
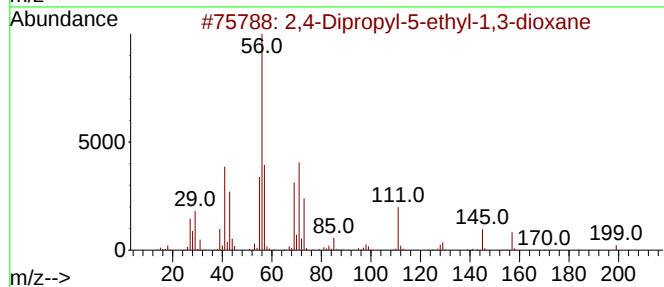
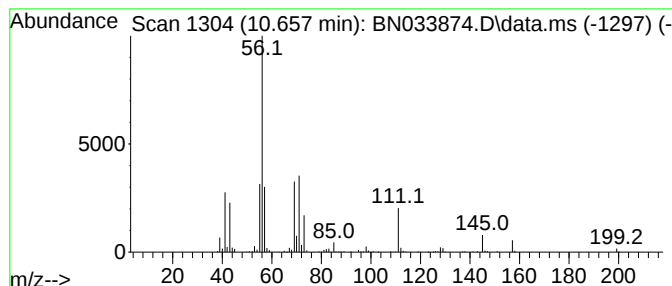
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Quant Title : SVOA CALIBRATION

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

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Peak Number 26 2,4-Dipropyl-5-ethyl-1,3-di... Concentration Rank 40

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 10.657 | 2.86 ng/ul | 618224 | Naphthalene-d8   | 10.257 |

| Hit# | of 5 | Tentative ID                      | MW  | MolForm  | CAS#        | Qual |
|------|------|-----------------------------------|-----|----------|-------------|------|
| 1    |      | 2,4-Dipropyl-5-ethyl-1,3-dioxane  | 200 | C12H24O2 | 006413-83-8 | 50   |
| 2    |      | Cyclohexylamine                   | 99  | C6H13N   | 000108-91-8 | 38   |
| 3    |      | Cyclohexylamine                   | 99  | C6H13N   | 000108-91-8 | 38   |
| 4    |      | Cyclobutane, 1,2,3,4-tetramethyl- | 112 | C8H16    | 069531-57-3 | 32   |
| 5    |      | 2-Methyl-1-nonene                 | 140 | C10H20   | 002980-71-4 | 32   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

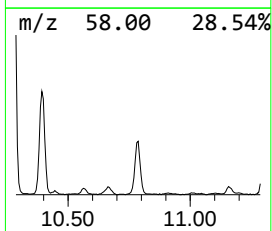
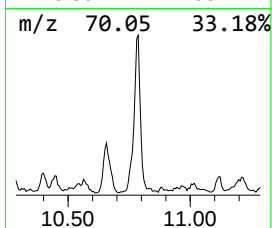
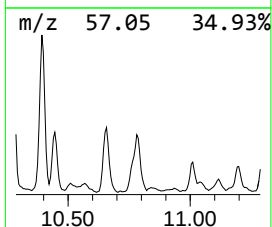
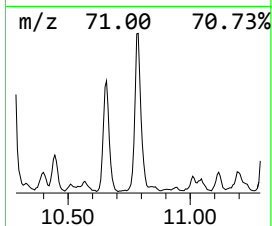
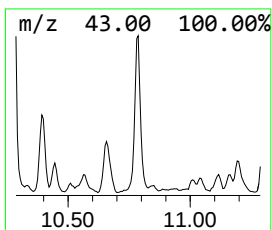
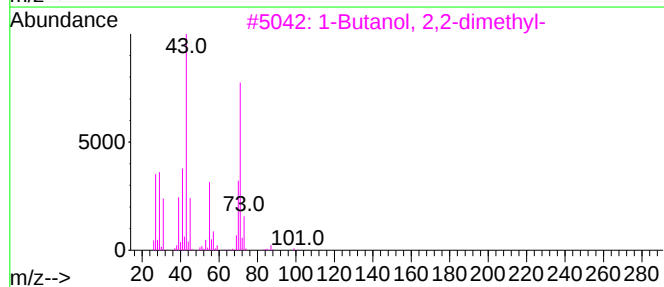
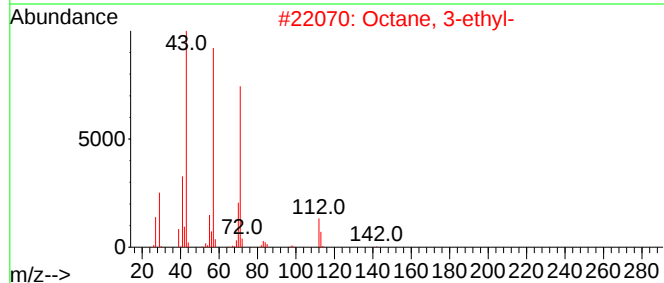
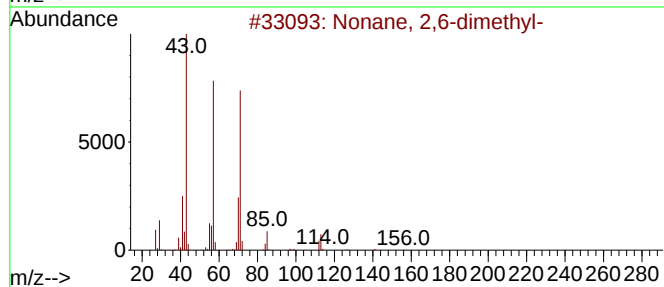
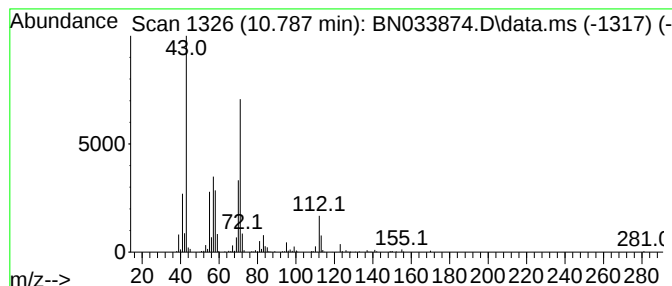
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TIC Integration Parameters: LSCINT.P

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Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 46

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 10.787 | 2.59 ng/ul | 560271 | Naphthalene-d8   | 10.257 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Nonane, 2,6-dimethyl-               | 156 | C11H24   | 017302-28-2  | 53   |
| 2    |    |   | Octane, 3-ethyl-                    | 142 | C10H22   | 005881-17-4  | 52   |
| 3    |    |   | 1-Butanol, 2,2-dimethyl-            | 102 | C6H14O   | 001185-33-7  | 50   |
| 4    |    |   | 2,4,4-Trimethyl-1-hexene            | 126 | C9H18    | 051174-12-0  | 49   |
| 5    |    |   | Oxalic acid, 2-ethylhexyl pentyl... | 272 | C15H28O4 | 1000309-38-7 | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

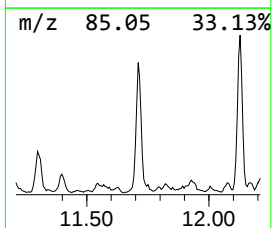
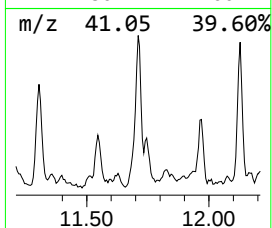
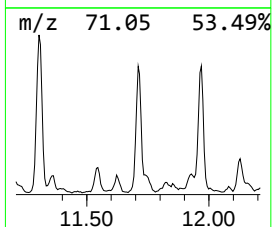
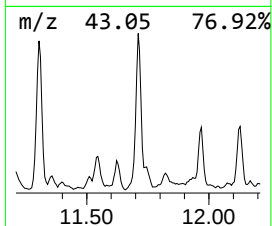
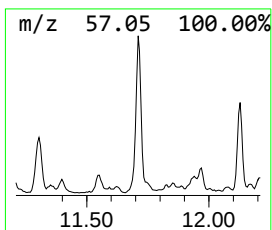
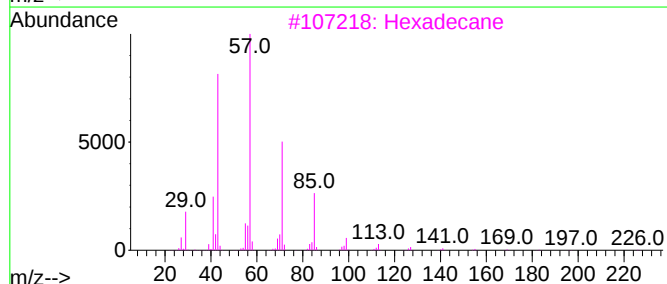
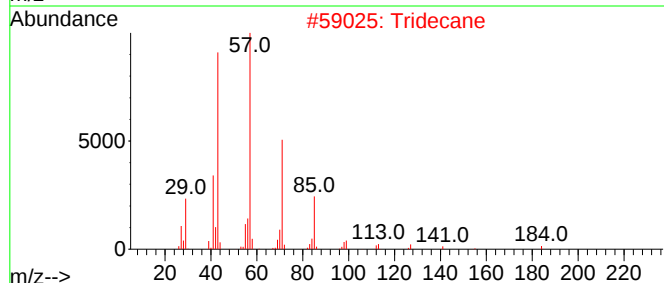
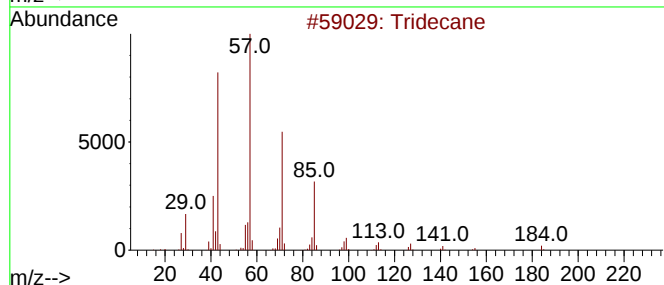
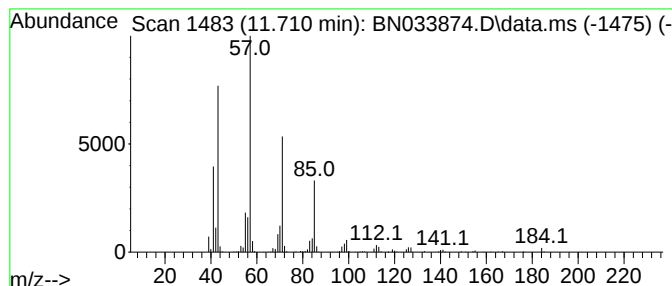
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TIC Integration Parameters: LSCINT.P

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Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 44

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 11.710 | 2.78 ng/ul | 599823 | Naphthalene-d8   | 10.257 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Tridecane    | 184 | C13H28  | 000629-50-5 | 94   |
| 2    |    |   | Tridecane    | 184 | C13H28  | 000629-50-5 | 90   |
| 3    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 90   |
| 4    |    |   | Tetradecane  | 198 | C14H30  | 000629-59-4 | 90   |
| 5    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

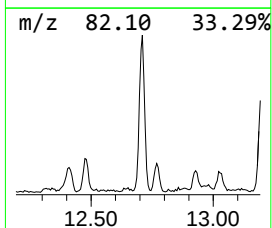
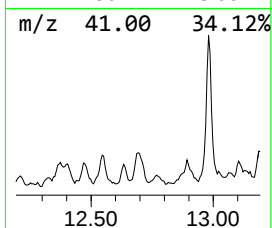
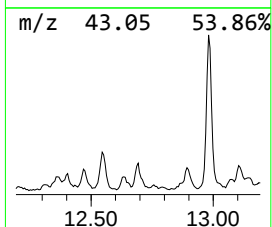
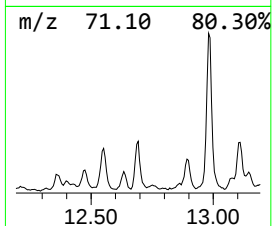
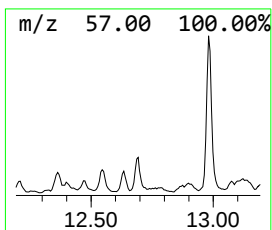
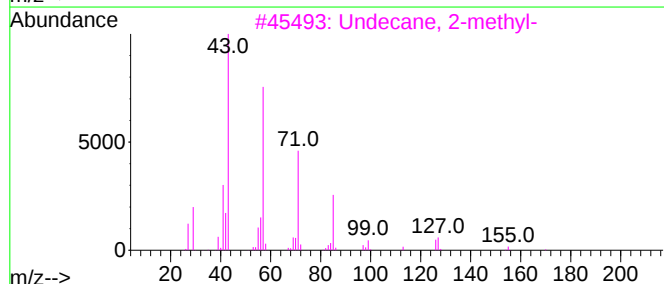
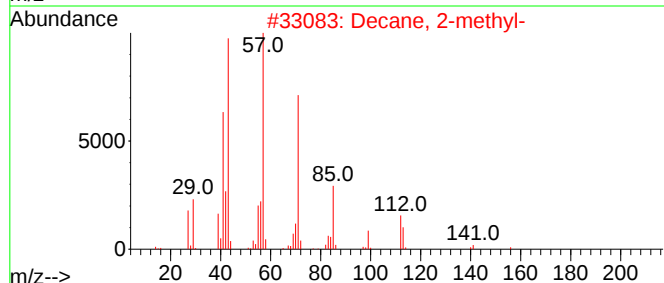
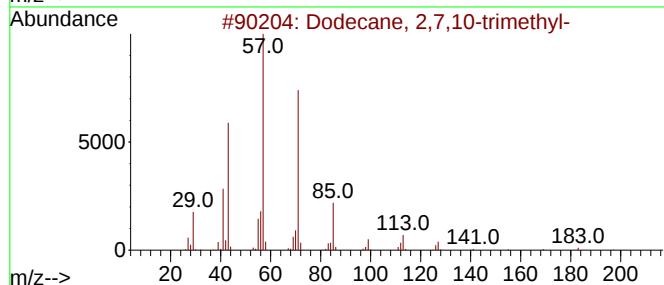
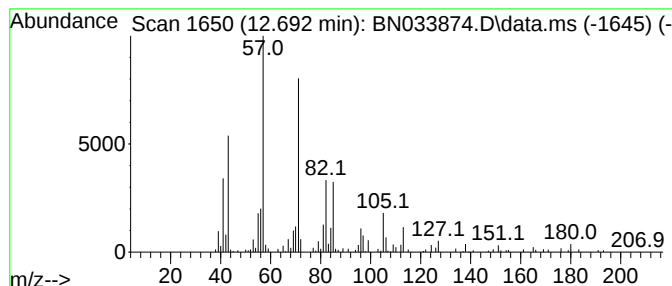
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 47

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 12.692 | 2.56 ng/ul | 283412 | Acenaphthene-d10 | 14.139 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Dodecane, 2,7,10-trimethyl-         | 212 | C15H32   | 074645-98-0  | 64   |
| 2    |    |   | Decane, 2-methyl-                   | 156 | C11H24   | 006975-98-0  | 59   |
| 3    |    |   | Undecane, 2-methyl-                 | 170 | C12H26   | 007045-71-8  | 53   |
| 4    |    |   | Oxalic acid, 2-ethylhexyl nonyl ... | 328 | C19H36O4 | 1000309-39-2 | 53   |
| 5    |    |   | Dodecane, 2,6,10-trimethyl-         | 212 | C15H32   | 003891-98-3  | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

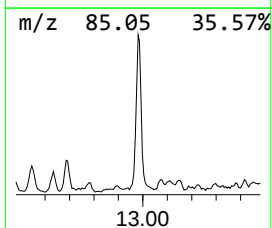
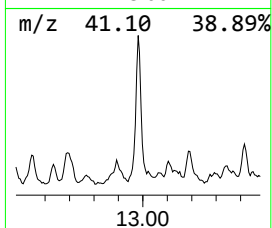
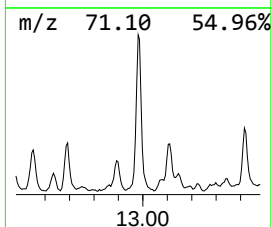
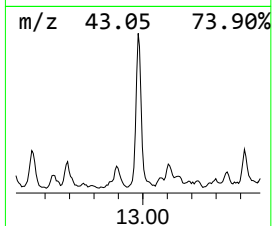
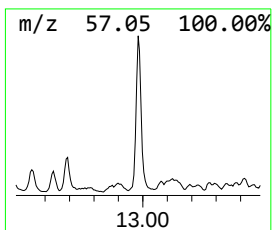
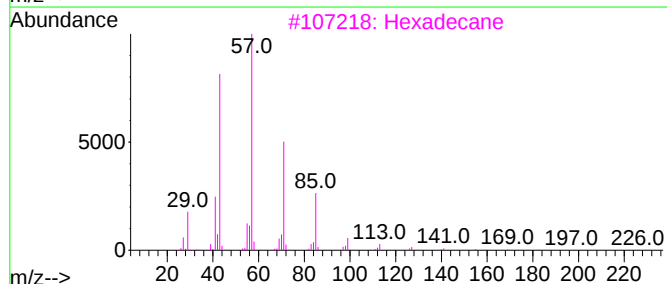
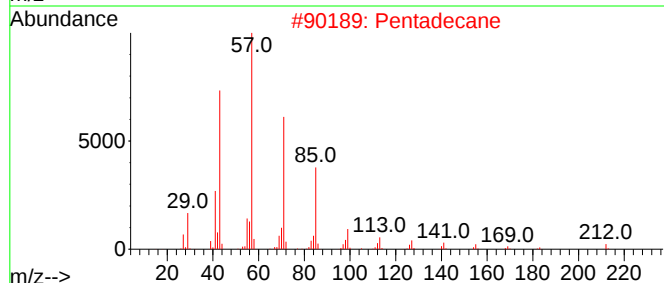
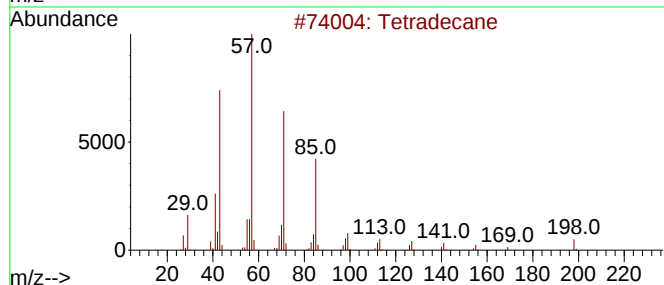
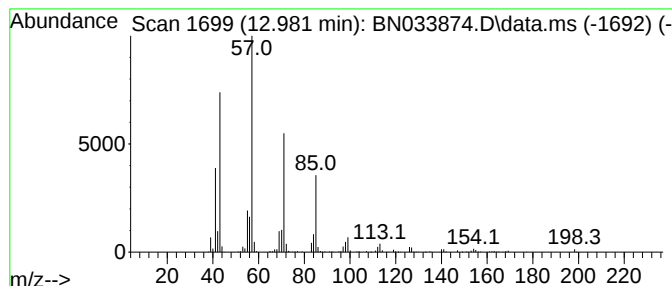
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 12.981 | 6.26 ng/ul | 692706 | Acenaphthene-d10 | 14.139 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Tetradecane  | 198 | C14H30  | 000629-59-4 | 94   |
| 2    |    |   | Pentadecane  | 212 | C15H32  | 000629-62-9 | 90   |
| 3    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 86   |
| 4    |    |   | Tridecane    | 184 | C13H28  | 000629-50-5 | 86   |
| 5    |    |   | Tetradecane  | 198 | C14H30  | 000629-59-4 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

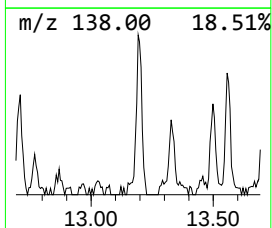
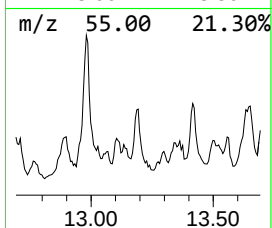
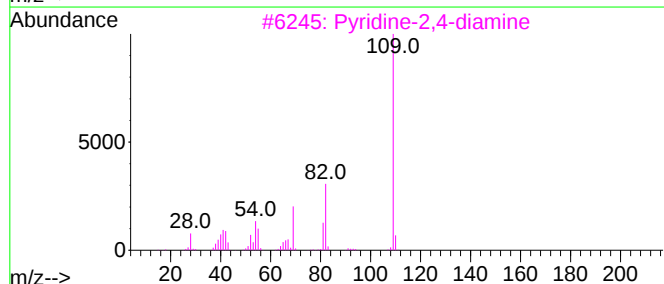
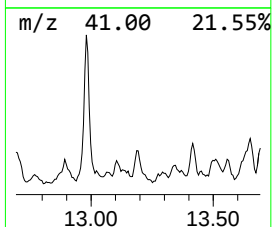
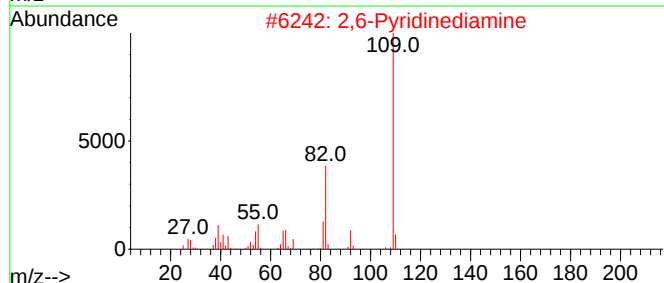
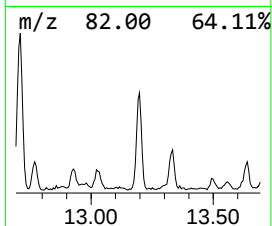
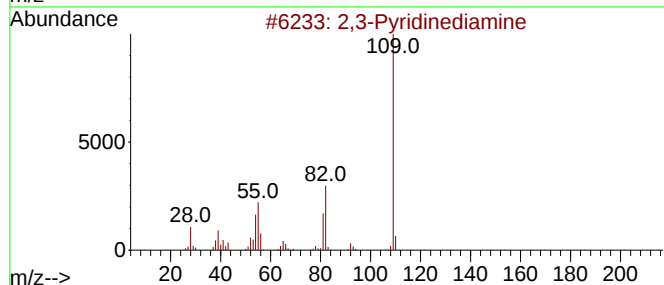
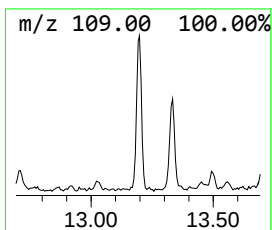
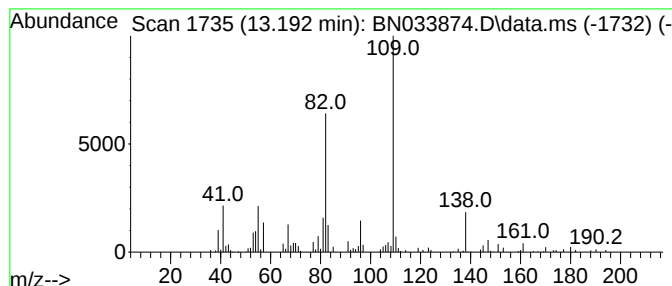
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Quant Title : SVOA CALIBRATION

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

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Peak Number 33 2,3-Pyridinediamine Concentration Rank 43

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.192 | 2.78 ng/ul | 307583 | Acenaphthene-d10 | 14.139 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | 2,3-Pyridinediamine  | 109 | C5H7N3  | 000452-58-4 | 59   |
| 2    |    |   | 2,6-Pyridinediamine  | 109 | C5H7N3  | 000141-86-6 | 47   |
| 3    |    |   | Pyridine-2,4-diamine | 109 | C5H7N3  | 000461-88-1 | 47   |
| 4    |    |   | m-Isopropoxyaniline  | 151 | C9H13NO | 041406-00-2 | 46   |
| 5    |    |   | Metacetamol          | 151 | C8H9NO2 | 000621-42-1 | 46   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

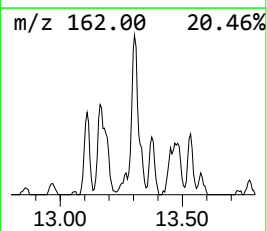
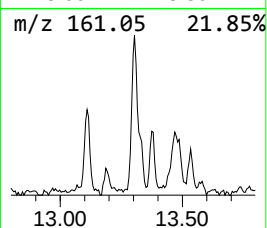
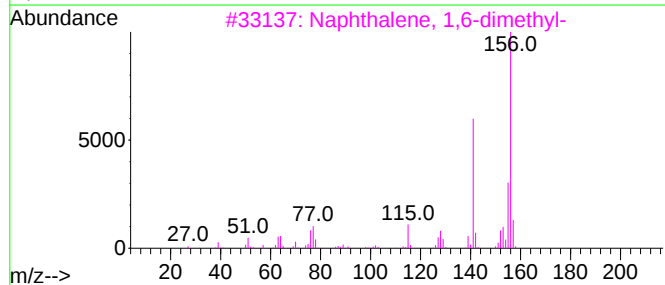
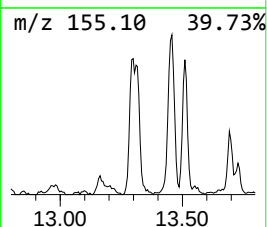
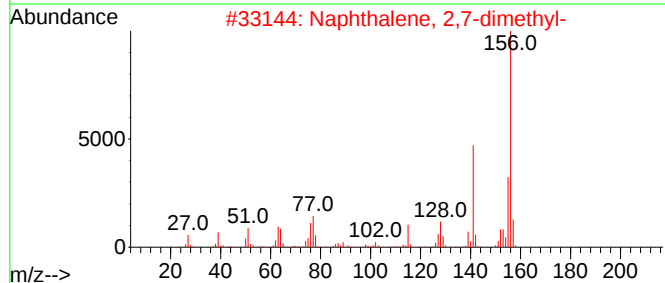
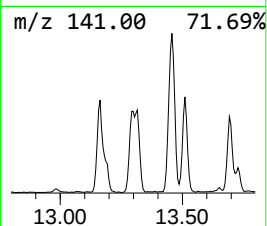
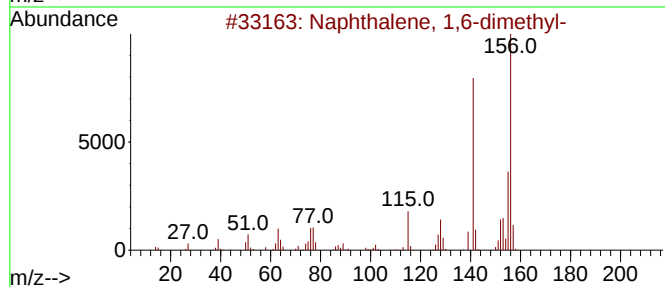
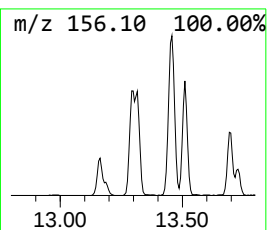
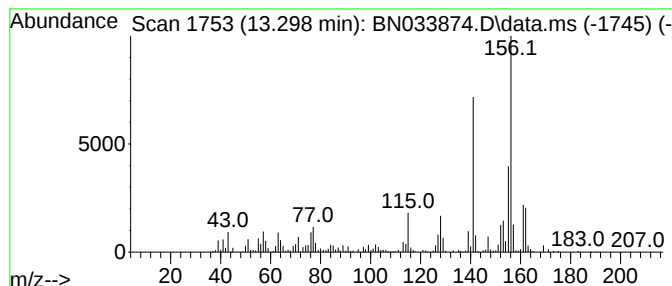
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Quant Title : SVOA CALIBRATION

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

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Peak Number 34 Naphthalene, 1,6-dimethyl- Concentration Rank 12

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.298 | 7.05 ng/ul | 779189 | Acenaphthene-d10 | 14.139 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 98   |
| 2    |    |   | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 97   |
| 3    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |
| 4    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 96   |
| 5    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

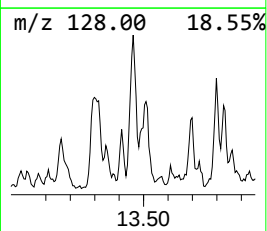
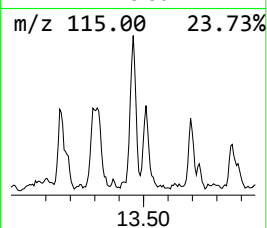
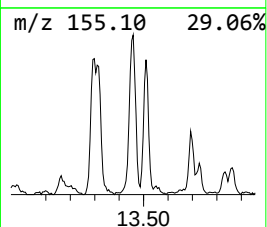
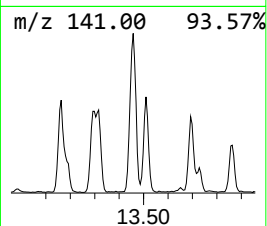
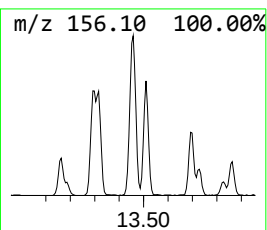
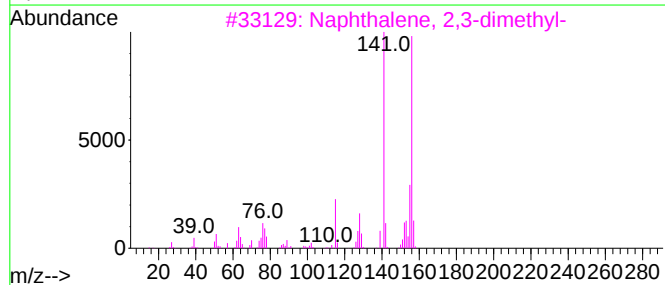
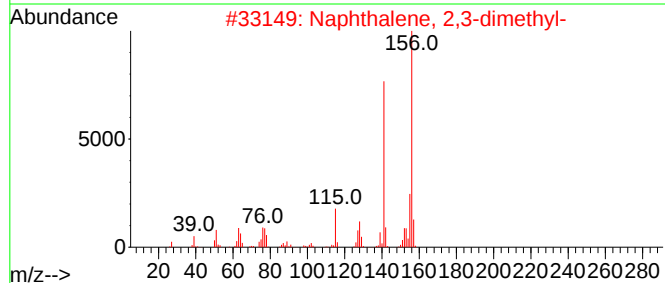
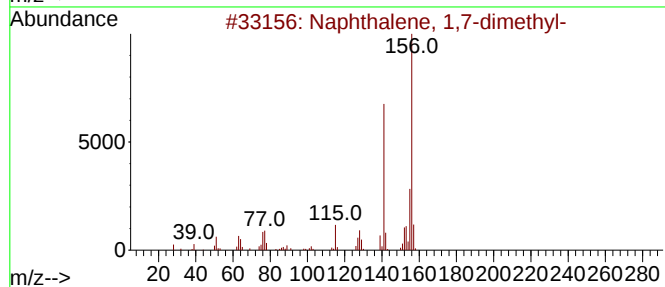
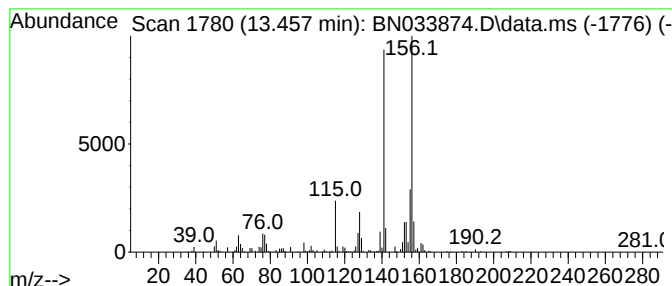
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 35 Naphthalene, 1,7-dimethyl- Concentration Rank 33

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.457 | 3.81 ng/ul | 421034 | Acenaphthene-d10 | 14.139 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 96   |
| 2    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 96   |
| 3    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 96   |
| 4    |    |   | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 96   |
| 5    |    |   | Naphthalene, 1,2-dimethyl- | 156 | C12H12  | 000573-98-8 | 95   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

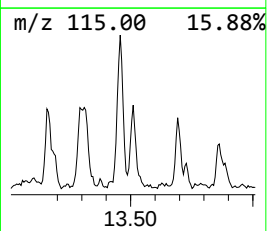
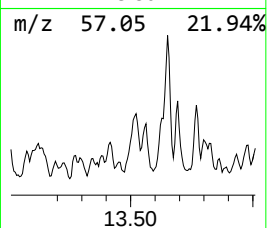
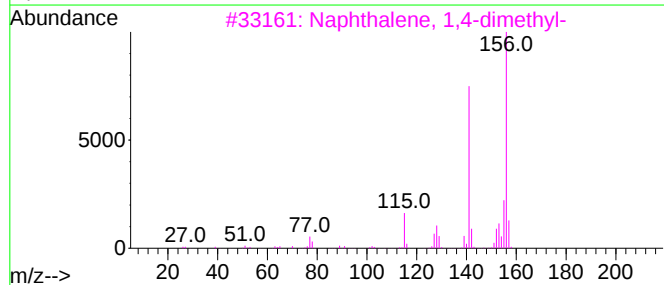
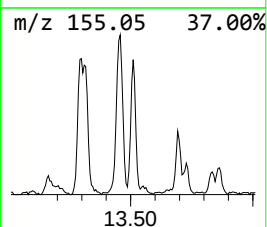
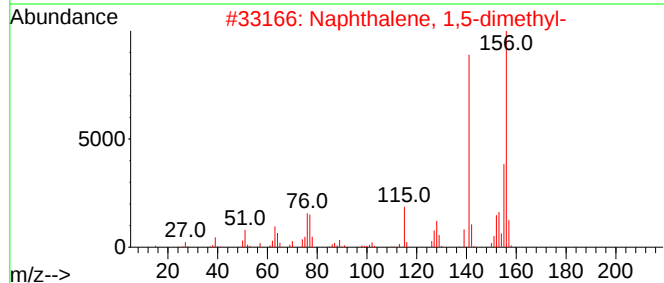
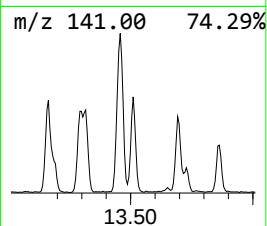
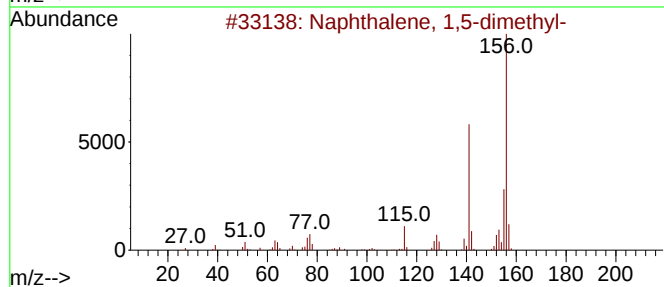
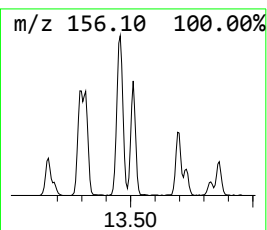
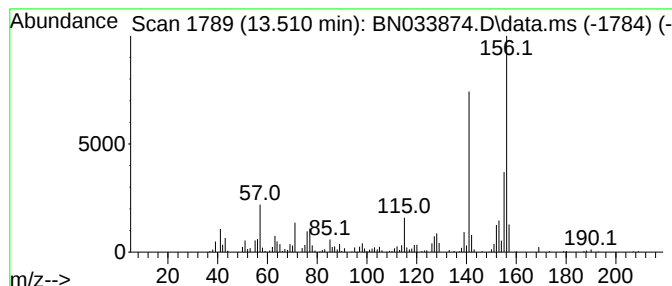
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Quant Title : SVOA CALIBRATION

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

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Peak Number 36 Naphthalene, 1,5-dimethyl- Concentration Rank 35

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.510 | 3.34 ng/ul | 368849 | Acenaphthene-d10 | 14.139 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 97   |
| 2    |    |   | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 96   |
| 3    |    |   | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 96   |
| 4    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |
| 5    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

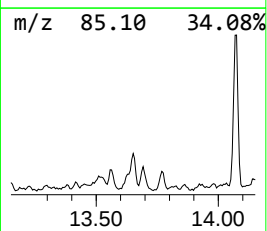
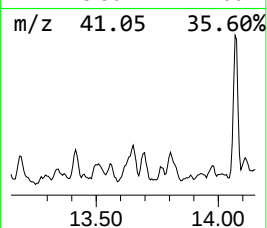
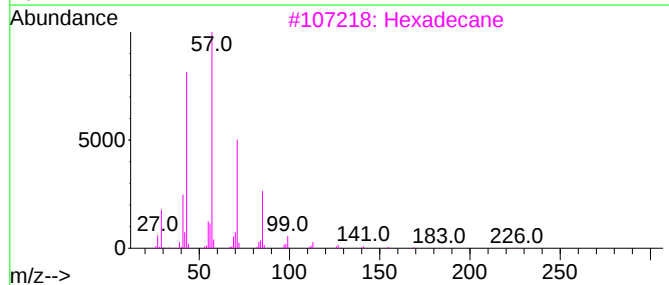
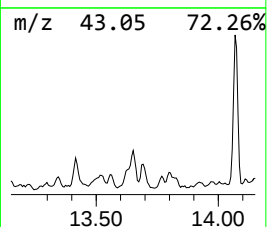
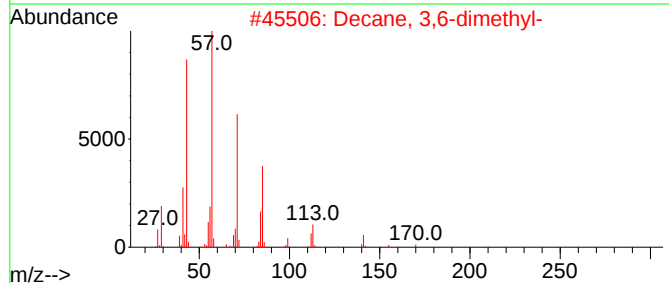
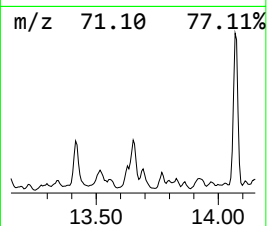
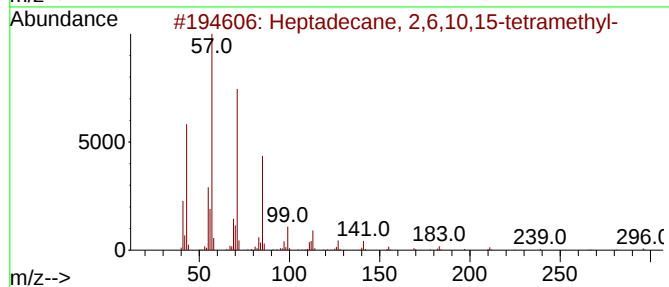
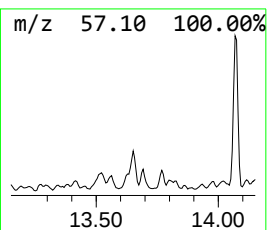
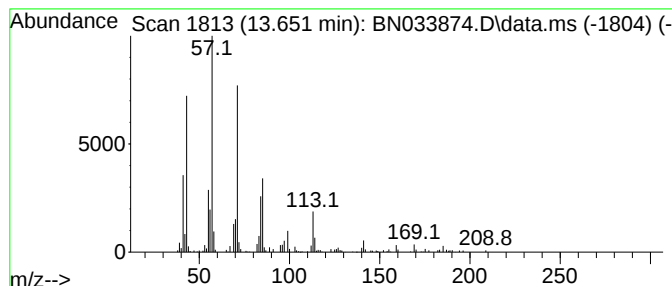
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 37 (DEL) Alkane: Straight-Chai... Concentration Rank 41

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.651 | 2.86 ng/ul | 316330 | Acenaphthene-d10 | 14.139 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 72   |
| 2    |      | Decane, 3,6-dimethyl-               | 170 | C12H26  | 017312-53-7 | 64   |
| 3    |      | Hexadecane                          | 226 | C16H34  | 000544-76-3 | 53   |
| 4    |      | Heptacosane                         | 380 | C27H56  | 000593-49-7 | 53   |
| 5    |      | Nonane, 3,7-dimethyl-               | 156 | C11H24  | 017302-32-8 | 49   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

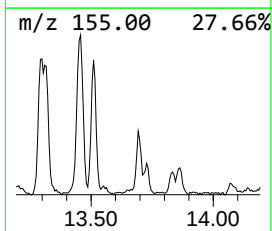
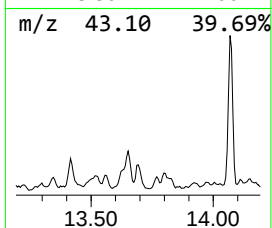
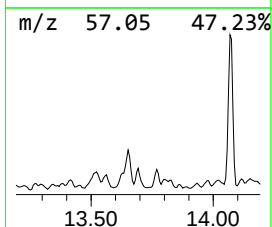
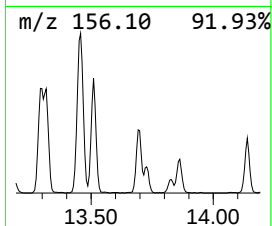
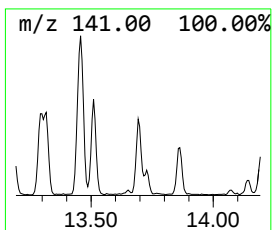
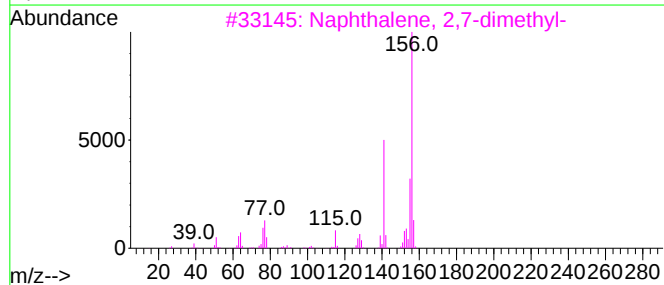
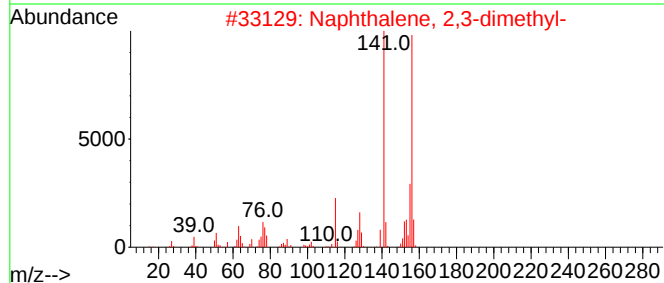
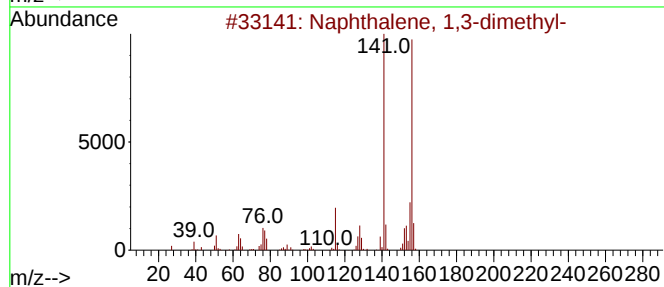
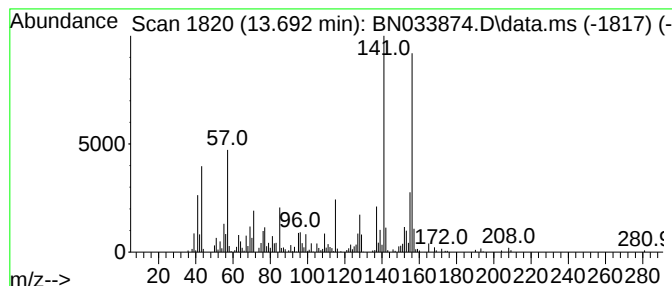
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Quant Title : SVOA CALIBRATION

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

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Peak Number 38 Naphthalene, 1,3-dimethyl- Concentration Rank 54

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.692 | 2.09 ng/ul | 231632 | Acenaphthene-d10 | 14.139 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 97   |
| 2    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |
| 3    |    |   | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 97   |
| 4    |    |   | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 97   |
| 5    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 97   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

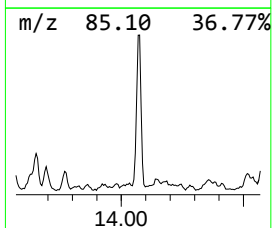
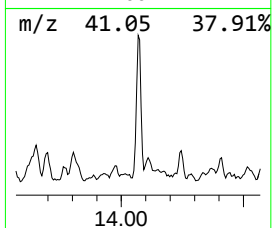
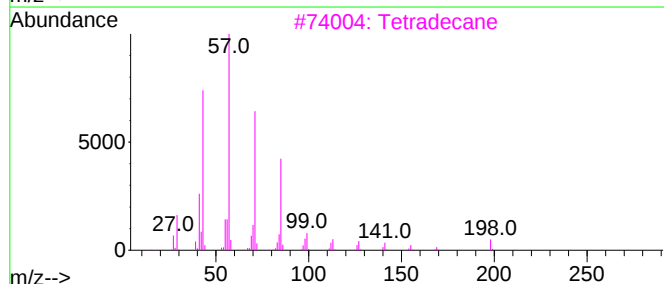
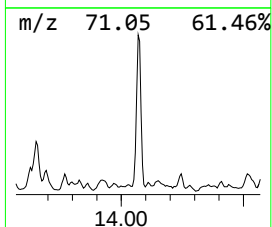
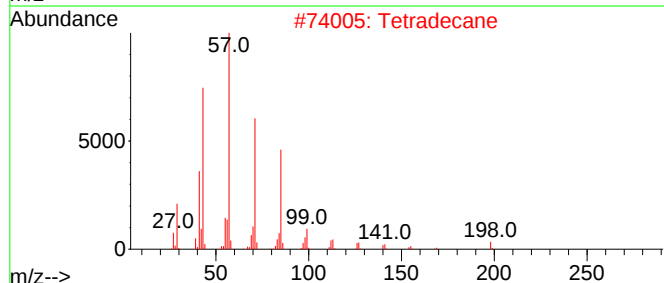
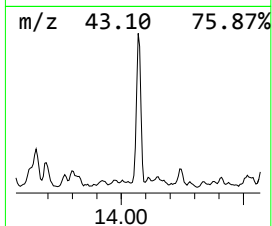
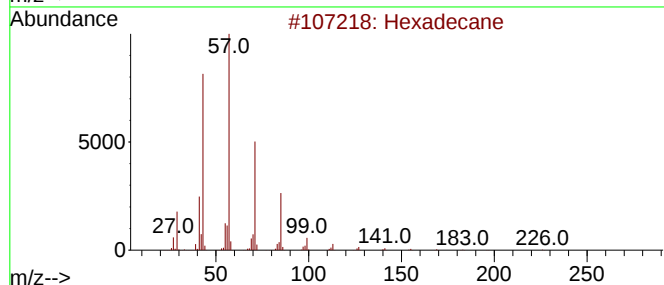
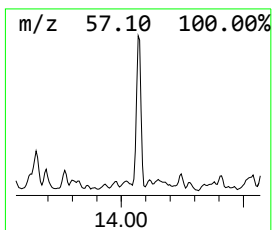
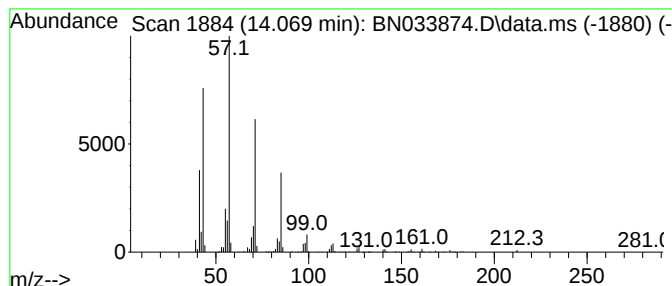
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 39 (DEL) Alkane: Straight-Chai... Concentration Rank 24

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 14.069 | 5.29 ng/ul | 585369 | Acenaphthene-d10 | 14.139 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane                   | 226 | C16H34  | 000544-76-3 | 90   |
| 2    |    |   | Tetradecane                  | 198 | C14H30  | 000629-59-4 | 90   |
| 3    |    |   | Tetradecane                  | 198 | C14H30  | 000629-59-4 | 90   |
| 4    |    |   | Hexadecane                   | 226 | C16H34  | 000544-76-3 | 90   |
| 5    |    |   | Dodecane, 2-methyl-6-propyl- | 226 | C16H34  | 055045-08-4 | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

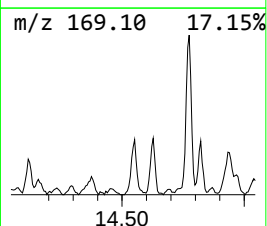
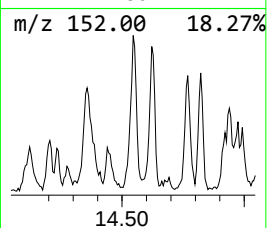
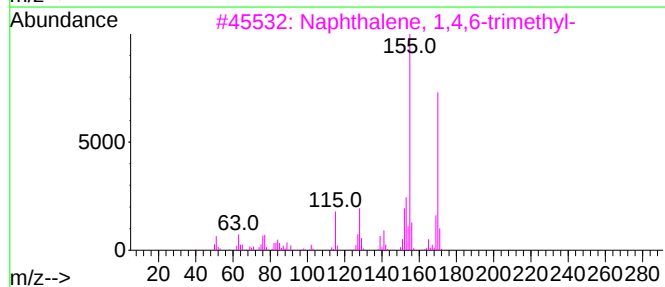
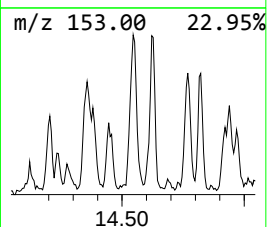
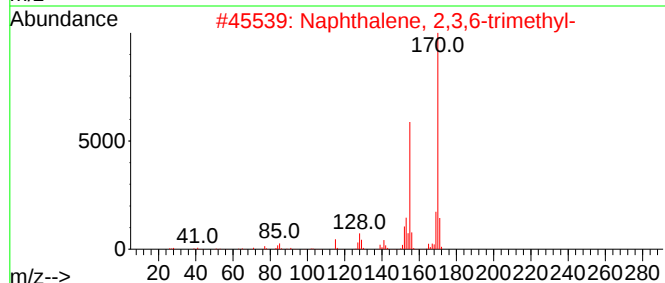
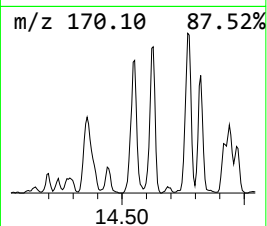
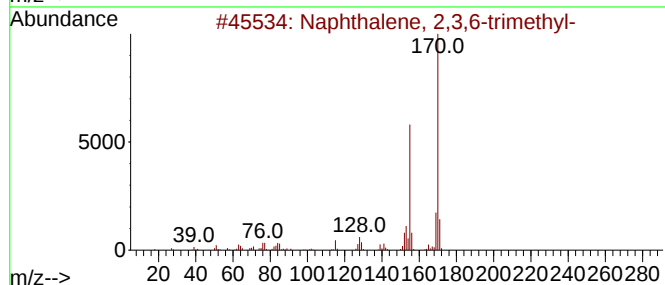
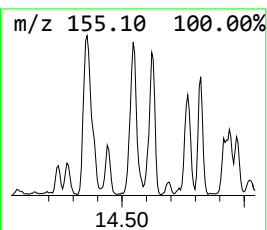
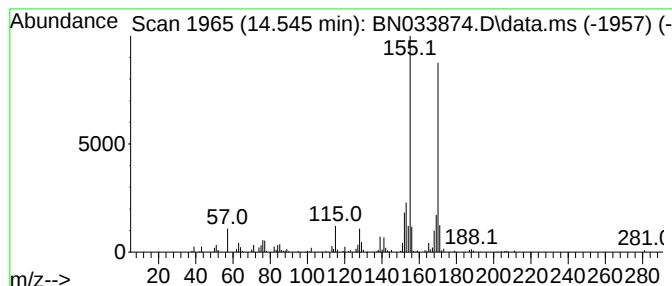
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 40 Naphthalene, 2,3,6-trimethyl- Concentration Rank 49

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 14.545 | 2.40 ng/ul | 265530 | Acenaphthene-d10 | 14.139 |

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



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

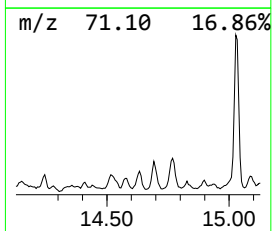
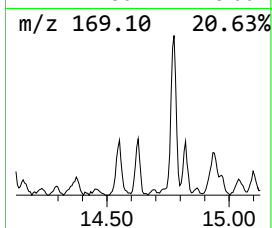
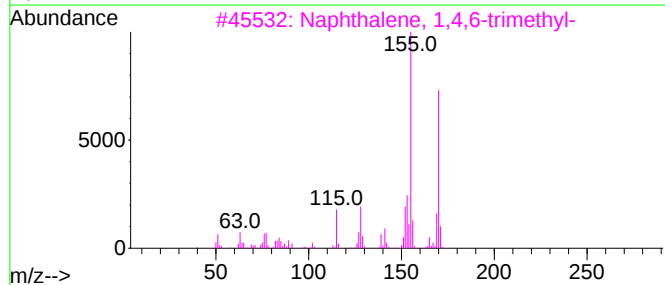
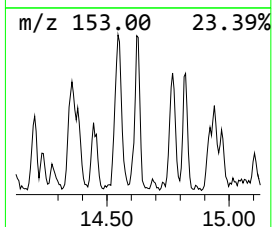
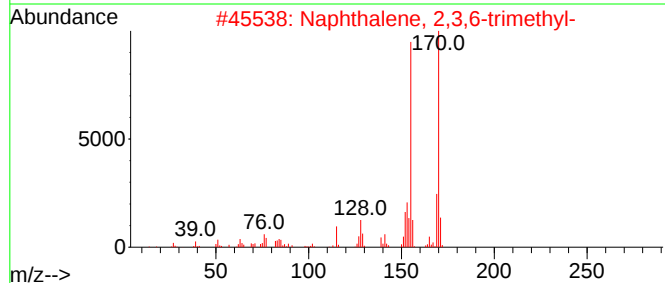
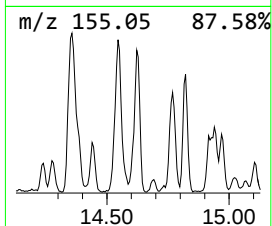
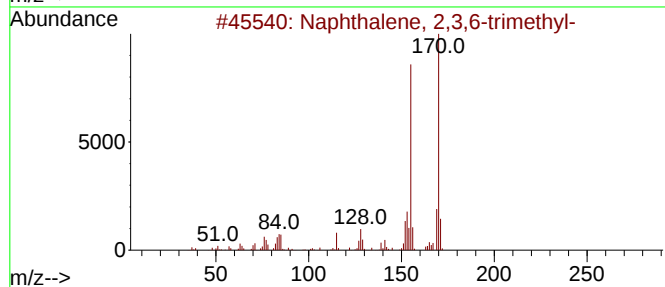
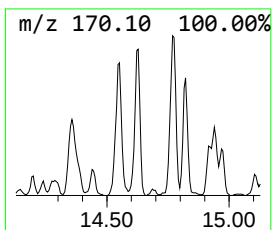
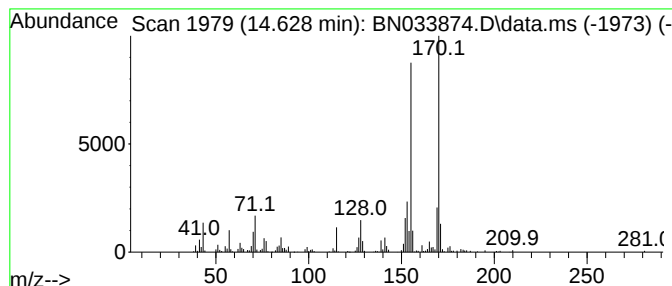
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Quant Title : SVOA CALIBRATION

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

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Peak Number 41 Naphthalene, 1,4,6-trimethyl- Concentration Rank 51

| R.T.      | EstConc                       | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------|--------|------------------|-------------|------|
| 14.628    | 2.29 ng/ul                    | 252800 | Acenaphthene-d10 | 14.139      |      |
| Hit# of 5 | Tentative ID                  | MW     | MolForm          | CAS#        | Qual |
| 1         | Naphthalene, 2,3,6-trimethyl- | 170    | C13H14           | 000829-26-5 | 97   |
| 2         | Naphthalene, 2,3,6-trimethyl- | 170    | C13H14           | 000829-26-5 | 96   |
| 3         | Naphthalene, 1,4,6-trimethyl- | 170    | C13H14           | 002131-42-2 | 96   |
| 4         | Naphthalene, 1,6,7-trimethyl- | 170    | C13H14           | 002245-38-7 | 95   |
| 5         | Naphthalene, 1,6,7-trimethyl- | 170    | C13H14           | 002245-38-7 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

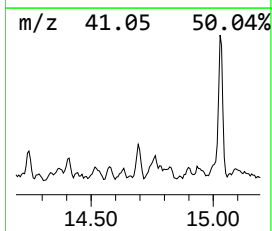
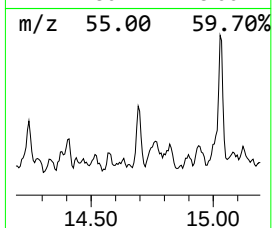
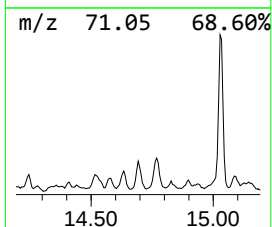
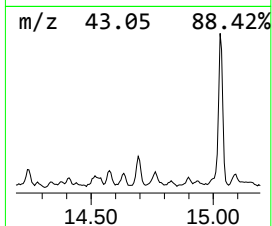
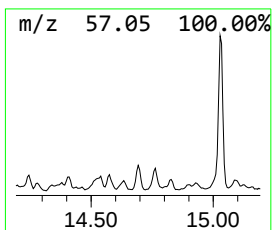
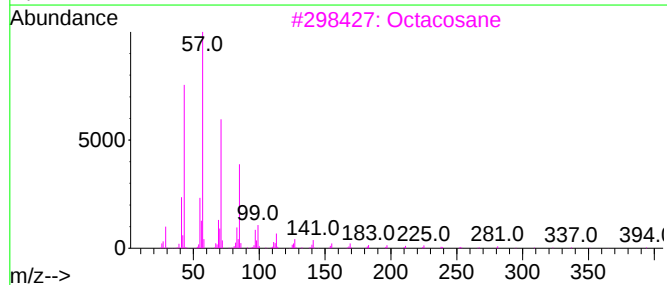
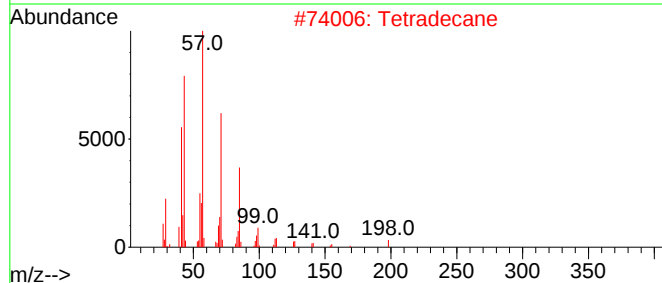
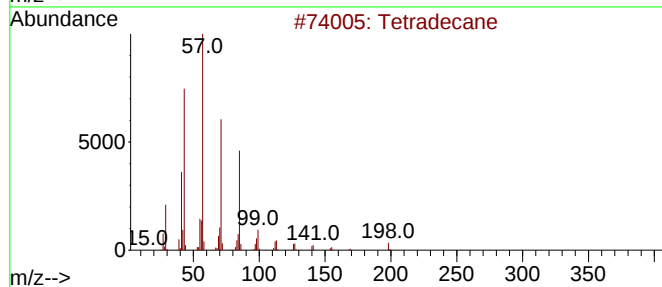
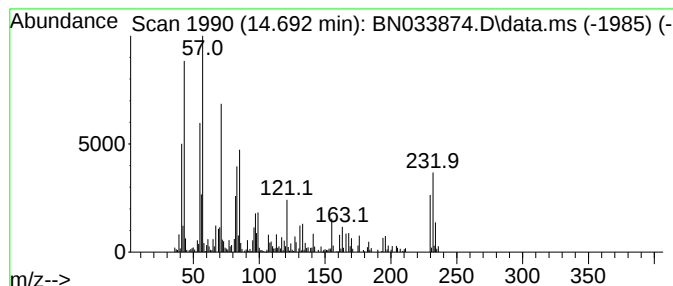
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TIC Integration Parameters: LSCINT.P

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Peak Number 42 (DEL) Alkane: Straight-Chai... Concentration Rank 50

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 14.692 | 2.36 ng/ul | 260889 | Acenaphthene-d10 | 14.139 |

| Hit# | of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------|--------------|-----|---------|-------------|------|
| 1    |      | Tetradecane  | 198 | C14H30  | 000629-59-4 | 55   |
| 2    |      | Tetradecane  | 198 | C14H30  | 000629-59-4 | 35   |
| 3    |      | Octacosane   | 394 | C28H58  | 000630-02-4 | 30   |
| 4    |      | Nonadecane   | 268 | C19H40  | 000629-92-5 | 30   |
| 5    |      | Octadecane   | 254 | C18H38  | 000593-45-3 | 30   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

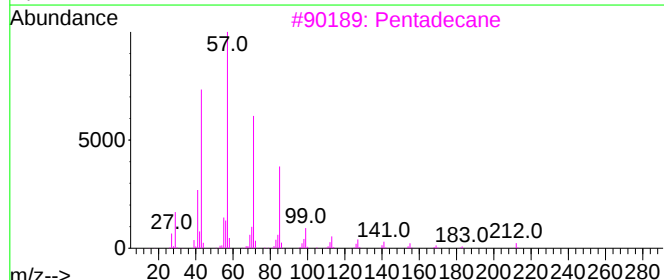
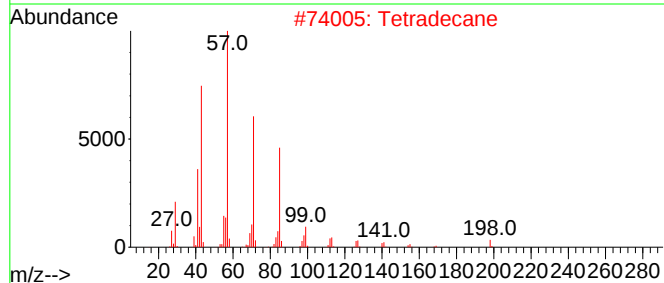
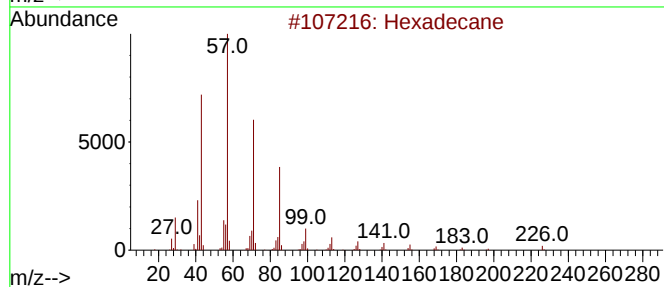
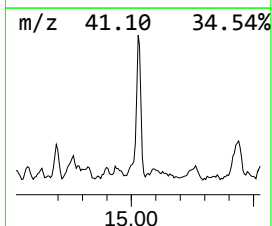
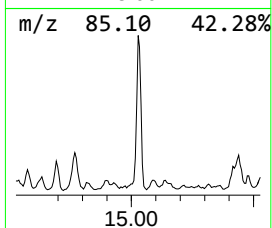
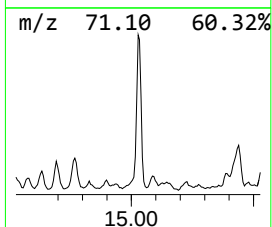
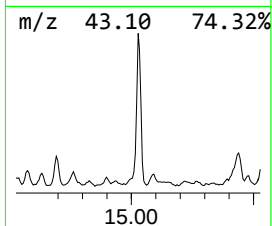
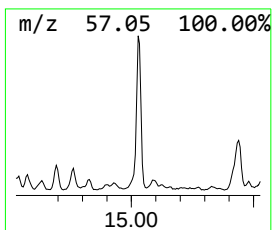
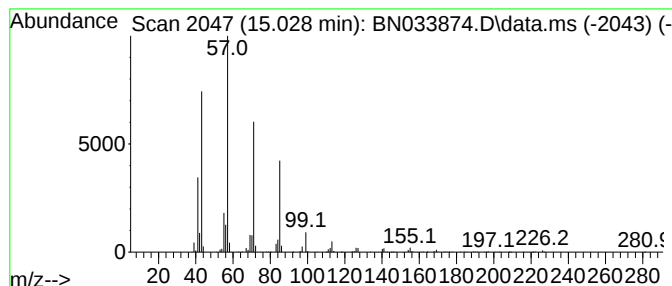
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 22

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 15.028 | 5.42 ng/ul | 599608 | Acenaphthene-d10 | 14.139 |

| Hit# | of | 5 | Tentative ID    | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane      | 226 | C16H34  | 000544-76-3 | 91   |
| 2    |    |   | Tetradecane     | 198 | C14H30  | 000629-59-4 | 90   |
| 3    |    |   | Pentadecane     | 212 | C15H32  | 000629-62-9 | 90   |
| 4    |    |   | Octadecane      | 254 | C18H38  | 000593-45-3 | 80   |
| 5    |    |   | Decane, 1-iodo- | 268 | C10H21I | 002050-77-3 | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

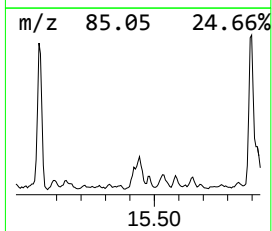
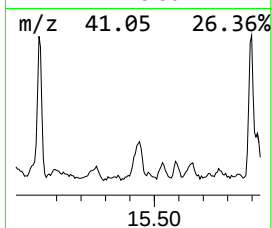
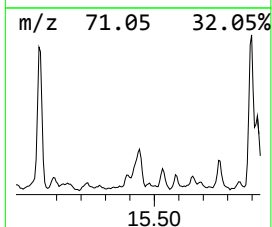
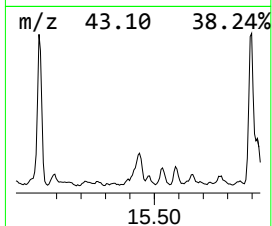
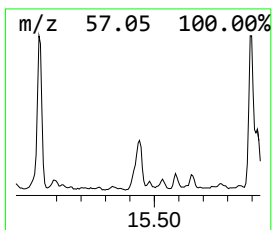
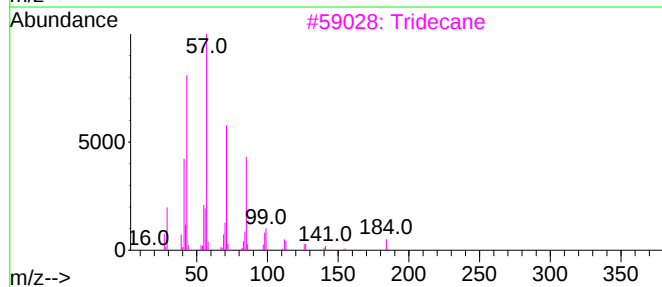
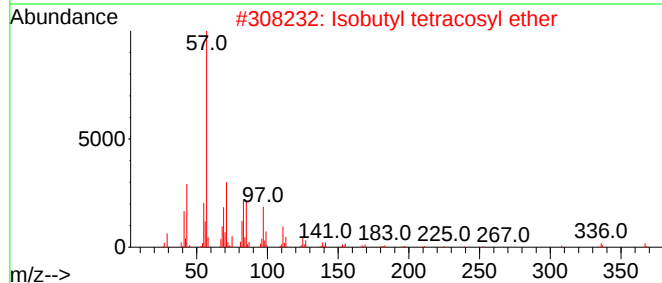
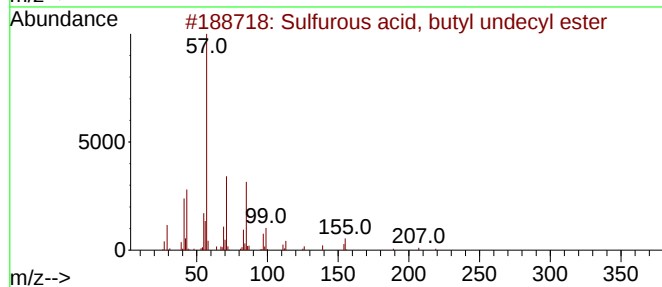
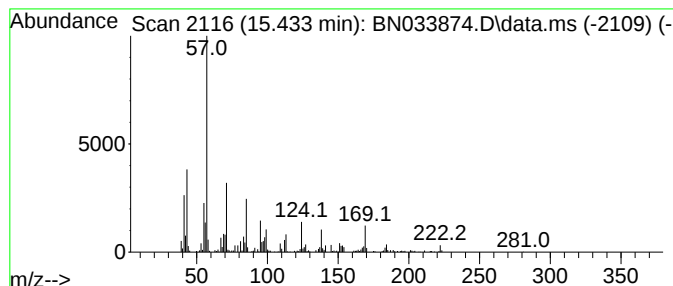
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 44 unknown-02 Concentration Rank 34

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 15.434 | 3.69 ng/ul | 407713 | Acenaphthene-d10 | 14.139 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Sulfurous acid, butyl undecyl ester | 292 | C15H32O3S | 1000309-17-8 | 49   |
| 2    |    |   | Isobutyl tetracosyl ether           | 410 | C28H58O   | 1000406-33-2 | 47   |
| 3    |    |   | Tridecane                           | 184 | C13H28    | 000629-50-5  | 45   |
| 4    |    |   | Tetracosane                         | 338 | C24H50    | 000646-31-1  | 43   |
| 5    |    |   | Octacosane                          | 394 | C28H58    | 000630-02-4  | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

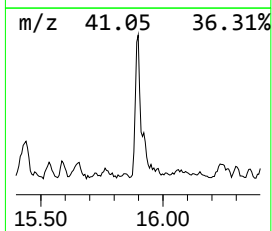
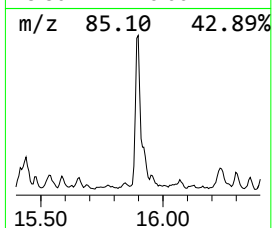
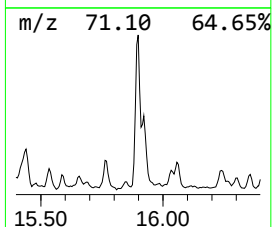
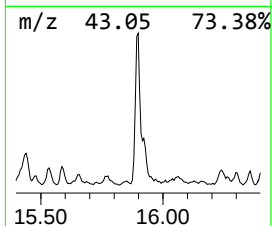
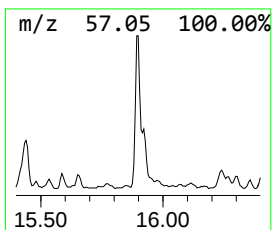
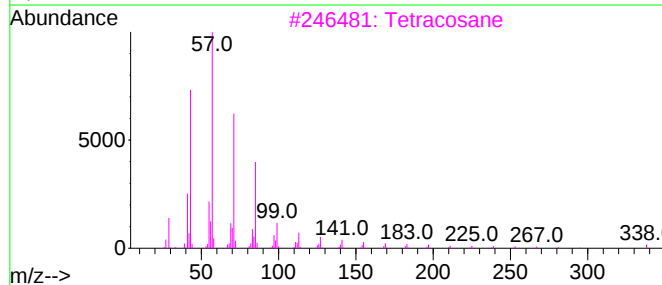
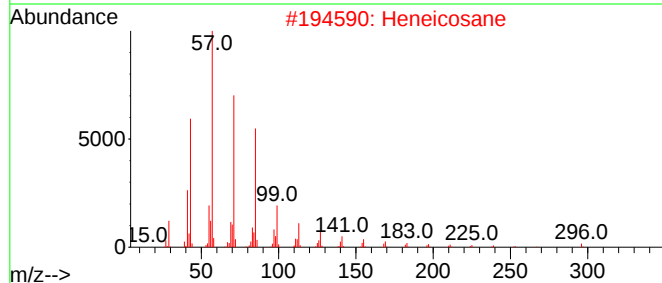
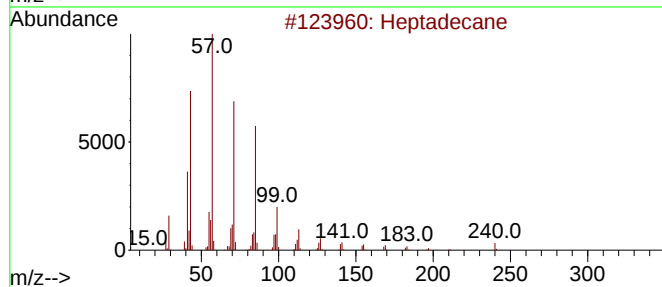
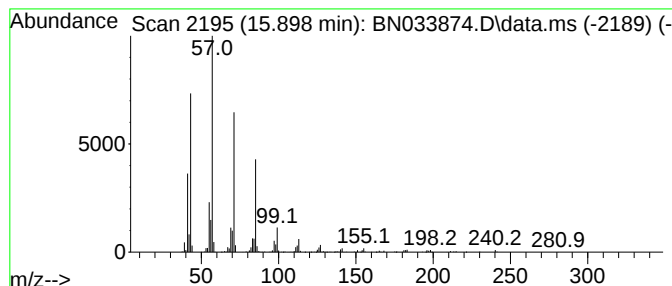
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TIC Integration Parameters: LSCINT.P

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Peak Number 45 (DEL) Alkane: Straight-Chai... Concentration Rank 16

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 15.898 | 6.06 ng/ul | 785461 | Phenanthrene-d10 | 16.892 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane  | 240 | C17H36  | 000629-78-7 | 97   |
| 2    |    |   | Heneicosane  | 296 | C21H44  | 000629-94-7 | 90   |
| 3    |    |   | Tetracosane  | 338 | C24H50  | 000646-31-1 | 90   |
| 4    |    |   | Heptadecane  | 240 | C17H36  | 000629-78-7 | 90   |
| 5    |    |   | Nonadecane   | 268 | C19H40  | 000629-92-5 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

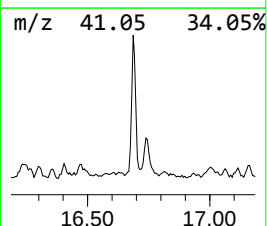
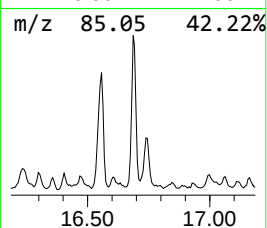
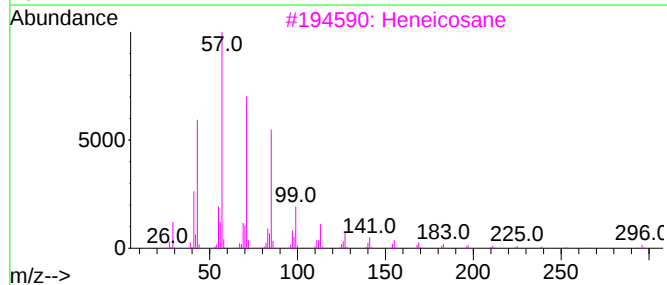
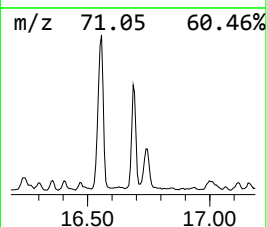
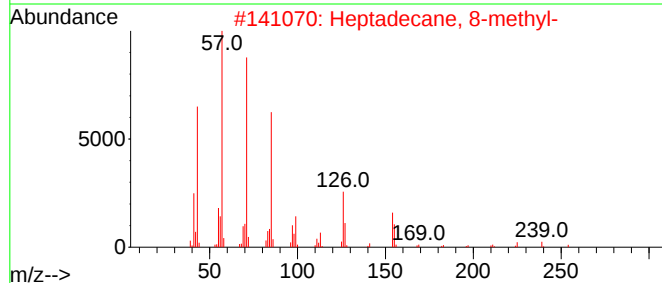
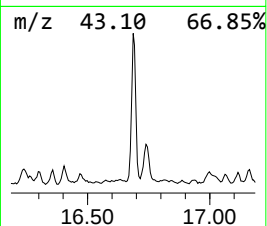
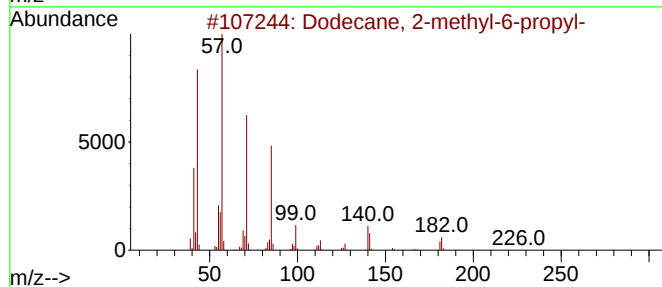
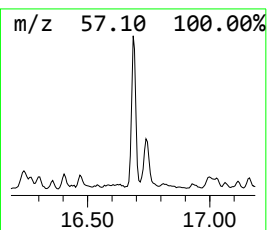
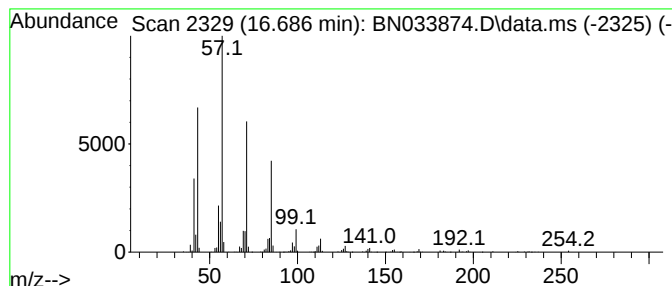
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 46 (DEL) Alkane: Straight-Chai... Concentration Rank 23

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.686 | 5.30 ng/ul | 686958 | Phenanthrene-d10 | 16.892 |

| Hit# | of 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|------|------------------------------|-----|---------|-------------|------|
| 1    |      | Dodecane, 2-methyl-6-propyl- | 226 | C16H34  | 055045-08-4 | 91   |
| 2    |      | Heptadecane, 8-methyl-       | 254 | C18H38  | 013287-23-5 | 91   |
| 3    |      | Heneicosane                  | 296 | C21H44  | 000629-94-7 | 90   |
| 4    |      | Hexadecane, 5-butyl-         | 282 | C20H42  | 006912-07-8 | 86   |
| 5    |      | Tetradecane                  | 198 | C14H30  | 000629-59-4 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

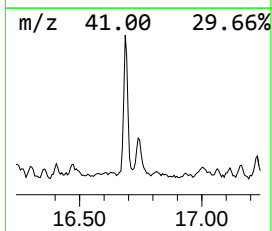
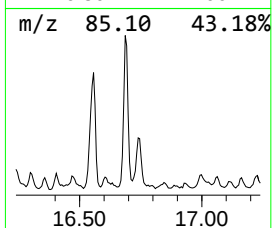
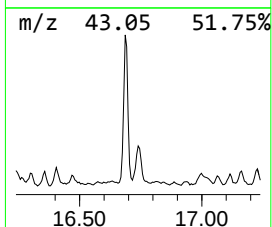
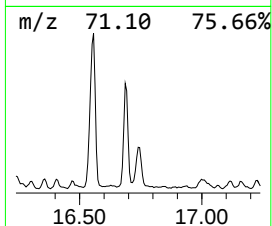
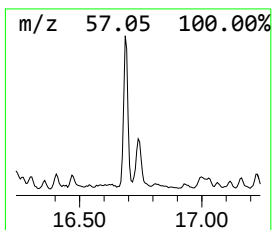
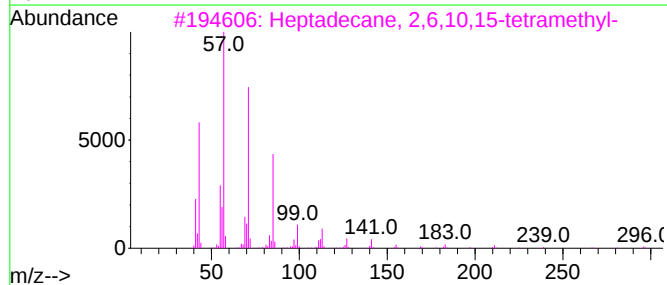
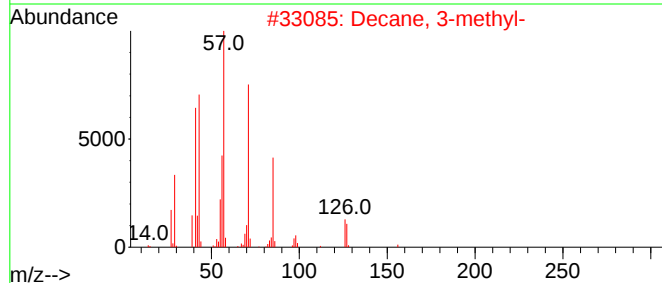
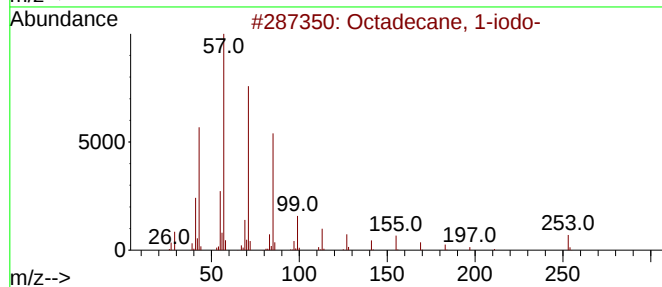
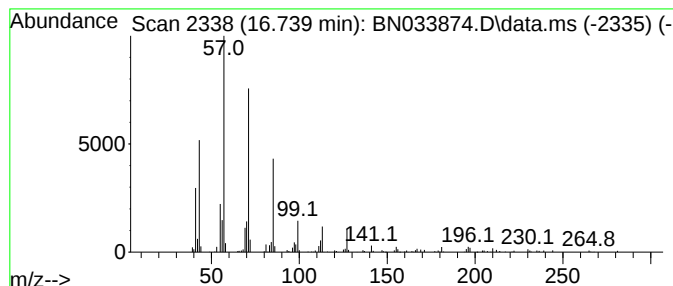
Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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 Octadecane, 1-iodo- Concentration Rank 48

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 16.739    | 2.50 ng/ul                          | 323486 | Phenanthrene-d10 | 16.892       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | Octadecane, 1-iodo-                 | 380    | C18H37I          | 000629-93-6  | 90   |
| 2         | Decane, 3-methyl-                   | 156    | C11H24           | 013151-34-3  | 87   |
| 3         | Heptadecane, 2,6,10,15-tetramethyl- | 296    | C21H44           | 054833-48-6  | 87   |
| 4         | Tetradecane, 1-iodo-                | 324    | C14H29I          | 019218-94-1  | 80   |
| 5         | Sulfurous acid, 2-ethylhexyl hex... | 278    | C14H30O3S        | 1000309-20-2 | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

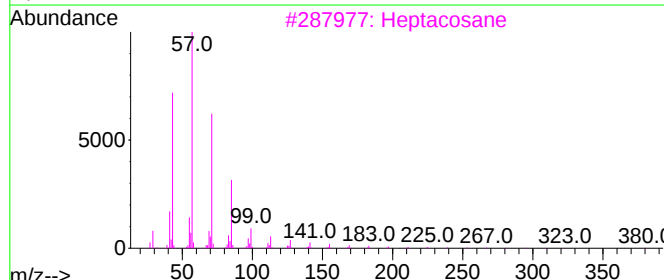
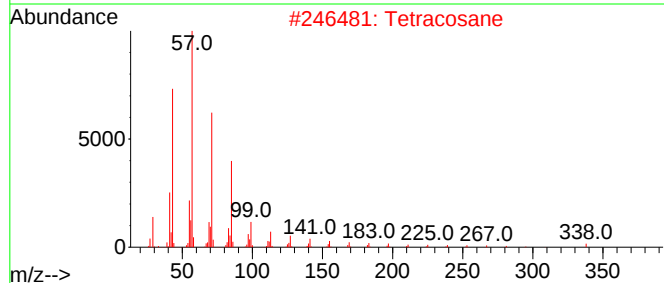
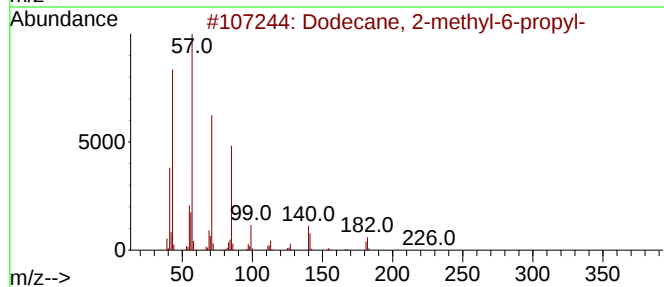
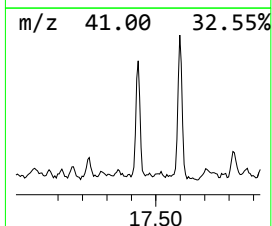
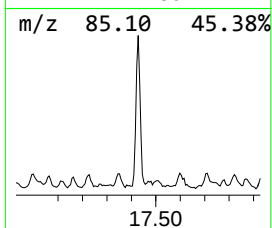
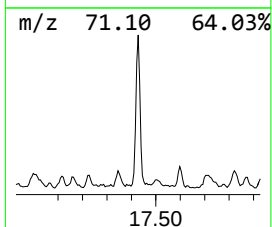
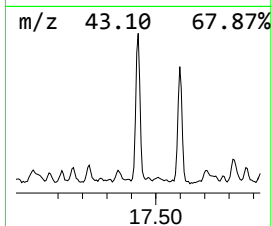
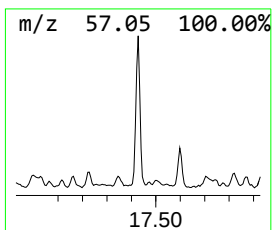
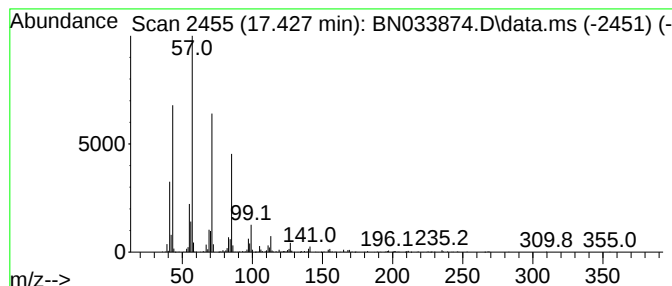
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TIC Integration Parameters: LSCINT.P

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Peak Number 48 (DEL) Alkane: Straight-Chai... Concentration Rank 29

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.427 | 4.54 ng/ul | 588639 | Phenanthrene-d10 | 16.892 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Dodecane, 2-methyl-6-propyl-        | 226 | C16H34  | 055045-08-4 | 93   |
| 2    |    |   | Tetracosane                         | 338 | C24H50  | 000646-31-1 | 91   |
| 3    |    |   | Heptacosane                         | 380 | C27H56  | 000593-49-7 | 91   |
| 4    |    |   | Pentadecane                         | 212 | C15H32  | 000629-62-9 | 91   |
| 5    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

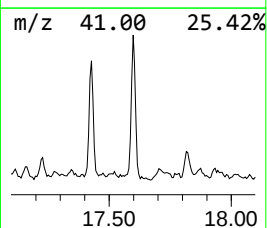
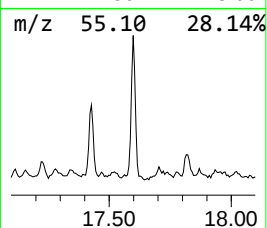
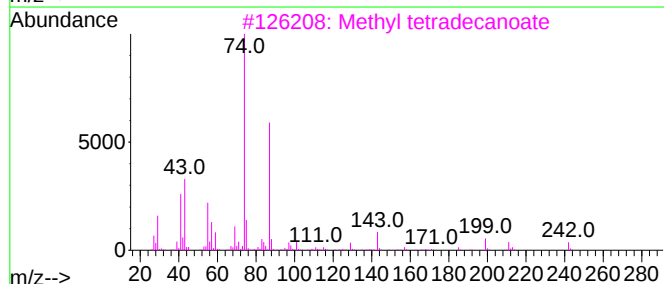
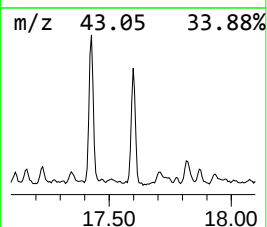
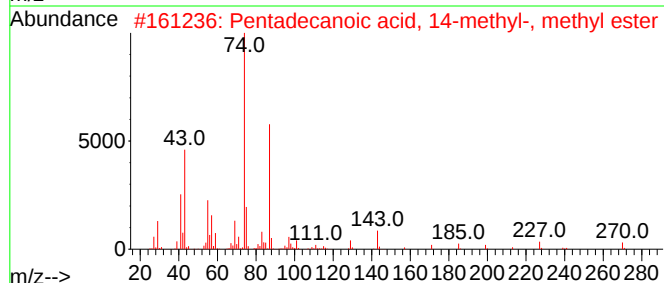
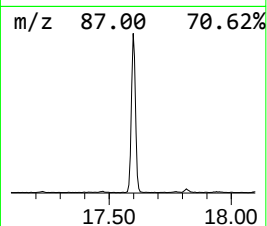
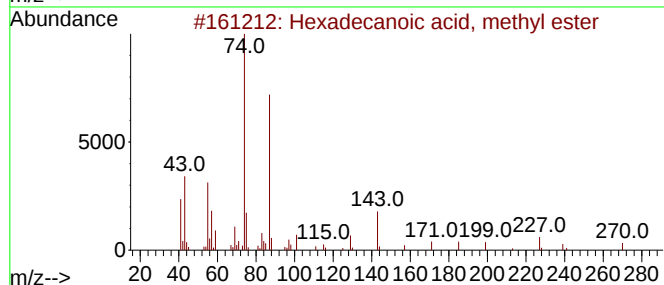
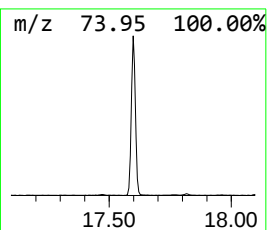
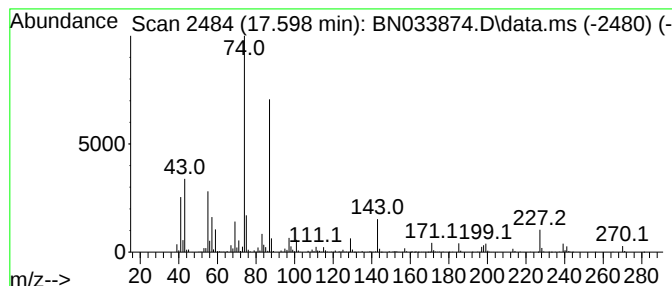
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Quant Title : SVOA CALIBRATION

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

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Peak Number 49 Hexadecanoic acid, methyl e... Concentration Rank 8

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 17.598 | 8.23 ng/ul | 1066180 | Phenanthrene-d10 | 16.892 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Hexadecanoic acid, methyl ester     | 270 | C17H34O2 | 000112-39-0 | 98   |
| 2    |    |   | Pentadecanoic acid, 14-methyl-, ... | 270 | C17H34O2 | 005129-60-2 | 96   |
| 3    |    |   | Methyl tetradecanoate               | 242 | C15H30O2 | 000124-10-7 | 94   |
| 4    |    |   | Hexadecanoic acid, methyl ester     | 270 | C17H34O2 | 000112-39-0 | 92   |
| 5    |    |   | Hexadecanoic acid, methyl ester     | 270 | C17H34O2 | 000112-39-0 | 92   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

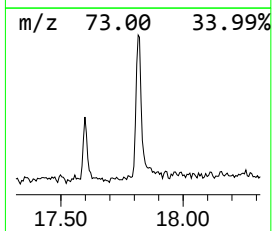
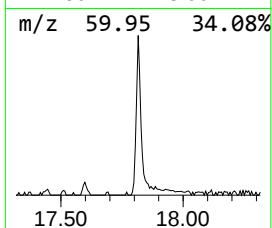
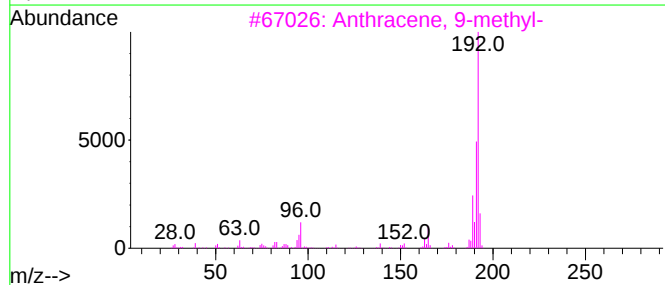
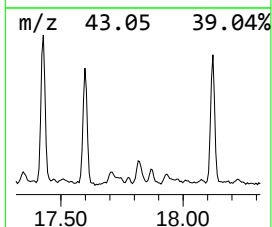
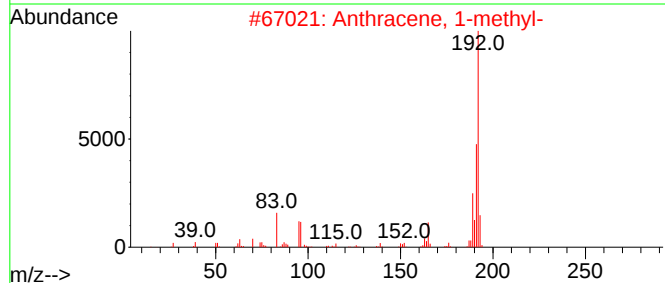
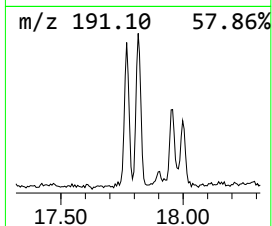
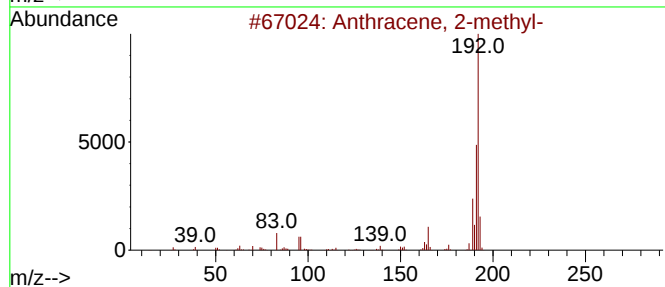
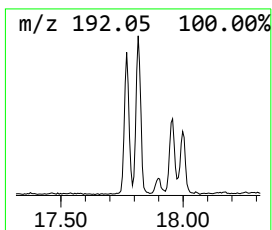
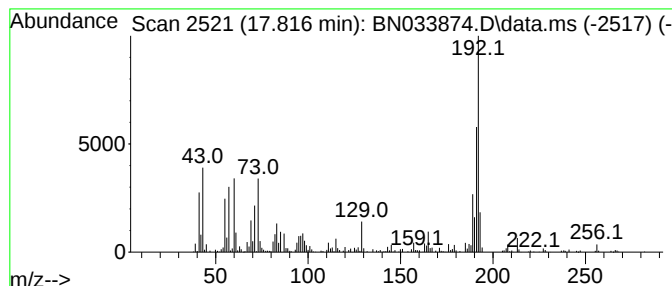
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 50 Anthracene, 2-methyl- Concentration Rank 52

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.816 | 2.28 ng/ul | 295038 | Phenanthrene-d10 | 16.892 |

| Hit# | of 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------|-----|---------|-------------|------|
| 1    |      | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 93   |
| 2    |      | Anthracene, 1-methyl-   | 192 | C15H12  | 000610-48-0 | 92   |
| 3    |      | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 91   |
| 4    |      | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 91   |
| 5    |      | Phenanthrene, 4-methyl- | 192 | C15H12  | 000832-64-4 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

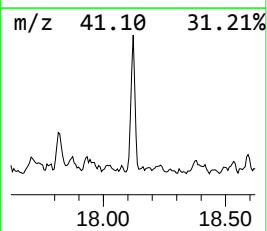
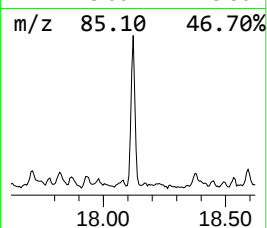
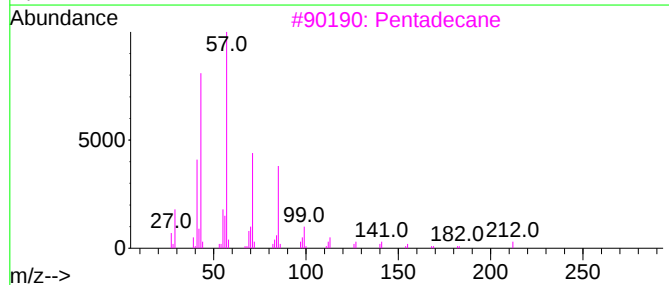
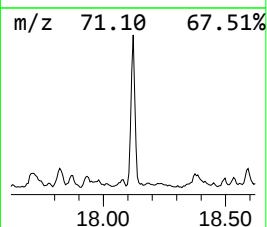
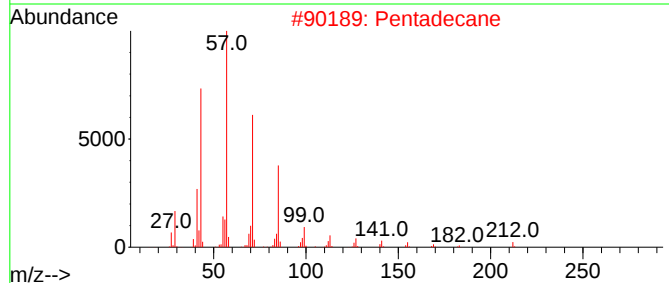
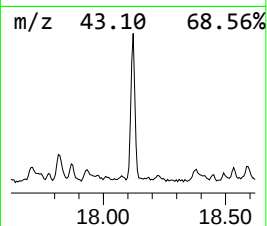
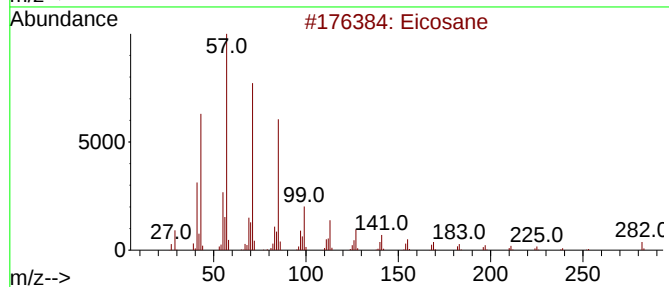
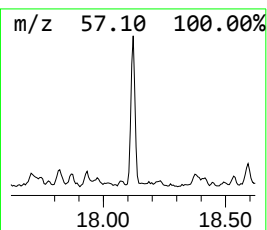
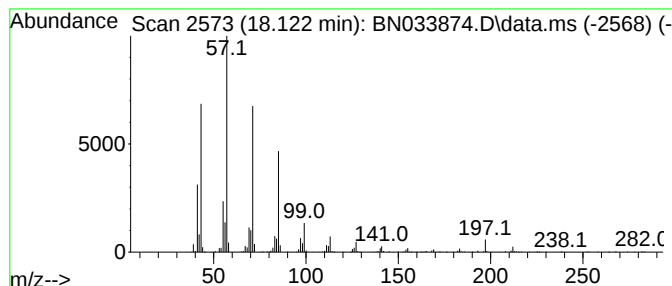
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 51 (DEL) Alkane: Straight-Chai... Concentration Rank 32

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.122 | 4.01 ng/ul | 519326 | Phenanthrene-d10 | 16.892 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Eicosane     | 282 | C20H42  | 000112-95-8 | 96   |
| 2    |    |   | Pentadecane  | 212 | C15H32  | 000629-62-9 | 96   |
| 3    |    |   | Pentadecane  | 212 | C15H32  | 000629-62-9 | 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\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

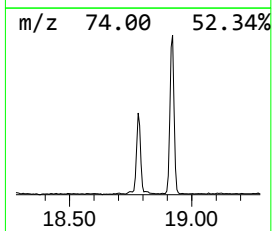
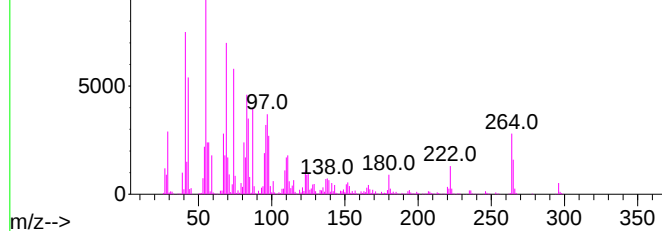
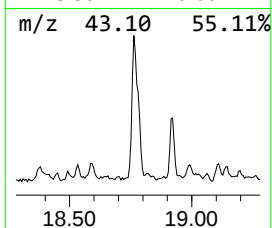
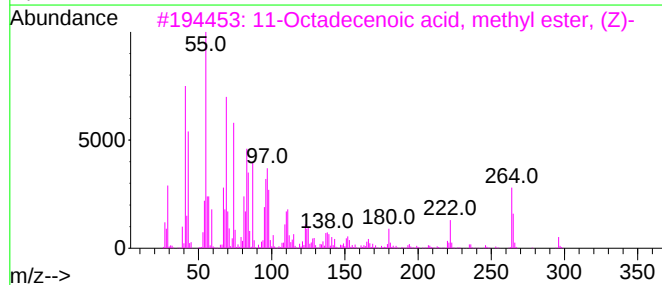
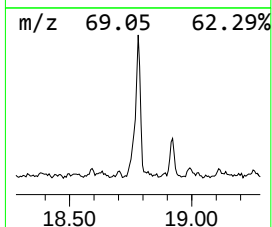
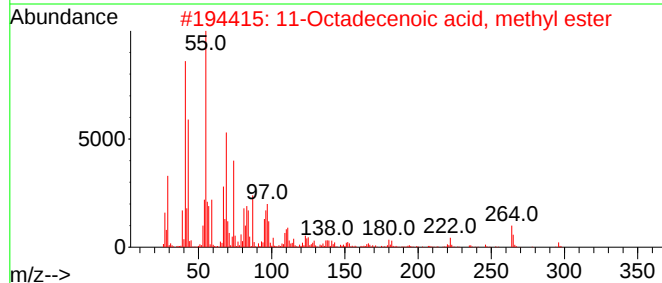
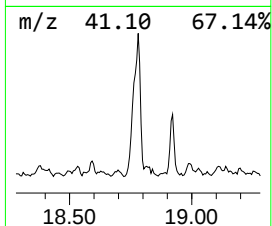
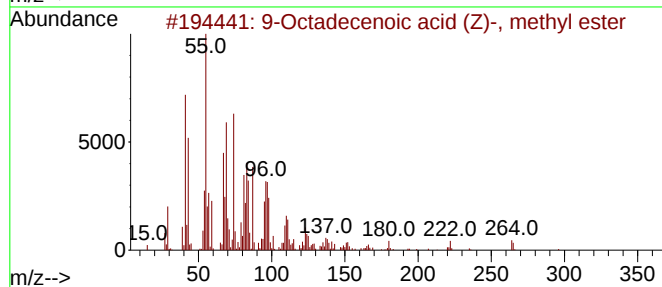
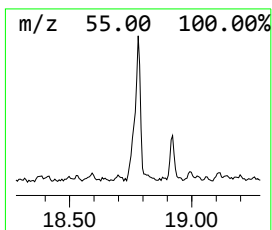
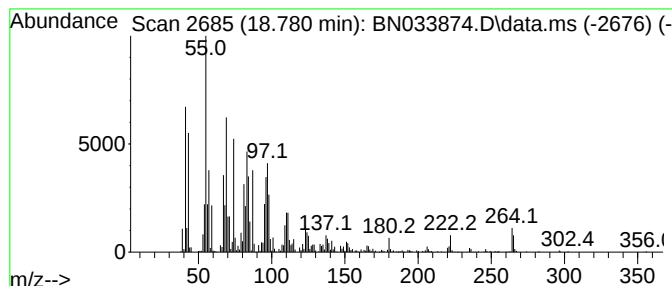
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TIC Integration Parameters: LSCINT.P

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Peak Number 52 9-Octadecenoic acid (Z)-, m... Concentration Rank 3

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 18.780 | 9.54 ng/ul | 1236050 | Phenanthrene-d10 | 16.892 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 9-Octadecenoic acid (Z)-, methyl... | 296 | C19H36O2 | 000112-62-9 | 99   |
| 2    |    |   | 11-Octadecenoic acid, methyl ester  | 296 | C19H36O2 | 052380-33-3 | 99   |
| 3    |    |   | 11-Octadecenoic acid, methyl est... | 296 | C19H36O2 | 001937-63-9 | 98   |
| 4    |    |   | 11-Octadecenoic acid, methyl ester  | 296 | C19H36O2 | 052380-33-3 | 98   |
| 5    |    |   | 16-Octadecenoic acid, methyl ester  | 296 | C19H36O2 | 056554-49-5 | 98   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

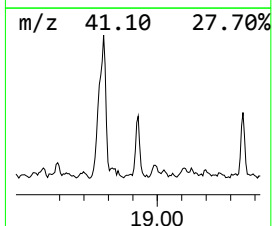
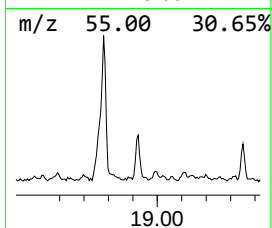
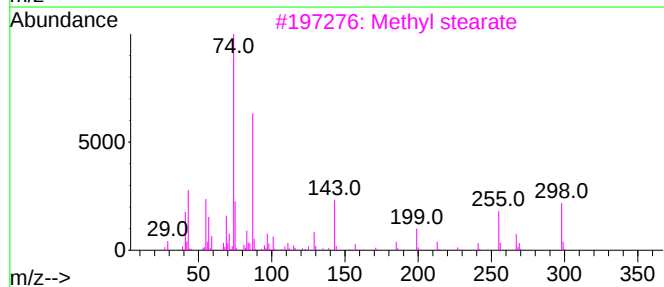
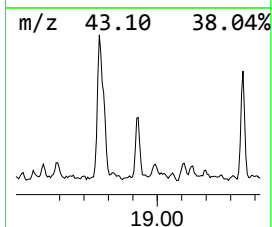
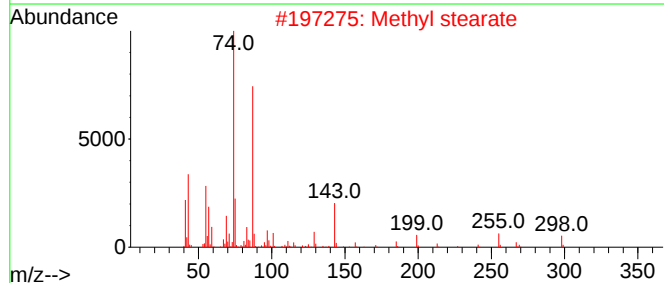
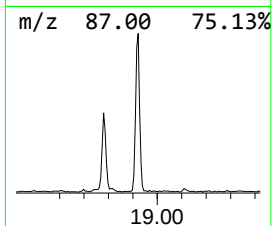
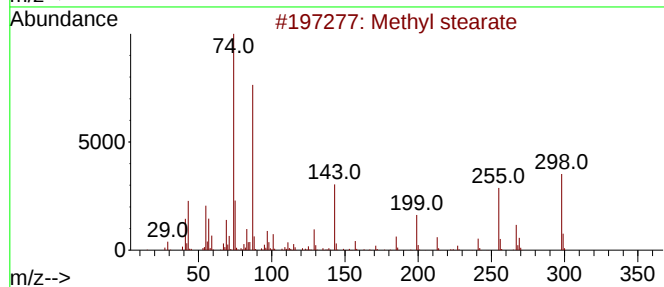
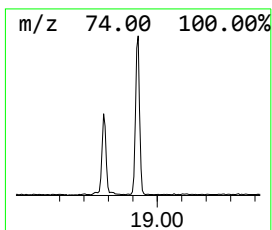
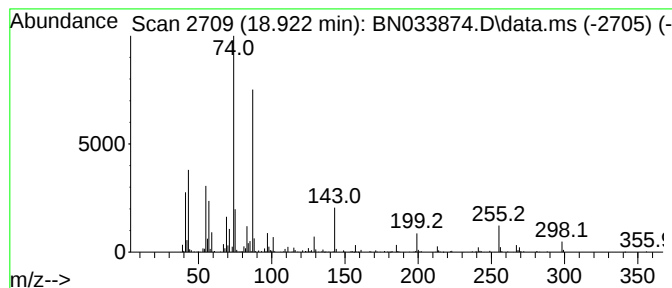
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Quant Title : SVOA CALIBRATION

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

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Peak Number 53 Methyl stearate Concentration Rank 42

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.922 | 2.85 ng/ul | 368842 | Phenanthrene-d10 | 16.892 |

| Hit# | of | 5 | Tentative ID    | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-----------------|-----|----------|-------------|------|
| 1    |    |   | Methyl stearate | 298 | C19H38O2 | 000112-61-8 | 98   |
| 2    |    |   | Methyl stearate | 298 | C19H38O2 | 000112-61-8 | 98   |
| 3    |    |   | Methyl stearate | 298 | C19H38O2 | 000112-61-8 | 98   |
| 4    |    |   | Methyl stearate | 298 | C19H38O2 | 000112-61-8 | 97   |
| 5    |    |   | Methyl stearate | 298 | C19H38O2 | 000112-61-8 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

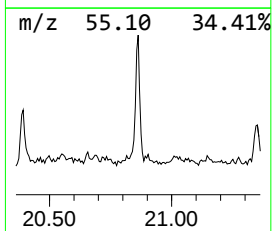
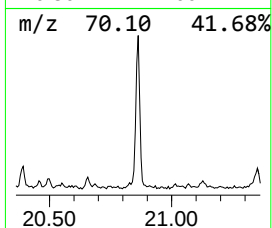
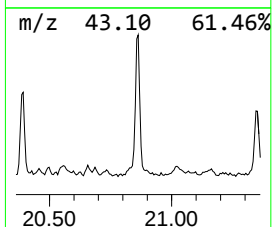
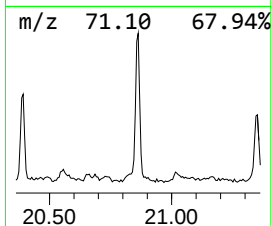
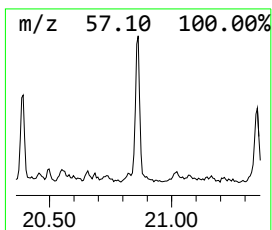
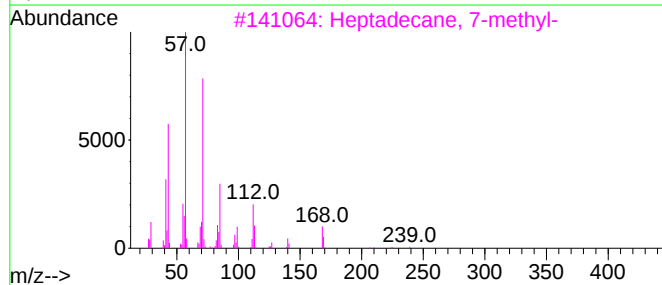
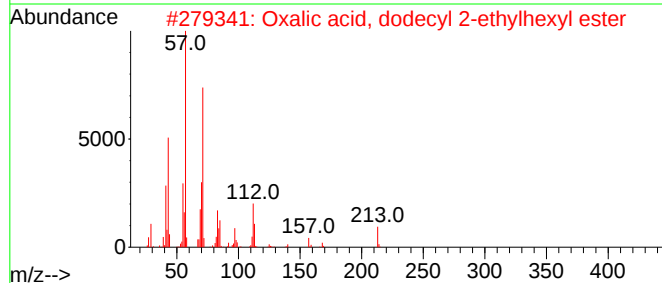
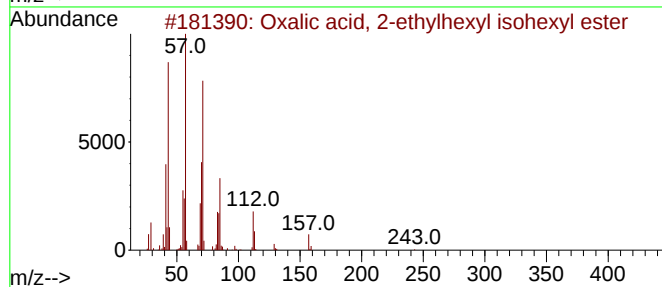
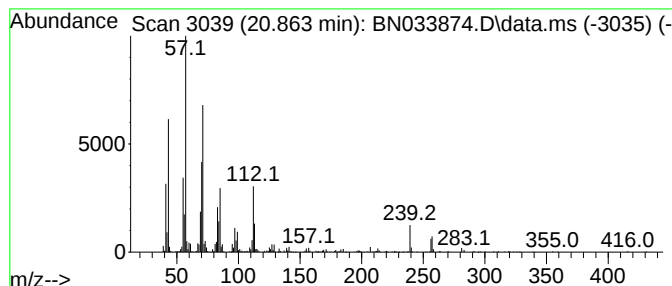
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 54 Oxalic acid, 2-ethylhexyl i... Concentration Rank 38

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 20.863 | 2.90 ng/ul | 505155 | Chrysene-d12     | 21.127 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Oxalic acid, 2-ethylhexyl isohex... | 286 | C16H30O4  | 1000309-38-8 | 64   |
| 2    |    |   | Oxalic acid, dodecyl 2-ethylhexy... | 370 | C22H42O4  | 1000309-39-5 | 58   |
| 3    |    |   | Heptadecane, 7-methyl-              | 254 | C18H38    | 020959-33-5  | 52   |
| 4    |    |   | Heneicosane                         | 296 | C21H44    | 000629-94-7  | 46   |
| 5    |    |   | Sulfurous acid, butyl dodecyl ester | 306 | C16H34O3S | 1000309-17-9 | 46   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
Data File : BN033874.D  
Acq On : 11 Sep 2024 23:28  
Operator : JU/RC  
Sample : P3878-02DL 20X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DDC46DL

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

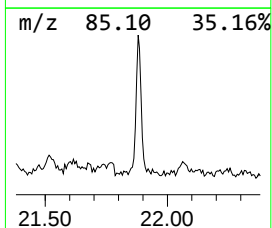
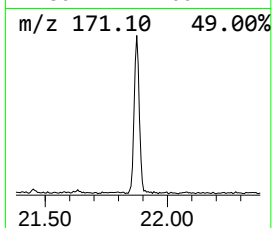
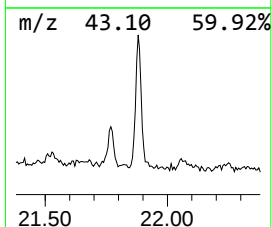
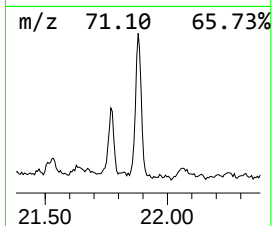
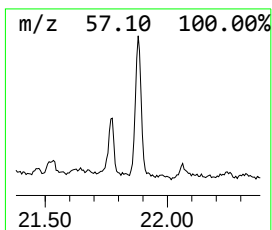
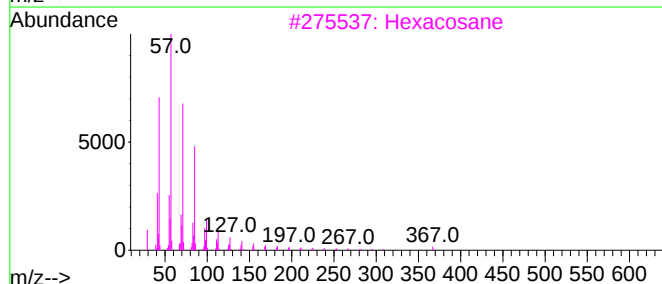
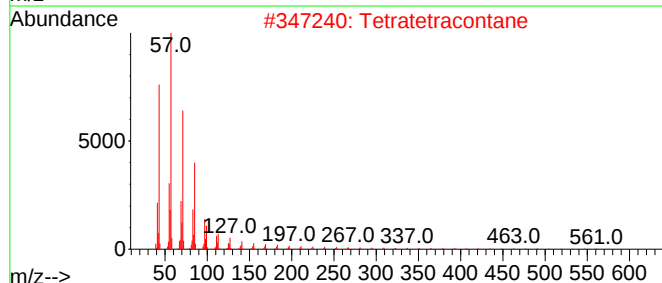
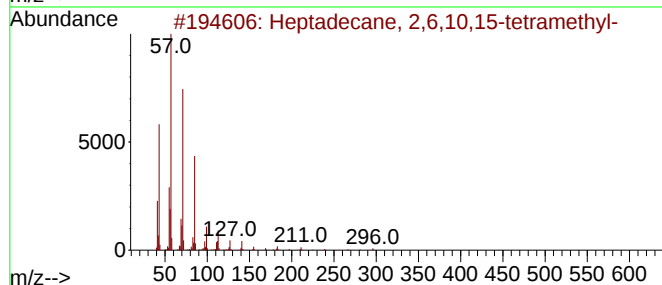
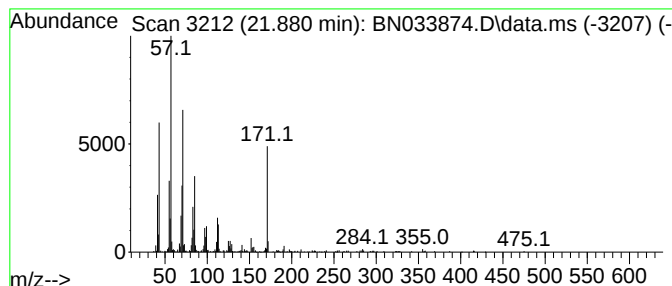
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 55 (DEL) Alkane: Straight-Chai... Concentration Rank 45

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 21.880 | 2.65 ng/ul | 460992 | Chrysene-d12     | 21.127 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44   | 054833-48-6  | 90   |
| 2    |    |   | Tetratetracontane                   | 619 | C44H90   | 007098-22-8  | 58   |
| 3    |    |   | Hexacosane                          | 366 | C26H54   | 000630-01-3  | 58   |
| 4    |    |   | Oxalic acid, decyl 2-ethylhexyl ... | 342 | C20H38O4 | 1000309-39-3 | 52   |
| 5    |    |   | Tritetracontane                     | 605 | C43H88   | 007098-21-7  | 52   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091124\  
 Data File : BN033874.D  
 Acq On : 11 Sep 2024 23:28  
 Operator : JU/RC  
 Sample : P3878-02DL 20X  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 DDC46DL

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

TIC Library : C:\Database\NIST0.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
|                    |        |         |       |          | #                     | RT     | Resp    | Conc |
| (DEL) Alkane: S... | 5.563  | 7.3     | ng/ul | 469729   | 1                     | 7.499  | 1285840 | 20.0 |
| (DEL) Alkane: S... | 6.093  | 2.9     | ng/ul | 186709   | 1                     | 7.499  | 1285840 | 20.0 |
| (DEL) Alkane: C... | 6.163  | 3.0     | ng/ul | 196035   | 1                     | 7.499  | 1285840 | 20.0 |
| Benzene, propyl-   | 6.493  | 3.0     | ng/ul | 195686   | 1                     | 7.499  | 1285840 | 20.0 |
| (DEL) Alkane: S... | 6.516  | 4.8     | ng/ul | 305845   | 1                     | 7.499  | 1285840 | 20.0 |
| (DEL) Alkane: S... | 6.575  | 5.5     | ng/ul | 353556   | 1                     | 7.499  | 1285840 | 20.0 |
| Benzene, 1-ethy... | 6.604  | 8.2     | ng/ul | 524610   | 1                     | 7.499  | 1285840 | 20.0 |
| Benzene, 1,2,3-... | 6.740  | 4.2     | ng/ul | 271028   | 1                     | 7.499  | 1285840 | 20.0 |
| Benzene, 1-ethy... | 6.899  | 5.6     | ng/ul | 360343   | 1                     | 7.499  | 1285840 | 20.0 |
| 2-Heptanone, 4,... | 6.946  | 9.3     | ng/ul | 598415   | 1                     | 7.499  | 1285840 | 20.0 |
| (DEL) Alkane: S... | 7.140  | 42.3    | ng/ul | 2717650  | 1                     | 7.499  | 1285840 | 20.0 |
| Butyl dodecyl e... | 7.575  | 5.2     | ng/ul | 331775   | 1                     | 7.499  | 1285840 | 20.0 |
| Mesitylene         | 7.610  | 6.1     | ng/ul | 391935   | 1                     | 7.499  | 1285840 | 20.0 |
| (DEL) Alkane: C... | 7.775  | 6.0     | ng/ul | 387163   | 1                     | 7.499  | 1285840 | 20.0 |
| Cyclohexanone, ... | 7.922  | 9.3     | ng/ul | 598564   | 1                     | 7.499  | 1285840 | 20.0 |
| unknown-01         | 7.981  | 6.5     | ng/ul | 420930   | 1                     | 7.499  | 1285840 | 20.0 |
| Benzene, 1-meth... | 8.040  | 9.4     | ng/ul | 602339   | 1                     | 7.499  | 1285840 | 20.0 |
| (DEL) Alkane: S... | 8.093  | 4.6     | ng/ul | 295505   | 1                     | 7.499  | 1285840 | 20.0 |
| Benzene, 1,4-di... | 8.134  | 8.9     | ng/ul | 575647   | 1                     | 7.499  | 1285840 | 20.0 |
| (DEL) Alkane: S... | 8.269  | 7.1     | ng/ul | 459184   | 1                     | 7.499  | 1285840 | 20.0 |
| Benzene, 1-meth... | 8.293  | 5.2     | ng/ul | 335687   | 1                     | 7.499  | 1285840 | 20.0 |
| Benzene, 2-ethy... | 8.440  | 4.1     | ng/ul | 262320   | 1                     | 7.499  | 1285840 | 20.0 |
| o-Cymene           | 8.493  | 5.7     | ng/ul | 368257   | 1                     | 7.499  | 1285840 | 20.0 |
| (DEL) Alkane: S... | 8.728  | 28.8    | ng/ul | 1849680  | 1                     | 7.499  | 1285840 | 20.0 |
| Benzene, 1-meth... | 8.840  | 5.7     | ng/ul | 366903   | 1                     | 7.499  | 1285840 | 20.0 |
| 2,4-Dipropyl-5-... | 10.657 | 2.9     | ng/ul | 618224   | 2                     | 10.257 | 4318740 | 20.0 |
| (DEL) Alkane: S... | 10.787 | 2.6     | ng/ul | 560271   | 2                     | 10.257 | 4318740 | 20.0 |
| (DEL) Alkane: S... | 11.710 | 2.8     | ng/ul | 599823   | 2                     | 10.257 | 4318740 | 20.0 |
| (DEL) Alkane: S... | 12.692 | 2.6     | ng/ul | 283412   | 3                     | 14.139 | 2211820 | 20.0 |
| (DEL) Alkane: S... | 12.981 | 6.3     | ng/ul | 692706   | 3                     | 14.139 | 2211820 | 20.0 |
| 2,3-Pyridinedia... | 13.192 | 2.8     | ng/ul | 307583   | 3                     | 14.139 | 2211820 | 20.0 |
| Naphthalene, 1,... | 13.298 | 7.0     | ng/ul | 779189   | 3                     | 14.139 | 2211820 | 20.0 |
| Naphthalene, 1,... | 13.457 | 3.8     | ng/ul | 421034   | 3                     | 14.139 | 2211820 | 20.0 |
| Naphthalene, 1,... | 13.510 | 3.3     | ng/ul | 368849   | 3                     | 14.139 | 2211820 | 20.0 |
| (DEL) Alkane: S... | 13.651 | 2.9     | ng/ul | 316330   | 3                     | 14.139 | 2211820 | 20.0 |
| Naphthalene, 1,... | 13.692 | 2.1     | ng/ul | 231632   | 3                     | 14.139 | 2211820 | 20.0 |
| (DEL) Alkane: S... | 14.069 | 5.3     | ng/ul | 585369   | 3                     | 14.139 | 2211820 | 20.0 |
| Naphthalene, 2,... | 14.545 | 2.4     | ng/ul | 265530   | 3                     | 14.139 | 2211820 | 20.0 |
| Naphthalene, 1,... | 14.628 | 2.3     | ng/ul | 252800   | 3                     | 14.139 | 2211820 | 20.0 |
| (DEL) Alkane: S... | 14.692 | 2.4     | ng/ul | 260889   | 3                     | 14.139 | 2211820 | 20.0 |
| (DEL) Alkane: S... | 15.028 | 5.4     | ng/ul | 599608   | 3                     | 14.139 | 2211820 | 20.0 |
| unknown-02         | 15.434 | 3.7     | ng/ul | 407713   | 3                     | 14.139 | 2211820 | 20.0 |
| (DEL) Alkane: S... | 15.898 | 6.1     | ng/ul | 785461   | 4                     | 16.892 | 2591340 | 20.0 |
| (DEL) Alkane: S... | 16.686 | 5.3     | ng/ul | 686958   | 4                     | 16.892 | 2591340 | 20.0 |
| Octadecane, 1-i... | 16.739 | 2.5     | ng/ul | 323486   | 4                     | 16.892 | 2591340 | 20.0 |
| (DEL) Alkane: S... | 17.427 | 4.5     | ng/ul | 588639   | 4                     | 16.892 | 2591340 | 20.0 |
| Hexadecanoic ac... | 17.598 | 8.2     | ng/ul | 1066180  | 4                     | 16.892 | 2591340 | 20.0 |
| Anthracene, 2-m... | 17.816 | 2.3     | ng/ul | 295038   | 4                     | 16.892 | 2591340 | 20.0 |
| (DEL) Alkane: S... | 18.122 | 4.0     | ng/ul | 519326   | 4                     | 16.892 | 2591340 | 20.0 |
| 9-Octadecenoic ... | 18.780 | 9.5     | ng/ul | 1236050  | 4                     | 16.892 | 2591340 | 20.0 |
| Methyl stearate    | 18.922 | 2.9     | ng/ul | 368842   | 4                     | 16.892 | 2591340 | 20.0 |
| Oxalic acid, 2-... | 20.863 | 2.9     | ng/ul | 505155   | 5                     | 21.127 | 3478570 | 20.0 |

**Instrument :**  
BNA\_N  
**ClientSampleId :**  
DDC46DL