

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
 Data File : BN031129.D  
 Acq On : 22 Apr 2024 19:35  
 Operator : MA/JU  
 Sample : P2037-23  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 E1CX8

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-BN041924.MA.M  
 Title : SVOA CALIBRATION

Signal : TIC: BN031129.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         | 3.417       | 69            | 73          | 86           | rBV      | 24243          | 49584         | 2.11%           | 0.126%        |
| 2         | 3.564       | 96            | 98          | 112          | rBV      | 53140          | 120257        | 5.12%           | 0.306%        |
| 3         | 4.228       | 206           | 211         | 220          | rVB6     | 12235          | 29708         | 1.27%           | 0.076%        |
| 4         | 4.481       | 250           | 254         | 259          | rVB      | 15554          | 21988         | 0.94%           | 0.056%        |
| 5         | 4.622       | 272           | 278         | 280          | rBV2     | 53979          | 100791        | 4.29%           | 0.256%        |
| 6         | 4.664       | 280           | 285         | 299          | rVB4     | 89938          | 253091        | 10.78%          | 0.644%        |
| 7         | 4.787       | 299           | 306         | 308          | rBV3     | 57234          | 106960        | 4.55%           | 0.272%        |
| 8         | 4.828       | 309           | 313         | 318          | rVB      | 62729          | 91167         | 3.88%           | 0.232%        |
| 9         | 4.911       | 318           | 327         | 339          | rVB3     | 181929         | 597155        | 25.43%          | 1.519%        |
| 10        | 5.011       | 340           | 344         | 346          | rBV2     | 36160          | 51346         | 2.19%           | 0.131%        |
| 11        | 5.069       | 351           | 354         | 360          | rVV      | 101099         | 189140        | 8.05%           | 0.481%        |
| 12        | 5.116       | 360           | 362         | 367          | rVV3     | 27329          | 42796         | 1.82%           | 0.109%        |
| 13        | 5.181       | 367           | 373         | 379          | rVV2     | 72251          | 167638        | 7.14%           | 0.426%        |
| 14        | 5.234       | 379           | 382         | 390          | rVB      | 47881          | 91341         | 3.89%           | 0.232%        |
| 15        | 5.316       | 390           | 396         | 402          | rBV2     | 34586          | 79910         | 3.40%           | 0.203%        |
| 16        | 5.381       | 402           | 407         | 416          | rVV3     | 46873          | 132775        | 5.65%           | 0.338%        |
| 17        | 5.469       | 416           | 422         | 433          | rVB6     | 22244          | 62915         | 2.68%           | 0.160%        |
| 18        | 5.622       | 444           | 448         | 457          | rVB      | 34368          | 67871         | 2.89%           | 0.173%        |
| 19        | 6.510       | 595           | 599         | 607          | rVB9     | 10875          | 23925         | 1.02%           | 0.061%        |
| 20        | 6.810       | 638           | 650         | 663          | rBV      | 150990         | 289330        | 12.32%          | 0.736%        |
| 21        | 6.975       | 670           | 678         | 688          | rBV      | 184218         | 296542        | 12.63%          | 0.754%        |
| 22        | 7.163       | 705           | 710         | 719          | rVB      | 99952          | 171485        | 7.30%           | 0.436%        |
| 23        | 7.252       | 721           | 725         | 733          | rVV2     | 10531          | 17647         | 0.75%           | 0.045%        |
| 24        | 7.522       | 766           | 771         | 779          | rBV8     | 6916           | 16882         | 0.72%           | 0.043%        |
| 25        | 7.634       | 779           | 790         | 796          | rBV      | 261105         | 439605        | 18.72%          | 1.118%        |
| 26        | 7.705       | 796           | 802         | 815          | rVV      | 50019          | 127099        | 5.41%           | 0.323%        |
| 27        | 8.022       | 852           | 856         | 860          | rBV4     | 15013          | 24183         | 1.03%           | 0.062%        |
| 28        | 8.281       | 891           | 900         | 904          | rBV3     | 30550          | 55542         | 2.37%           | 0.141%        |
| 29        | 8.340       | 905           | 910         | 915          | rBV      | 69087          | 131049        | 5.58%           | 0.333%        |
| 30        | 8.452       | 924           | 929         | 939          | rVB2     | 85616          | 140770        | 5.99%           | 0.358%        |
| 31        | 8.587       | 946           | 952         | 960          | rBV3     | 40058          | 66835         | 2.85%           | 0.170%        |
| 32        | 8.728       | 965           | 976         | 981          | rBV5     | 45664          | 131129        | 5.58%           | 0.334%        |
| 33        | 8.787       | 981           | 986         | 993          | rVV      | 232766         | 397936        | 16.94%          | 1.012%        |
| 34        | 8.846       | 993           | 996         | 1001         | rVB      | 48786          | 67972         | 2.89%           | 0.173%        |
| 35        | 8.963       | 1006          | 1016        | 1022         | rBV7     | 43282          | 117131        | 4.99%           | 0.298%        |
| 36        | 9.075       | 1028          | 1035        | 1039         | rBV10    | 14395          | 33873         | 1.44%           | 0.086%        |

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 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 E1CX8

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-BN041924.MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 9.128  | 1039 | 1044 | 1050 | rVB9 | 18863   | 42729   | 1.82%  | 0.109% |
| 38 | 9.263  | 1056 | 1067 | 1071 | rBV6 | 33700   | 110273  | 4.70%  | 0.281% |
| 39 | 9.304  | 1071 | 1074 | 1080 | rVV3 | 21639   | 48378   | 2.06%  | 0.123% |
| 40 | 9.357  | 1080 | 1083 | 1093 | rVB5 | 18741   | 36962   | 1.57%  | 0.094% |
| 41 | 9.504  | 1101 | 1108 | 1113 | rBV2 | 60682   | 112572  | 4.79%  | 0.286% |
| 42 | 9.587  | 1119 | 1122 | 1125 | rBV5 | 13433   | 15425   | 0.66%  | 0.039% |
| 43 | 9.763  | 1144 | 1152 | 1156 | rBV9 | 13444   | 25413   | 1.08%  | 0.065% |
| 44 | 9.822  | 1156 | 1162 | 1167 | rVV5 | 18765   | 44809   | 1.91%  | 0.114% |
| 45 | 9.881  | 1168 | 1172 | 1179 | rVB4 | 12680   | 26128   | 1.11%  | 0.066% |
| 46 | 10.040 | 1195 | 1199 | 1206 | rVB  | 66826   | 107820  | 4.59%  | 0.274% |
| 47 | 10.110 | 1206 | 1211 | 1218 | rBV  | 17683   | 31371   | 1.34%  | 0.080% |
| 48 | 10.246 | 1228 | 1234 | 1243 | rBV  | 237526  | 419287  | 17.85% | 1.067% |
| 49 | 10.410 | 1252 | 1262 | 1271 | rBV  | 413506  | 831719  | 35.42% | 2.116% |
| 50 | 10.557 | 1278 | 1287 | 1295 | rBV2 | 240669  | 462108  | 19.68% | 1.176% |
| 51 | 10.628 | 1297 | 1299 | 1307 | rVB6 | 17002   | 28121   | 1.20%  | 0.072% |
| 52 | 10.834 | 1328 | 1334 | 1338 | rBV6 | 14339   | 27976   | 1.19%  | 0.071% |
| 53 | 10.963 | 1348 | 1356 | 1363 | rBV8 | 16941   | 51667   | 2.20%  | 0.131% |
| 54 | 11.157 | 1384 | 1389 | 1397 | rVB2 | 47114   | 86339   | 3.68%  | 0.220% |
| 55 | 11.240 | 1400 | 1403 | 1411 | rVV7 | 17112   | 32217   | 1.37%  | 0.082% |
| 56 | 11.322 | 1413 | 1417 | 1424 | rVB7 | 26252   | 53629   | 2.28%  | 0.136% |
| 57 | 11.434 | 1430 | 1436 | 1445 | rBV  | 61577   | 128481  | 5.47%  | 0.327% |
| 58 | 11.610 | 1457 | 1466 | 1471 | rBV3 | 42727   | 108594  | 4.62%  | 0.276% |
| 59 | 11.687 | 1475 | 1479 | 1485 | rVB7 | 29386   | 55728   | 2.37%  | 0.142% |
| 60 | 11.840 | 1499 | 1505 | 1515 | rBV  | 153878  | 264237  | 11.25% | 0.672% |
| 61 | 12.151 | 1553 | 1558 | 1564 | rBV6 | 27732   | 60512   | 2.58%  | 0.154% |
| 62 | 12.240 | 1565 | 1573 | 1578 | rBV  | 181916  | 328797  | 14.00% | 0.836% |
| 63 | 12.304 | 1580 | 1584 | 1590 | rVV2 | 51147   | 91569   | 3.90%  | 0.233% |
| 64 | 12.387 | 1592 | 1598 | 1602 | rVV  | 392813  | 654450  | 27.87% | 1.665% |
| 65 | 12.440 | 1602 | 1607 | 1616 | rVV  | 620523  | 1448986 | 61.70% | 3.686% |
| 66 | 12.516 | 1617 | 1620 | 1626 | rVB3 | 90954   | 146319  | 6.23%  | 0.372% |
| 67 | 12.640 | 1635 | 1641 | 1646 | rBV2 | 333222  | 758685  | 32.31% | 1.930% |
| 68 | 12.693 | 1646 | 1650 | 1656 | rVB2 | 208898  | 406895  | 17.33% | 1.035% |
| 69 | 12.751 | 1657 | 1660 | 1664 | rBV2 | 53475   | 70267   | 2.99%  | 0.179% |
| 70 | 12.804 | 1665 | 1669 | 1673 | rBV  | 234235  | 307432  | 13.09% | 0.782% |
| 71 | 12.857 | 1673 | 1678 | 1680 | rBV2 | 278345  | 465994  | 19.84% | 1.185% |
| 72 | 12.957 | 1691 | 1695 | 1698 | rBV  | 254772  | 392459  | 16.71% | 0.998% |
| 73 | 13.034 | 1704 | 1708 | 1714 | rVB3 | 199272  | 341813  | 14.55% | 0.870% |
| 74 | 13.098 | 1715 | 1719 | 1725 | rVB  | 239289  | 335283  | 14.28% | 0.853% |
| 75 | 13.587 | 1798 | 1802 | 1807 | rBV5 | 105561  | 188549  | 8.03%  | 0.480% |
| 76 | 13.692 | 1815 | 1820 | 1825 | rBV  | 1120108 | 1534217 | 65.33% | 3.903% |

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Instrument :  
 BNA\_N  
 ClientSampleId :  
 E1CX8

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

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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-BN041924.MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |        |
|-----|--------|------|------|------|------|---------|---------|---------|--------|
| 77  | 13.763 | 1828 | 1832 | 1835 | rVV2 | 238991  | 327548  | 13.95%  | 0.833% |
| 78  | 13.804 | 1836 | 1839 | 1847 | rVB3 | 170726  | 309941  | 13.20%  | 0.788% |
| 79  | 13.928 | 1856 | 1860 | 1862 | rBV4 | 65071   | 98631   | 4.20%   | 0.251% |
| 80  | 13.969 | 1862 | 1867 | 1871 | rVV  | 976751  | 1406132 | 59.88%  | 3.577% |
| 81  | 14.010 | 1871 | 1874 | 1880 | rVB5 | 178797  | 308767  | 13.15%  | 0.785% |
| 82  | 14.104 | 1885 | 1890 | 1897 | rVB2 | 690021  | 1089008 | 46.37%  | 2.770% |
| 83  | 14.181 | 1898 | 1903 | 1907 | rBV  | 481775  | 715416  | 30.46%  | 1.820% |
| 84  | 14.239 | 1908 | 1913 | 1916 | rVV2 | 298686  | 587164  | 25.00%  | 1.494% |
| 85  | 14.281 | 1916 | 1920 | 1927 | rVB  | 868747  | 1319967 | 56.21%  | 3.358% |
| 86  | 14.687 | 1986 | 1989 | 1992 | rVB3 | 119763  | 144667  | 6.16%   | 0.368% |
| 87  | 14.804 | 2005 | 2009 | 2016 | rBV5 | 307500  | 591427  | 25.18%  | 1.505% |
| 88  | 14.904 | 2023 | 2026 | 2028 | rBV3 | 170160  | 247429  | 10.54%  | 0.629% |
| 89  | 14.998 | 2037 | 2042 | 2048 | rBV6 | 268646  | 597512  | 25.44%  | 1.520% |
| 90  | 15.063 | 2048 | 2053 | 2058 | rBV3 | 825670  | 1337723 | 56.96%  | 3.403% |
| 91  | 15.139 | 2063 | 2066 | 2069 | rBV  | 341801  | 405733  | 17.28%  | 1.032% |
| 92  | 15.275 | 2085 | 2089 | 2094 | rVB  | 1328903 | 1848197 | 78.70%  | 4.702% |
| 93  | 15.322 | 2094 | 2097 | 2100 | rBV  | 200955  | 263379  | 11.22%  | 0.670% |
| 94  | 15.545 | 2132 | 2135 | 2141 | rVB  | 475408  | 701365  | 29.87%  | 1.784% |
| 95  | 16.010 | 2209 | 2214 | 2215 | rBV  | 1280836 | 1879176 | 80.02%  | 4.781% |
| 96  | 16.851 | 2353 | 2357 | 2366 | rVB  | 1454485 | 2348445 | 100.00% | 5.974% |
| 97  | 17.033 | 2384 | 2388 | 2392 | rBV  | 1249144 | 2047795 | 87.20%  | 5.209% |
| 98  | 17.133 | 2401 | 2405 | 2410 | rVB  | 1397756 | 2006102 | 85.42%  | 5.103% |
| 99  | 17.539 | 2471 | 2474 | 2479 | rVB  | 1470949 | 1865785 | 79.45%  | 4.746% |
| 100 | 18.233 | 2590 | 2592 | 2597 | rBV  | 1109796 | 1372133 | 58.43%  | 3.491% |

Sum of corrected areas: 39309020



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Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

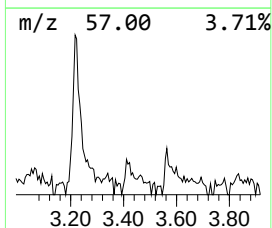
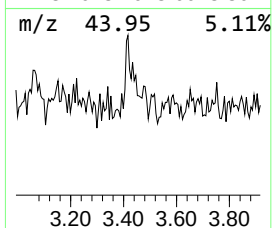
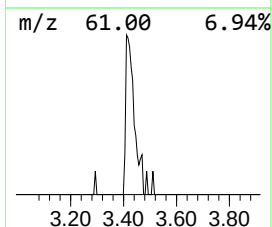
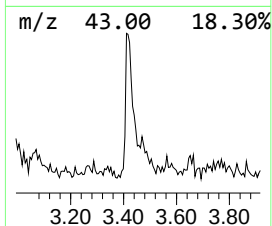
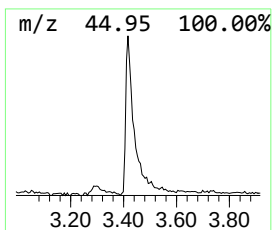
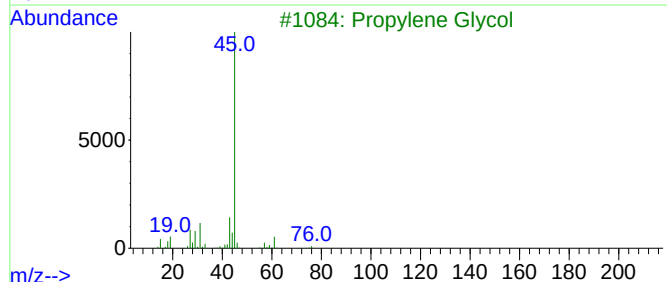
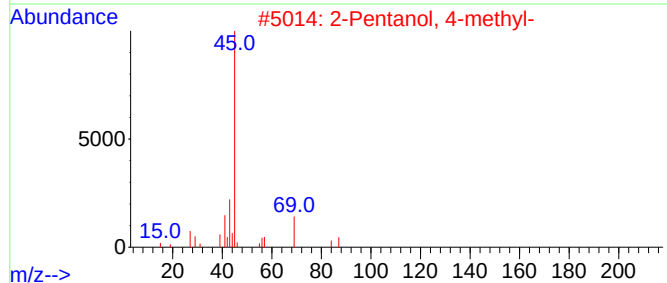
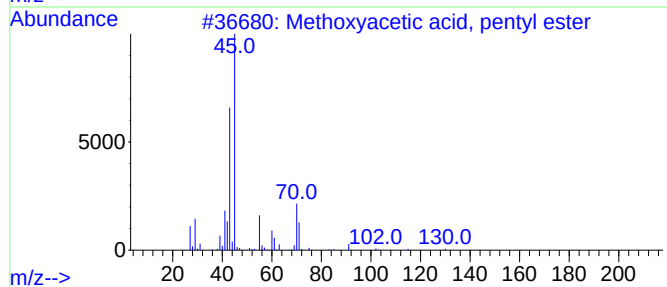
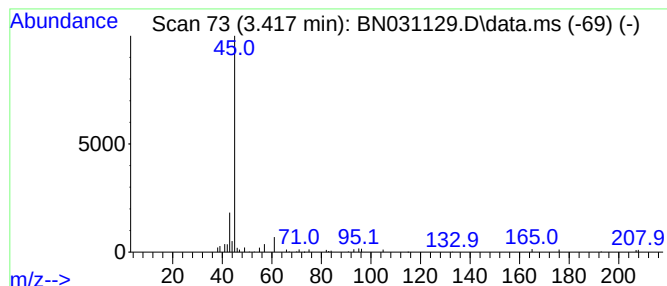
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 1 unknown-01 Concentration Rank 50

| R.T.  | EstConc    | Area  | Relative to ISTD       | R.T.  |
|-------|------------|-------|------------------------|-------|
| 3.417 | 2.26 ng/ul | 49584 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Methoxyacetic acid, pentyl ester | 160 | C8H16O3 | 168920-35-2 | 40   |
| 2    |    |   | 2-Pentanol, 4-methyl-            | 102 | C6H14O  | 000108-11-2 | 38   |
| 3    |    |   | Propylene Glycol                 | 76  | C3H8O2  | 000057-55-6 | 9    |
| 4    |    |   | 2-Hexanol                        | 102 | C6H14O  | 000626-93-7 | 9    |
| 5    |    |   | Lactic acid                      | 90  | C3H6O3  | 000050-21-5 | 9    |



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Quant Title : SVOA CALIBRATION

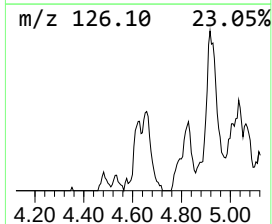
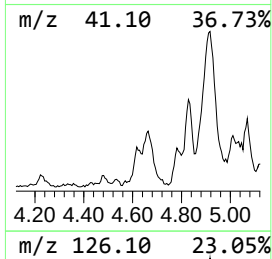
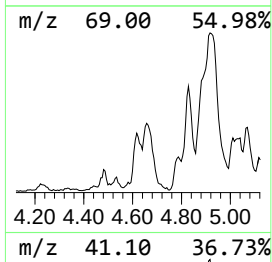
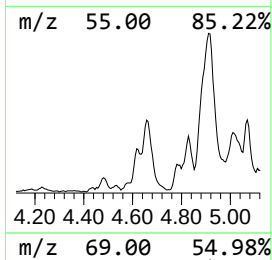
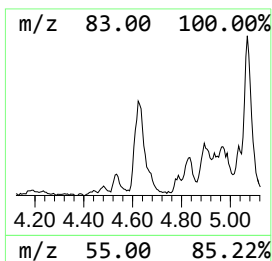
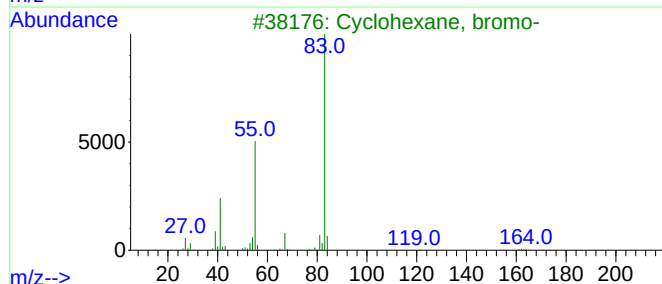
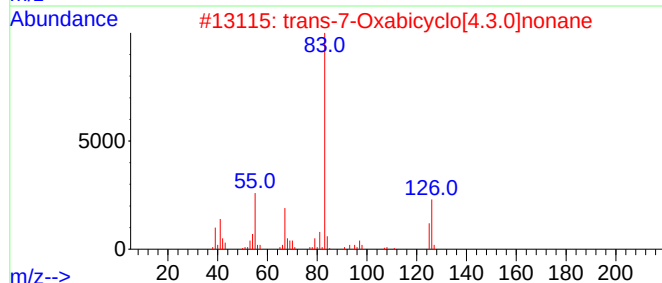
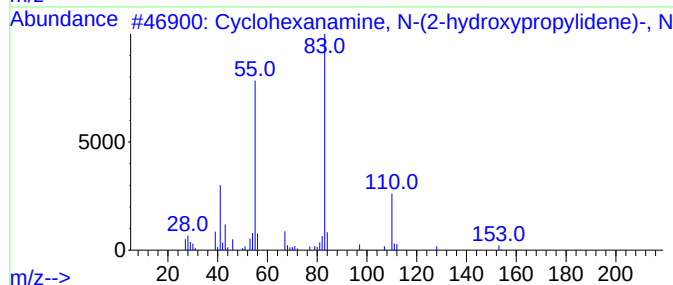
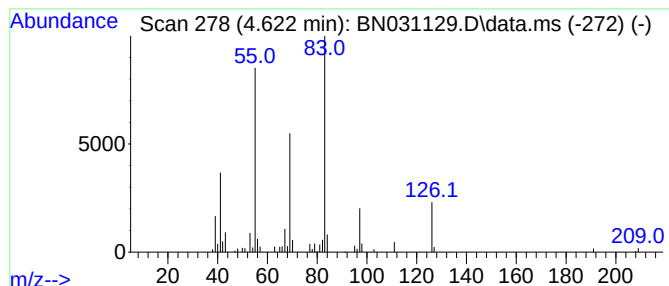
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Cyclohexanamine, N-(2-hydro... Concentration Rank 35

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 4.622 | 4.59 ng/ul | 100791 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Cyclohexanamine, N-(2-hydroxypro... | 171 | C9H17NO2 | 160423-87-0  | 53   |
| 2    |    |   | trans-7-Oxabicyclo[4.3.0]nonane     | 126 | C8H14O   | 027345-70-6  | 52   |
| 3    |    |   | Cyclohexane, bromo-                 | 162 | C6H11Br  | 000108-85-0  | 50   |
| 4    |    |   | Oxalic acid, dicyclohexyl ester     | 254 | C14H22O4 | 1000309-30-9 | 50   |
| 5    |    |   | 2-Pentene, 2,4-dimethyl-            | 98  | C7H14    | 000625-65-0  | 50   |



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E1CX8

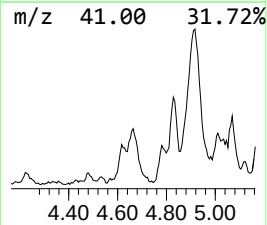
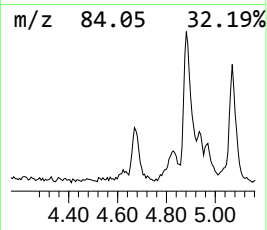
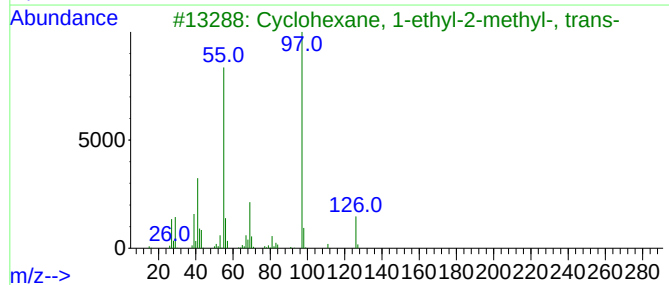
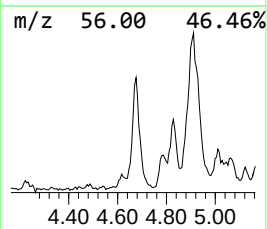
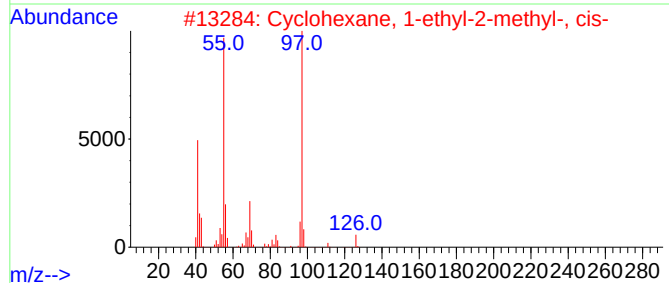
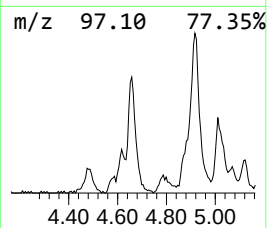
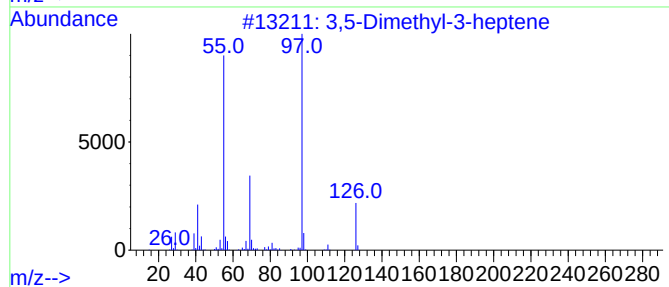
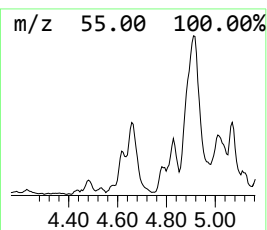
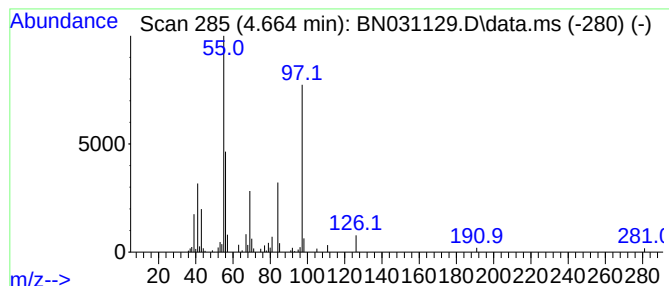
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 4.664 | 11.51 ng/ul | 253091 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | 3,5-Dimethyl-3-heptene              | 126 | C9H18   | 059643-68-4 | 52   |
| 2    |      | Cyclohexane, 1-ethyl-2-methyl-, ... | 126 | C9H18   | 004923-77-7 | 50   |
| 3    |      | Cyclohexane, 1-ethyl-2-methyl-, ... | 126 | C9H18   | 004923-78-8 | 50   |
| 4    |      | 3,5-Dimethyl-3-heptene              | 126 | C9H18   | 059643-68-4 | 50   |
| 5    |      | 1-Ethyl-4-methylcyclohexane         | 126 | C9H18   | 003728-56-1 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

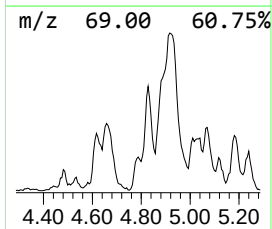
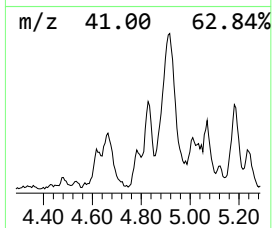
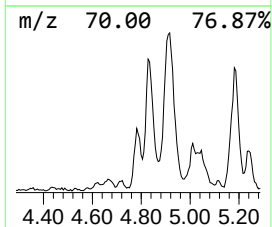
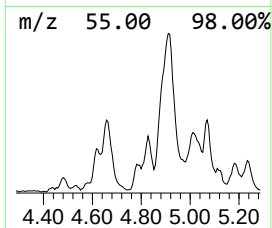
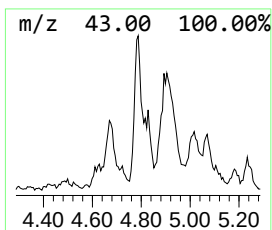
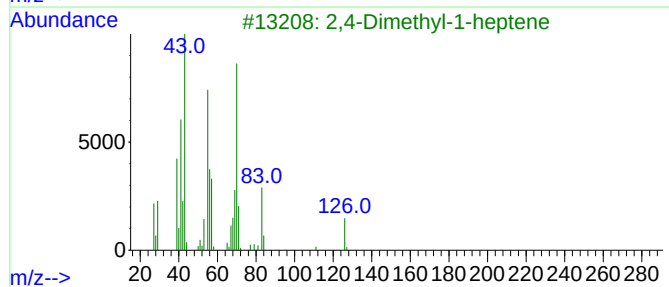
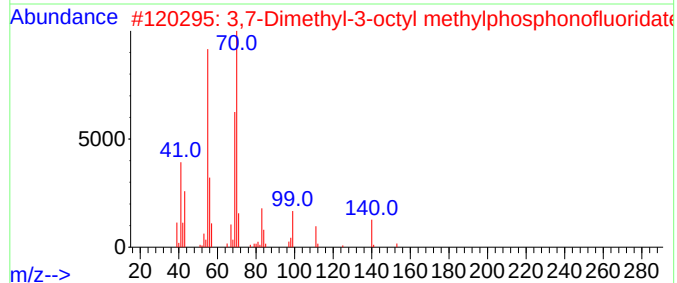
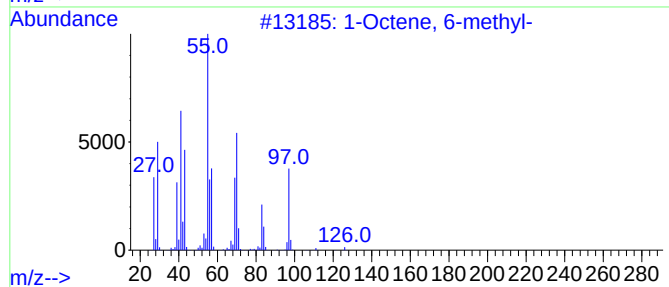
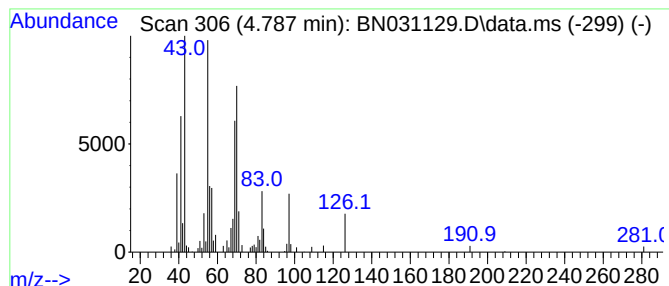
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 32

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 4.787 | 4.87 ng/ul | 106960 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|------|-------------------------------------|-----|------------|-------------|------|
| 1    |      | 1-Octene, 6-methyl-                 | 126 | C9H18      | 013151-10-5 | 78   |
| 2    |      | 3,7-Dimethyl-3-octyl methylphosp... | 238 | C11H24F02P | 159395-79-6 | 59   |
| 3    |      | 2,4-Dimethyl-1-heptene              | 126 | C9H18      | 019549-87-2 | 58   |
| 4    |      | 1.alpha.,2.beta.,3.alpha.,4.beta... | 126 | C9H18      | 002532-67-4 | 58   |
| 5    |      | Cyclohexane, 1,2,3-trimethyl-       | 126 | C9H18      | 001678-97-3 | 49   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

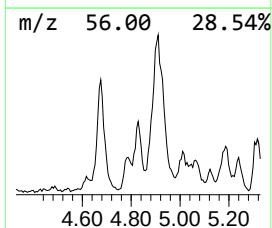
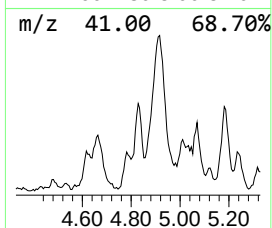
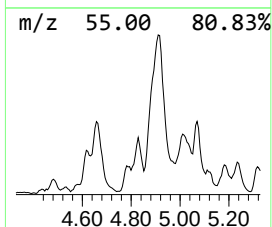
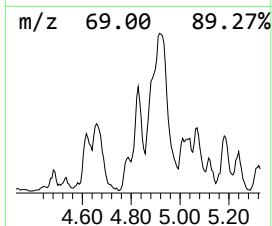
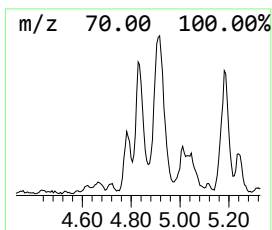
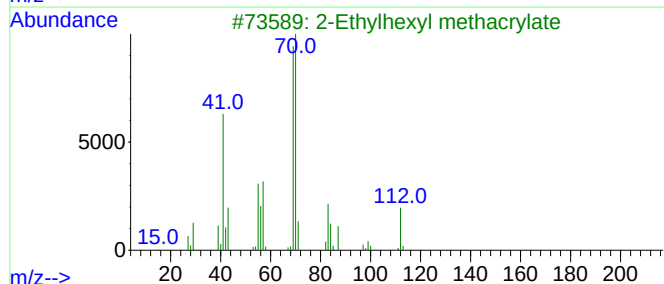
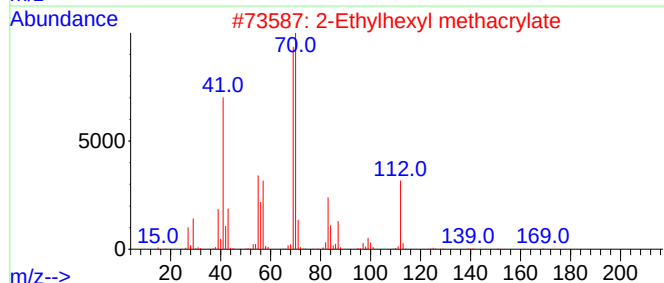
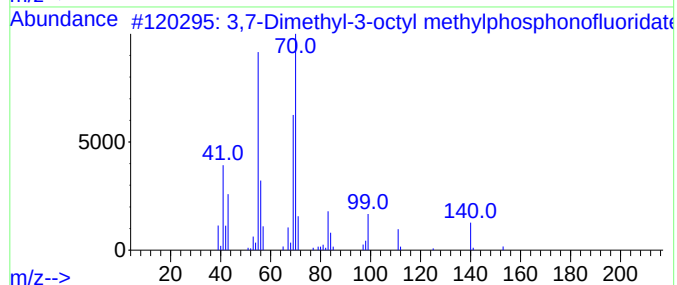
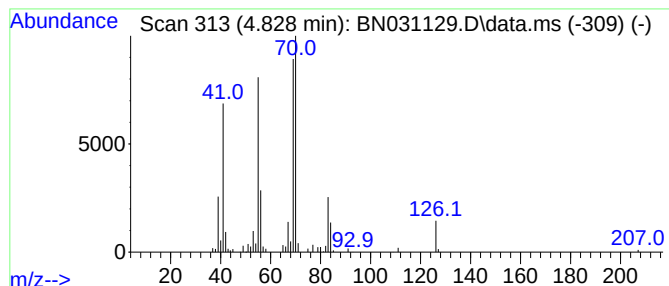
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 3,7-Dimethyl-3-octyl methyl... Concentration Rank 37

| R.T.  | EstConc    | Area  | Relative to ISTD       | R.T.  |
|-------|------------|-------|------------------------|-------|
| 4.828 | 4.15 ng/ul | 91167 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|------------|-------------|------|
| 1    |    |   | 3,7-Dimethyl-3-octyl methylphosp... | 238 | C11H24F02P | 159395-79-6 | 78   |
| 2    |    |   | 2-Ethylhexyl methacrylate           | 198 | C12H22O2   | 000688-84-6 | 64   |
| 3    |    |   | 2-Ethylhexyl methacrylate           | 198 | C12H22O2   | 000688-84-6 | 59   |
| 4    |    |   | Butane, 2-cyclopropyl-              | 98  | C7H14      | 005750-02-7 | 47   |
| 5    |    |   | Pentane, 3-methylene-               | 84  | C6H12      | 000760-21-4 | 46   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

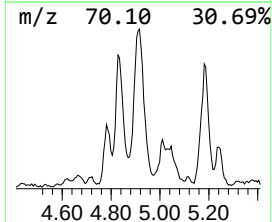
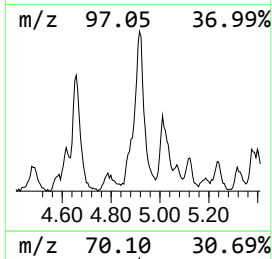
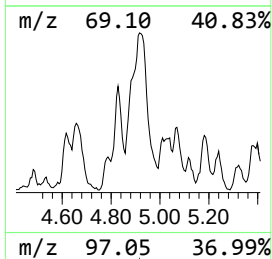
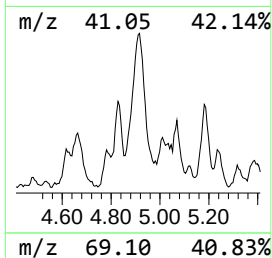
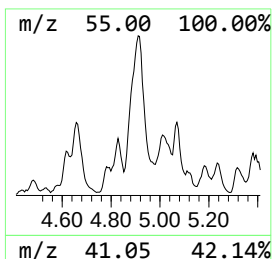
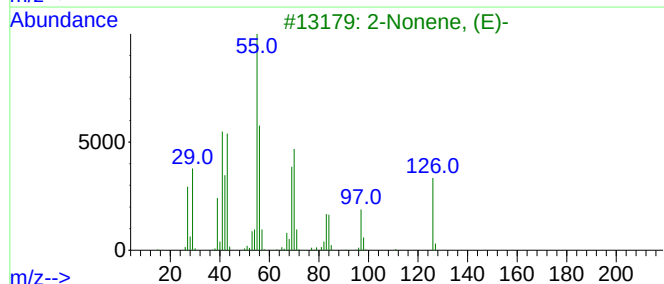
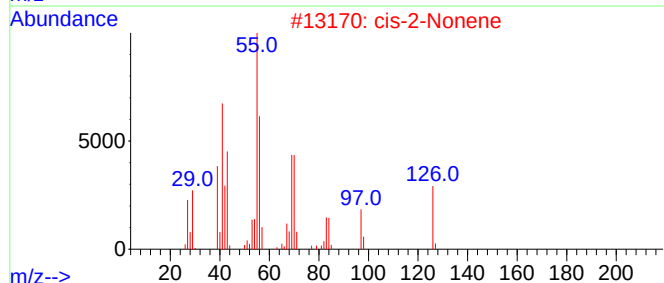
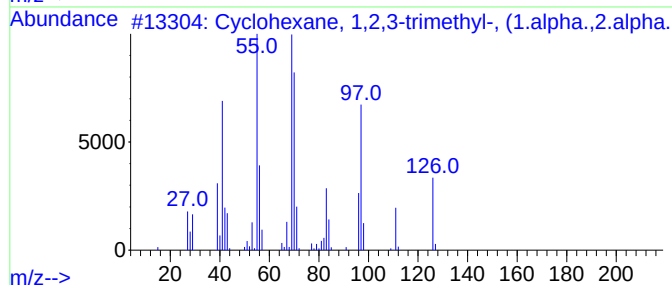
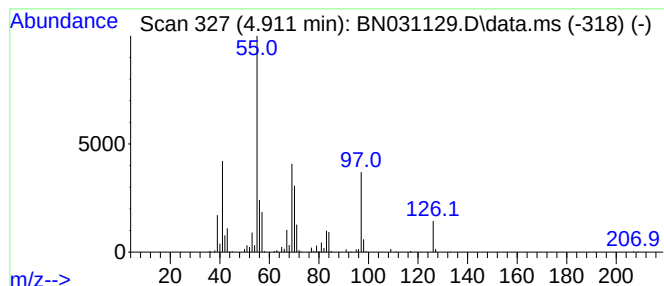
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Quant Title : SVOA CALIBRATION

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

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Peak Number 6 (DEL) Alkane: Cyclic4.911 Concentration Rank 1

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 4.911 | 27.17 ng/ul | 597155 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, 1,2,3-trimethyl-, (... | 126 | C9H18   | 001839-88-9 | 72   |
| 2    |    |   | cis-2-Nonene                        | 126 | C9H18   | 006434-77-1 | 59   |
| 3    |    |   | 2-Nonene, (E)-                      | 126 | C9H18   | 006434-78-2 | 58   |
| 4    |    |   | 2-Nonene                            | 126 | C9H18   | 002216-38-8 | 58   |
| 5    |    |   | trans-1,2-Diethyl cyclopentane      | 126 | C9H18   | 000932-40-1 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
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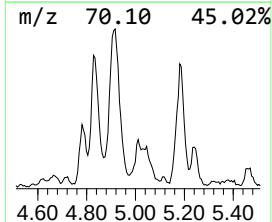
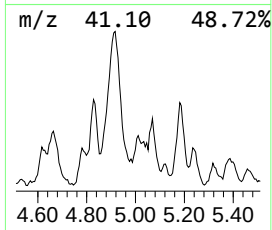
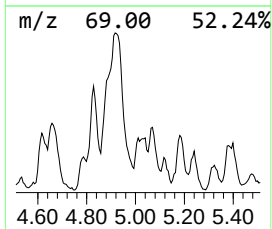
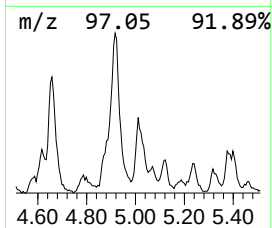
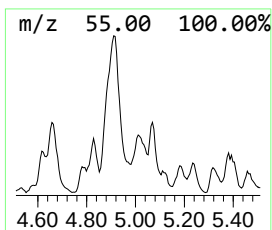
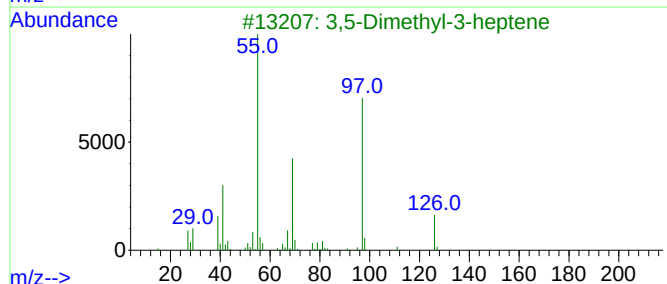
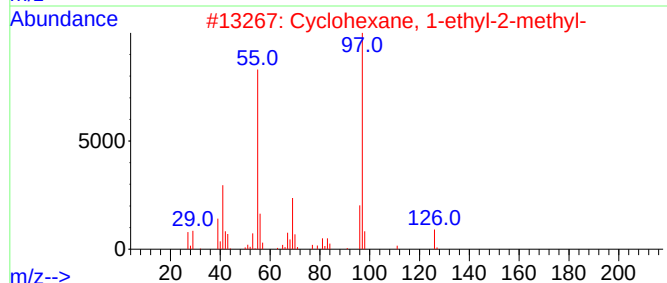
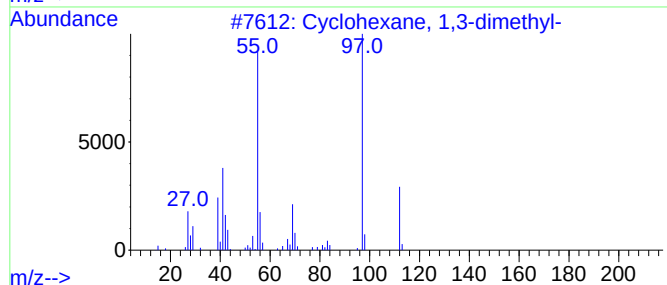
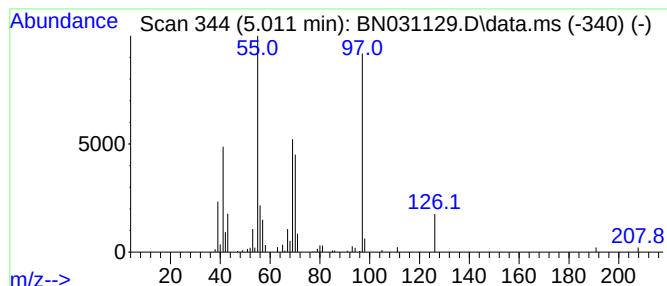
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Quant Title : SVOA CALIBRATION

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

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Peak Number 7 (DEL) Alkane: Cyclic5.011 Concentration Rank 49

| R.T.  | EstConc    | Area  | Relative to ISTD       | R.T.  |
|-------|------------|-------|------------------------|-------|
| 5.011 | 2.34 ng/ul | 51346 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, 1,3-dimethyl-          | 112 | C8H16   | 000591-21-9 | 59   |
| 2    |    |   | Cyclohexane, 1-ethyl-2-methyl-      | 126 | C9H18   | 003728-54-9 | 58   |
| 3    |    |   | 3,5-Dimethyl-3-heptene              | 126 | C9H18   | 059643-68-4 | 58   |
| 4    |    |   | Cyclohexane, 1-ethyl-4-methyl-, ... | 126 | C9H18   | 004926-78-7 | 58   |
| 5    |    |   | 3,5-Dimethyl-3-heptene              | 126 | C9H18   | 059643-68-4 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

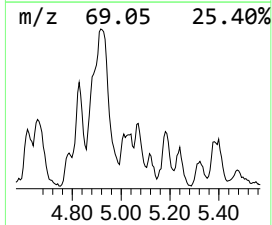
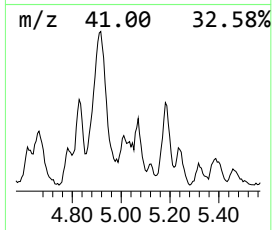
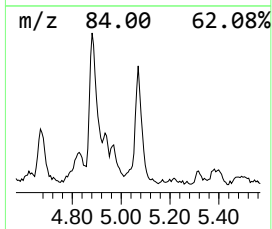
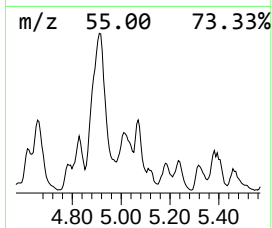
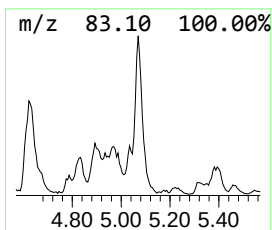
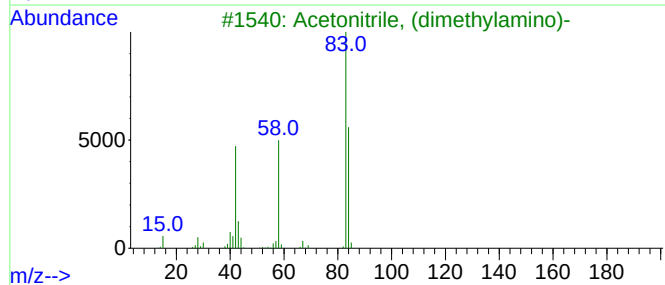
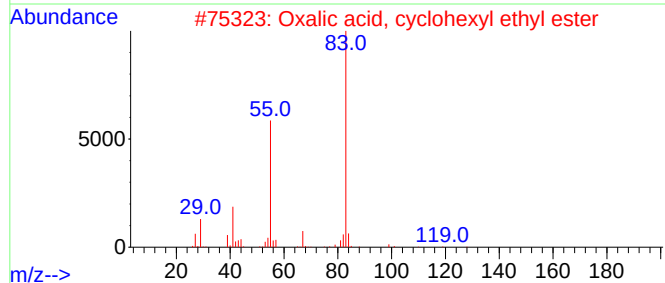
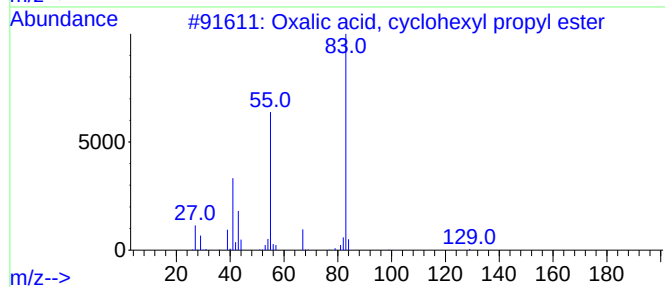
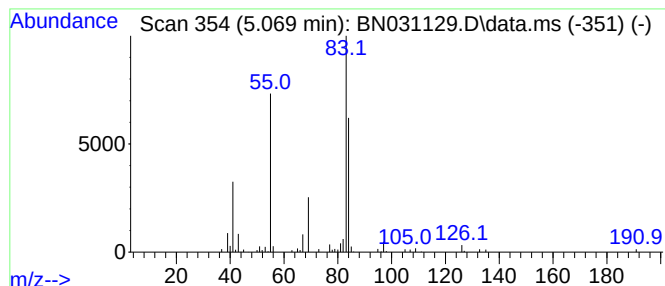
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Quant Title : SVOA CALIBRATION

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

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Peak Number 8 Oxalic acid, cyclohexyl pro... Concentration Rank 17

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 5.069 | 8.61 ng/ul | 189140 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Oxalic acid, cyclohexyl propyl e... | 214 | C11H18O4 | 1000309-30-3 | 50   |
| 2    |    |   | Oxalic acid, cyclohexyl ethyl ester | 200 | C10H16O4 | 1000309-30-2 | 50   |
| 3    |    |   | Acetonitrile, (dimethylamino)-      | 84  | C4H8N2   | 000926-64-7  | 45   |
| 4    |    |   | Cyclobutanecarboxylic acid, 6-et... | 240 | C15H28O2 | 1000282-61-1 | 43   |
| 5    |    |   | Oxalic acid, cyclohexyl butyl ester | 228 | C12H20O4 | 1000309-30-5 | 40   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

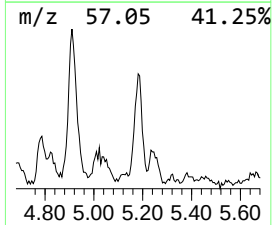
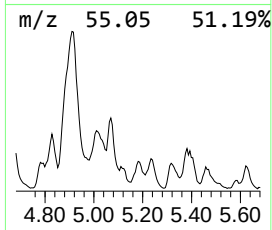
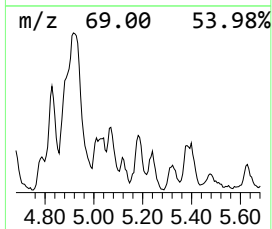
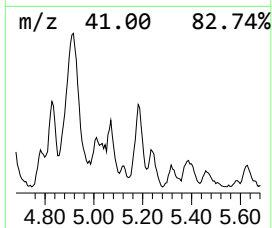
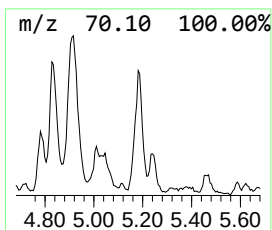
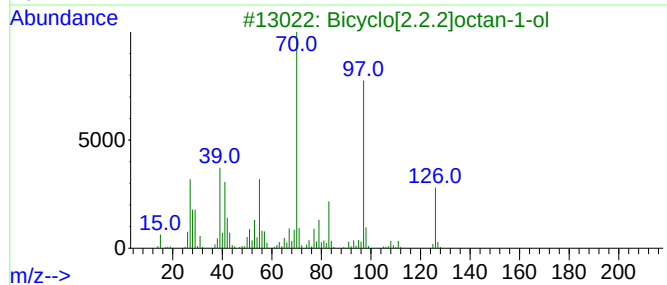
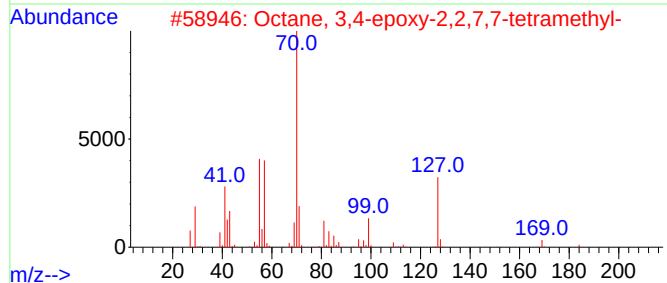
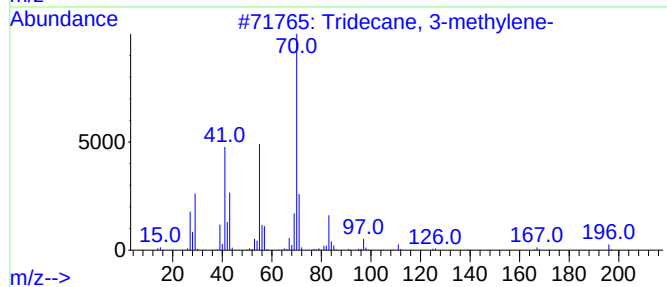
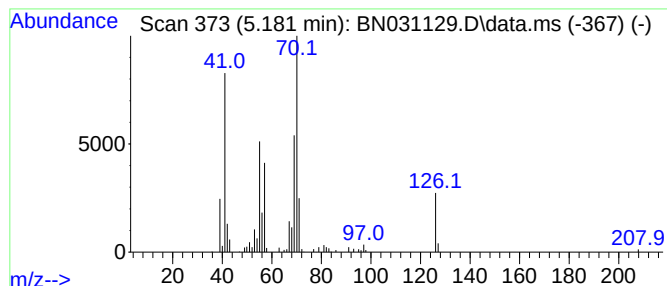
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Quant Title : SVOA CALIBRATION

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

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Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 19

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 5.181 | 7.63 ng/ul | 167638 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Tridecane, 3-methylene-             | 196 | C14H28   | 019780-34-8  | 53   |
| 2    |    |   | Octane, 3,4-epoxy-2,2,7,7-tetram... | 184 | C12H24O  | 1000163-73-3 | 45   |
| 3    |    |   | Bicyclo[2.2.2]octan-1-ol            | 126 | C8H14O   | 020534-58-1  | 42   |
| 4    |    |   | 1-Cyclopropylpropan-1-amine         | 99  | C6H13N   | 1000436-93-8 | 42   |
| 5    |    |   | Butanoic acid, 3-methyl-, 3-meth... | 172 | C10H20O2 | 000659-70-1  | 40   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

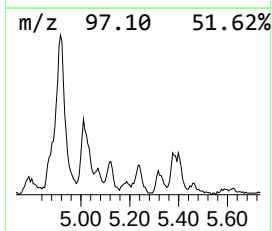
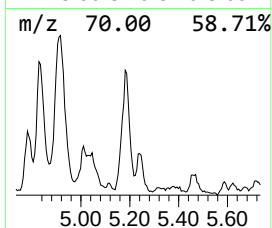
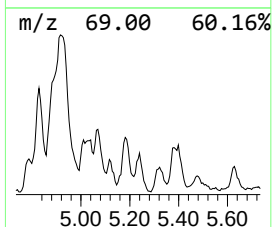
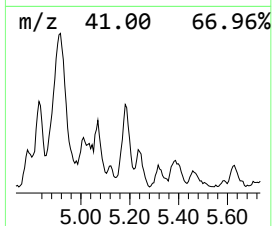
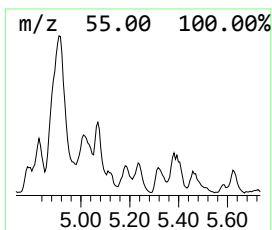
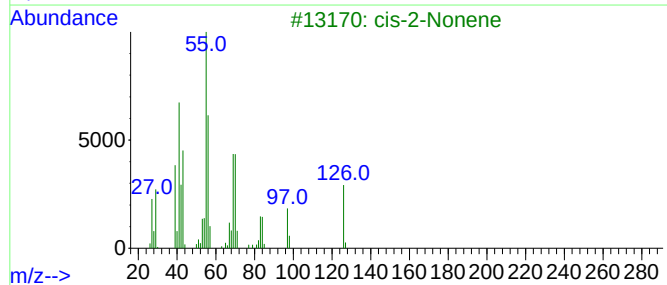
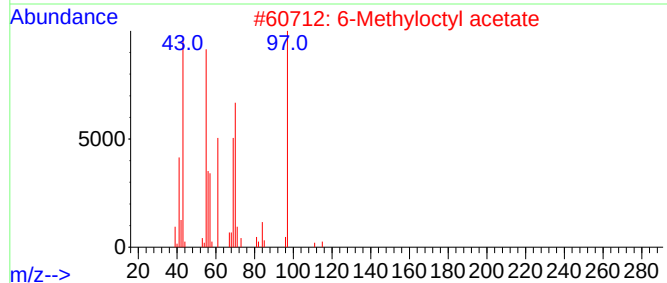
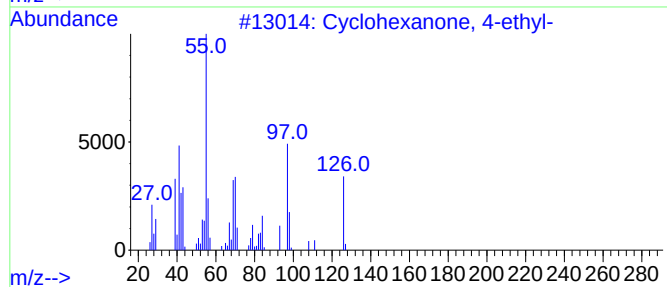
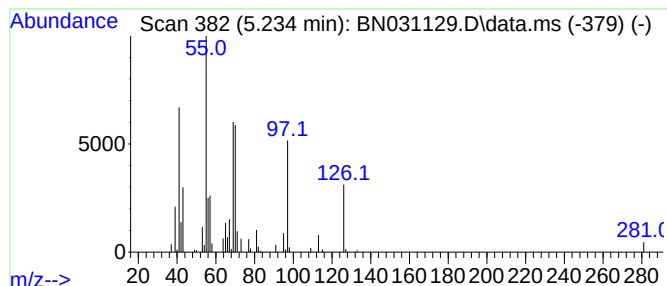
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TIC Integration Parameters: LSCINT.P

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Peak Number 10 Cyclohexanone, 4-ethyl- Concentration Rank 36

| R.T.  | EstConc    | Area  | Relative to ISTD       | R.T.  |
|-------|------------|-------|------------------------|-------|
| 5.234 | 4.16 ng/ul | 91341 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------|-----|----------|--------------|------|
| 1    |    |   | Cyclohexanone, 4-ethyl- | 126 | C8H14O   | 005441-51-0  | 53   |
| 2    |    |   | 6-Methyloctyl acetate   | 186 | C11H22O2 | 1000439-66-5 | 53   |
| 3    |    |   | cis-2-Nonene            | 126 | C9H18    | 006434-77-1  | 43   |
| 4    |    |   | cis-3-Nonene            | 126 | C9H18    | 020237-46-1  | 43   |
| 5    |    |   | cis-2-Nonene            | 126 | C9H18    | 006434-77-1  | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

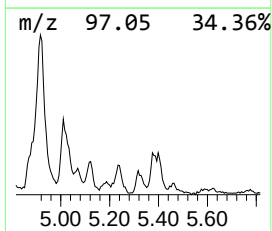
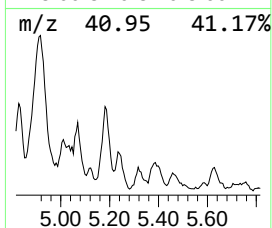
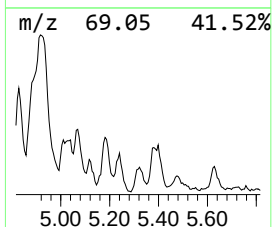
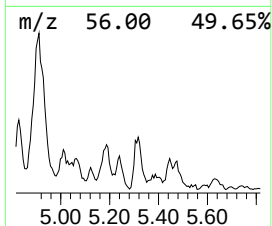
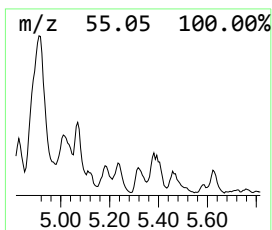
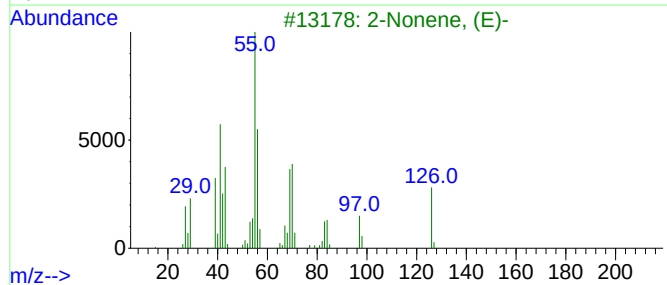
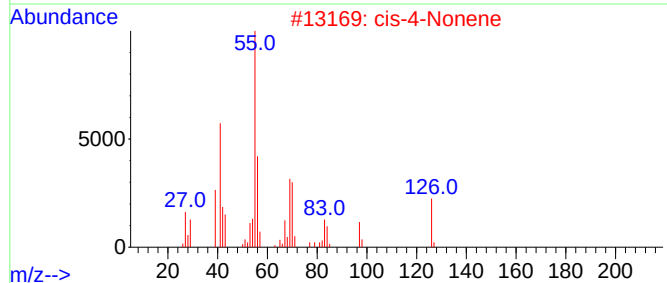
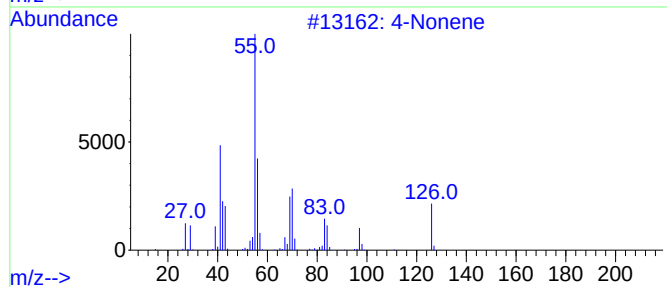
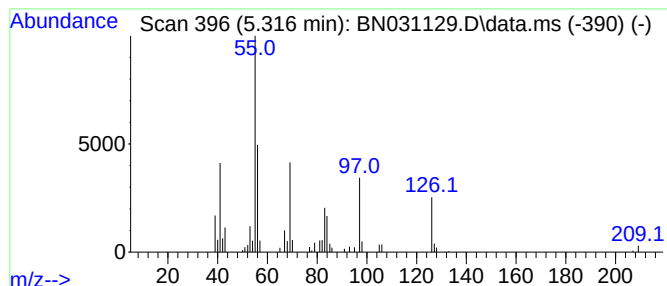
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Quant Title : SVOA CALIBRATION

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

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Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 39

| R.T.  | EstConc    | Area  | Relative to ISTD       | R.T.  |
|-------|------------|-------|------------------------|-------|
| 5.316 | 3.64 ng/ul | 79910 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of 5 | Tentative ID   | MW  | MolForm | CAS#        | Qual |
|------|------|----------------|-----|---------|-------------|------|
| 1    |      | 4-Nonene       | 126 | C9H18   | 002198-23-4 | 86   |
| 2    |      | cis-4-Nonene   | 126 | C9H18   | 010405-84-2 | 72   |
| 3    |      | 2-Nonene, (E)- | 126 | C9H18   | 006434-78-2 | 72   |
| 4    |      | 4-Nonene       | 126 | C9H18   | 002198-23-4 | 72   |
| 5    |      | cis-2-Nonene   | 126 | C9H18   | 006434-77-1 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

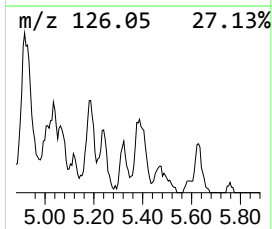
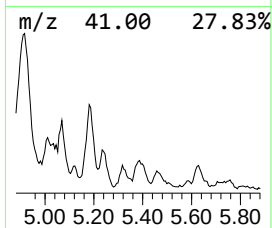
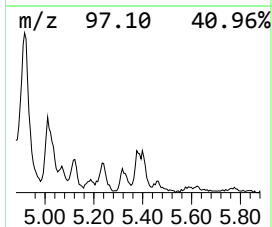
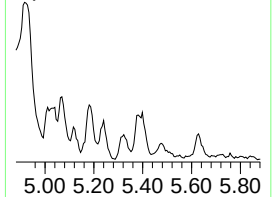
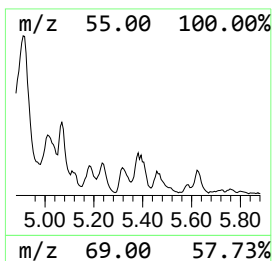
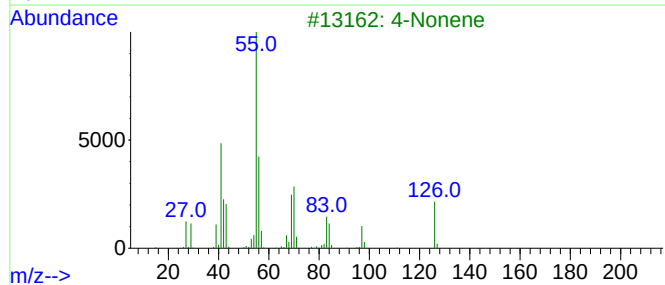
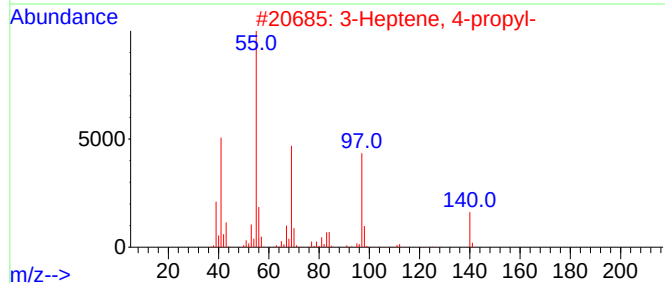
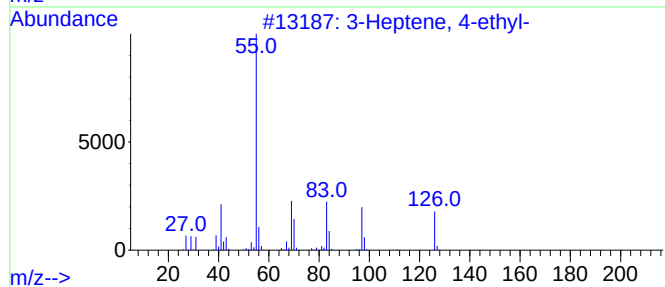
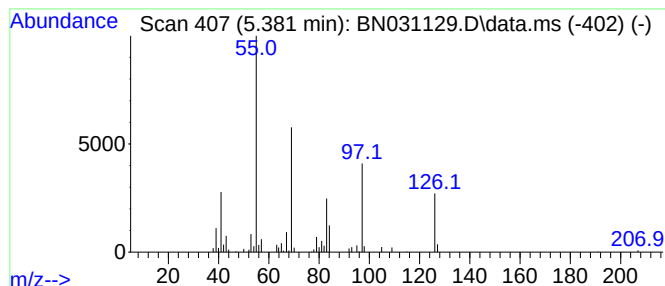
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 24

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 5.381 | 6.04 ng/ul | 132775 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|------|--------------------------|-----|---------|-------------|------|
| 1    |      | 3-Heptene, 4-ethyl-      | 126 | C9H18   | 033933-74-3 | 59   |
| 2    |      | 3-Heptene, 4-propyl-     | 140 | C10H20  | 004485-13-6 | 50   |
| 3    |      | 4-Nonene                 | 126 | C9H18   | 002198-23-4 | 50   |
| 4    |      | Cyclohexanone, 4-ethyl-  | 126 | C8H14O  | 005441-51-0 | 50   |
| 5    |      | 2,2,4-Trimethyl-3-hexene | 126 | C9H18   | 059643-72-0 | 50   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

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Quant Title : SVOA CALIBRATION

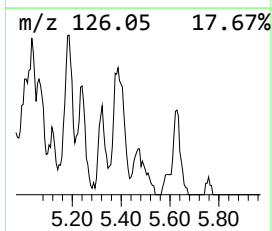
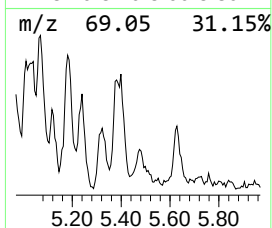
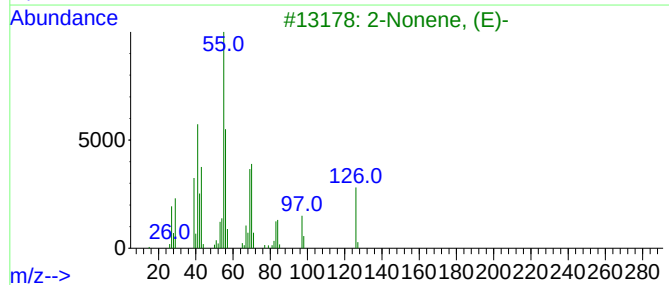
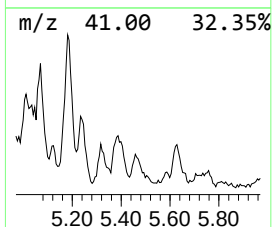
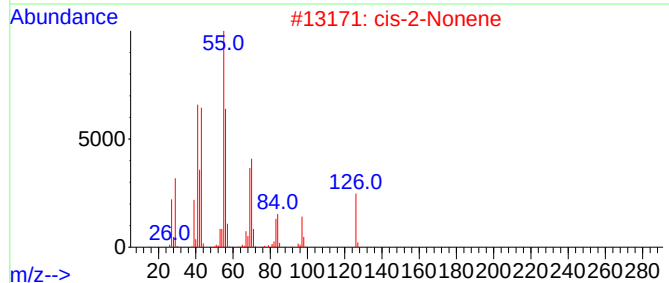
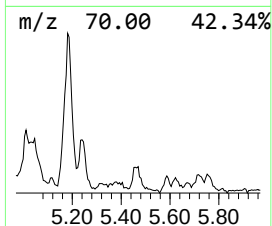
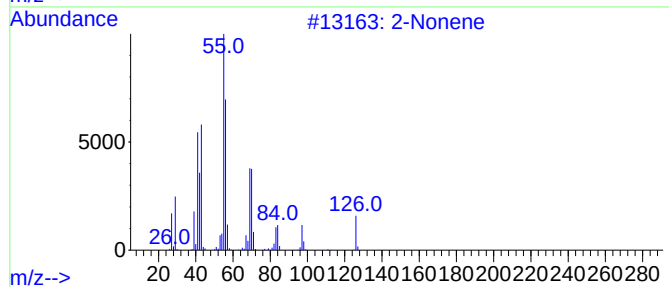
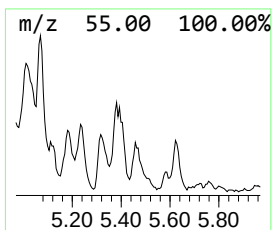
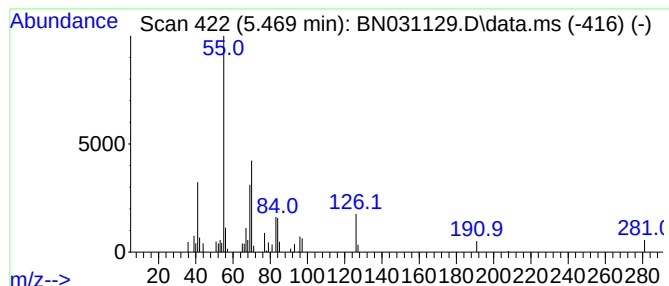
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TIC Integration Parameters: LSCINT.P

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Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 44

| R.T.  | EstConc    | Area  | Relative to ISTD       | R.T.  |
|-------|------------|-------|------------------------|-------|
| 5.469 | 2.86 ng/ul | 62915 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of              | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|-----------------|---|--------------|-----|---------|-------------|------|
| 1    | 2-Nonene        |   |              | 126 | C9H18   | 002216-38-8 | 72   |
| 2    | cis-2-Nonene    |   |              | 126 | C9H18   | 006434-77-1 | 64   |
| 3    | 2-Nonene, (E)-  |   |              | 126 | C9H18   | 006434-78-2 | 64   |
| 4    | trans--4-Nonene |   |              | 126 | C9H18   | 010405-85-3 | 64   |
| 5    | 4-Nonene        |   |              | 126 | C9H18   | 002198-23-4 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

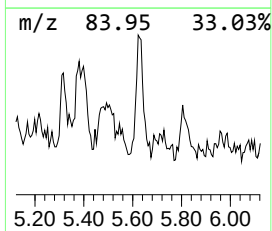
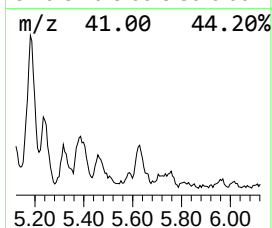
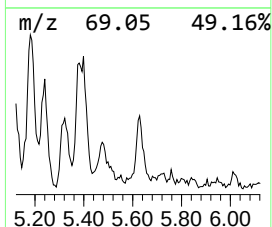
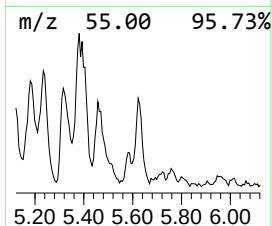
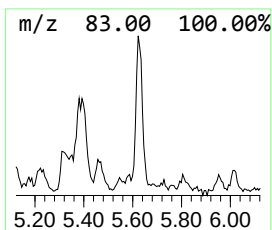
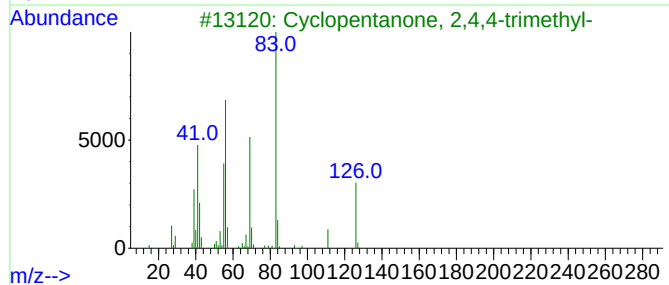
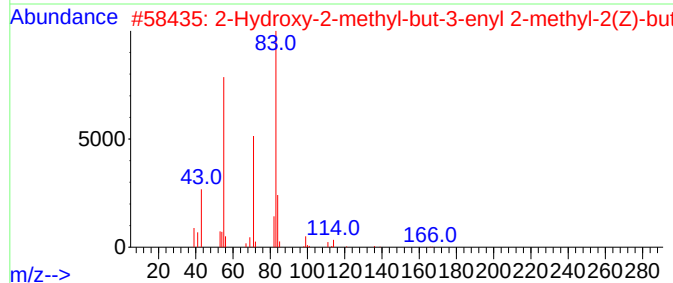
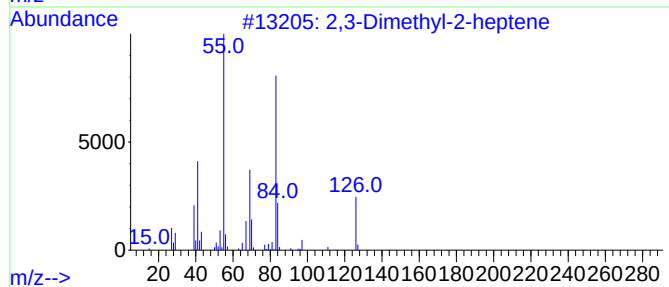
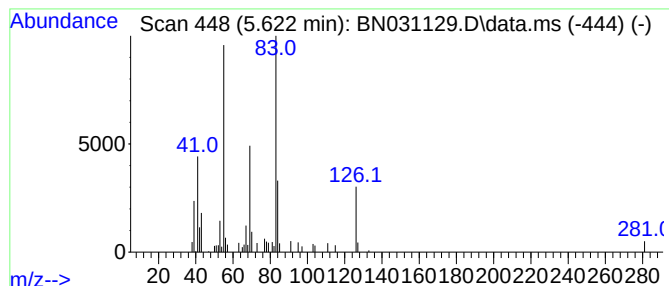
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 42

| R.T.  | EstConc    | Area  | Relative to ISTD       | R.T.  |
|-------|------------|-------|------------------------|-------|
| 5.622 | 3.09 ng/ul | 67871 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 2,3-Dimethyl-2-heptene              | 126 | C9H18     | 003074-64-4  | 83   |
| 2    |    |   | 2-Hydroxy-2-methyl-but-3-enyl 2-... | 184 | C10H16O3  | 080758-67-4  | 59   |
| 3    |    |   | Cyclopentanone, 2,4,4-trimethyl-    | 126 | C8H14O    | 004694-12-6  | 53   |
| 4    |    |   | 5-Oxa-6-azaspiro[3.4]oct-6-ene      | 111 | C6H9NO    | 1000362-18-0 | 50   |
| 5    |    |   | Phosphonous dibromide, cyclohexyl-  | 272 | C6H11Br2P | 086012-32-0  | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

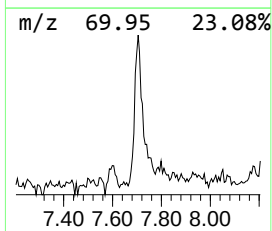
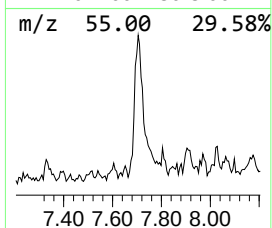
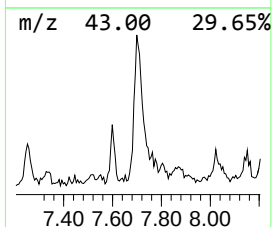
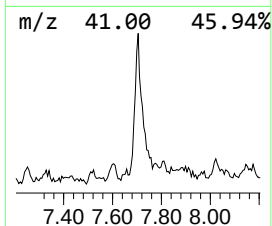
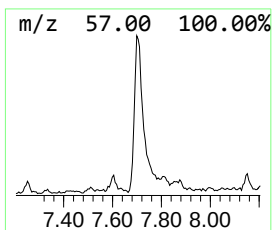
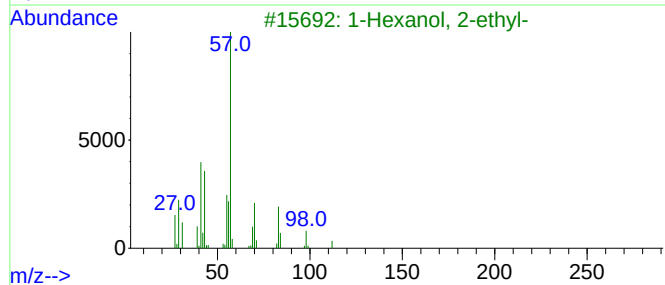
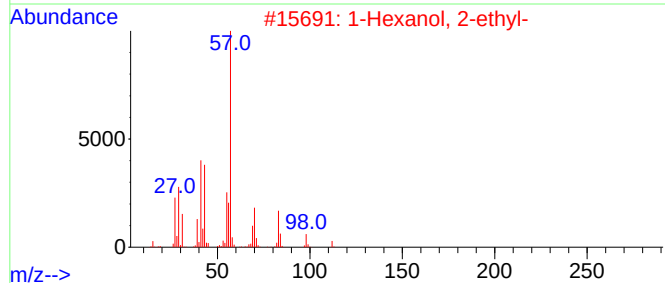
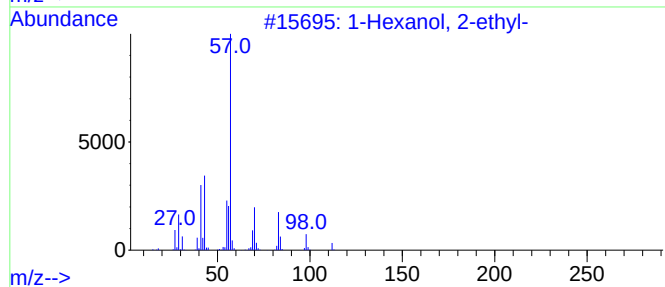
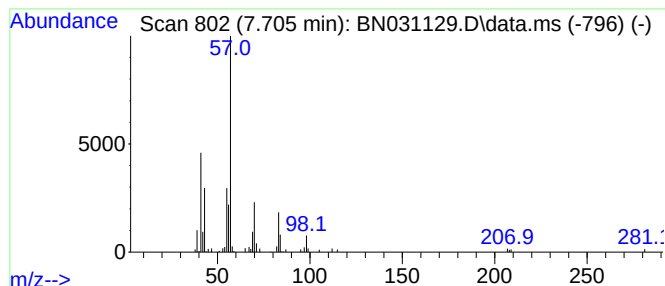
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Quant Title : SVOA CALIBRATION

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

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Peak Number 15 1-Hexanol, 2-ethyl- Concentration Rank 27

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 7.705 | 5.78 ng/ul | 127099 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 1-Hexanol, 2-ethyl-                 | 130 | C8H18O     | 000104-76-7  | 78   |
| 2    |    |   | 1-Hexanol, 2-ethyl-                 | 130 | C8H18O     | 000104-76-7  | 64   |
| 3    |    |   | 1-Hexanol, 2-ethyl-                 | 130 | C8H18O     | 000104-76-7  | 64   |
| 4    |    |   | 2-Propyl-1-pentanol                 | 130 | C8H18O     | 058175-57-8  | 53   |
| 5    |    |   | 2-Propyl-1-pentanol, pentafluoro... | 276 | C11H17F5O2 | 1000365-51-2 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

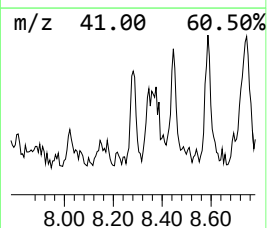
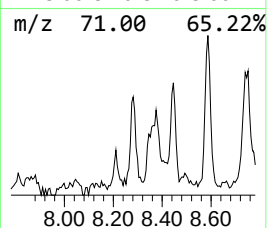
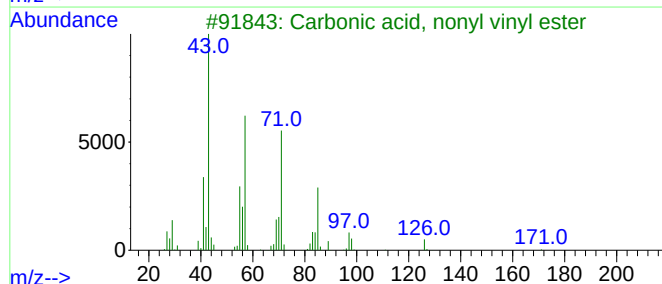
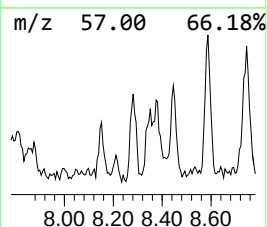
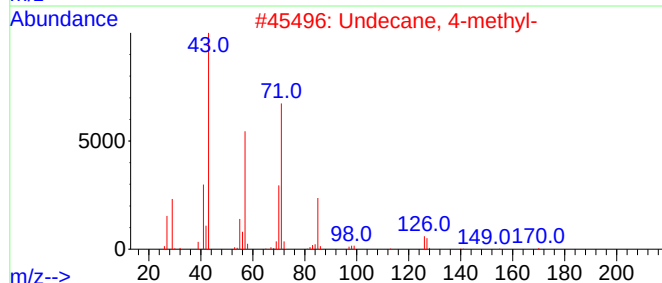
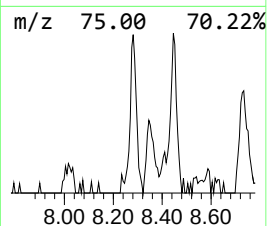
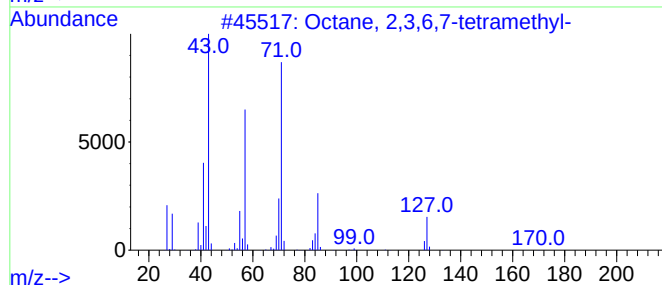
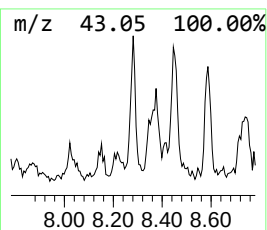
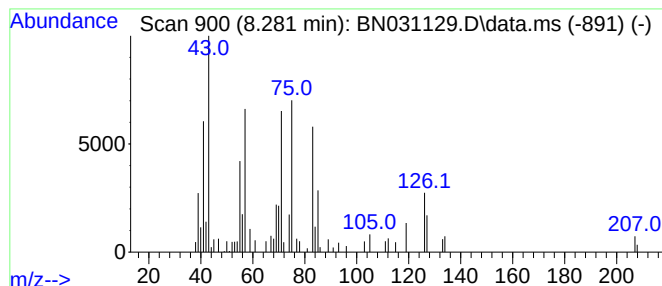
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TIC Integration Parameters: LSCINT.P

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Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 48

| R.T.  | EstConc    | Area  | Relative to ISTD       | R.T.  |
|-------|------------|-------|------------------------|-------|
| 8.281 | 2.53 ng/ul | 55542 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm  | CAS#         | Qual |
|------|----|---|----------------------------------|-----|----------|--------------|------|
| 1    |    |   | Octane, 2,3,6,7-tetramethyl-     | 170 | C12H26   | 052670-34-5  | 38   |
| 2    |    |   | Undecane, 4-methyl-              | 170 | C12H26   | 002980-69-0  | 35   |
| 3    |    |   | Carbonic acid, nonyl vinyl ester | 214 | C12H22O3 | 1000383-25-6 | 27   |
| 4    |    |   | Undecane                         | 156 | C11H24   | 001120-21-4  | 27   |
| 5    |    |   | Hexane, 3,3-dimethyl-            | 114 | C8H18    | 000563-16-6  | 27   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

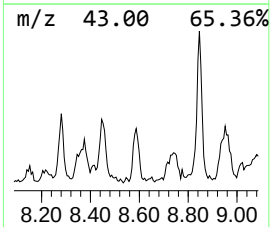
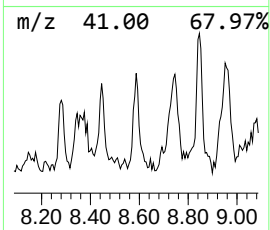
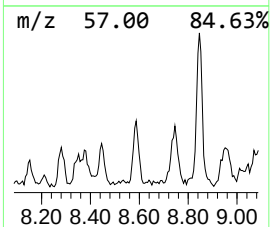
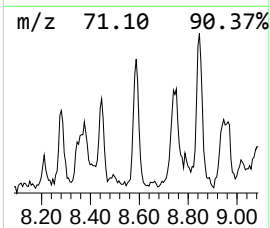
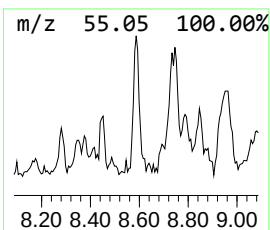
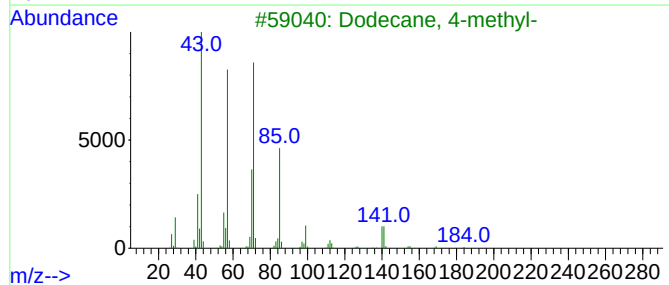
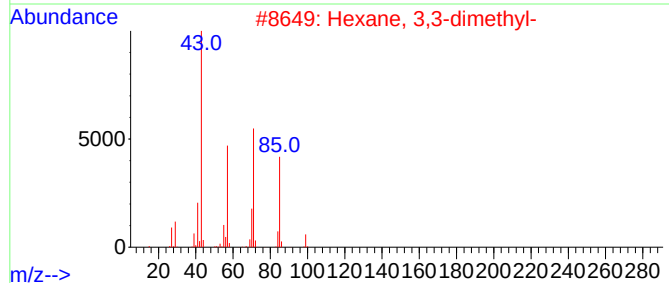
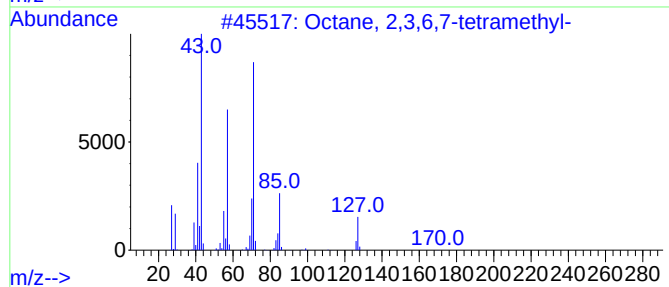
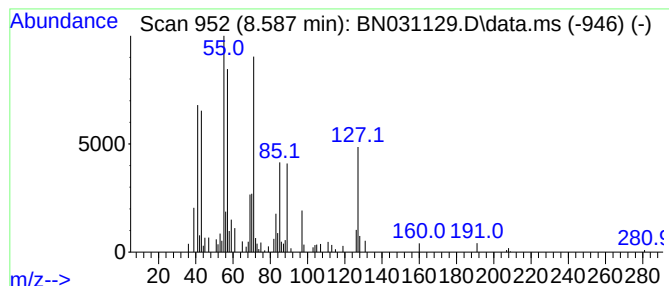
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TIC Integration Parameters: LSCINT.P

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Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 43

| R.T.  | EstConc    | Area  | Relative to ISTD       | R.T.  |
|-------|------------|-------|------------------------|-------|
| 8.587 | 3.04 ng/ul | 66835 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|------|-------------------------------------|-----|-----------|--------------|------|
| 1    |      | Octane, 2,3,6,7-tetramethyl-        | 170 | C12H26    | 052670-34-5  | 43   |
| 2    |      | Hexane, 3,3-dimethyl-               | 114 | C8H18     | 000563-16-6  | 43   |
| 3    |      | Dodecane, 4-methyl-                 | 184 | C13H28    | 006117-97-1  | 38   |
| 4    |      | Decane, 5-propyl-                   | 184 | C13H28    | 017312-62-8  | 38   |
| 5    |      | Sulfurous acid, hexyl 2-pentyl e... | 236 | C11H24O3S | 1000309-15-6 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

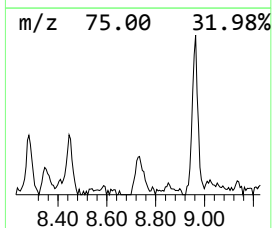
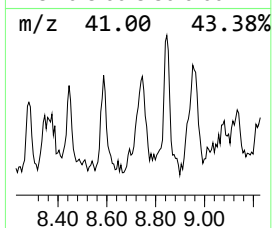
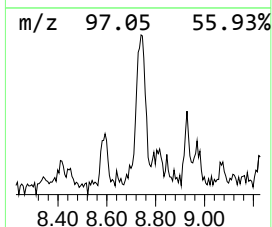
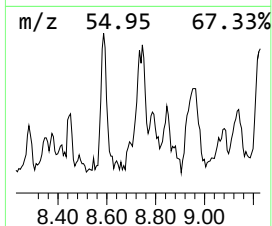
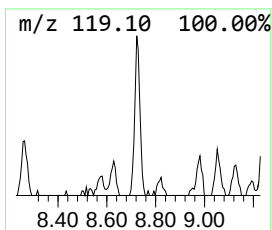
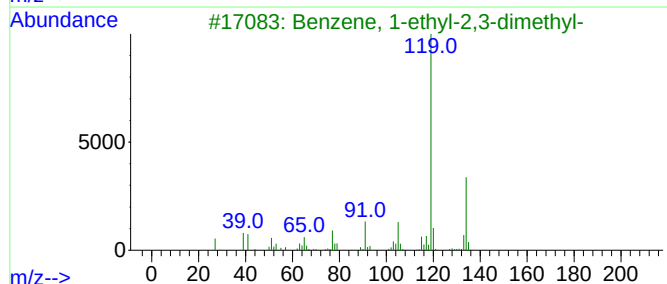
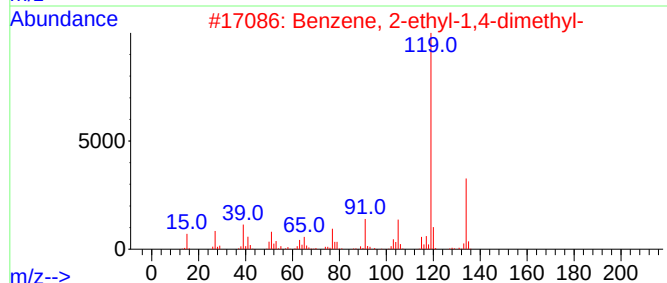
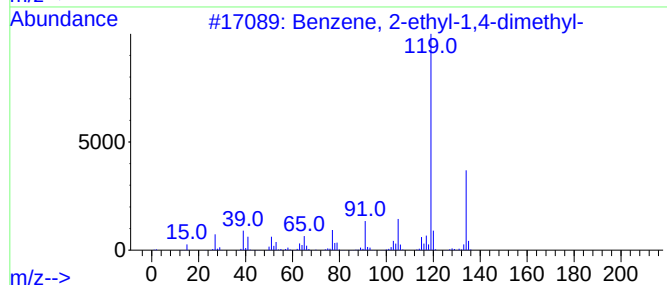
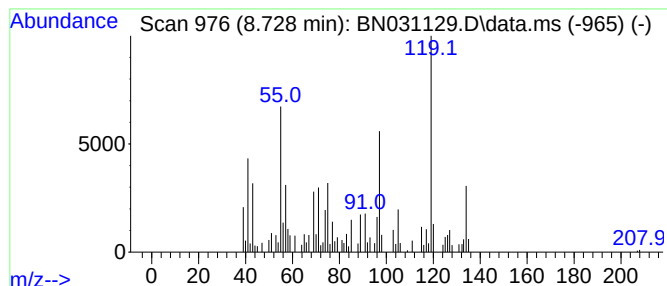
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TIC Integration Parameters: LSCINT.P

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Peak Number 18 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 25

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.728 | 5.97 ng/ul | 131129 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 91   |
| 2    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 83   |
| 3    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 83   |
| 4    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 64   |
| 5    |    |   | Benzene, 1,2,4,5-tetramethyl-  | 134 | C10H14  | 000095-93-2 | 60   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

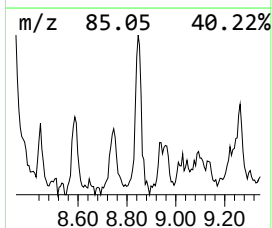
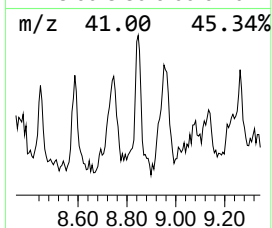
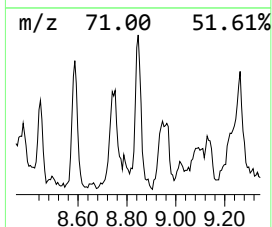
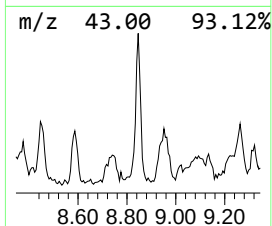
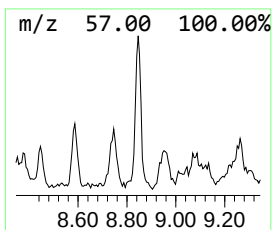
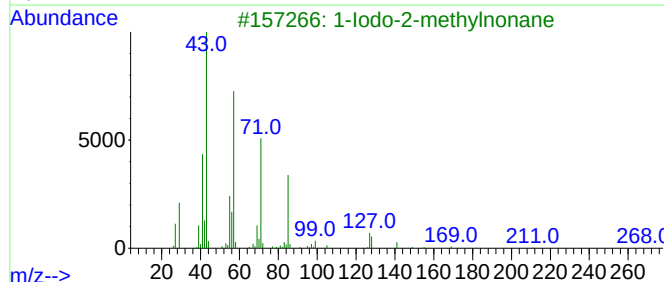
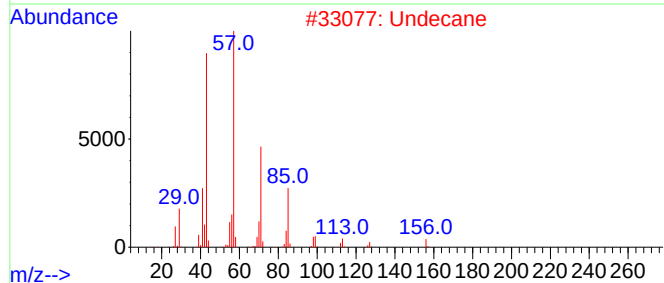
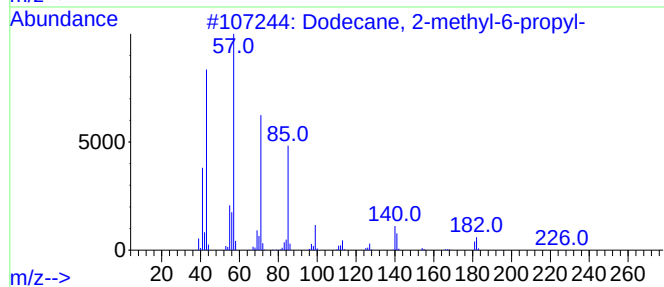
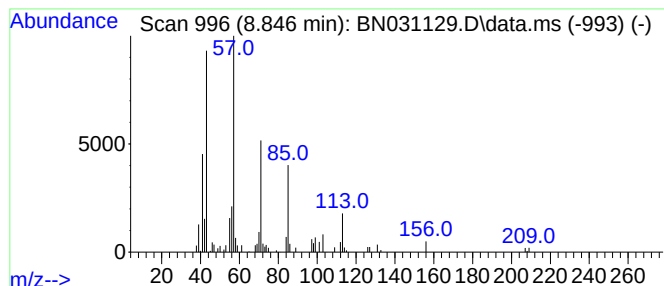
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Quant Title : SVOA CALIBRATION

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

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Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 40

| R.T.  | EstConc    | Area  | Relative to ISTD       | R.T.  |
|-------|------------|-------|------------------------|-------|
| 8.846 | 3.09 ng/ul | 67972 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of 5 | Tentative ID                 | MW  | MolForm | CAS#         | Qual |
|------|------|------------------------------|-----|---------|--------------|------|
| 1    |      | Dodecane, 2-methyl-6-propyl- | 226 | C16H34  | 055045-08-4  | 64   |
| 2    |      | Undecane                     | 156 | C11H24  | 001120-21-4  | 60   |
| 3    |      | 1-Iodo-2-methylnonane        | 268 | C10H21I | 1000101-47-9 | 59   |
| 4    |      | Hexadecane                   | 226 | C16H34  | 000544-76-3  | 58   |
| 5    |      | Undecane                     | 156 | C11H24  | 001120-21-4  | 55   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

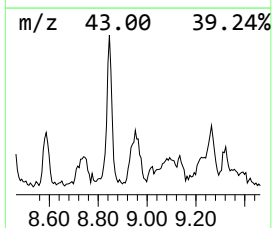
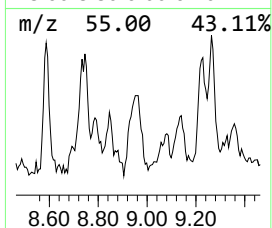
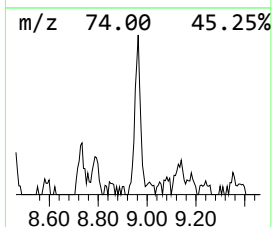
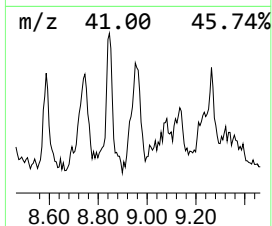
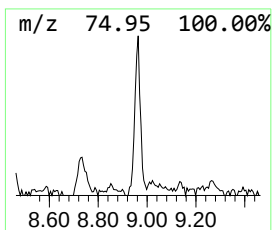
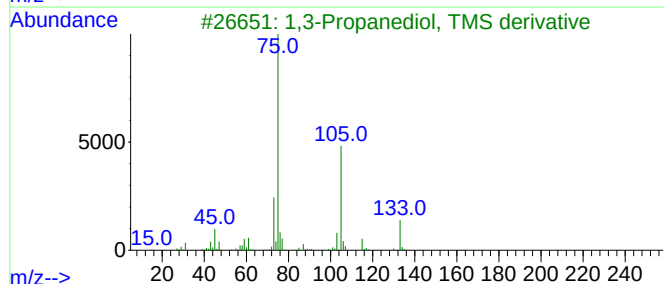
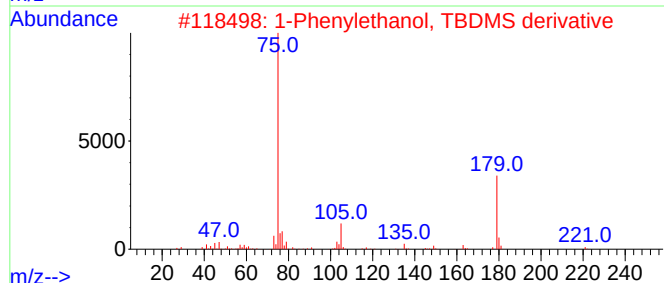
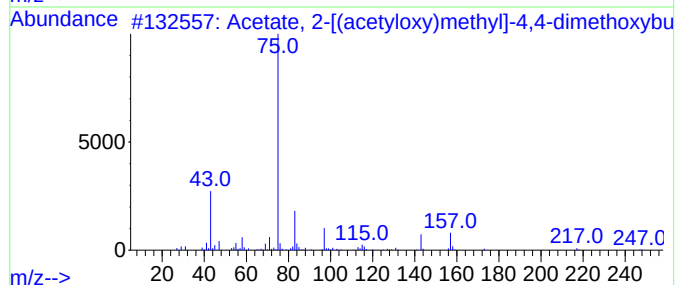
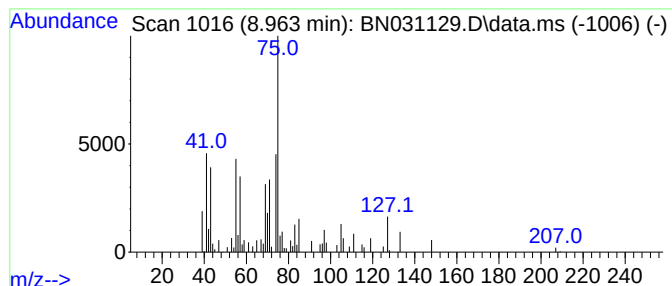
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Quant Title : SVOA CALIBRATION

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

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Peak Number 20 unknown-02 Concentration Rank 28

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.963 | 5.33 ng/ul | 117131 | 1,4-Dichlorobenzene-d4 | 7.634 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Acetate, 2-[(acetyloxy)methyl]-4... | 248 | C11H20O6  | 119117-34-9  | 32   |
| 2    |    |   | 1-Phenylethanol, TBDMS derivative   | 236 | C14H24OSi | 1000364-39-7 | 25   |
| 3    |    |   | 1,3-Propanediol, TMS derivative     | 148 | C6H16O2Si | 1000333-05-2 | 25   |
| 4    |    |   | Isobutanol, TMS derivative          | 146 | C7H18OSi  | 018269-50-6  | 25   |
| 5    |    |   | Cyclohexanol, 2-(trimethylsilyl)... | 172 | C9H20OSi  | 020584-43-4  | 25   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

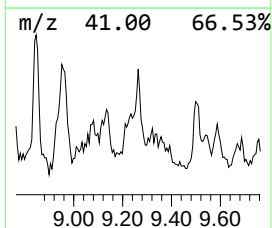
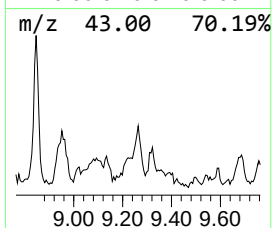
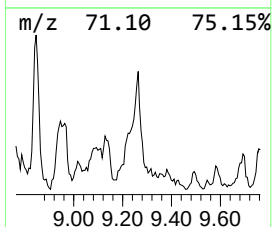
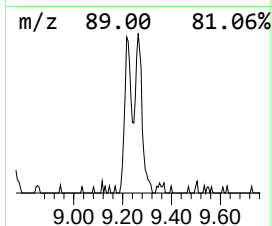
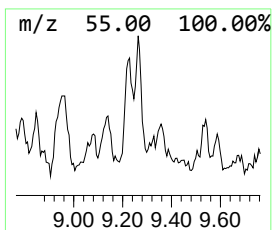
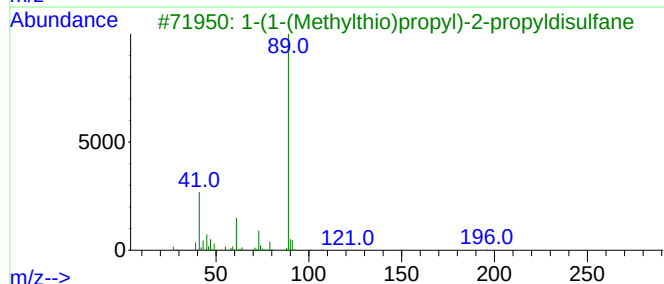
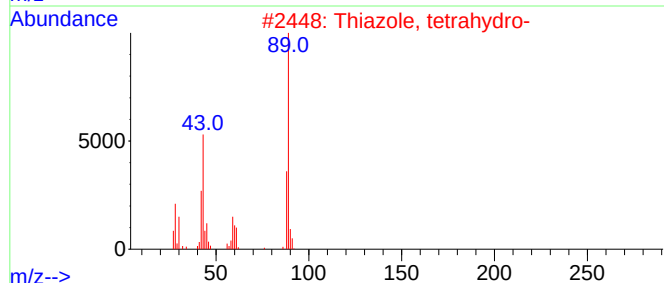
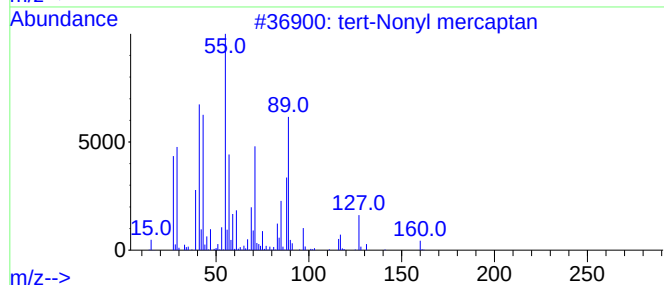
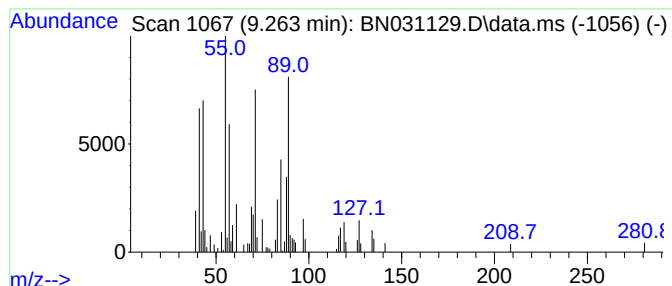
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Quant Title : SVOA CALIBRATION

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

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Peak Number 21 tert-Nonyl mercaptan Concentration Rank 46

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.   |
|-------|------------|--------|------------------|--------|
| 9.263 | 2.65 ng/ul | 110273 | Naphthalene-d8   | 10.410 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | tert-Nonyl mercaptan                | 160 | C9H20S   | 025360-10-5 | 64   |
| 2    |    |   | Thiazole, tetrahydro-               | 89  | C3H7NS   | 000504-78-9 | 50   |
| 3    |    |   | 1-(1-(Methylthio)propyl)-2-propy... | 196 | C7H16S3  | 126876-22-0 | 50   |
| 4    |    |   | Isobutyric acid, tetradecyl ester   | 284 | C18H36O2 | 167871-30-9 | 42   |
| 5    |    |   | Isopropyl butyrate                  | 130 | C7H14O2  | 000638-11-9 | 40   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

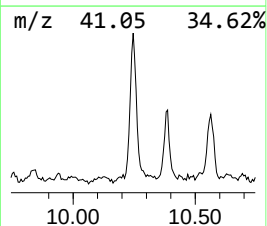
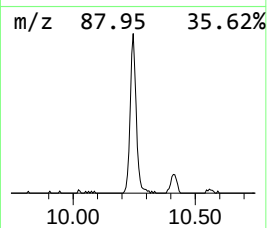
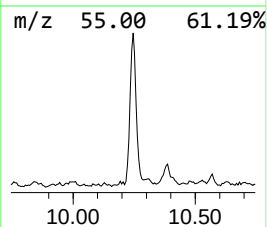
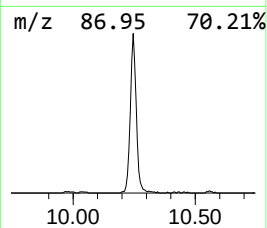
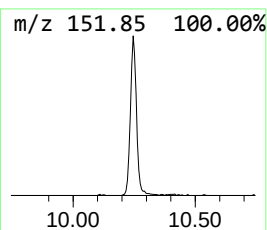
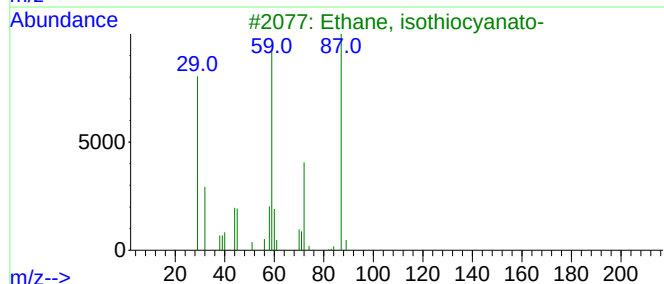
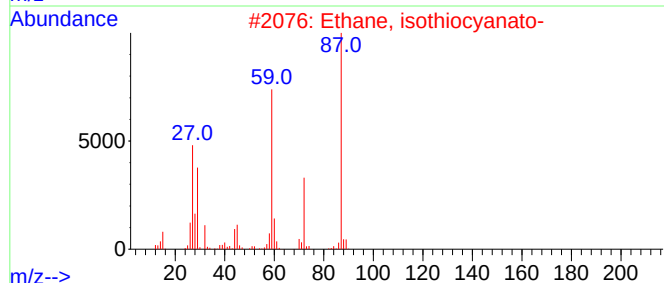
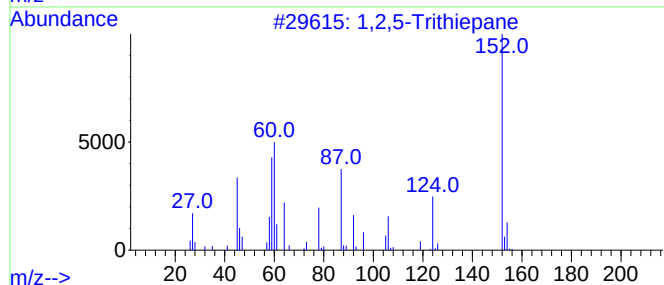
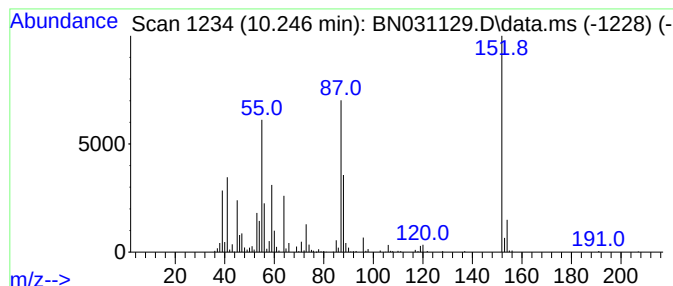
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TIC Integration Parameters: LSCINT.P

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Peak Number 22 unknown-03 Concentration Rank 13

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 10.246 | 10.08 ng/ul | 419287 | Naphthalene-d8   | 10.410 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1,2,5-Trithiepane                   | 152 | C4H8S3  | 006576-93-8 | 28   |
| 2    |    |   | Ethane, isothiocyanato-             | 87  | C3H5NS  | 000542-85-8 | 14   |
| 3    |    |   | Ethane, isothiocyanato-             | 87  | C3H5NS  | 000542-85-8 | 14   |
| 4    |    |   | 2-Heptenoic acid, methyl ester, ... | 142 | C8H14O2 | 038693-91-3 | 12   |
| 5    |    |   | 2-Heptenoic acid, methyl ester      | 142 | C8H14O2 | 022104-69-4 | 12   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

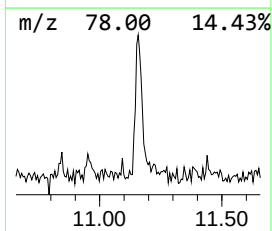
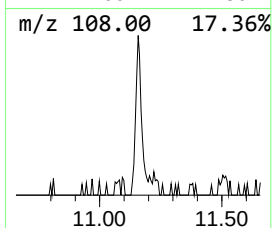
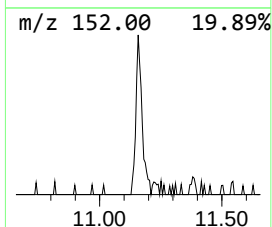
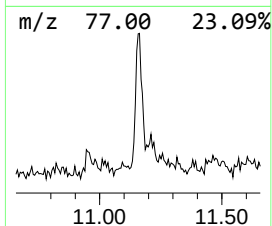
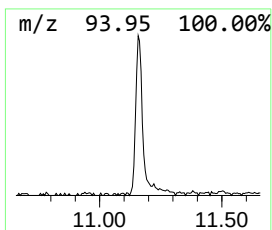
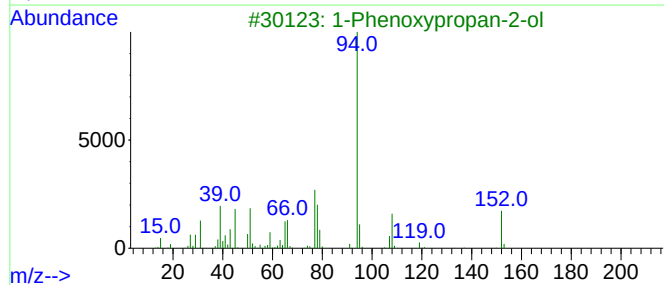
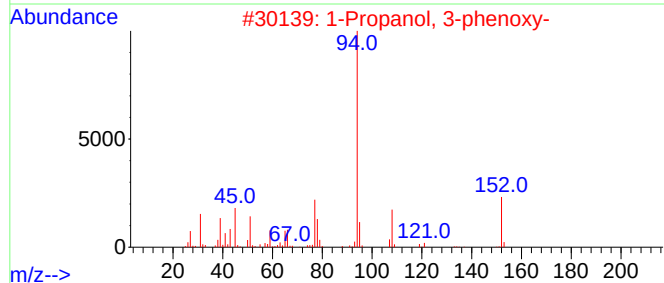
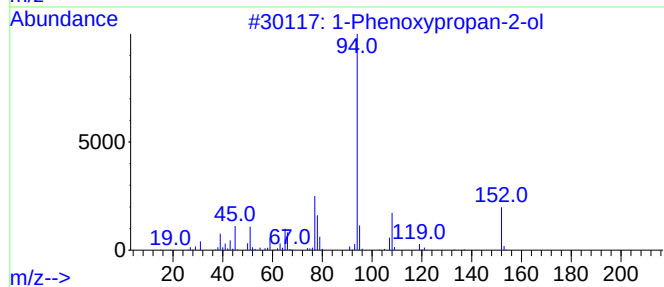
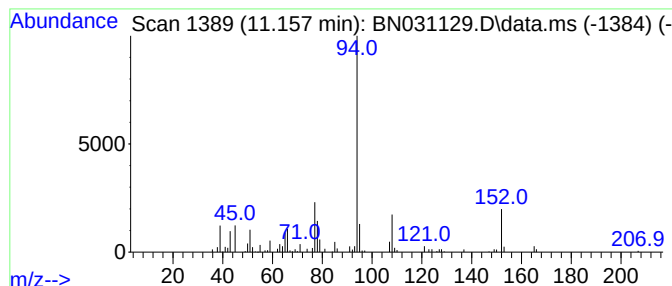
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Quant Title : SVOA CALIBRATION

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

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Peak Number 23 1-Phenoxypropan-2-ol Concentration Rank 54

| R.T.   | EstConc    | Area  | Relative to ISTD | R.T.   |
|--------|------------|-------|------------------|--------|
| 11.157 | 2.08 ng/ul | 86339 | Naphthalene-d8   | 10.410 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | 1-Phenoxypropan-2-ol   | 152 | C9H12O2 | 000770-35-4 | 97   |
| 2    |    |   | 1-Propanol, 3-phenoxy- | 152 | C9H12O2 | 006180-61-6 | 93   |
| 3    |    |   | 1-Phenoxypropan-2-ol   | 152 | C9H12O2 | 000770-35-4 | 91   |
| 4    |    |   | 1-Phenoxypropan-2-ol   | 152 | C9H12O2 | 000770-35-4 | 90   |
| 5    |    |   | 1-Phenoxypropan-2-ol   | 152 | C9H12O2 | 000770-35-4 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

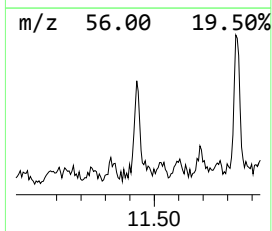
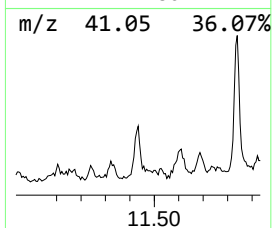
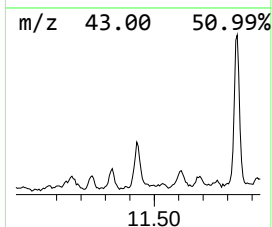
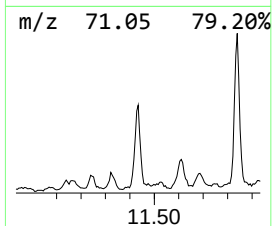
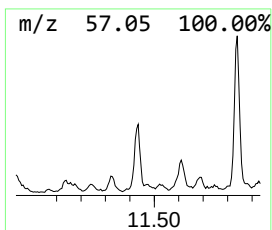
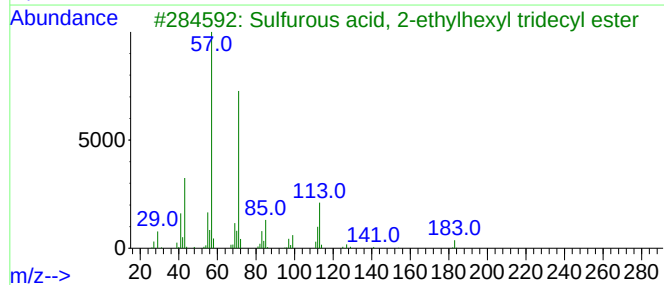
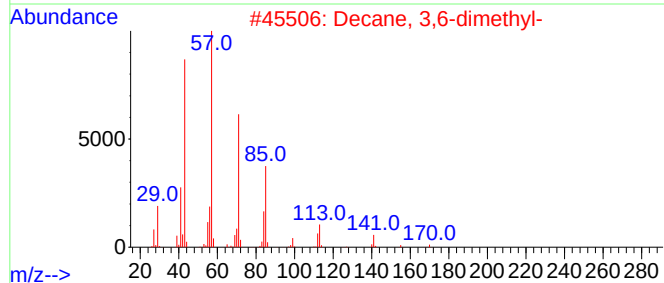
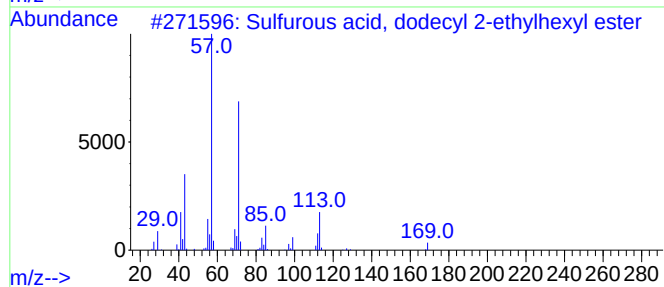
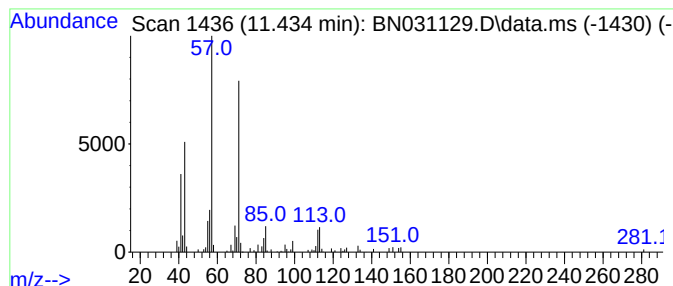
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 24 Sulfurous acid, dodecyl 2-e... Concentration Rank 41

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 11.434 | 3.09 ng/ul | 128481 | Naphthalene-d8   | 10.410 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Sulfurous acid, dodecyl 2-ethylh... | 362 | C20H42O3S | 1000309-19-5 | 72   |
| 2    |    |   | Decane, 3,6-dimethyl-               | 170 | C12H26    | 017312-53-7  | 72   |
| 3    |    |   | Sulfurous acid, 2-ethylhexyl tri... | 376 | C21H44O3S | 1000309-19-6 | 64   |
| 4    |    |   | Dodecane, 2,6,10-trimethyl-         | 212 | C15H32    | 003891-98-3  | 64   |
| 5    |    |   | Dodecane, 2-methyl-6-propyl-        | 226 | C16H34    | 055045-08-4  | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

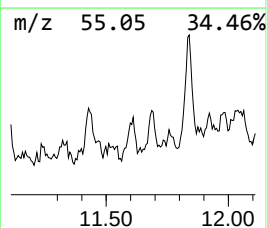
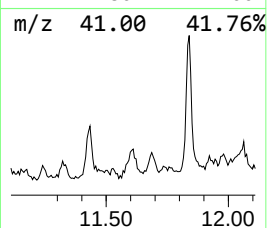
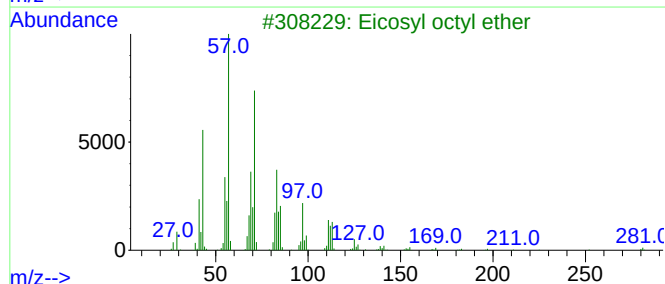
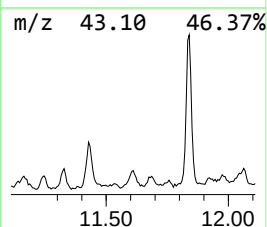
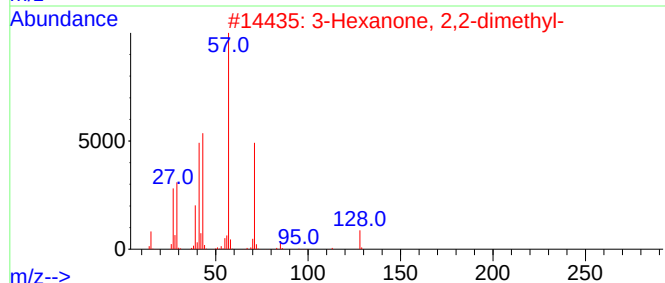
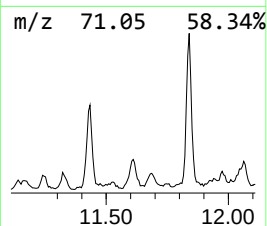
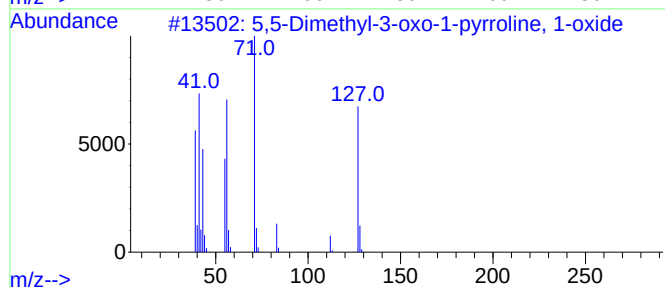
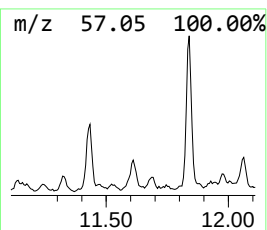
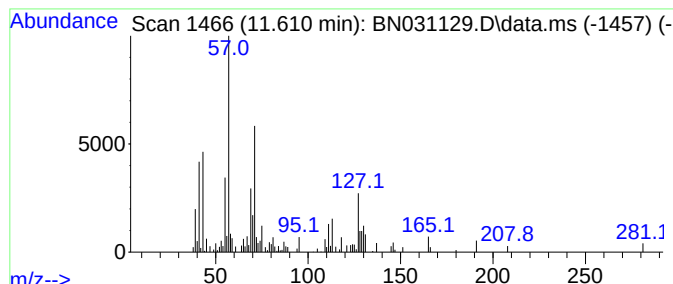
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Quant Title : SVOA CALIBRATION

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

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Peak Number 25 unknown-04 Concentration Rank 47

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 11.610 | 2.61 ng/ul | 108594 | Naphthalene-d8   | 10.410 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 5,5-Dimethyl-3-oxo-1-pyrroline, ... | 127 | C6H9NO2   | 1000305-98-3 | 43   |
| 2    |    |   | 3-Hexanone, 2,2-dimethyl-           | 128 | C8H16O    | 005405-79-8  | 38   |
| 3    |    |   | Eicosyl octyl ether                 | 410 | C28H58O   | 1000406-38-8 | 37   |
| 4    |    |   | Sulfurous acid, 2-ethylhexyl non... | 320 | C17H36O3S | 1010309-19-2 | 37   |
| 5    |    |   | Sulfurous acid, di(2-ethylhexyl)... | 306 | C16H34O3S | 1000309-19-1 | 32   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

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

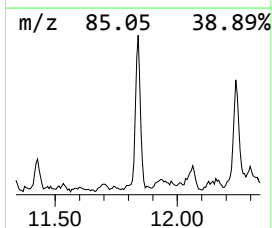
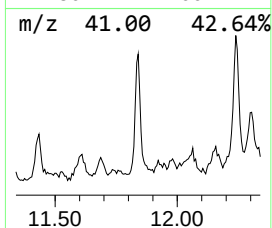
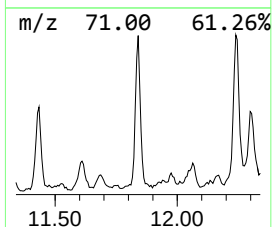
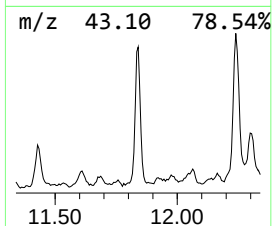
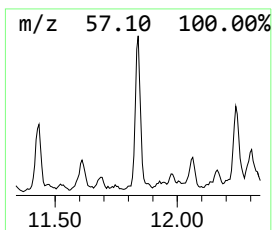
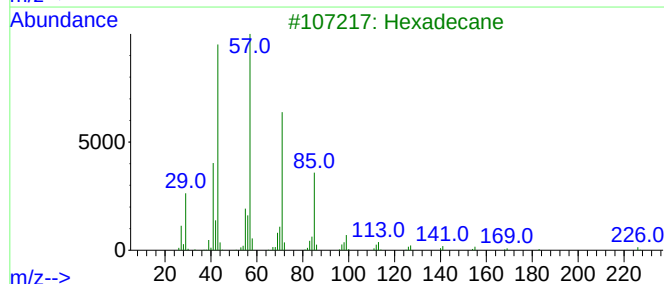
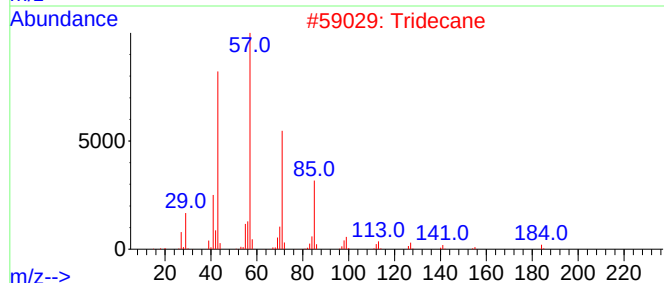
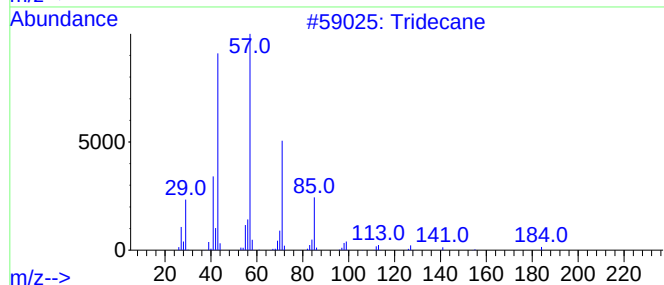
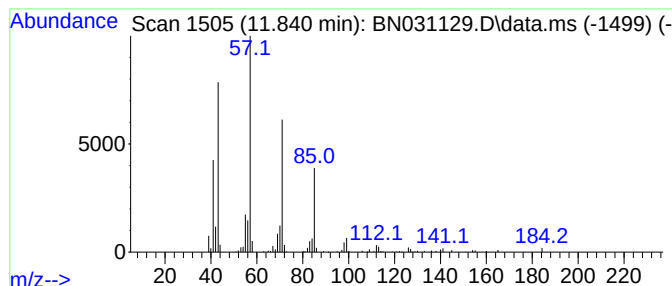
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Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 21

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 11.840 | 6.35 ng/ul | 264237 | Naphthalene-d8   | 10.410 |

| Hit# | of 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|------|----------------------|-----|---------|-------------|------|
| 1    |      | Tridecane            | 184 | C13H28  | 000629-50-5 | 93   |
| 2    |      | Tridecane            | 184 | C13H28  | 000629-50-5 | 91   |
| 3    |      | Hexadecane           | 226 | C16H34  | 000544-76-3 | 90   |
| 4    |      | Tetradecane          | 198 | C14H30  | 000629-59-4 | 90   |
| 5    |      | Eicosane, 10-methyl- | 296 | C21H44  | 054833-23-7 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
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Quant Title : SVOA CALIBRATION

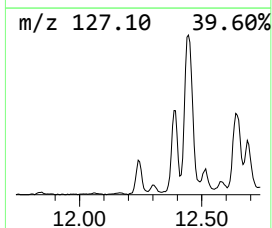
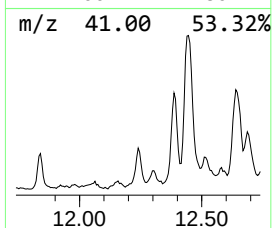
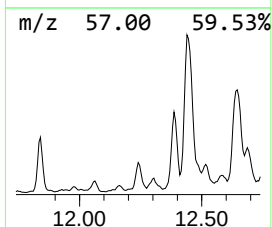
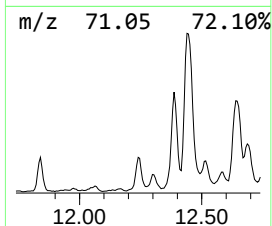
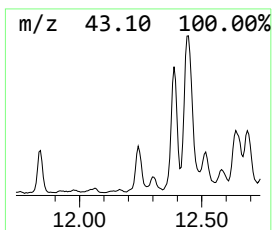
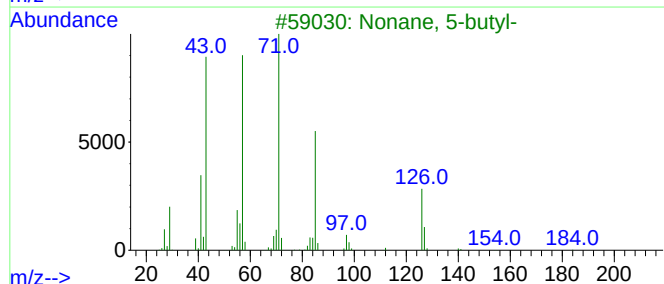
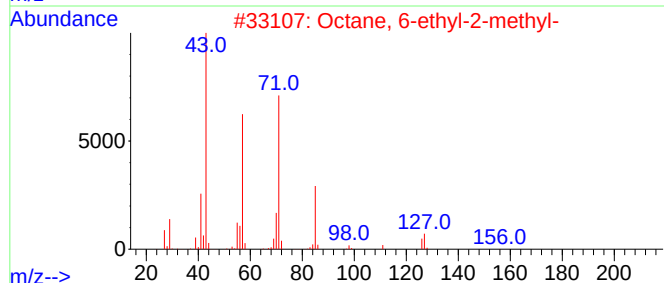
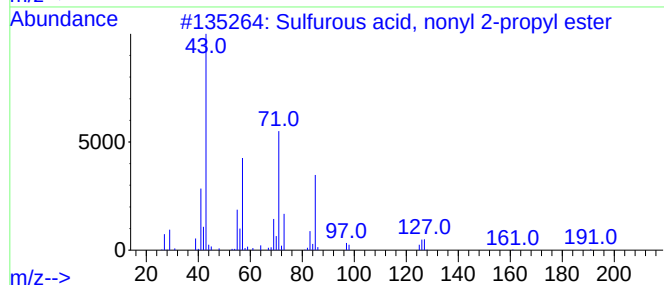
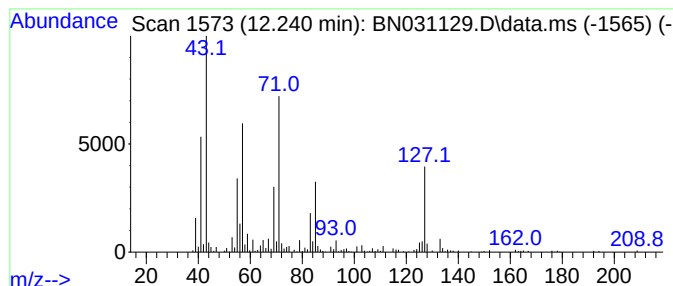
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TIC Integration Parameters: LSCINT.P

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Peak Number 27 Sulfurous acid, nonyl 2-pro... Concentration Rank 18

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 12.240 | 7.91 ng/ul | 328797 | Naphthalene-d8   | 10.410 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Sulfurous acid, nonyl 2-propyl e... | 250 | C12H26O3S | 1000309-12-0 | 50   |
| 2    |    |   | Octane, 6-ethyl-2-methyl-           | 156 | C11H24    | 062016-19-7  | 43   |
| 3    |    |   | Nonane, 5-butyl-                    | 184 | C13H28    | 017312-63-9  | 38   |
| 4    |    |   | Dodecane, 2,7,10-trimethyl-         | 212 | C15H32    | 074645-98-0  | 38   |
| 5    |    |   | Pentane, 2,3,3-trimethyl-           | 114 | C8H18     | 000560-21-4  | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

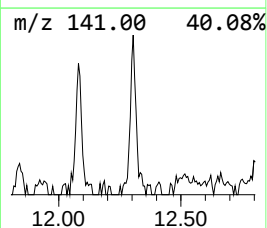
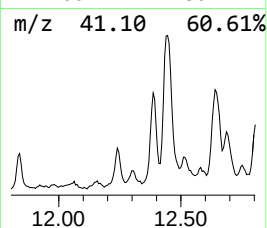
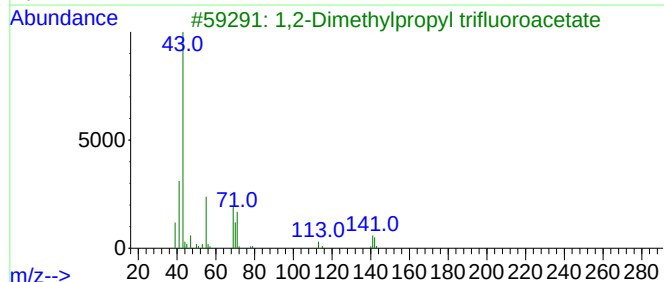
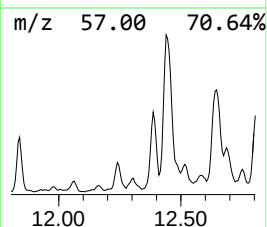
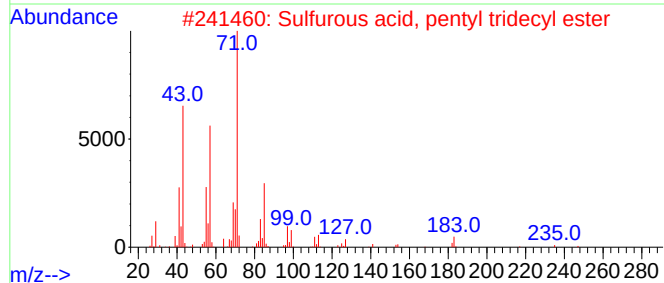
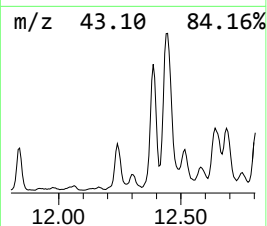
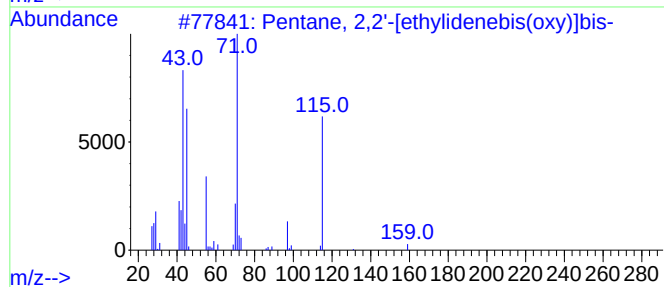
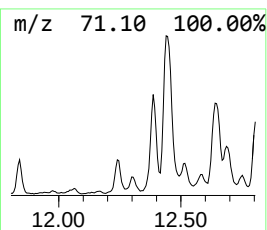
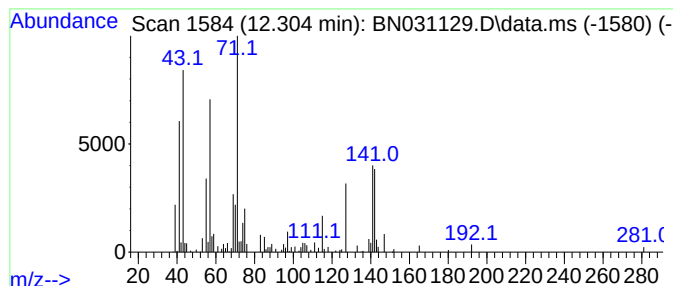
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 28 unknown-05 Concentration Rank 52

| R.T.   | EstConc    | Area  | Relative to ISTD | R.T.   |
|--------|------------|-------|------------------|--------|
| 12.304 | 2.20 ng/ul | 91569 | Naphthalene-d8   | 10.410 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm   | CAS#         | Qual |
|------|----|---|--------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Pentane, 2,2'-[ethylidenebis(oxy...] | 202 | C12H26O2  | 054889-47-3  | 35   |
| 2    |    |   | Sulfurous acid, pentyl tridecyl ...  | 334 | C18H38O3S | 1000309-14-6 | 35   |
| 3    |    |   | 1,2-Dimethylpropyl trifluoroacetate  | 184 | C7H11F3O2 | 000691-75-8  | 32   |
| 4    |    |   | Propanoic acid, 2-methyl-, 1-met...  | 158 | C9H18O2   | 054340-93-1  | 27   |
| 5    |    |   | Pentane, 3,3-dimethyl-               | 100 | C7H16     | 000562-49-2  | 27   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

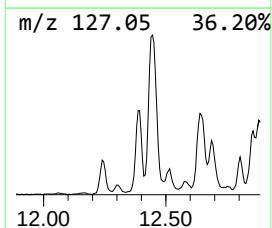
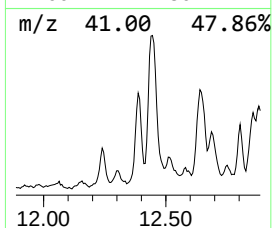
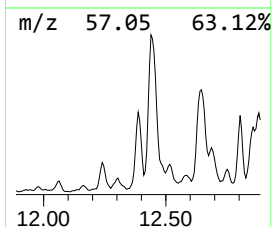
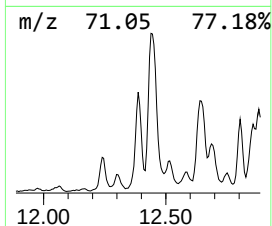
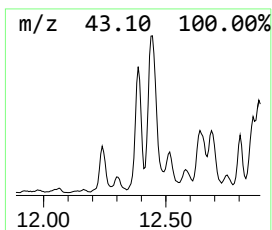
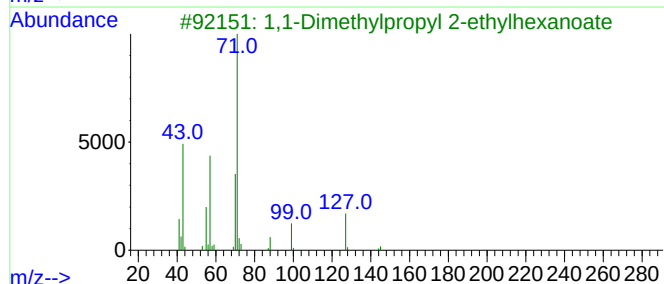
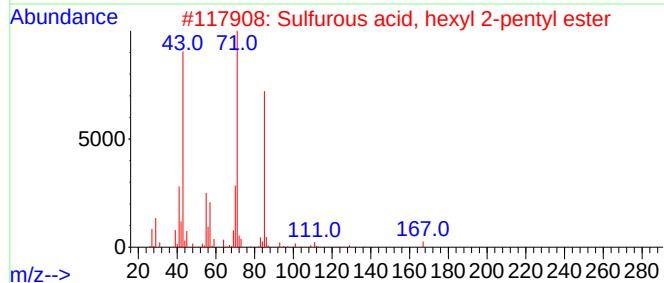
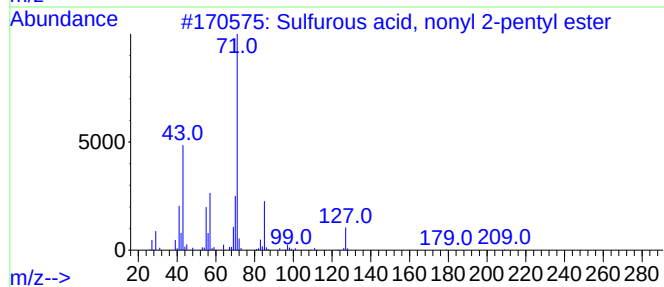
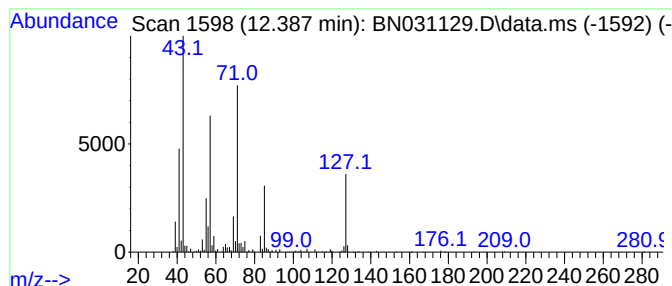
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TIC Integration Parameters: LSCINT.P

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Peak Number 29 Sulfurous acid, nonyl 2-pen... Concentration Rank 14

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 12.387 | 9.92 ng/ul | 654450 | Acenaphthene-d10 | 14.281 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Sulfurous acid, nonyl 2-pentyl e... | 278 | C14H30O3S | 1010309-15-8 | 72   |
| 2    |    |   | Sulfurous acid, hexyl 2-pentyl e... | 236 | C11H24O3S | 1000309-15-6 | 64   |
| 3    |    |   | 1,1-Dimethylpropyl 2-ethylhexanoate | 214 | C13H26O2  | 1000099-99-2 | 59   |
| 4    |    |   | Undecane, 2,8-dimethyl-             | 184 | C13H28    | 017301-25-6  | 59   |
| 5    |    |   | 1-Methylpentyl isobutyrate          | 172 | C10H20O2  | 1000429-31-7 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

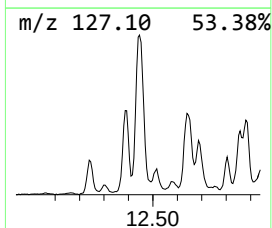
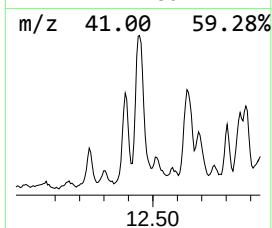
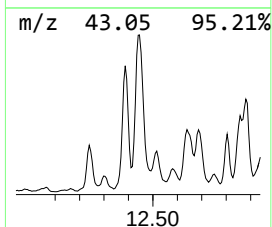
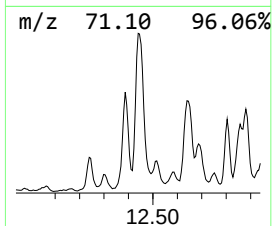
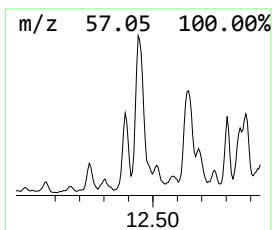
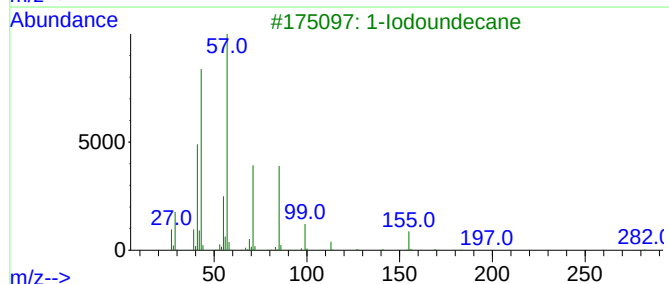
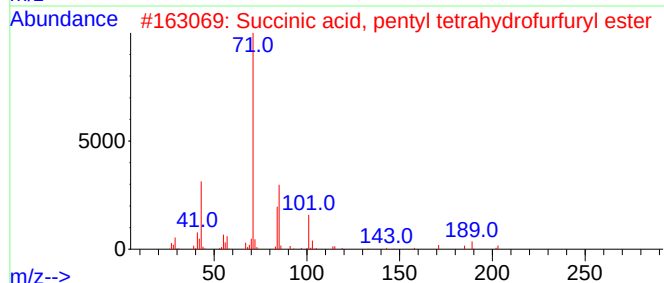
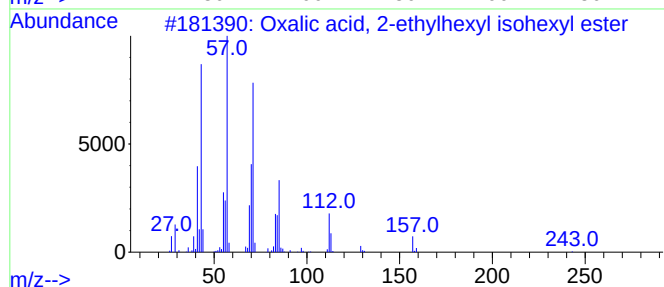
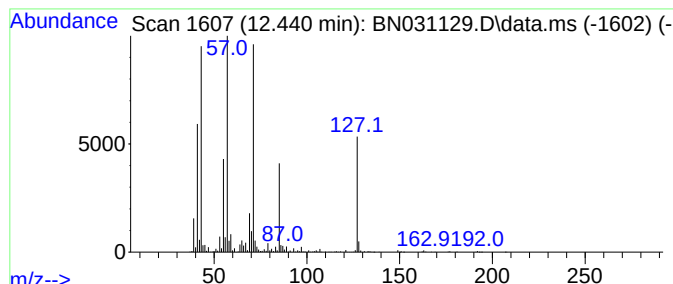
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Quant Title : SVOA CALIBRATION

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

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Peak Number 30 Oxalic acid, 2-ethylhexyl i... Concentration Rank 3

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 12.440 | 21.95 ng/ul | 1448990 | Acenaphthene-d10 | 14.281 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Oxalic acid, 2-ethylhexyl isohex... | 286 | C16H30O4 | 1000309-38-8 | 53   |
| 2    |    |   | Succinic acid, pentyl tetrahydro... | 272 | C14H24O5 | 1000329-86-3 | 50   |
| 3    |    |   | 1-Iodoundecane                      | 282 | C11H23I  | 004282-44-4  | 38   |
| 4    |    |   | Oxalic acid, neopentyl propyl ester | 202 | C10H18O4 | 1000309-72-5 | 38   |
| 5    |    |   | Dodecane, 1-iodo-                   | 296 | C12H25I  | 004292-19-7  | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

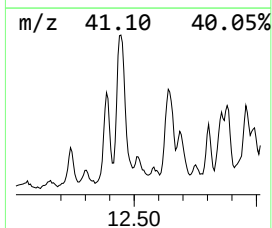
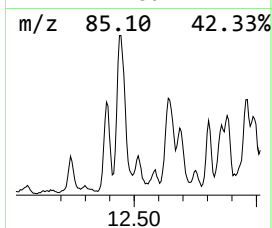
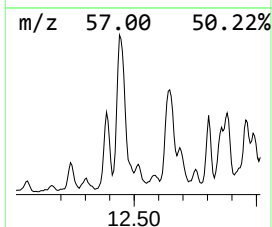
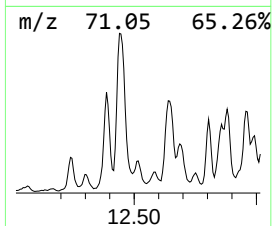
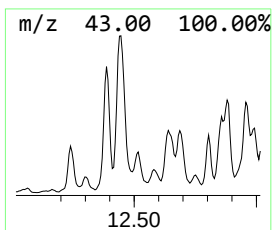
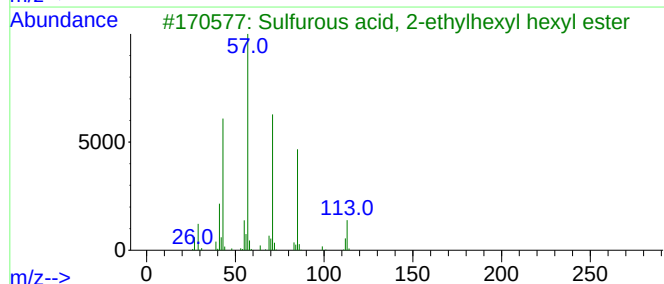
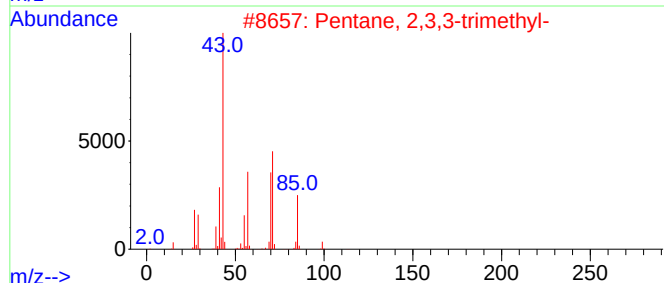
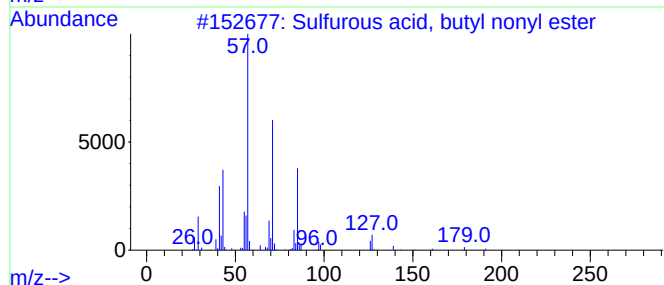
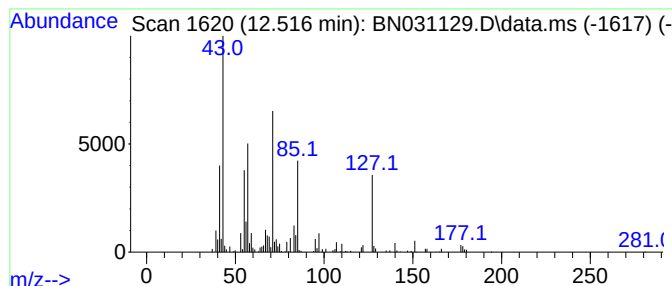
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TIC Integration Parameters: LSCINT.P

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Peak Number 31 Sulfurous acid, butyl nonyl... Concentration Rank 51

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 12.516 | 2.22 ng/ul | 146319 | Acenaphthene-d10 | 14.281 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Sulfurous acid, butyl nonyl ester   | 264 | C13H28O3S | 1000309-17-6 | 53   |
| 2    |    |   | Pentane, 2,3,3-trimethyl-           | 114 | C8H18     | 000560-21-4  | 47   |
| 3    |    |   | Sulfurous acid, 2-ethylhexyl hex... | 278 | C14H30O3S | 1000309-20-2 | 47   |
| 4    |    |   | Hexane, 2,3,4-trimethyl-            | 128 | C9H20     | 000921-47-1  | 47   |
| 5    |    |   | Tetradecane, 1-iodo-                | 324 | C14H29I   | 019218-94-1  | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

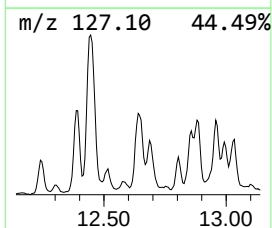
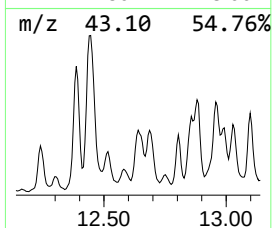
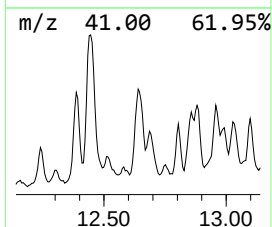
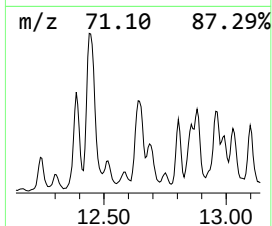
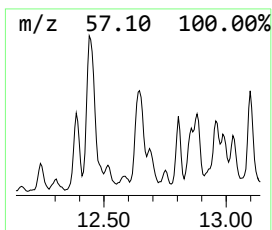
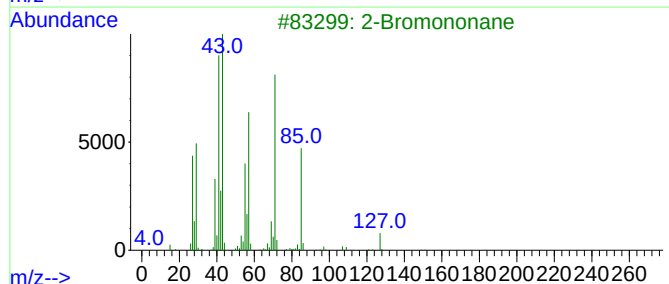
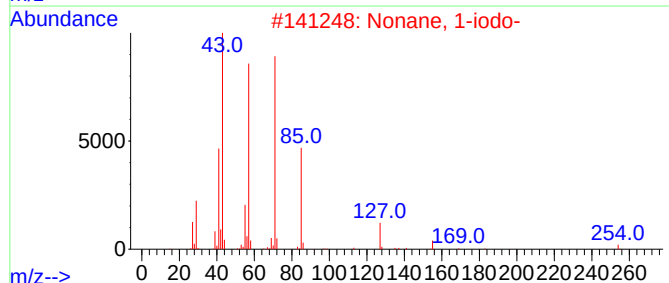
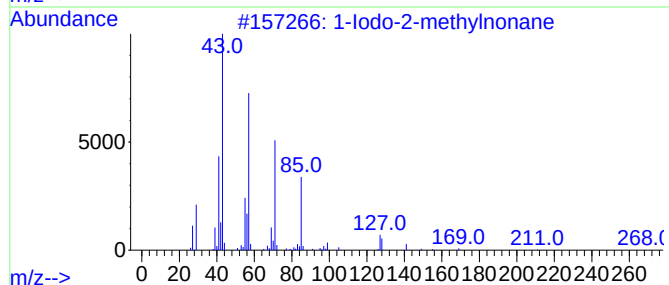
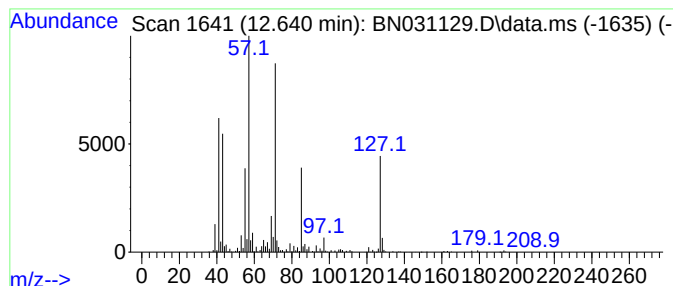
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Quant Title : SVOA CALIBRATION

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

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Peak Number 32 1-Iodo-2-methylnonane Concentration Rank 10

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 12.640 | 11.50 ng/ul | 758685 | Acenaphthene-d10 | 14.281 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 1-Iodo-2-methylnonane               | 268 | C10H21I  | 1000101-47-9 | 59   |
| 2    |    |   | Nonane, 1-iodo-                     | 254 | C9H19I   | 004282-42-2  | 50   |
| 3    |    |   | 2-Bromononane                       | 206 | C9H19Br  | 002216-35-5  | 50   |
| 4    |    |   | Succinic acid, pentyl tetrahydro... | 272 | C14H24O5 | 1000329-86-3 | 47   |
| 5    |    |   | Dodecane, 4-methyl-                 | 184 | C13H28   | 006117-97-1  | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

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

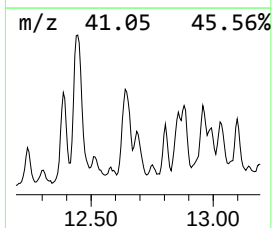
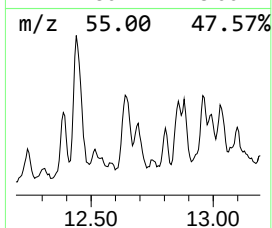
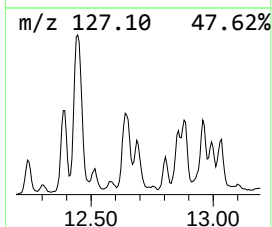
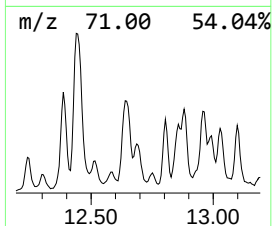
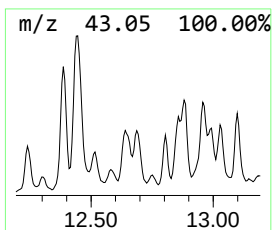
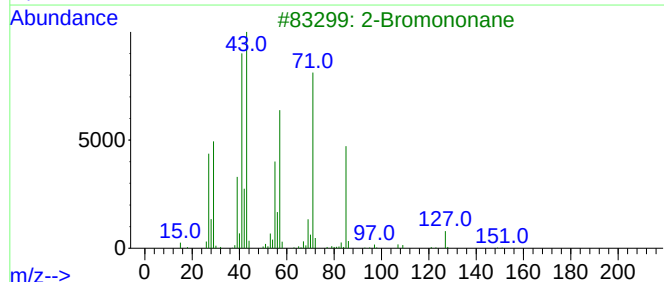
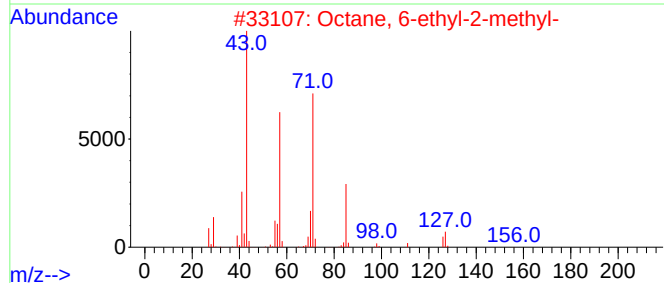
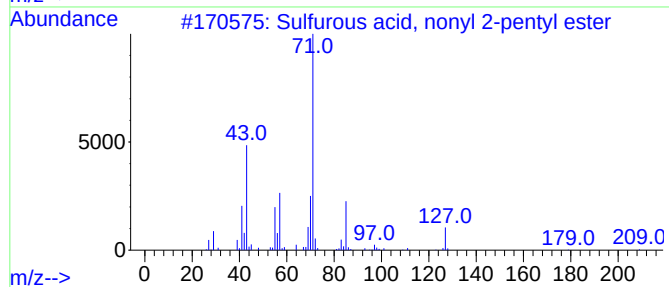
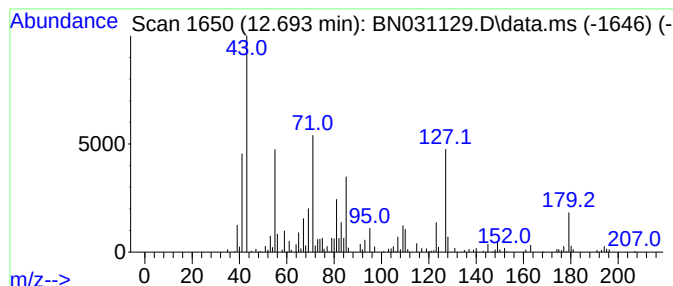
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Peak Number 33 unknown-06

Concentration Rank 22

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 12.693 | 6.17 ng/ul | 406895 | Acenaphthene-d10 | 14.281 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Sulfurous acid, nonyl 2-pentyl e... | 278 | C14H30O3S | 1010309-15-8 | 43   |
| 2    |    |   | Octane, 6-ethyl-2-methyl-           | 156 | C11H24    | 062016-19-7  | 38   |
| 3    |    |   | 2-Bromononane                       | 206 | C9H19Br   | 002216-35-5  | 35   |
| 4    |    |   | 1,3-Dimethylcyclopentanol           | 114 | C7H14O    | 019550-46-0  | 27   |
| 5    |    |   | (1R,2S)-2-Acetyl-1-methylcyclobu... | 170 | C9H14O3   | 068225-41-2  | 25   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

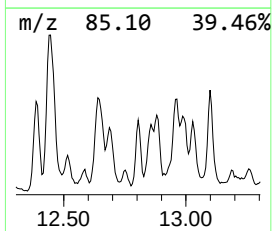
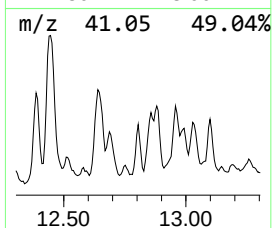
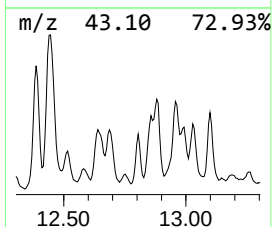
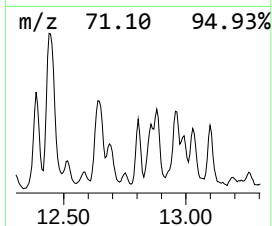
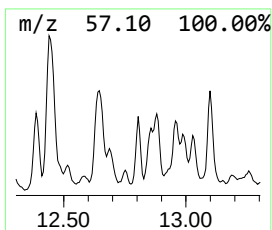
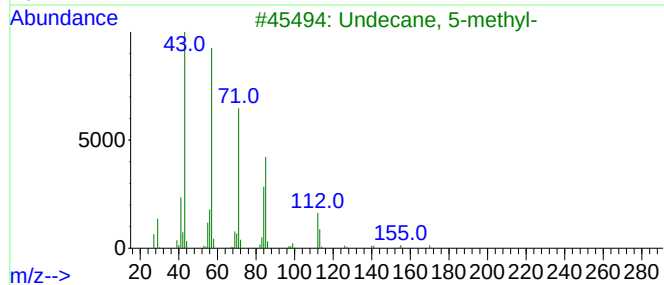
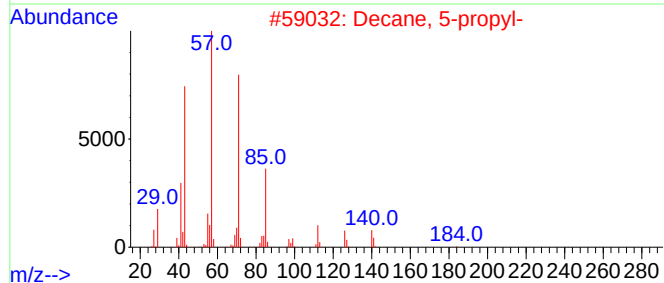
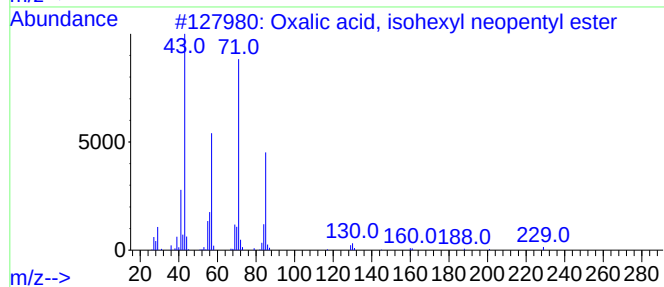
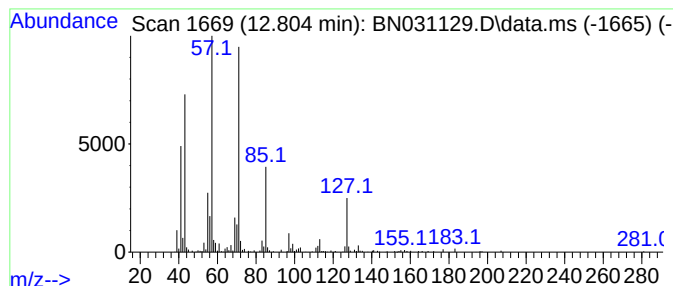
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 34 Oxalic acid, isohexyl neope... Concentration Rank 34

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 12.804 | 4.66 ng/ul | 307432 | Acenaphthene-d10 | 14.281 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Oxalic acid, isohexyl neopentyl ... | 244 | C13H24O4  | 1000309-73-0 | 64   |
| 2    |    |   | Decane, 5-propyl-                   | 184 | C13H28    | 017312-62-8  | 64   |
| 3    |    |   | Undecane, 5-methyl-                 | 170 | C12H26    | 001632-70-8  | 53   |
| 4    |    |   | Sulfurous acid, 2-ethylhexyl hex... | 418 | C24H50O3S | 1000309-19-9 | 53   |
| 5    |    |   | Sulfurous acid, di(2-ethylhexyl)... | 306 | C16H34O3S | 1000309-19-1 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

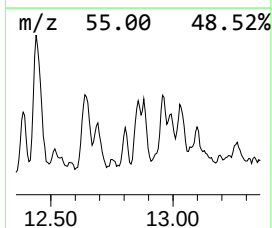
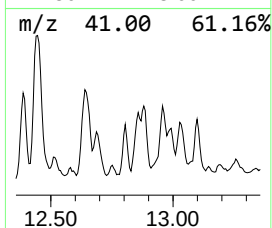
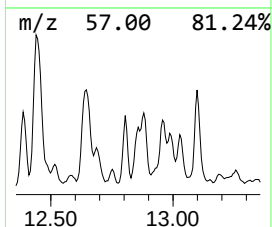
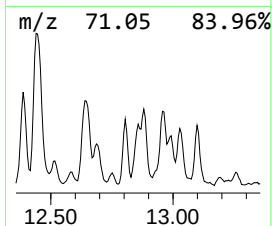
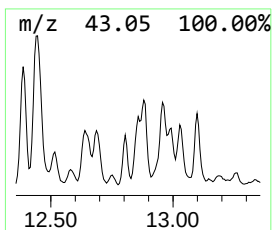
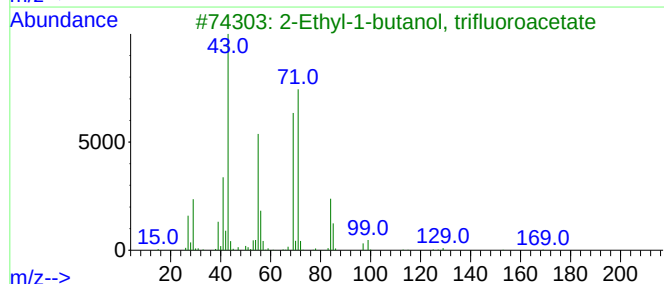
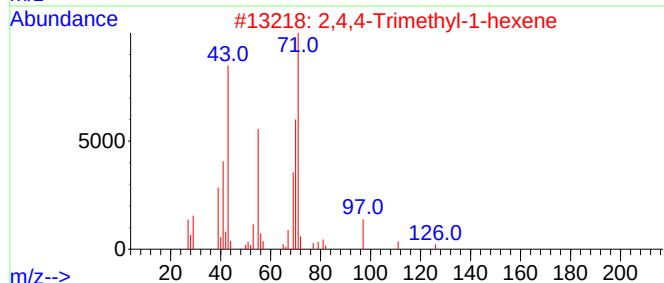
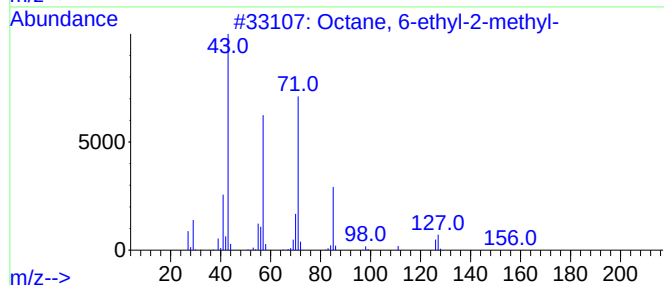
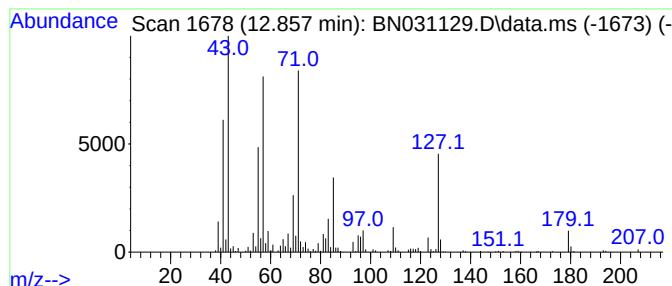
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Quant Title : SVOA CALIBRATION

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

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Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 20

| R.T.      | EstConc                             | Area   | Relative to ISTD |           |             | R.T.   |
|-----------|-------------------------------------|--------|------------------|-----------|-------------|--------|
| 12.857    | 7.06 ng/ul                          | 465994 | Acenaphthene-d10 |           |             | 14.281 |
| Hit# of 5 | Tentative ID                        |        | MW               | MolForm   | CAS#        | Qual   |
| 1         | Octane, 6-ethyl-2-methyl-           |        | 156              | C11H24    | 062016-19-7 | 47     |
| 2         | 2,4,4-Trimethyl-1-hexene            |        | 126              | C9H18     | 051174-12-0 | 38     |
| 3         | 2-Ethyl-1-butanol, trifluoroacetate |        | 198              | C8H13F3O2 | 116464-98-3 | 35     |
| 4         | Butane, 2,2-dimethyl-               |        | 86               | C6H14     | 000075-83-2 | 35     |
| 5         | 2,4,4-Trimethyl-1-hexene            |        | 126              | C9H18     | 051174-12-0 | 30     |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

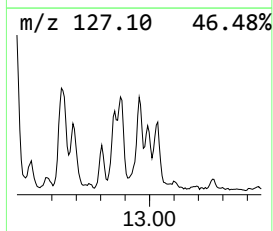
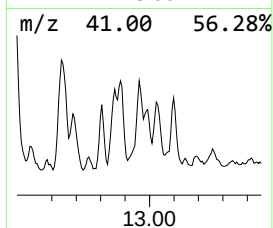
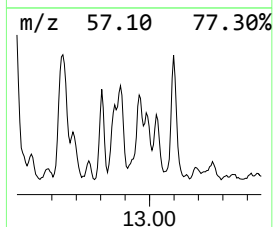
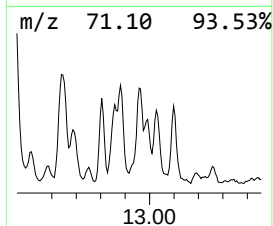
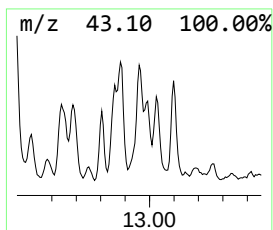
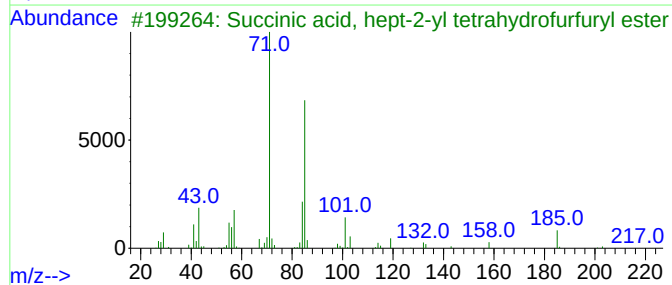
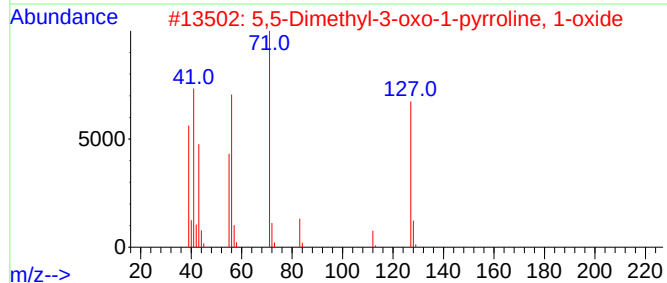
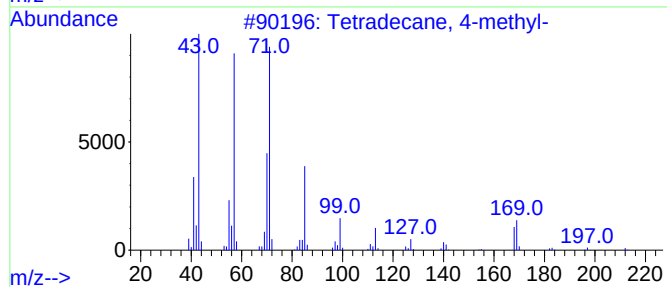
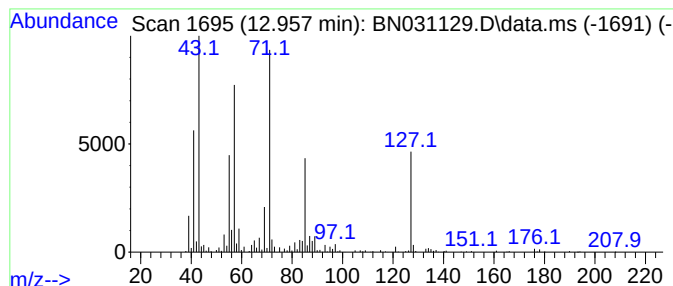
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TIC Integration Parameters: LSCINT.P

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Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 26

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 12.957 | 5.95 ng/ul | 392459 | Acenaphthene-d10 | 14.281 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Tetradecane, 4-methyl-              | 212 | C15H32    | 025117-24-2  | 53   |
| 2    |    |   | 5,5-Dimethyl-3-oxo-1-pyrroline, ... | 127 | C6H9NO2   | 1000305-98-3 | 49   |
| 3    |    |   | Succinic acid, hept-2-yl tetrahy... | 300 | C16H28O5  | 1000390-71-8 | 47   |
| 4    |    |   | Succinic acid, isohexyl tetrahyd... | 286 | C15H26O5  | 1000329-86-4 | 43   |
| 5    |    |   | Sulfurous acid, butyl nonyl ester   | 264 | C13H28O3S | 1000309-17-6 | 43   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

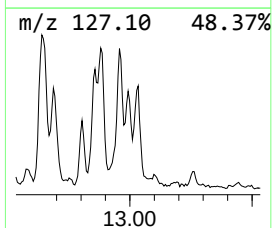
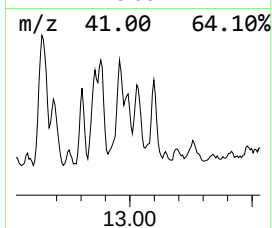
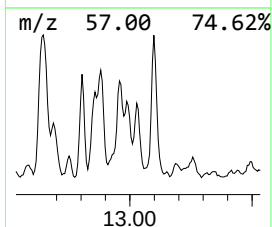
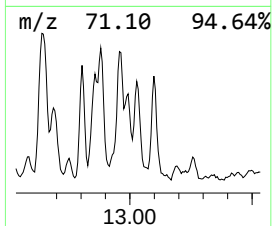
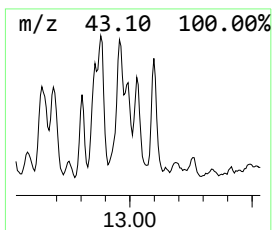
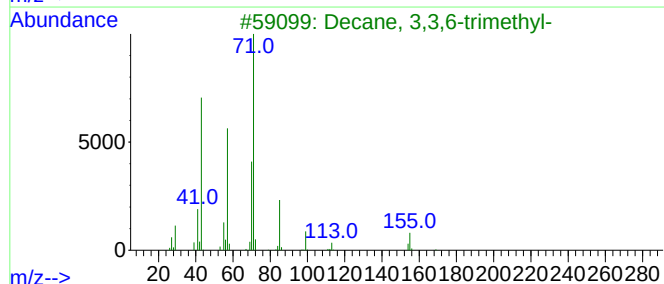
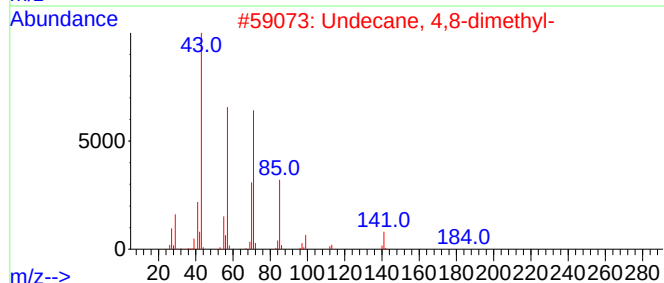
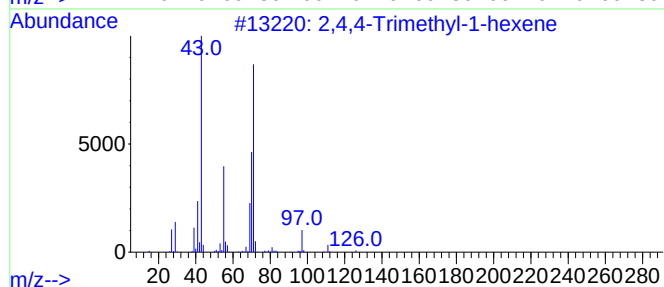
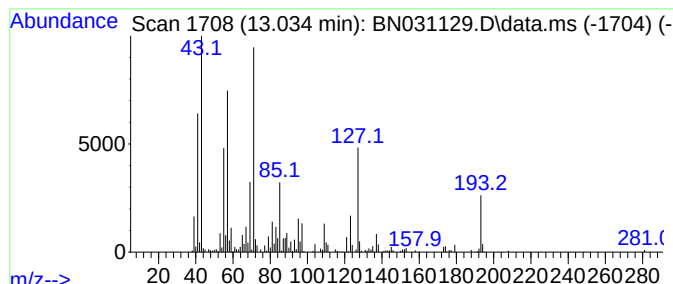
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ClientSampleId :  
E1CX8

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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 37 (DEL) Alkane: Straight-Chai... Concentration Rank 29

| R.T.      | EstConc                  | Area   | Relative to ISTD | R.T.        |      |
|-----------|--------------------------|--------|------------------|-------------|------|
| 13.034    | 5.18 ng/ul               | 341813 | Acenaphthene-d10 | 14.281      |      |
| Hit# of 5 | Tentative ID             | MW     | MolForm          | CAS#        | Qual |
| 1         | 2,4,4-Trimethyl-1-hexene | 126    | C9H18            | 051174-12-0 | 42   |
| 2         | Undecane, 4,8-dimethyl-  | 184    | C13H28           | 017301-33-6 | 35   |
| 3         | Decane, 3,3,6-trimethyl- | 184    | C13H28           | 062338-14-1 | 35   |
| 4         | Undecane, 6-ethyl-       | 184    | C13H28           | 017312-60-6 | 35   |
| 5         | Tridecane, 4-methyl-     | 198    | C14H30           | 026730-12-1 | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

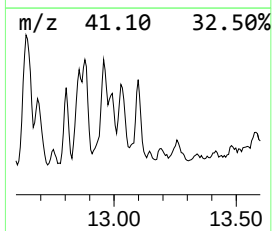
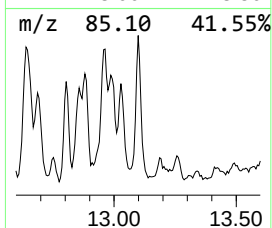
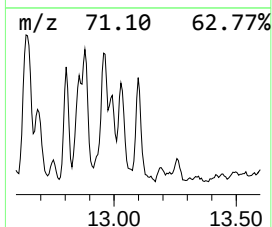
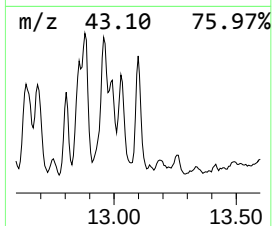
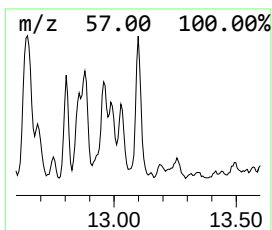
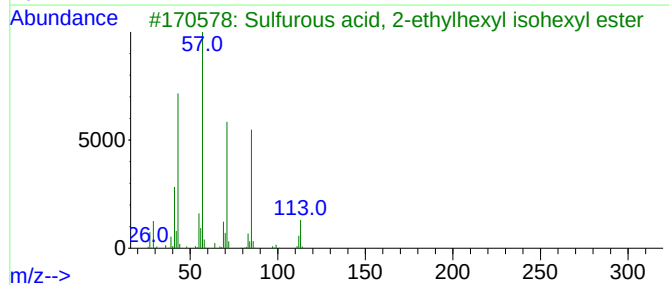
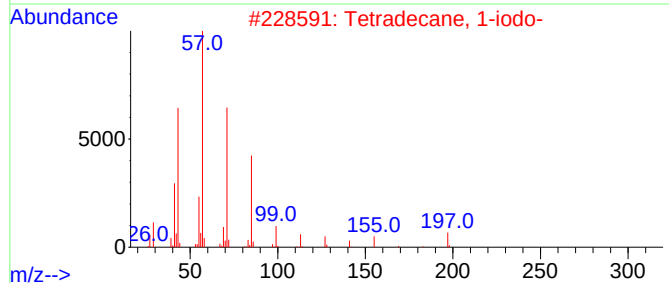
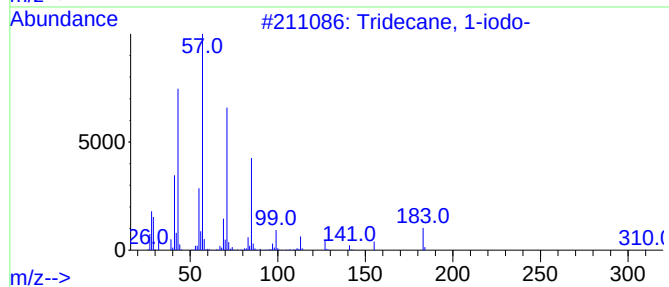
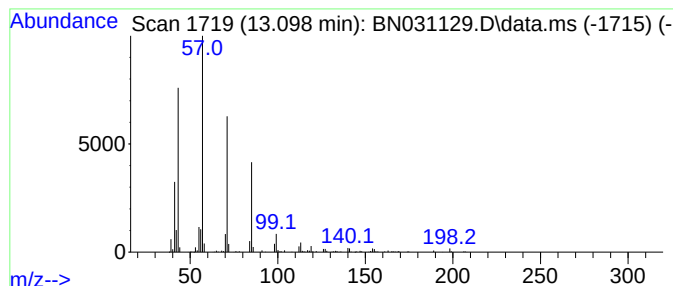
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Quant Title : SVOA CALIBRATION

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

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Peak Number 38 Tridecane, 1-iodo- Concentration Rank 30

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.098 | 5.08 ng/ul | 335283 | Acenaphthene-d10 | 14.281 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Tridecane, 1-iodo-                  | 310 | C13H27I   | 035599-77-0  | 78   |
| 2    |    |   | Tetradecane, 1-iodo-                | 324 | C14H29I   | 019218-94-1  | 78   |
| 3    |    |   | Sulfurous acid, 2-ethylhexyl iso... | 278 | C14H30O3S | 1000309-19-0 | 78   |
| 4    |    |   | Tridecane                           | 184 | C13H28    | 000629-50-5  | 72   |
| 5    |    |   | Hexadecane                          | 226 | C16H34    | 000544-76-3  | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

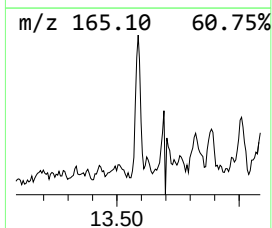
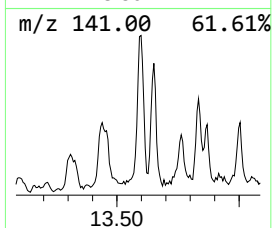
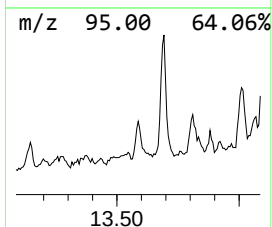
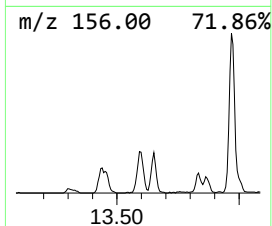
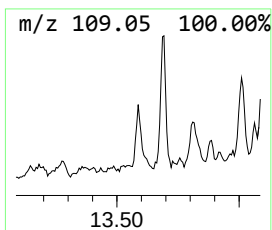
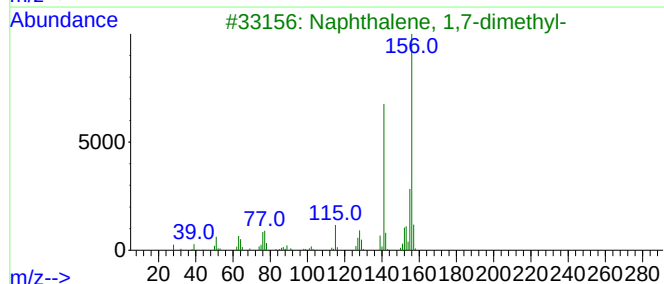
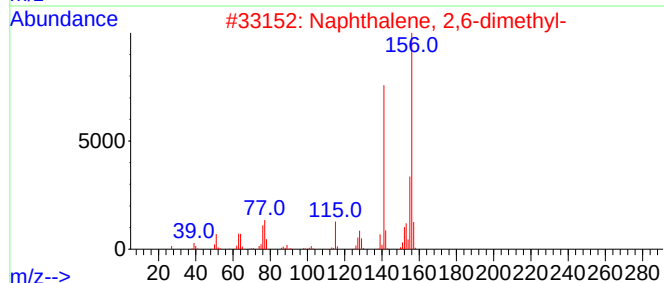
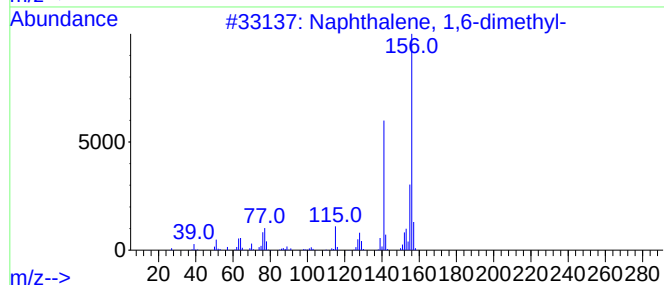
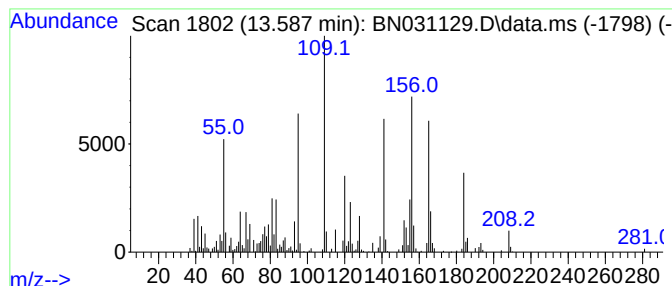
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Quant Title : SVOA CALIBRATION

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

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Peak Number 39 unknown-07 Concentration Rank 45

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.587 | 2.86 ng/ul | 188549 | Acenaphthene-d10 | 14.281 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 48   |
| 2    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 48   |
| 3    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 48   |
| 4    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 48   |
| 5    |    |   | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 48   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

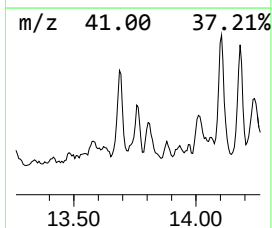
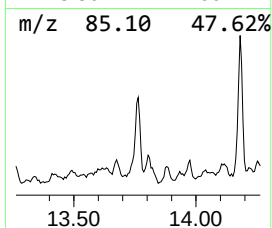
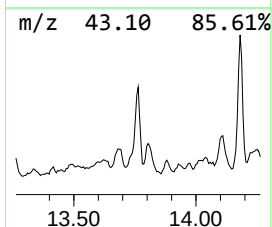
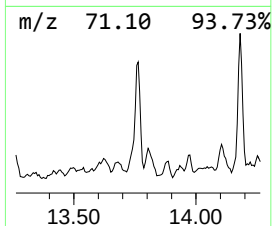
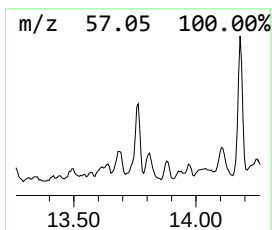
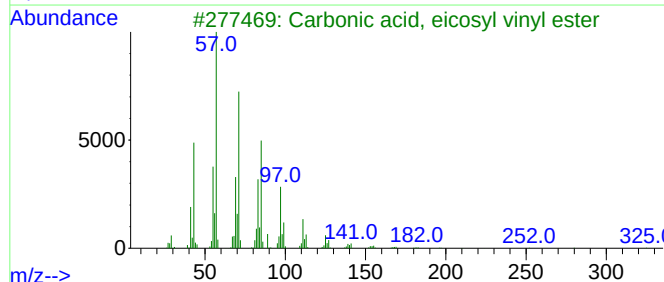
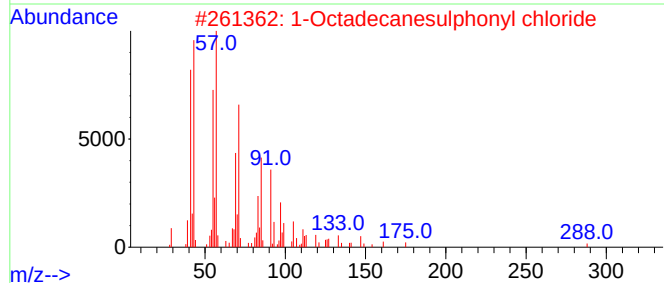
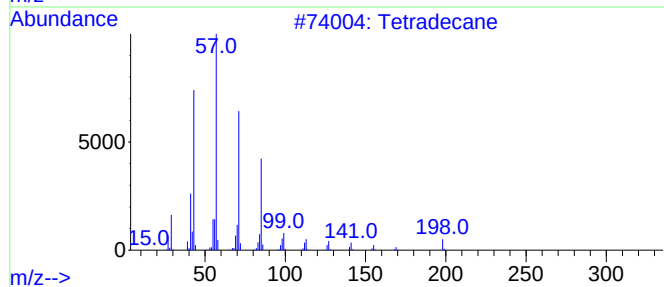
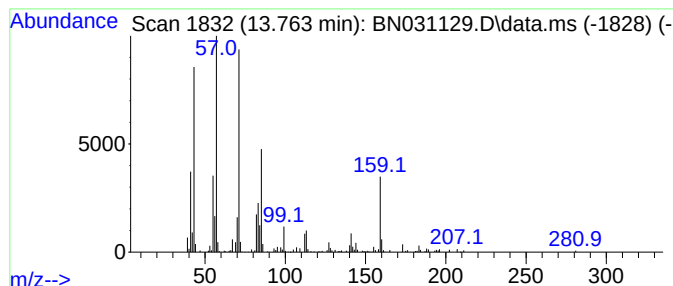
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Quant Title : SVOA CALIBRATION

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

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Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 31

| R.T.      | EstConc                            | Area   | Relative to ISTD | R.T.         |      |
|-----------|------------------------------------|--------|------------------|--------------|------|
| 13.763    | 4.96 ng/ul                         | 327548 | Acenaphthene-d10 | 14.281       |      |
| Hit# of 5 | Tentative ID                       | MW     | MolForm          | CAS#         | Qual |
| 1         | Tetradecane                        | 198    | C14H30           | 000629-59-4  | 72   |
| 2         | 1-Octadecanesulphonyl chloride     | 352    | C18H37ClO2S      | 1000342-70-4 | 72   |
| 3         | Carbonic acid, eicosyl vinyl ester | 368    | C23H44O3         | 1000382-54-3 | 72   |
| 4         | Dodecane, 4-methyl-                | 184    | C13H28           | 006117-97-1  | 70   |
| 5         | Undecane, 3,8-dimethyl-            | 184    | C13H28           | 017301-30-3  | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

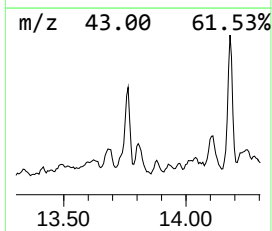
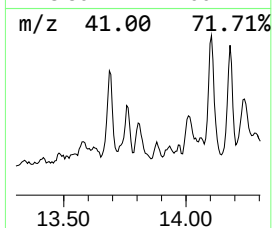
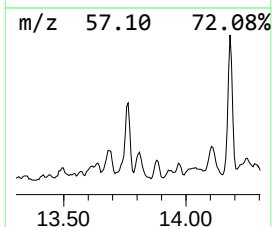
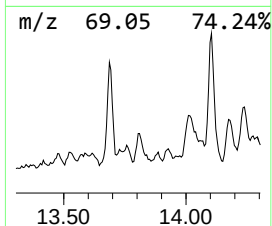
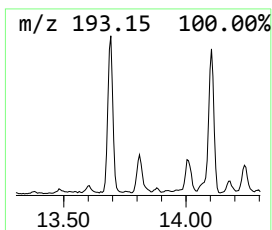
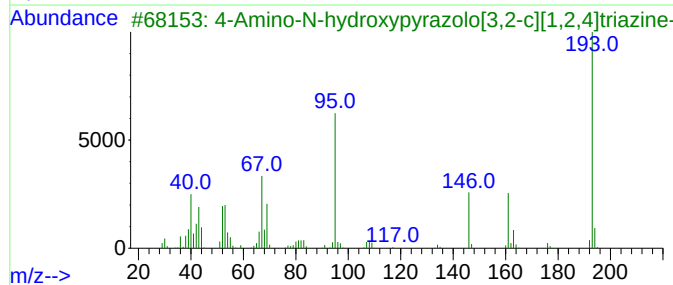
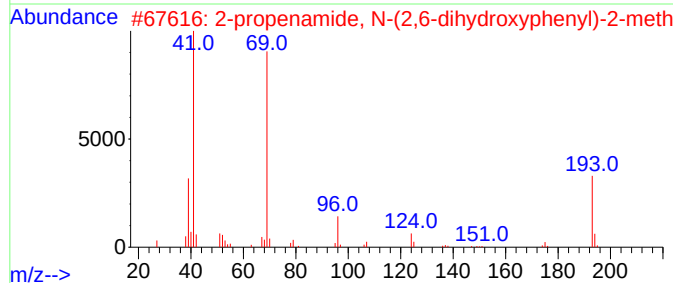
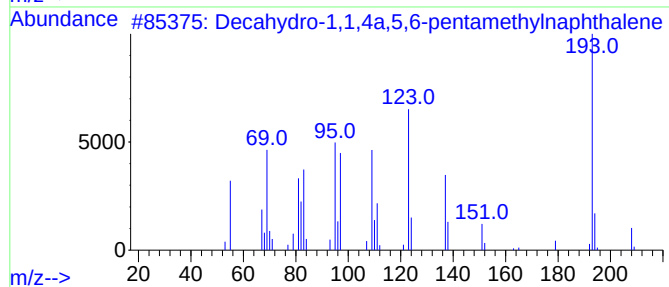
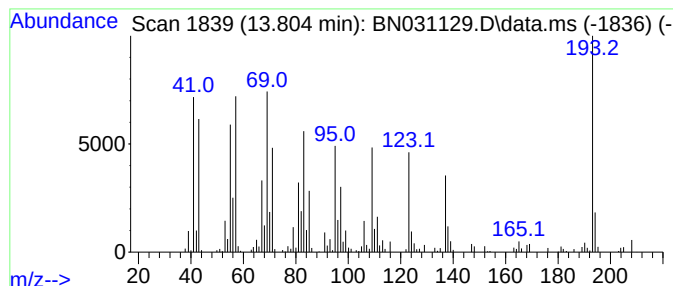
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Quant Title : SVOA CALIBRATION

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

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Peak Number 41 unknown-08 Concentration Rank 33

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.804 | 4.70 ng/ul | 309941 | Acenaphthene-d10 | 14.281 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Decahydro-1,1,4a,5,6-pentamethyl... | 208 | C15H28    | 080655-44-3  | 46   |
| 2    |    |   | 2-propenamide, N-(2,6-dihydroxyp... | 193 | C10H11NO3 | 1000395-98-7 | 35   |
| 3    |    |   | 4-Amino-N-hydroxypyrazolo[3,2-c]... | 193 | C6H7N7O   | 1000443-50-2 | 14   |
| 4    |    |   | 3-Decene, 2,2-dimethyl-, (E)-       | 168 | C12H24    | 055499-02-0  | 14   |
| 5    |    |   | Iminostilbene                       | 193 | C14H11N   | 000256-96-2  | 14   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

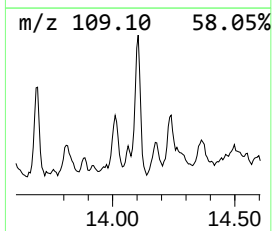
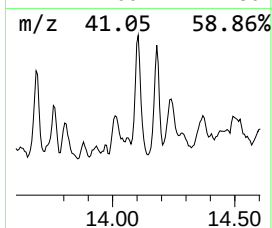
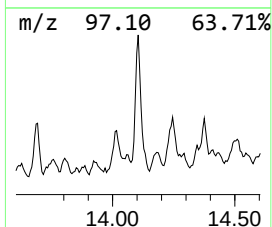
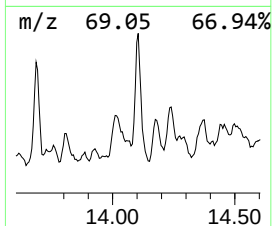
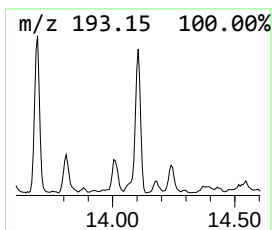
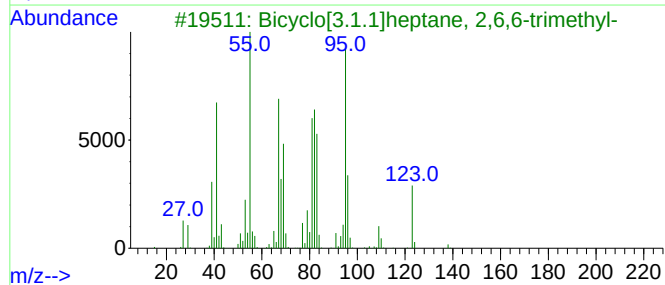
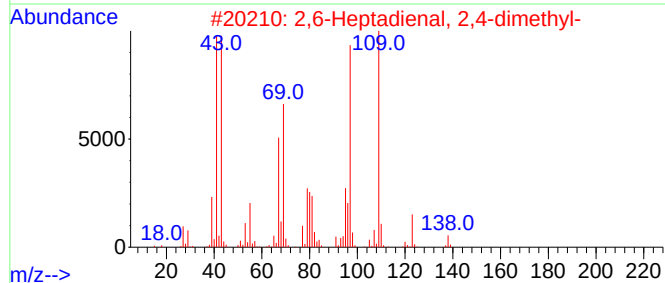
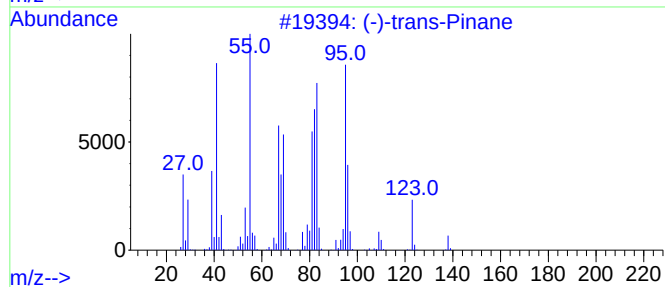
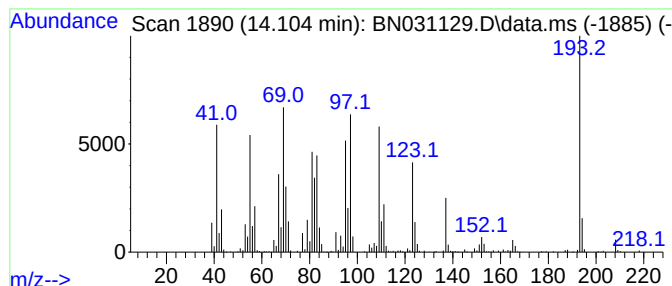
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Quant Title : SVOA CALIBRATION

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

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Peak Number 42 unknown-09 Concentration Rank 7

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 14.104 | 16.50 ng/ul | 1089010 | Acenaphthene-d10 | 14.281 |

| Hit# | of                                  | 5 | Tentative ID | MW  | MolForm  | CAS#        | Qual |
|------|-------------------------------------|---|--------------|-----|----------|-------------|------|
| 1    | (-)-trans-Pinane                    |   |              | 138 | C10H18   | 033626-25-4 | 35   |
| 2    | 2,6-Heptadienal, 2,4-dimethyl-      |   |              | 138 | C9H14O   | 085136-08-9 | 35   |
| 3    | Bicyclo[3.1.1]heptane, 2,6,6-tri... |   |              | 138 | C10H18   | 000473-55-2 | 30   |
| 4    | Hexadecanedinitrile                 |   |              | 248 | C16H28N2 | 006812-44-8 | 22   |
| 5    | Cyclohexane, 1,2-dimethyl-3-pent... |   |              | 224 | C16H32   | 062376-17-4 | 18   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

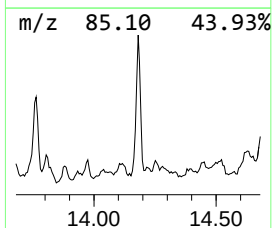
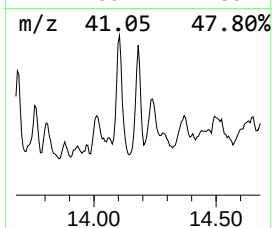
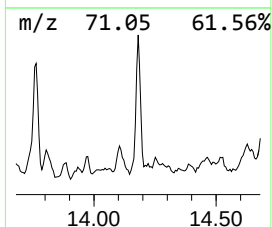
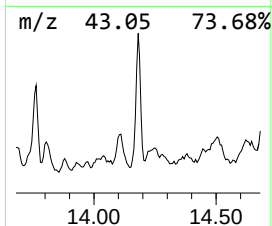
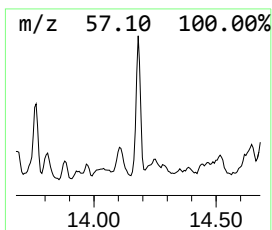
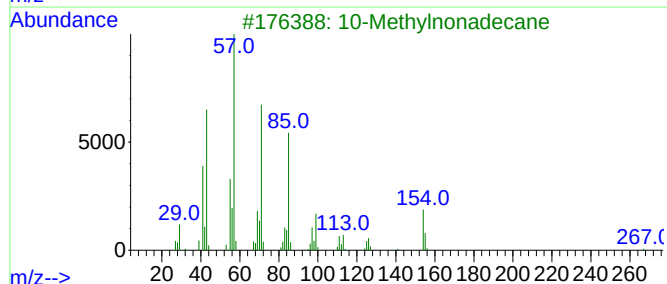
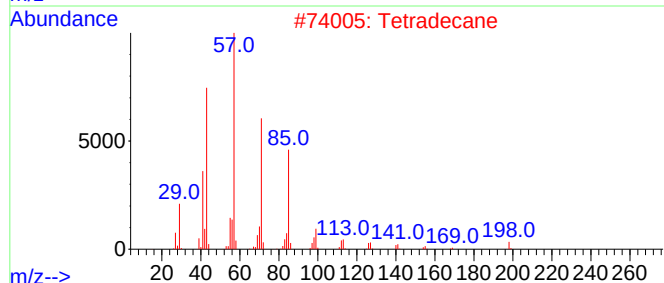
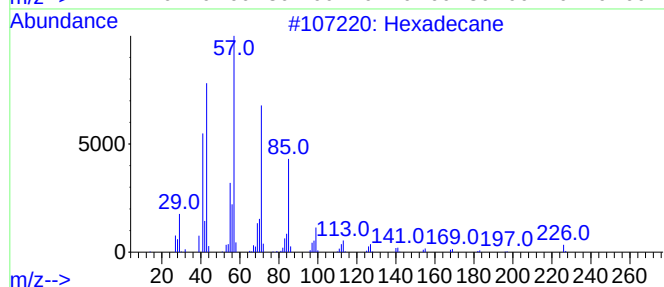
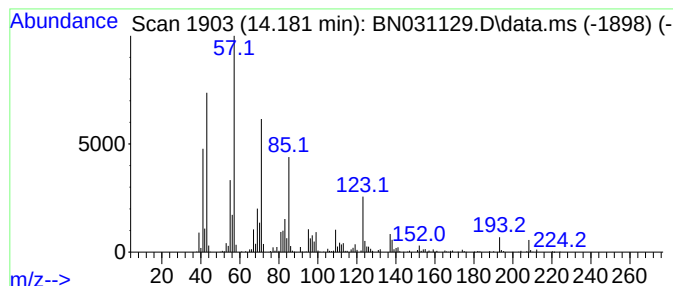
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Quant Title : SVOA CALIBRATION

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

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Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T.      | EstConc             | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|---------------------|--------|------------------|---------|-------------|------|
| 14.181    | 10.84 ng/ul         | 715416 | Acenaphthene-d10 |         | 14.281      |      |
| Hit# of 5 | Tentative ID        |        | MW               | MolForm | CAS#        | Qual |
| 1         | Hexadecane          |        | 226              | C16H34  | 000544-76-3 | 58   |
| 2         | Tetradecane         |        | 198              | C14H30  | 000629-59-4 | 58   |
| 3         | 10-Methylnonadecane |        | 282              | C20H42  | 056862-62-5 | 58   |
| 4         | Nonadecane          |        | 268              | C19H40  | 000629-92-5 | 58   |
| 5         | Pentadecane         |        | 212              | C15H32  | 000629-62-9 | 55   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

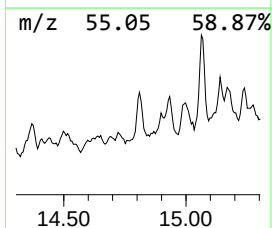
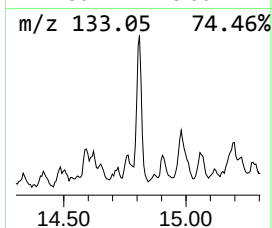
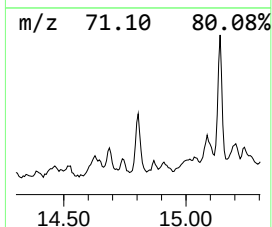
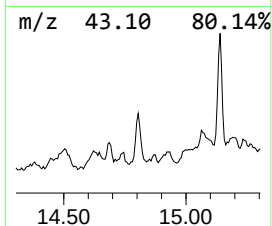
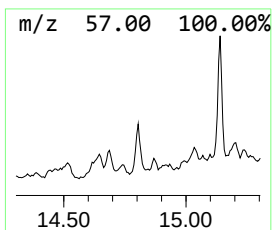
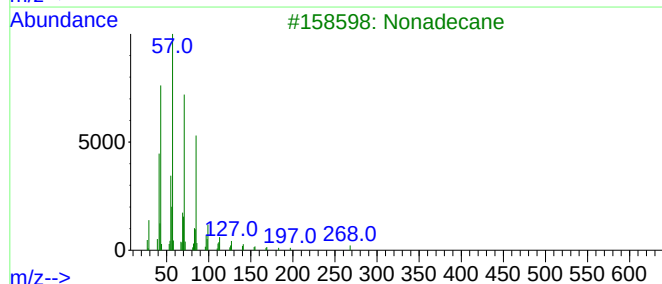
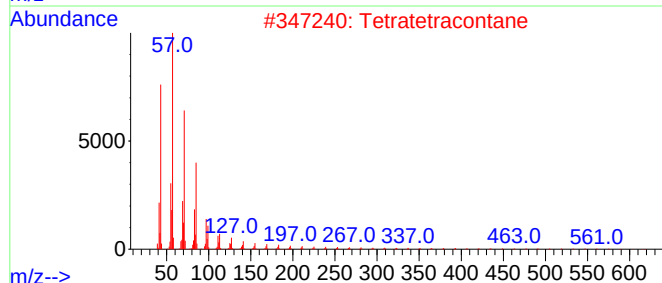
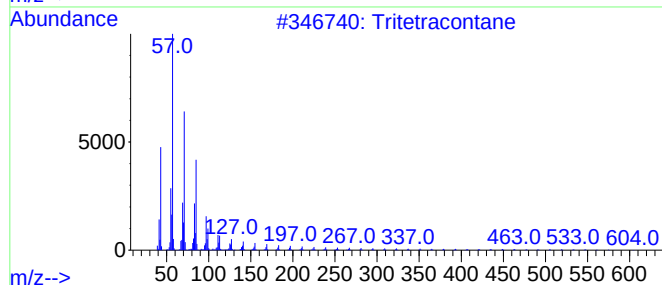
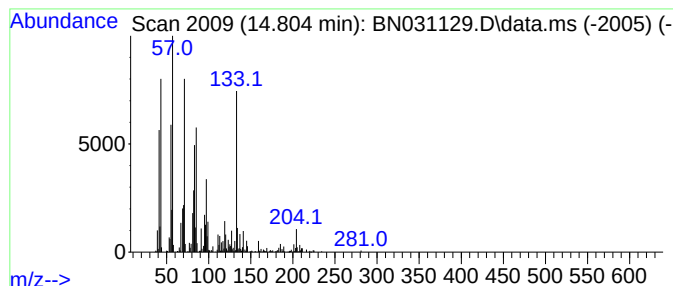
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Quant Title : SVOA CALIBRATION

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

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

| R.T.      | EstConc              | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|----------------------|--------|------------------|---------|-------------|------|
| 14.804    | 8.96 ng/ul           | 591427 | Acenaphthene-d10 |         | 14.281      |      |
| Hit# of 5 | Tentative ID         |        | MW               | MolForm | CAS#        | Qual |
| 1         | Tritetracontane      |        | 605              | C43H88  | 007098-21-7 | 43   |
| 2         | Tetratetracontane    |        | 619              | C44H90  | 007098-22-8 | 43   |
| 3         | Nonadecane           |        | 268              | C19H40  | 000629-92-5 | 41   |
| 4         | Tetradecane, 1-iodo- |        | 324              | C14H29I | 019218-94-1 | 38   |
| 5         | 10-Methylnonadecane  |        | 282              | C20H42  | 056862-62-5 | 38   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

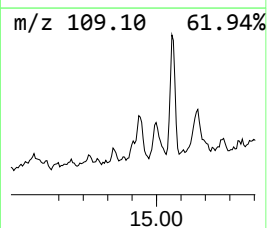
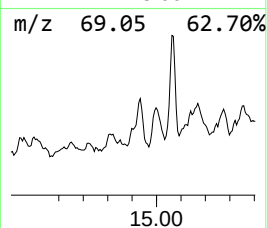
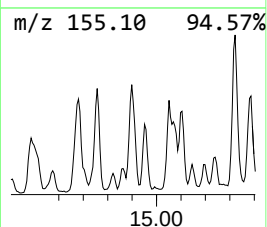
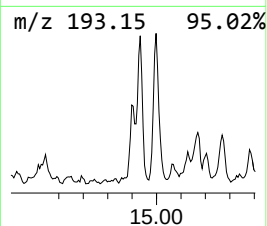
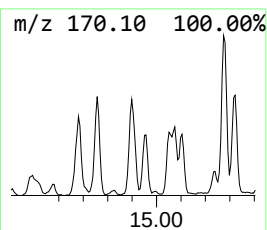
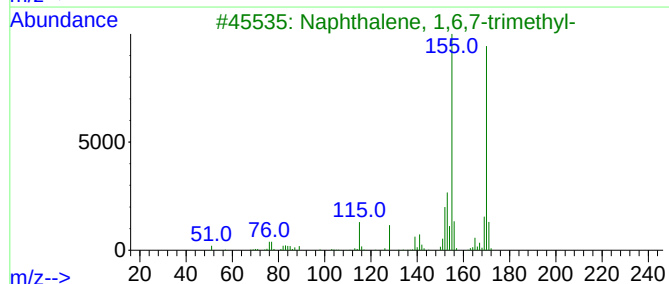
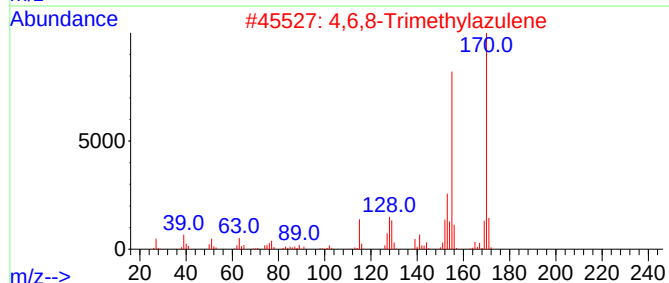
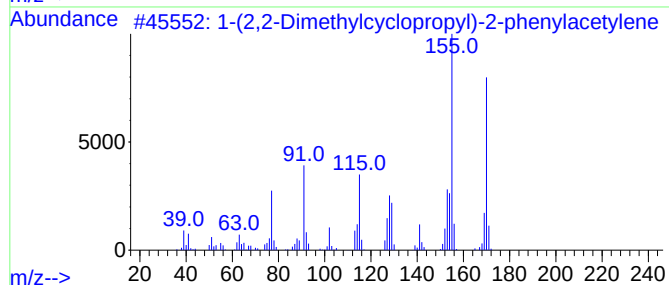
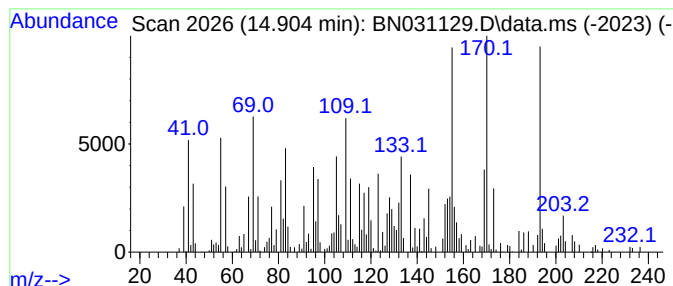
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Quant Title : SVOA CALIBRATION

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

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Peak Number 46 1-(2,2-Dimethylcyclopropyl)... Concentration Rank 38

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 14.904 | 3.75 ng/ul | 247429 | Acenaphthene-d10 | 14.281 |

| Hit# | of                                  | 5 | Tentative ID | MW     | MolForm      | CAS# | Qual |
|------|-------------------------------------|---|--------------|--------|--------------|------|------|
| 1    | 1-(2,2-Dimethylcyclopropyl)-2-ph... |   | 170          | C13H14 | 1000294-91-1 | 55   |      |
| 2    | 4,6,8-Trimethylazulene              |   | 170          | C13H14 | 000941-81-1  | 42   |      |
| 3    | Naphthalene, 1,6,7-trimethyl-       |   | 170          | C13H14 | 002245-38-7  | 42   |      |
| 4    | 4,6,8-Trimethylazulene              |   | 170          | C13H14 | 000941-81-1  | 42   |      |
| 5    | Naphthalene, 1,4,6-trimethyl-       |   | 170          | C13H14 | 002131-42-2  | 42   |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

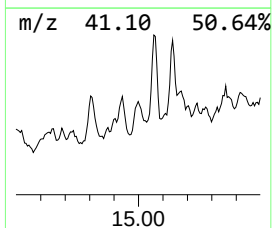
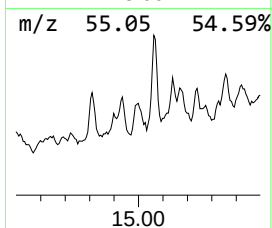
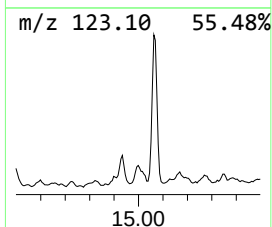
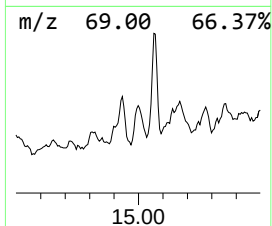
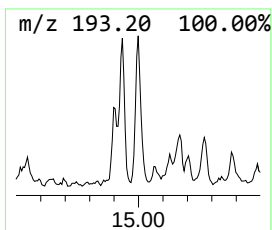
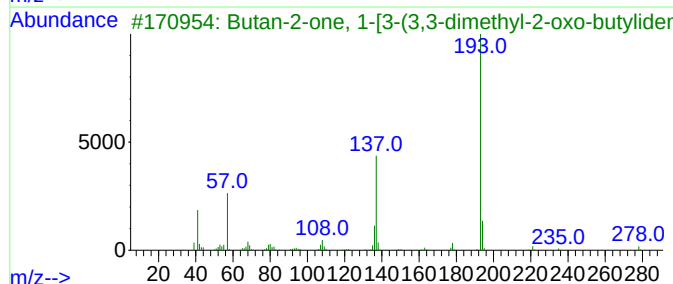
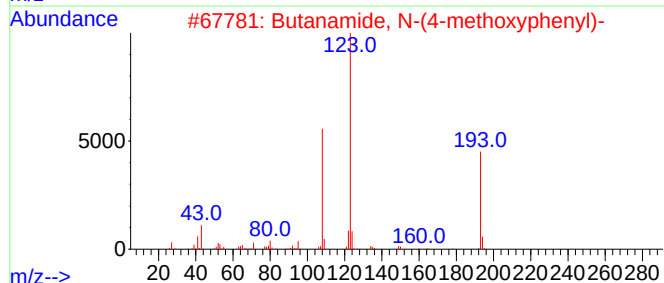
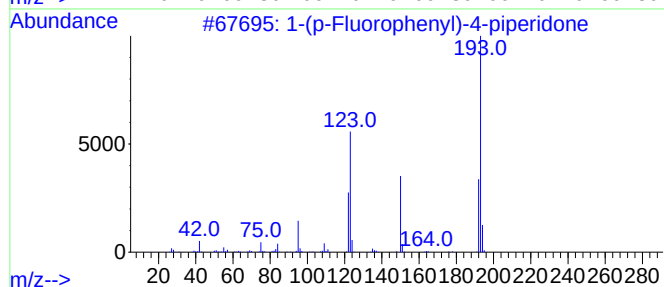
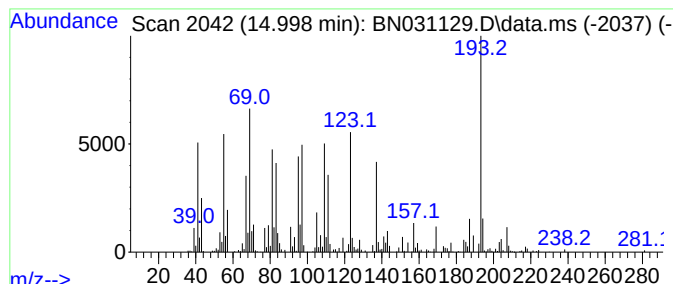
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Quant Title : SVOA CALIBRATION

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

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Peak Number 47 unknown-10 Concentration Rank 15

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 14.998    | 9.05 ng/ul                          | 597512 | Acenaphthene-d10 | 14.281       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 1-(p-Fluorophenyl)-4-piperidone     | 193    | C11H12FNO        | 1000238-56-7 | 27   |
| 2         | Butanamide, N-(4-methoxyphenyl)-    | 193    | C11H15NO2        | 1000306-91-5 | 22   |
| 3         | Butan-2-one, 1-[3-(3,3-dimethyl-... | 278    | C16H26N2O2       | 1000303-34-2 | 22   |
| 4         | Cyclohexane, 1,1',1''-(1-ethanyl... | 276    | C20H36           | 055682-86-5  | 22   |
| 5         | 6-Nitroundec-5-ene                  | 199    | C11H21NO2        | 1000192-40-3 | 14   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

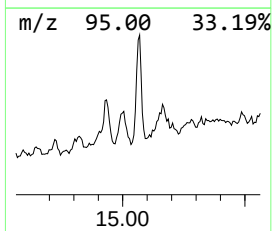
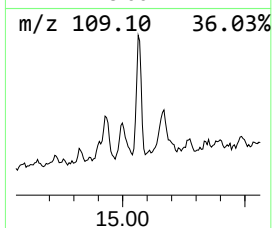
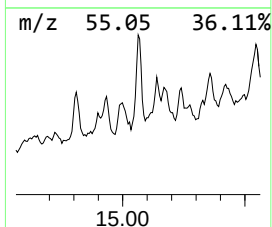
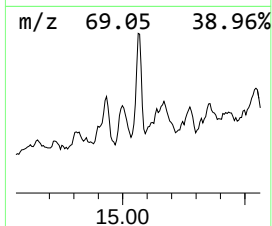
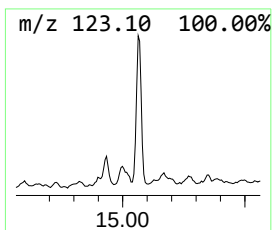
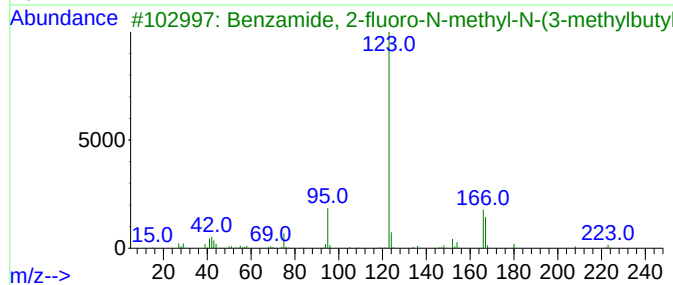
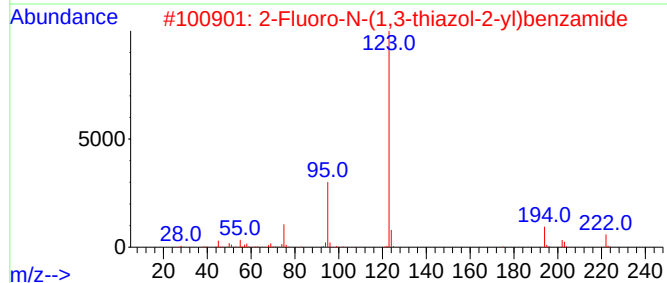
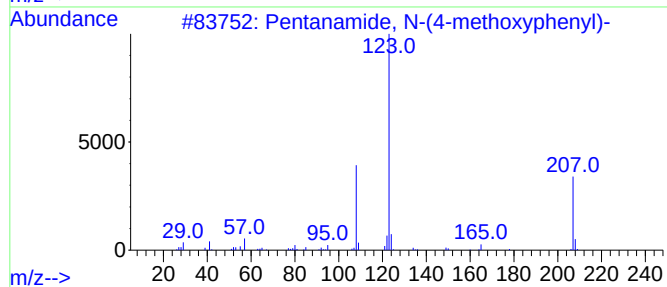
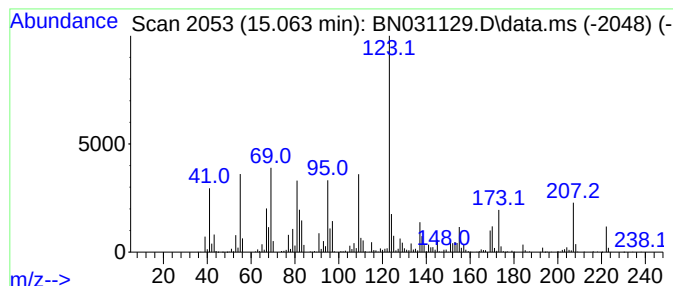
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Quant Title : SVOA CALIBRATION

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

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Peak Number 48 unknown-11 Concentration Rank 4

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 15.063 | 20.27 ng/ul | 1337720 | Acenaphthene-d10 | 14.281 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Pentanamide, N-(4-methoxyphenyl)-   | 207 | C12H17NO2  | 1000306-89-3 | 43   |
| 2    |    |   | 2-Fluoro-N-(1,3-thiazol-2-yl)ben... | 222 | C10H7FN2OS | 000319-38-0  | 43   |
| 3    |    |   | Benzamide, 2-fluoro-N-methyl-N-(... | 223 | C13H18FNO  | 1000421-27-0 | 38   |
| 4    |    |   | 4-Fluorobenzoic acid, 3-methylbu... | 208 | C12H13FO2  | 1000299-15-1 | 38   |
| 5    |    |   | 3-Fluorobenzoic acid, 3-methylbu... | 208 | C12H13FO2  | 154559-38-3  | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

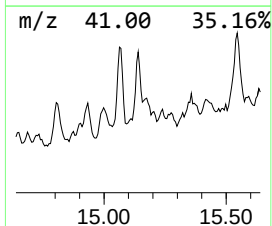
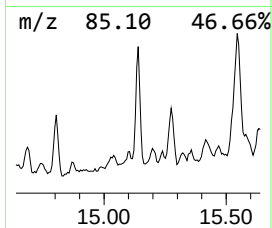
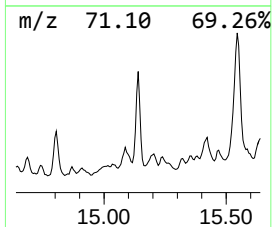
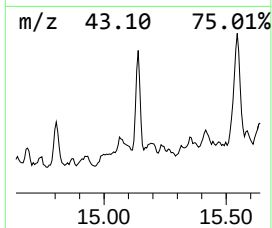
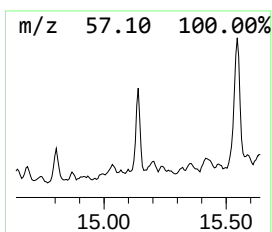
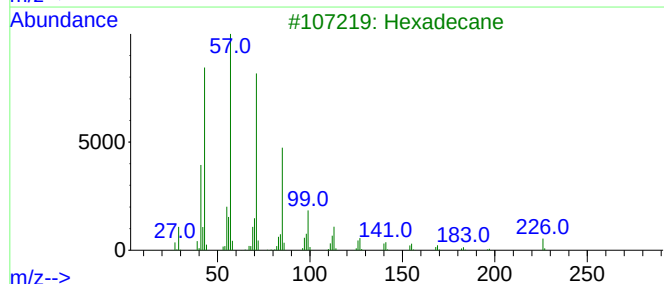
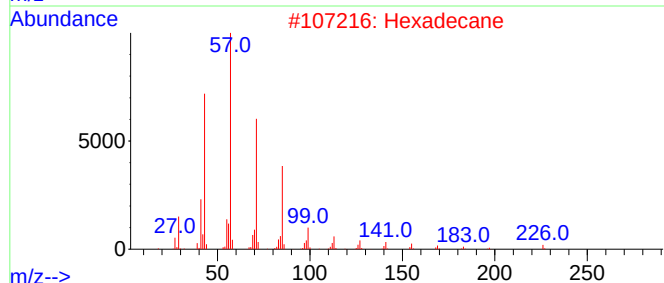
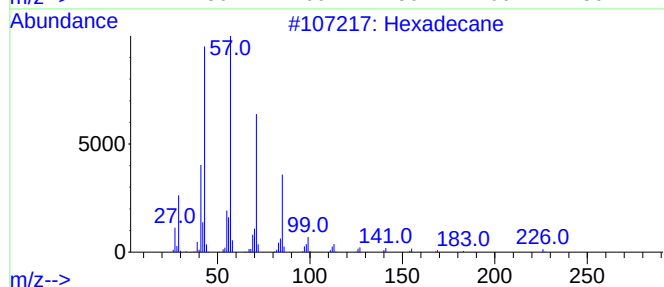
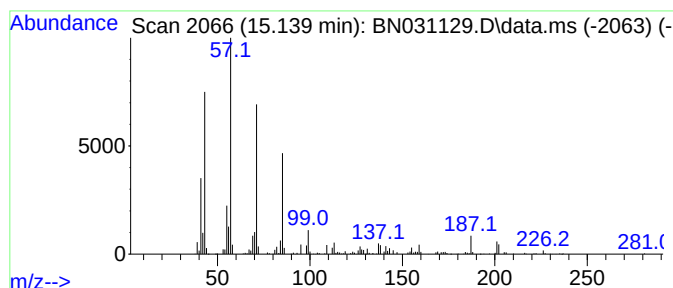
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 49 (DEL) Alkane: Straight-Chai... Concentration Rank 23

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 15.139 | 6.15 ng/ul | 405733 | Acenaphthene-d10 | 14.281 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 90   |
| 2    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 90   |
| 3    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 83   |
| 4    |    |   | Tetradecane  | 198 | C14H30  | 000629-59-4 | 74   |
| 5    |    |   | Pentadecane  | 212 | C15H32  | 000629-62-9 | 74   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

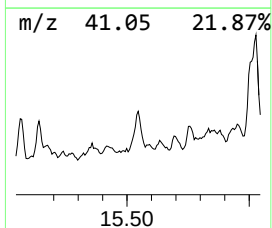
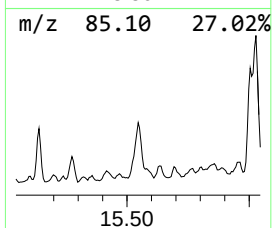
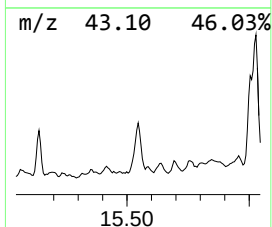
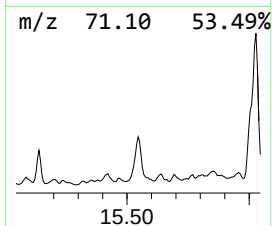
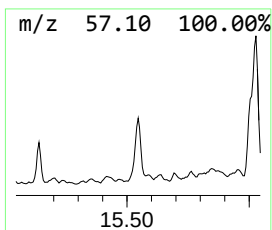
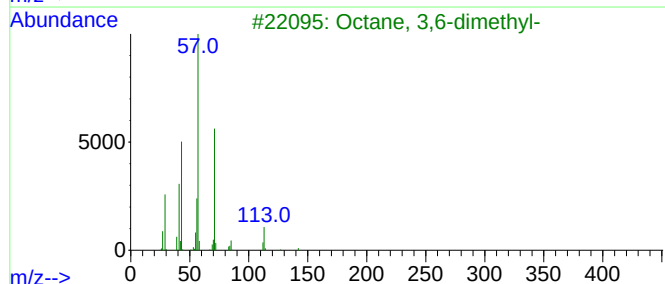
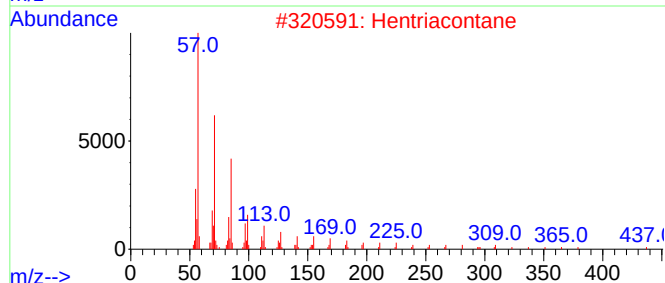
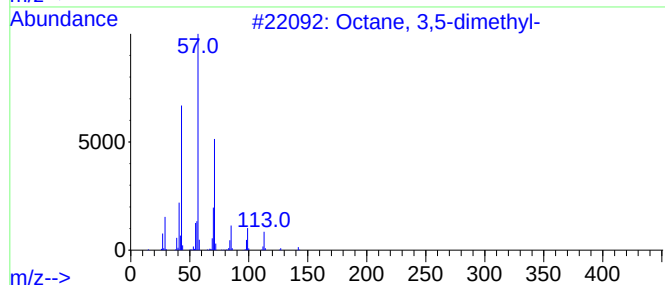
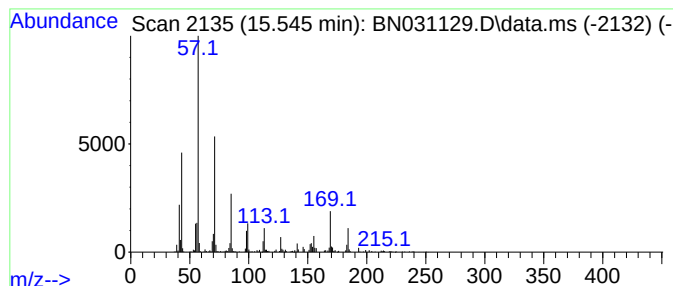
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Quant Title : SVOA CALIBRATION

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

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Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 12

| R.T.      | EstConc                            | Area   | Relative to ISTD | R.T.        |      |
|-----------|------------------------------------|--------|------------------|-------------|------|
| 15.545    | 10.63 ng/ul                        | 701365 | Acenaphthene-d10 | 14.281      |      |
| Hit# of 5 | Tentative ID                       | MW     | MolForm          | CAS#        | Qual |
| 1         | Octane, 3,5-dimethyl-              | 142    | C10H22           | 015869-93-9 | 62   |
| 2         | Hentriacontane                     | 437    | C31H64           | 000630-04-6 | 58   |
| 3         | Octane, 3,6-dimethyl-              | 142    | C10H22           | 015869-94-0 | 55   |
| 4         | Octane, 2,6-dimethyl-              | 142    | C10H22           | 002051-30-1 | 55   |
| 5         | Hexadecane, 2,6,10,14-tetramethyl- | 282    | C20H42           | 000638-36-8 | 52   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

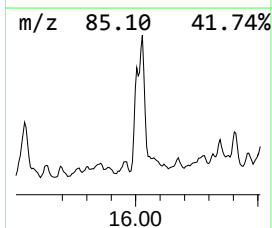
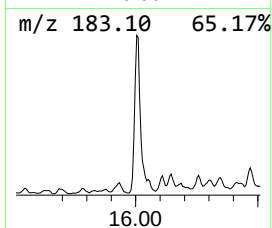
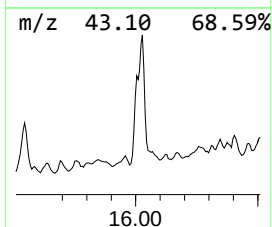
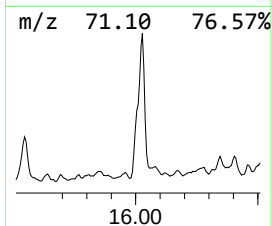
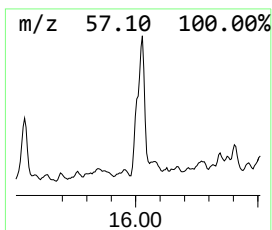
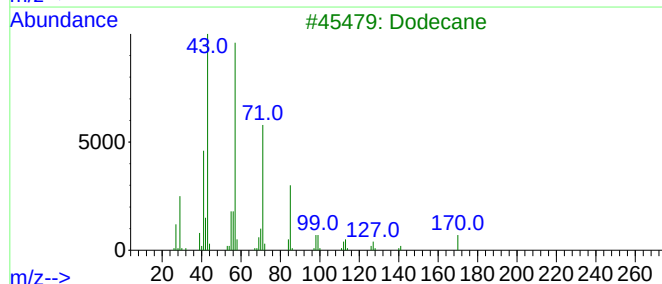
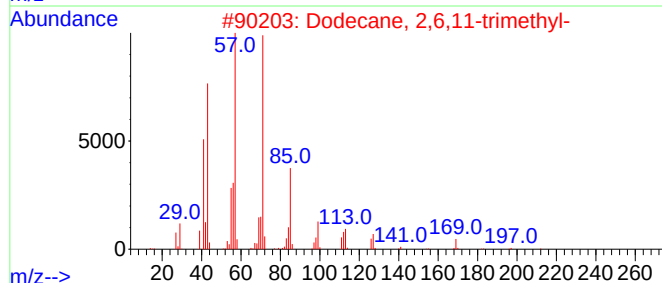
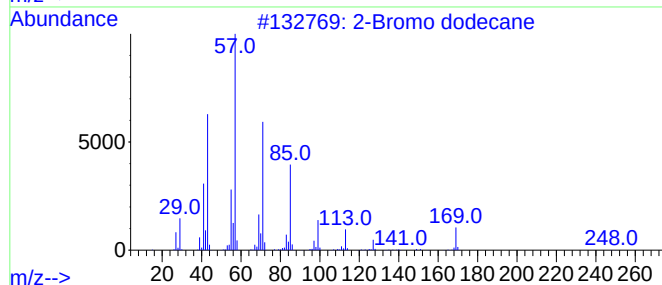
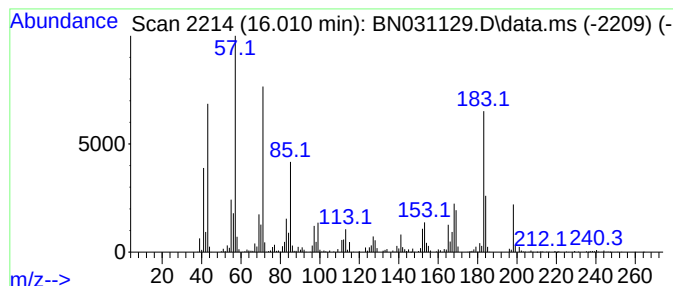
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Quant Title : SVOA CALIBRATION

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

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Peak Number 51 2-Bromo dodecane Concentration Rank 5

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 16.010 | 18.35 ng/ul | 1879180 | Phenanthrene-d10 | 17.033 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Bromo dodecane                    | 248 | C12H25Br | 013187-99-0 | 89   |
| 2    |    |   | Dodecane, 2,6,11-trimethyl-         | 212 | C15H32   | 031295-56-4 | 78   |
| 3    |    |   | Dodecane                            | 170 | C12H26   | 000112-40-3 | 60   |
| 4    |    |   | 6-Isopropyl-1,4-dimethylnaphthalene | 198 | C15H18   | 000489-77-0 | 60   |
| 5    |    |   | Naphthalene, 1,6-dimethyl-4-(1-m... | 198 | C15H18   | 000483-78-3 | 52   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

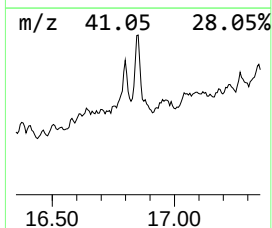
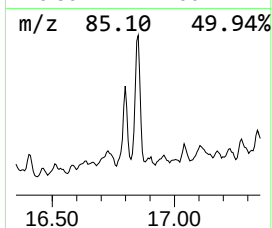
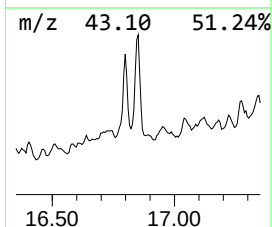
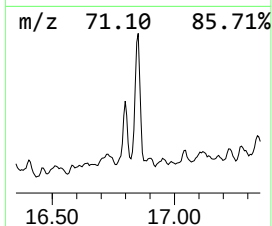
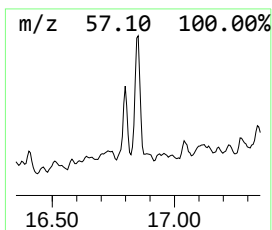
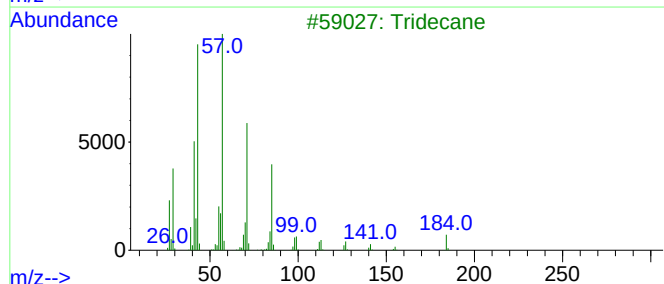
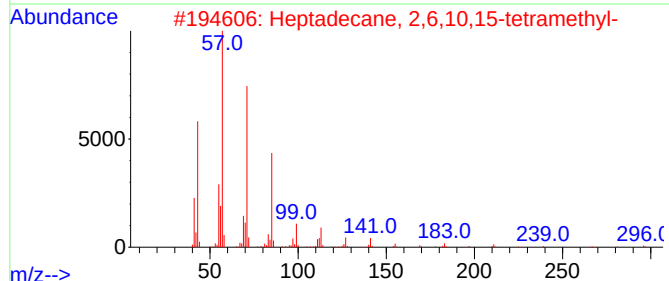
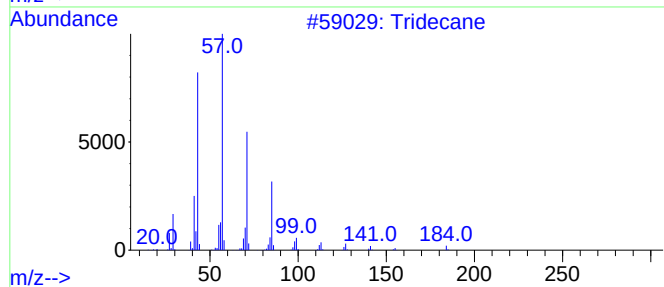
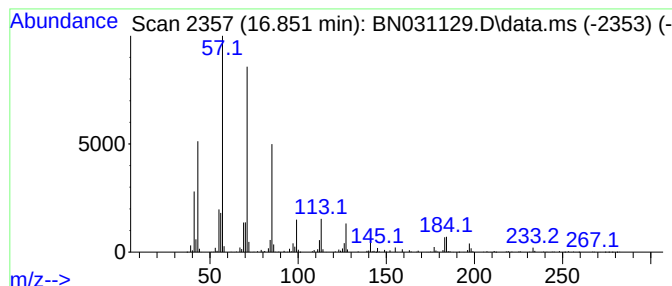
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Quant Title : SVOA CALIBRATION

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

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Peak Number 52 (DEL) Alkane: Straight-Chai... Concentration Rank 2

| R.T.      | EstConc                             | Area    | Relative to ISTD |          | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|----------|-------------|------|
| 16.851    | 22.94 ng/ul                         | 2348450 | Phenanthrene-d10 |          | 17.033      |      |
| Hit# of 5 | Tentative ID                        |         | MW               | MolForm  | CAS#        | Qual |
| 1         | Tridecane                           |         | 184              | C13H28   | 000629-50-5 | 91   |
| 2         | Heptadecane, 2,6,10,15-tetramethyl- |         | 296              | C21H44   | 054833-48-6 | 87   |
| 3         | Tridecane                           |         | 184              | C13H28   | 000629-50-5 | 87   |
| 4         | Octadecane, 1-iodo-                 |         | 380              | C18H37I  | 000629-93-6 | 86   |
| 5         | 2-Bromotetradecane                  |         | 276              | C14H29Br | 074036-95-6 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION

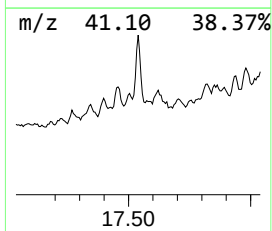
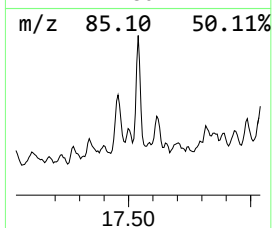
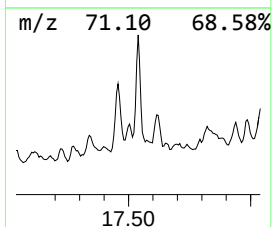
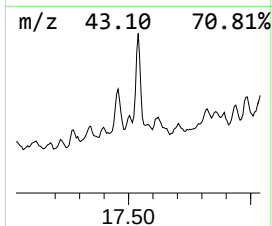
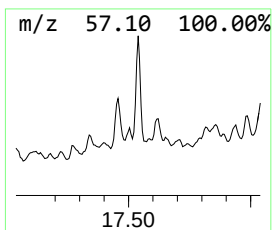
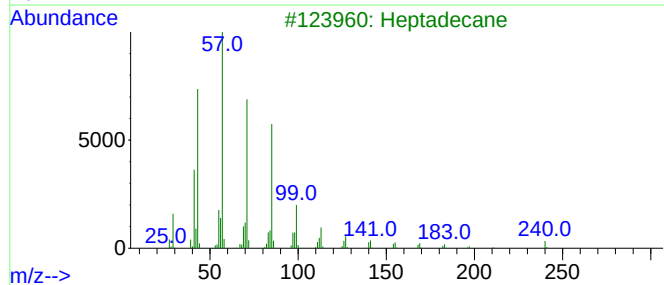
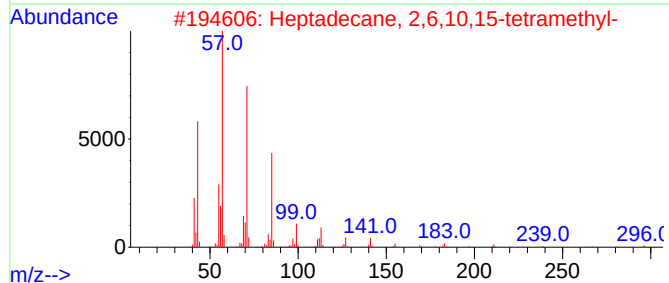
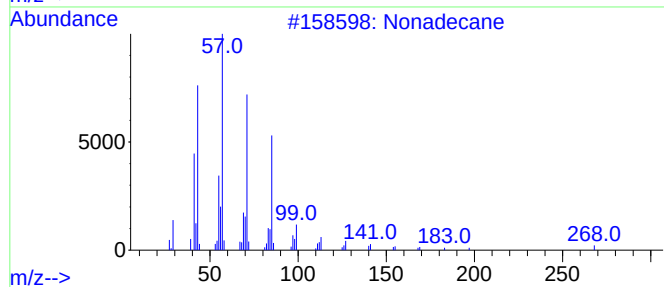
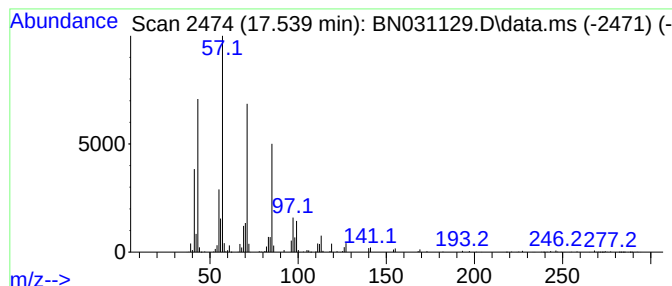
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TIC Integration Parameters: LSCINT.P

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Peak Number 53 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 17.539 | 18.22 ng/ul | 1865790 | Phenanthrene-d10 | 17.033 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Nonadecane                          | 268 | C19H40  | 000629-92-5 | 90   |
| 2    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 87   |
| 3    |    |   | Heptadecane                         | 240 | C17H36  | 000629-78-7 | 87   |
| 4    |    |   | Tritetracontane                     | 605 | C43H88  | 007098-21-7 | 86   |
| 5    |    |   | Tetradecane                         | 198 | C14H30  | 000629-59-4 | 83   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
Data File : BN031129.D  
Acq On : 22 Apr 2024 19:35  
Operator : MA/JU  
Sample : P2037-23  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8

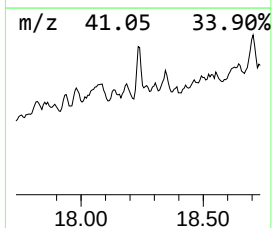
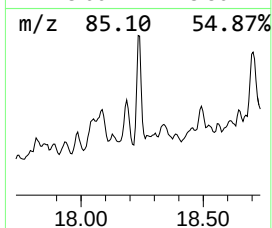
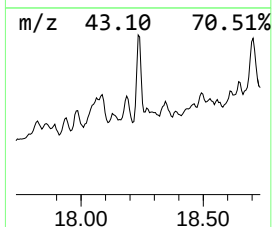
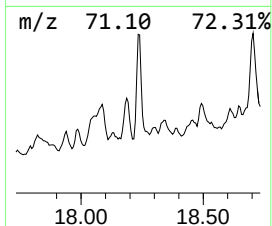
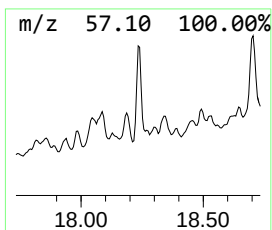
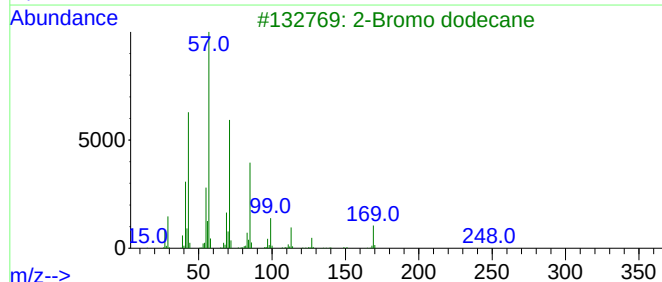
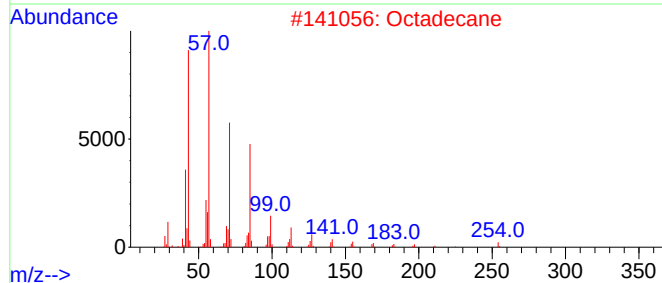
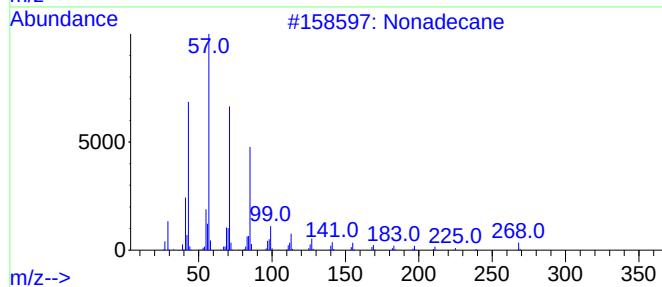
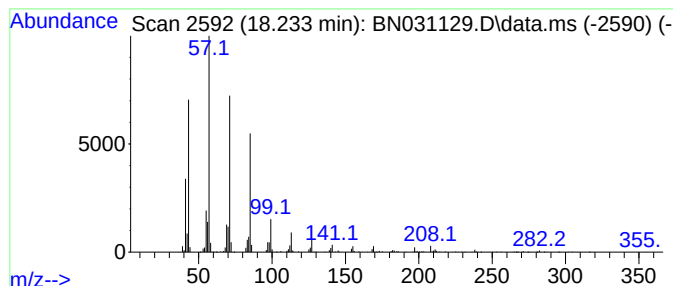
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 54 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 18.233 | 13.40 ng/ul | 1372130 | Phenanthrene-d10 | 17.033 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|---------------------|-----|----------|-------------|------|
| 1    |    |   | Nonadecane          | 268 | C19H40   | 000629-92-5 | 91   |
| 2    |    |   | Octadecane          | 254 | C18H38   | 000593-45-3 | 91   |
| 3    |    |   | 2-Bromo dodecane    | 248 | C12H25Br | 013187-99-0 | 90   |
| 4    |    |   | Nonadecane          | 268 | C19H40   | 000629-92-5 | 90   |
| 5    |    |   | Octadecane, 1-iodo- | 380 | C18H37I  | 000629-93-6 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042224\  
 Data File : BN031129.D  
 Acq On : 22 Apr 2024 19:35  
 Operator : MA/JU  
 Sample : P2037-23  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 E1CX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
 Quant Title : SVOA CALIBRATION

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

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
|                    |        |         |       |          | #                     | RT     | Resp    | Conc |
| unknown-01         | 3.417  | 2.3     | ng/ul | 49584    | 1                     | 7.634  | 439605  | 20.0 |
| Cyclohexanamine... | 4.622  | 4.6     | ng/ul | 100791   | 1                     | 7.634  | 439605  | 20.0 |
| (DEL) Alkane: S... | 4.664  | 11.5    | ng/ul | 253091   | 1                     | 7.634  | 439605  | 20.0 |
| (DEL) Alkane: S... | 4.787  | 4.9     | ng/ul | 106960   | 1                     | 7.634  | 439605  | 20.0 |
| 3,7-Dimethyl-3-... | 4.828  | 4.2     | ng/ul | 91167    | 1                     | 7.634  | 439605  | 20.0 |
| (DEL) Alkane: C... | 4.911  | 27.2    | ng/ul | 597155   | 1                     | 7.634  | 439605  | 20.0 |
| (DEL) Alkane: C... | 5.011  | 2.3     | ng/ul | 51346    | 1                     | 7.634  | 439605  | 20.0 |
| Oxalic acid, cy... | 5.069  | 8.6     | ng/ul | 189140   | 1                     | 7.634  | 439605  | 20.0 |
| (DEL) Alkane: S... | 5.181  | 7.6     | ng/ul | 167638   | 1                     | 7.634  | 439605  | 20.0 |
| Cyclohexanone, ... | 5.234  | 4.2     | ng/ul | 91341    | 1                     | 7.634  | 439605  | 20.0 |
| (DEL) Alkane: S... | 5.316  | 3.6     | ng/ul | 79910    | 1                     | 7.634  | 439605  | 20.0 |
| (DEL) Alkane: S... | 5.381  | 6.0     | ng/ul | 132775   | 1                     | 7.634  | 439605  | 20.0 |
| (DEL) Alkane: S... | 5.469  | 2.9     | ng/ul | 62915    | 1                     | 7.634  | 439605  | 20.0 |
| (DEL) Alkane: S... | 5.622  | 3.1     | ng/ul | 67871    | 1                     | 7.634  | 439605  | 20.0 |
| 1-Hexanol, 2-et... | 7.705  | 5.8     | ng/ul | 127099   | 1                     | 7.634  | 439605  | 20.0 |
| (DEL) Alkane: S... | 8.281  | 2.5     | ng/ul | 55542    | 1                     | 7.634  | 439605  | 20.0 |
| (DEL) Alkane: S... | 8.587  | 3.0     | ng/ul | 66835    | 1                     | 7.634  | 439605  | 20.0 |
| Benzene, 2-ethy... | 8.728  | 6.0     | ng/ul | 131129   | 1                     | 7.634  | 439605  | 20.0 |
| (DEL) Alkane: S... | 8.846  | 3.1     | ng/ul | 67972    | 1                     | 7.634  | 439605  | 20.0 |
| unknown-02         | 8.963  | 5.3     | ng/ul | 117131   | 1                     | 7.634  | 439605  | 20.0 |
| tert-Nonyl merc... | 9.263  | 2.6     | ng/ul | 110273   | 2                     | 10.410 | 831719  | 20.0 |
| unknown-03         | 10.246 | 10.1    | ng/ul | 419287   | 2                     | 10.410 | 831719  | 20.0 |
| 1-Phenoxypropan... | 11.157 | 2.1     | ng/ul | 86339    | 2                     | 10.410 | 831719  | 20.0 |
| Sulfurous acid,... | 11.434 | 3.1     | ng/ul | 128481   | 2                     | 10.410 | 831719  | 20.0 |
| unknown-04         | 11.610 | 2.6     | ng/ul | 108594   | 2                     | 10.410 | 831719  | 20.0 |
| (DEL) Alkane: S... | 11.840 | 6.3     | ng/ul | 264237   | 2                     | 10.410 | 831719  | 20.0 |
| Sulfurous acid,... | 12.240 | 7.9     | ng/ul | 328797   | 2                     | 10.410 | 831719  | 20.0 |
| unknown-05         | 12.304 | 2.2     | ng/ul | 91569    | 2                     | 10.410 | 831719  | 20.0 |
| Sulfurous acid,... | 12.387 | 9.9     | ng/ul | 654450   | 3                     | 14.281 | 1319970 | 20.0 |
| Oxalic acid, 2-... | 12.440 | 21.9    | ng/ul | 1448990  | 3                     | 14.281 | 1319970 | 20.0 |
| Sulfurous acid,... | 12.516 | 2.2     | ng/ul | 146319   | 3                     | 14.281 | 1319970 | 20.0 |
| 1-Iodo-2-methyl... | 12.640 | 11.5    | ng/ul | 758685   | 3                     | 14.281 | 1319970 | 20.0 |
| unknown-06         | 12.693 | 6.2     | ng/ul | 406895   | 3                     | 14.281 | 1319970 | 20.0 |
| Oxalic acid, is... | 12.804 | 4.7     | ng/ul | 307432   | 3                     | 14.281 | 1319970 | 20.0 |
| (DEL) Alkane: S... | 12.857 | 7.1     | ng/ul | 465994   | 3                     | 14.281 | 1319970 | 20.0 |
| (DEL) Alkane: S... | 12.957 | 6.0     | ng/ul | 392459   | 3                     | 14.281 | 1319970 | 20.0 |
| (DEL) Alkane: S... | 13.034 | 5.2     | ng/ul | 341813   | 3                     | 14.281 | 1319970 | 20.0 |
| Tridecane, 1-iodo- | 13.098 | 5.1     | ng/ul | 335283   | 3                     | 14.281 | 1319970 | 20.0 |
| unknown-07         | 13.587 | 2.9     | ng/ul | 188549   | 3                     | 14.281 | 1319970 | 20.0 |
| (DEL) Alkane: S... | 13.763 | 5.0     | ng/ul | 327548   | 3                     | 14.281 | 1319970 | 20.0 |
| unknown-08         | 13.804 | 4.7     | ng/ul | 309941   | 3                     | 14.281 | 1319970 | 20.0 |
| unknown-09         | 14.104 | 16.5    | ng/ul | 1089010  | 3                     | 14.281 | 1319970 | 20.0 |
| (DEL) Alkane: S... | 14.181 | 10.8    | ng/ul | 715416   | 3                     | 14.281 | 1319970 | 20.0 |
| (DEL) Alkane: S... | 14.804 | 9.0     | ng/ul | 591427   | 3                     | 14.281 | 1319970 | 20.0 |
| 1-(2,2-Dimethyl... | 14.904 | 3.8     | ng/ul | 247429   | 3                     | 14.281 | 1319970 | 20.0 |
| unknown-10         | 14.998 | 9.1     | ng/ul | 597512   | 3                     | 14.281 | 1319970 | 20.0 |
| unknown-11         | 15.063 | 20.3    | ng/ul | 1337720  | 3                     | 14.281 | 1319970 | 20.0 |
| (DEL) Alkane: S... | 15.139 | 6.2     | ng/ul | 405733   | 3                     | 14.281 | 1319970 | 20.0 |
| (DEL) Alkane: S... | 15.545 | 10.6    | ng/ul | 701365   | 3                     | 14.281 | 1319970 | 20.0 |
| 2-Bromo dodecane   | 16.010 | 18.4    | ng/ul | 1879180  | 4                     | 17.033 | 2047800 | 20.0 |
| (DEL) Alkane: S... | 16.851 | 22.9    | ng/ul | 2348450  | 4                     | 17.033 | 2047800 | 20.0 |
| (DEL) Alkane: S... | 17.539 | 18.2    | ng/ul | 1865790  | 4                     | 17.033 | 2047800 | 20.0 |

Instrument :  
BNA\_N  
ClientSampleId :  
E1CX8