

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
 Data File : BP022172.D  
 Acq On : 02 Oct 2024 22:14  
 Operator : RC/JU  
 Sample : P4016-09  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DCVW0

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\_P\Methods\SFAM-EPA-BP092424.MA.M  
 Title : SVOA CALIBRATION

Signal : TIC: BP022172.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         | 3.628       | 105           | 109         | 124          | rBV      | 390597         | 603264        | 8.65%           | 0.910%        |
| 2         | 5.752       | 464           | 470         | 477          | rVB      | 181395         | 258221        | 3.70%           | 0.389%        |
| 3         | 6.005       | 506           | 513         | 520          | rBV4     | 38273          | 81598         | 1.17%           | 0.123%        |
| 4         | 6.122       | 526           | 533         | 540          | rBV2     | 31316          | 65125         | 0.93%           | 0.098%        |
| 5         | 6.211       | 543           | 548         | 555          | rBV4     | 33404          | 68887         | 0.99%           | 0.104%        |
| 6         | 6.281       | 555           | 560         | 565          | rBV      | 78659          | 119038        | 1.71%           | 0.179%        |
| 7         | 6.393       | 576           | 579         | 585          | rVB2     | 56948          | 87370         | 1.25%           | 0.132%        |
| 8         | 6.710       | 629           | 633         | 637          | rVV      | 106638         | 162568        | 2.33%           | 0.245%        |
| 9         | 6.763       | 637           | 642         | 646          | rVV      | 126884         | 183808        | 2.63%           | 0.277%        |
| 10        | 6.881       | 653           | 662         | 677          | rVB2     | 628164         | 1188112       | 17.03%          | 1.791%        |
| 11        | 7.040       | 683           | 689         | 695          | rVV      | 392806         | 624059        | 8.95%           | 0.941%        |
| 12        | 7.240       | 718           | 723         | 732          | rVV      | 590310         | 961813        | 13.79%          | 1.450%        |
| 13        | 7.328       | 732           | 738         | 746          | rVB      | 639726         | 963486        | 13.81%          | 1.453%        |
| 14        | 7.534       | 765           | 773         | 777          | rVB4     | 36739          | 81714         | 1.17%           | 0.123%        |
| 15        | 7.616       | 778           | 787         | 792          | rBV5     | 56529          | 161676        | 2.32%           | 0.244%        |
| 16        | 7.693       | 792           | 800         | 807          | rVB2     | 458586         | 914383        | 13.11%          | 1.379%        |
| 17        | 7.763       | 807           | 812         | 818          | rVB2     | 105993         | 190227        | 2.73%           | 0.287%        |
| 18        | 7.887       | 828           | 833         | 836          | rBV2     | 68495          | 93340         | 1.34%           | 0.141%        |
| 19        | 7.981       | 846           | 849         | 856          | rVB3     | 67109          | 111964        | 1.60%           | 0.169%        |
| 20        | 8.122       | 868           | 873         | 878          | rBV4     | 45219          | 89154         | 1.28%           | 0.134%        |
| 21        | 8.222       | 883           | 890         | 896          | rBV2     | 102726         | 186983        | 2.68%           | 0.282%        |
| 22        | 8.281       | 896           | 900         | 904          | rVB      | 111280         | 158904        | 2.28%           | 0.240%        |
| 23        | 8.346       | 904           | 911         | 915          | rBV2     | 177031         | 285793        | 4.10%           | 0.431%        |
| 24        | 8.399       | 915           | 920         | 925          | rVB      | 674312         | 1055443       | 15.13%          | 1.591%        |
| 25        | 8.452       | 925           | 929         | 933          | rBV      | 128664         | 173718        | 2.49%           | 0.262%        |
| 26        | 8.681       | 962           | 968         | 976          | rVV8     | 32139          | 70976         | 1.02%           | 0.107%        |
| 27        | 8.793       | 983           | 987         | 989          | rVV2     | 55573          | 93088         | 1.33%           | 0.140%        |
| 28        | 8.834       | 989           | 994         | 1000         | rVV      | 491371         | 832843        | 11.94%          | 1.256%        |
| 29        | 8.910       | 1000          | 1007        | 1012         | rVB      | 656374         | 1042161       | 14.94%          | 1.571%        |
| 30        | 8.999       | 1017          | 1022        | 1023         | rBV2     | 43811          | 66915         | 0.96%           | 0.101%        |
| 31        | 9.028       | 1024          | 1027        | 1033         | rVV4     | 61762          | 110809        | 1.59%           | 0.167%        |
| 32        | 9.152       | 1042          | 1048        | 1058         | rVV5     | 42255          | 147223        | 2.11%           | 0.222%        |
| 33        | 9.263       | 1058          | 1067        | 1071         | rVV7     | 28987          | 71240         | 1.02%           | 0.107%        |
| 34        | 9.316       | 1071          | 1076        | 1083         | rVV2     | 44195          | 82737         | 1.19%           | 0.125%        |
| 35        | 9.393       | 1084          | 1089        | 1093         | rVV4     | 42971          | 72145         | 1.03%           | 0.109%        |
| 36        | 9.546       | 1108          | 1115        | 1121         | rBV      | 497316         | 842584        | 12.08%          | 1.270%        |

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 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 9.599  | 1121 | 1124 | 1129 | rVV5 | 48809   | 81436   | 1.17%  | 0.123% |
| 38 | 9.746  | 1140 | 1149 | 1154 | rBV2 | 72506   | 158722  | 2.28%  | 0.239% |
| 39 | 9.816  | 1154 | 1161 | 1166 | rVB3 | 61032   | 97658   | 1.40%  | 0.147% |
| 40 | 9.893  | 1167 | 1174 | 1183 | rBV3 | 76061   | 155120  | 2.22%  | 0.234% |
| 41 | 10.087 | 1197 | 1207 | 1215 | rVV  | 745491  | 1231155 | 17.65% | 1.856% |
| 42 | 10.451 | 1259 | 1269 | 1277 | rBV3 | 883928  | 1986014 | 28.47% | 2.994% |
| 43 | 10.587 | 1283 | 1292 | 1303 | rVV2 | 669529  | 1569922 | 22.50% | 2.367% |
| 44 | 10.834 | 1328 | 1334 | 1343 | rBV3 | 127579  | 230419  | 3.30%  | 0.347% |
| 45 | 10.957 | 1348 | 1355 | 1358 | rBV3 | 45955   | 89703   | 1.29%  | 0.135% |
| 46 | 11.151 | 1379 | 1388 | 1391 | rBV6 | 31954   | 78112   | 1.12%  | 0.118% |
| 47 | 11.346 | 1415 | 1421 | 1434 | rVV6 | 68276   | 187229  | 2.68%  | 0.282% |
| 48 | 11.481 | 1438 | 1444 | 1453 | rVV2 | 109588  | 224301  | 3.22%  | 0.338% |
| 49 | 11.710 | 1475 | 1483 | 1492 | rBV5 | 71902   | 159353  | 2.28%  | 0.240% |
| 50 | 11.869 | 1504 | 1510 | 1515 | rBV2 | 249095  | 424657  | 6.09%  | 0.640% |
| 51 | 11.910 | 1515 | 1517 | 1522 | rVB  | 71724   | 80825   | 1.16%  | 0.122% |
| 52 | 12.110 | 1546 | 1551 | 1558 | rVB4 | 37232   | 89964   | 1.29%  | 0.136% |
| 53 | 12.281 | 1576 | 1580 | 1586 | rVV7 | 63701   | 120230  | 1.72%  | 0.181% |
| 54 | 12.510 | 1610 | 1619 | 1623 | rBV5 | 31224   | 81891   | 1.17%  | 0.123% |
| 55 | 12.834 | 1670 | 1674 | 1678 | rBV  | 70936   | 96202   | 1.38%  | 0.145% |
| 56 | 13.122 | 1719 | 1723 | 1731 | rVB  | 249718  | 372123  | 5.33%  | 0.561% |
| 57 | 13.345 | 1757 | 1761 | 1769 | rVB8 | 33169   | 67187   | 0.96%  | 0.101% |
| 58 | 13.704 | 1817 | 1822 | 1829 | rVB  | 1294174 | 1971492 | 28.26% | 2.972% |
| 59 | 13.787 | 1829 | 1836 | 1840 | rVB3 | 98376   | 177829  | 2.55%  | 0.268% |
| 60 | 13.834 | 1840 | 1844 | 1852 | rVB6 | 43836   | 88799   | 1.27%  | 0.134% |
| 61 | 13.987 | 1864 | 1870 | 1878 | rVB  | 1298665 | 2248491 | 32.23% | 3.390% |
| 62 | 14.140 | 1893 | 1896 | 1903 | rVB7 | 35554   | 65501   | 0.94%  | 0.099% |
| 63 | 14.210 | 1903 | 1908 | 1913 | rBV  | 603853  | 764283  | 10.96% | 1.152% |
| 64 | 14.304 | 1917 | 1924 | 1932 | rVV  | 1022599 | 1594554 | 22.86% | 2.404% |
| 65 | 14.381 | 1932 | 1937 | 1944 | rVB4 | 92062   | 138738  | 1.99%  | 0.209% |
| 66 | 14.522 | 1955 | 1961 | 1969 | rBV  | 363104  | 617302  | 8.85%  | 0.931% |
| 67 | 14.710 | 1989 | 1993 | 1999 | rVB7 | 41458   | 82373   | 1.18%  | 0.124% |
| 68 | 14.834 | 2011 | 2014 | 2020 | rBV4 | 52968   | 82287   | 1.18%  | 0.124% |
| 69 | 14.898 | 2021 | 2025 | 2028 | rBV5 | 55929   | 95541   | 1.37%  | 0.144% |
| 70 | 14.934 | 2028 | 2031 | 2036 | rVB  | 124496  | 172115  | 2.47%  | 0.259% |
| 71 | 15.169 | 2064 | 2071 | 2076 | rVB  | 273763  | 463386  | 6.64%  | 0.699% |
| 72 | 15.310 | 2088 | 2095 | 2101 | rVB  | 1964762 | 2900235 | 41.57% | 4.373% |
| 73 | 15.428 | 2107 | 2115 | 2129 | rVB2 | 347298  | 640163  | 9.18%  | 0.965% |
| 74 | 15.586 | 2135 | 2142 | 2146 | rBV  | 114550  | 192510  | 2.76%  | 0.290% |
| 75 | 15.675 | 2152 | 2157 | 2164 | rVB  | 915841  | 1312985 | 18.82% | 1.980% |
| 76 | 15.804 | 2175 | 2179 | 2188 | rVB5 | 43915   | 85524   | 1.23%  | 0.129% |

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Integrator: RTE

Smoothing : OFF

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 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |         |
|-----|--------|------|------|------|------|---------|---------|---------|---------|
| 77  | 16.051 | 2216 | 2221 | 2224 | rBV  | 301063  | 477618  | 6.85%   | 0.720%  |
| 78  | 16.075 | 2224 | 2225 | 2233 | rVB  | 162310  | 169506  | 2.43%   | 0.256%  |
| 79  | 16.257 | 2252 | 2256 | 2264 | rVB  | 185377  | 260625  | 3.74%   | 0.393%  |
| 80  | 16.739 | 2332 | 2338 | 2348 | rVB  | 4670823 | 6975953 | 100.00% | 10.518% |
| 81  | 16.863 | 2355 | 2359 | 2363 | rVV  | 272562  | 344453  | 4.94%   | 0.519%  |
| 82  | 16.916 | 2363 | 2368 | 2376 | rVB2 | 143395  | 250963  | 3.60%   | 0.378%  |
| 83  | 17.092 | 2392 | 2398 | 2404 | rBV  | 1377251 | 2038212 | 29.22%  | 3.073%  |
| 84  | 17.198 | 2410 | 2416 | 2422 | rBV2 | 2332536 | 3406108 | 48.83%  | 5.135%  |
| 85  | 17.622 | 2484 | 2488 | 2493 | rBV  | 270536  | 336979  | 4.83%   | 0.508%  |
| 86  | 17.692 | 2497 | 2500 | 2503 | rVV  | 44729   | 73312   | 1.05%   | 0.111%  |
| 87  | 17.733 | 2503 | 2507 | 2511 | rVB  | 239113  | 319261  | 4.58%   | 0.481%  |
| 88  | 18.033 | 2553 | 2558 | 2564 | rBV  | 694383  | 937543  | 13.44%  | 1.414%  |
| 89  | 18.345 | 2606 | 2611 | 2617 | rBV  | 268118  | 346202  | 4.96%   | 0.522%  |
| 90  | 19.004 | 2720 | 2723 | 2730 | rVB  | 178412  | 244336  | 3.50%   | 0.368%  |
| 91  | 19.369 | 2781 | 2785 | 2796 | rBV4 | 113410  | 231177  | 3.31%   | 0.349%  |
| 92  | 19.580 | 2814 | 2821 | 2825 | rBV  | 2939835 | 4314145 | 61.84%  | 6.504%  |
| 93  | 20.174 | 2918 | 2922 | 2926 | rBV  | 161794  | 191374  | 2.74%   | 0.289%  |
| 94  | 20.686 | 3006 | 3009 | 3013 | rVB  | 136939  | 144986  | 2.08%   | 0.219%  |
| 95  | 21.204 | 3093 | 3097 | 3105 | rVB2 | 602797  | 926949  | 13.29%  | 1.398%  |
| 96  | 21.533 | 3147 | 3153 | 3162 | rBV  | 1772460 | 2564492 | 36.76%  | 3.866%  |
| 97  | 21.780 | 3192 | 3195 | 3205 | rVB2 | 132287  | 208006  | 2.98%   | 0.314%  |
| 98  | 22.410 | 3297 | 3302 | 3310 | rVB4 | 156196  | 363769  | 5.21%   | 0.548%  |
| 99  | 24.598 | 3665 | 3674 | 3686 | rVB  | 1890047 | 4835193 | 69.31%  | 7.290%  |
| 100 | 24.821 | 3704 | 3712 | 3724 | rVB  | 1076998 | 2785143 | 39.92%  | 4.199%  |

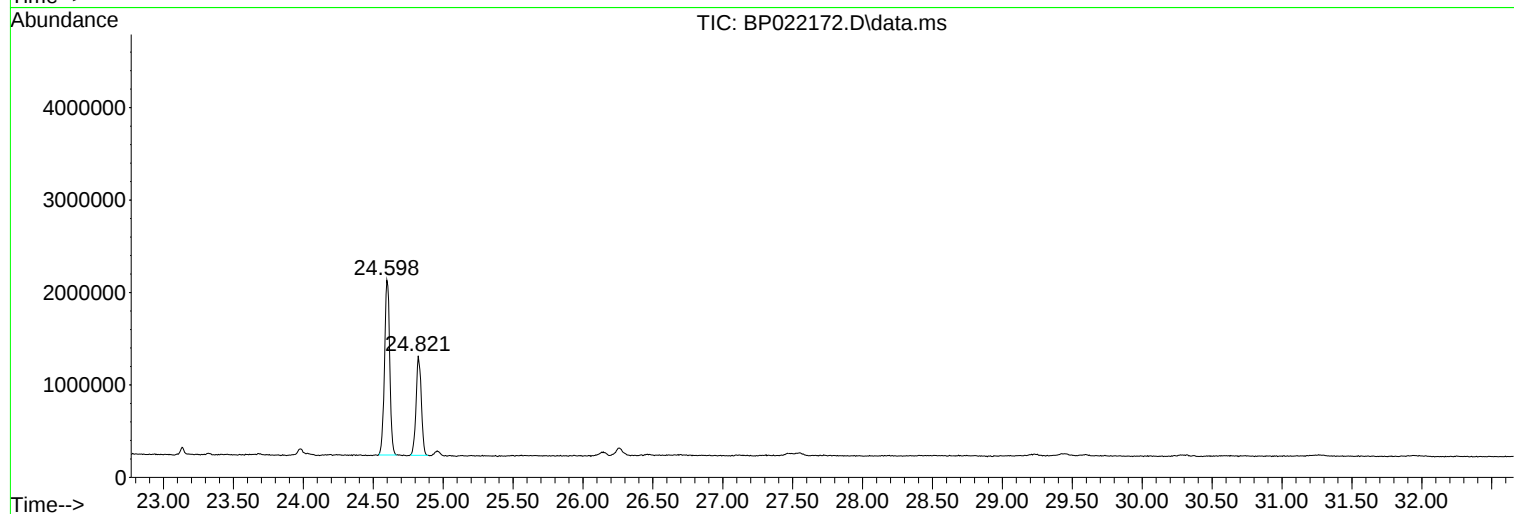
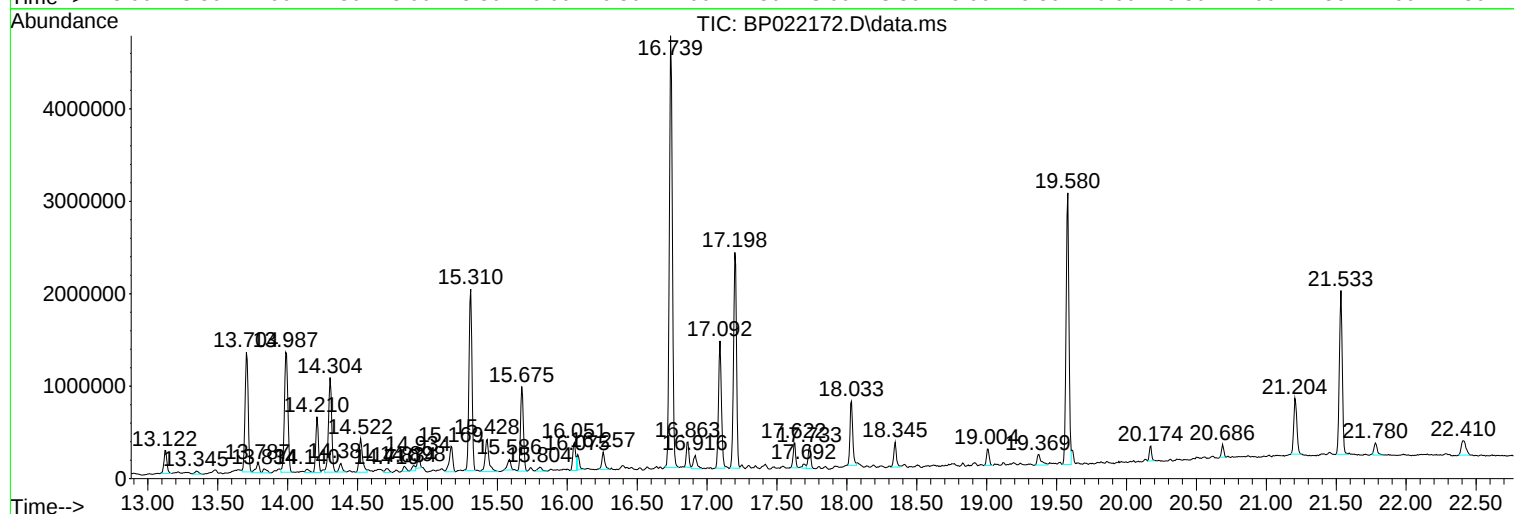
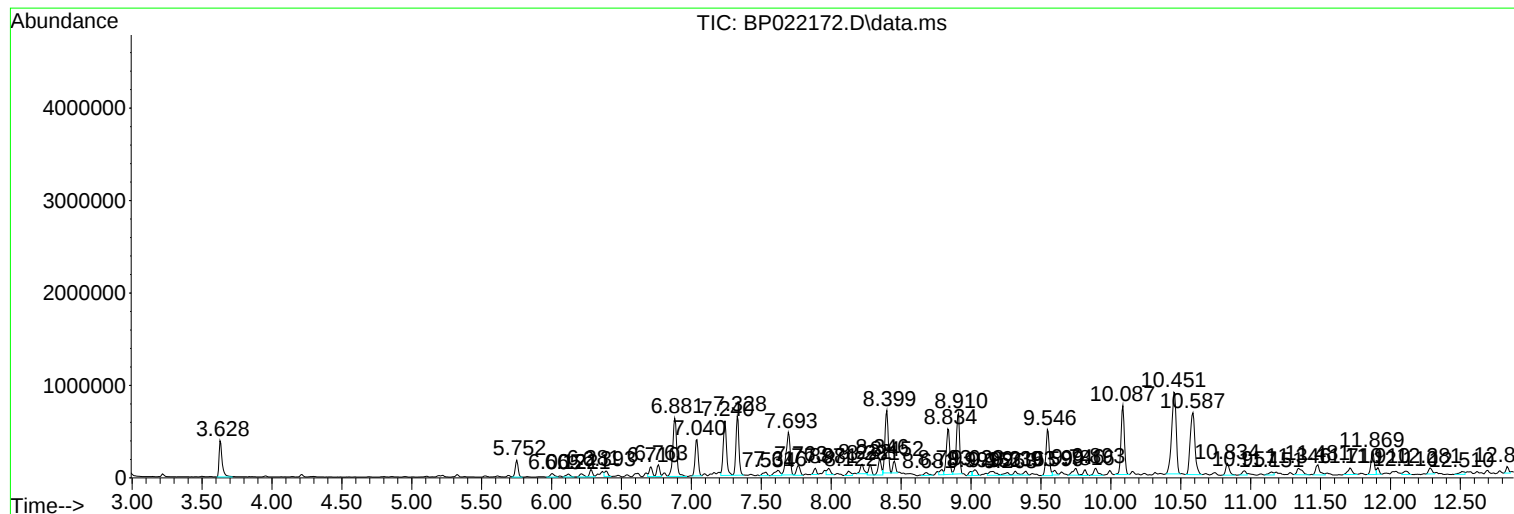
Sum of corrected areas: 66326005

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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

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



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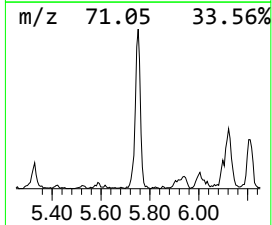
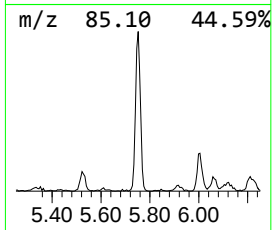
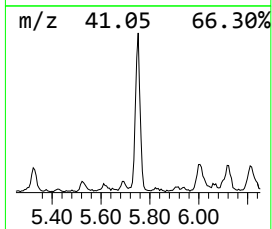
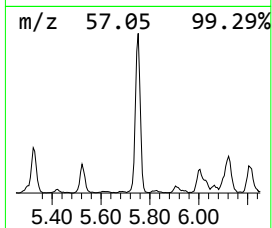
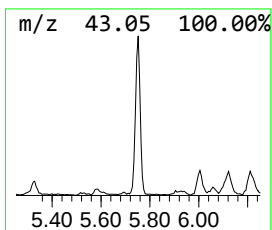
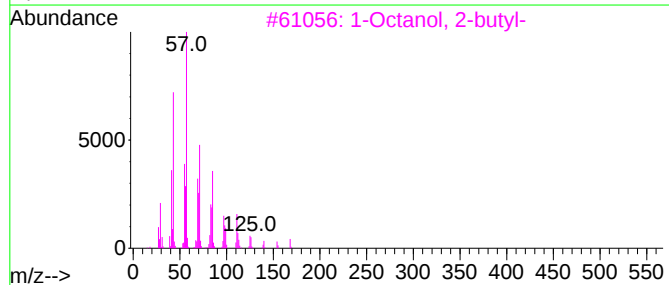
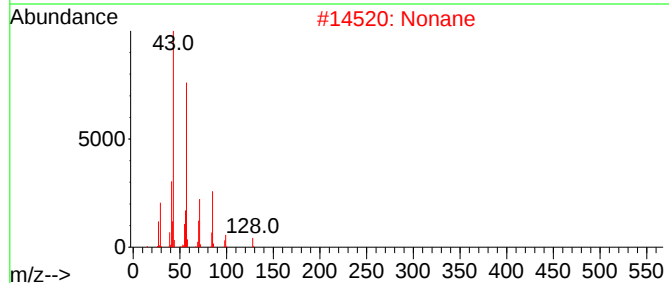
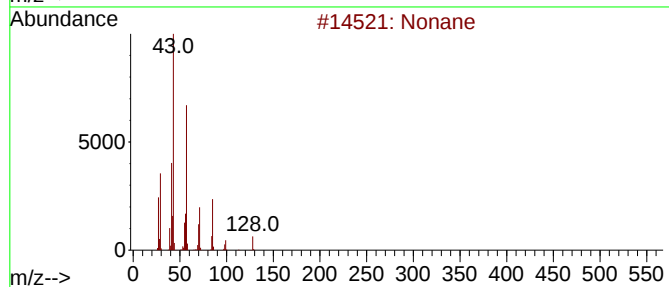
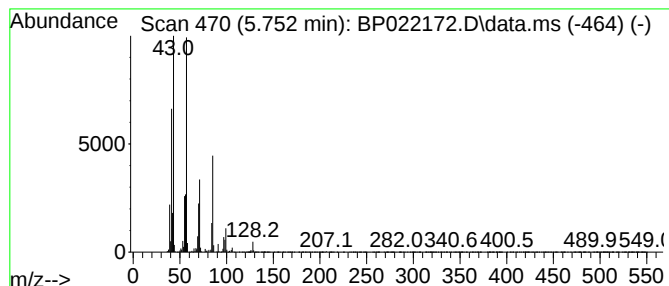
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 5.752 | 5.65 ng/ul | 258221 | 1,4-Dichlorobenzene-d4 | 7.699 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Nonane                  | 128 | C9H20   | 000111-84-2 | 94   |
| 2    |    |   | Nonane                  | 128 | C9H20   | 000111-84-2 | 91   |
| 3    |    |   | 1-Octanol, 2-butyl-     | 186 | C12H26O | 003913-02-8 | 64   |
| 4    |    |   | Decane                  | 142 | C10H22  | 000124-18-5 | 64   |
| 5    |    |   | 1-Iodo-2-methylundecane | 296 | C12H25I | 073105-67-6 | 64   |



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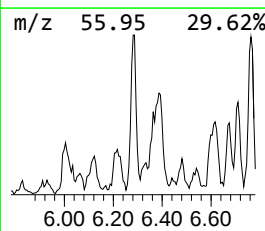
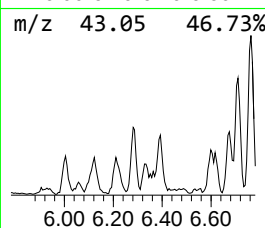
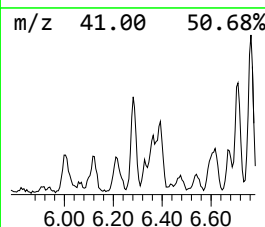
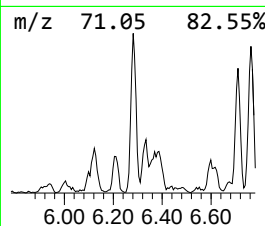
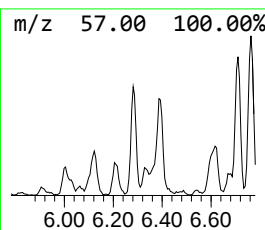
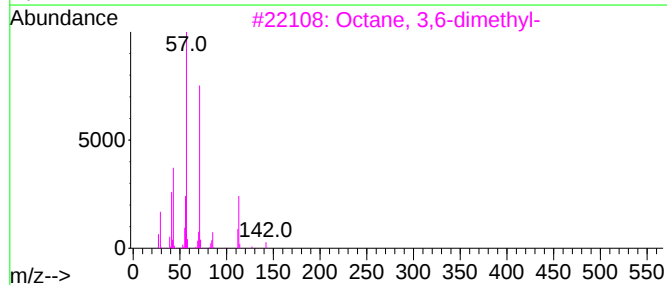
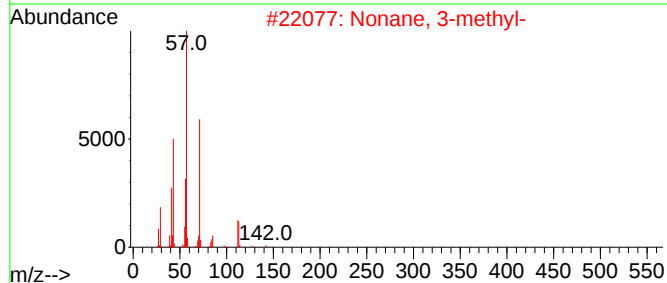
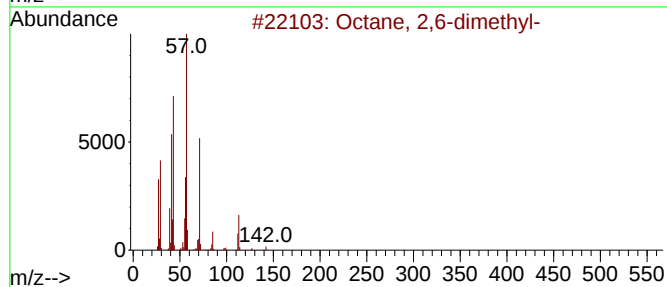
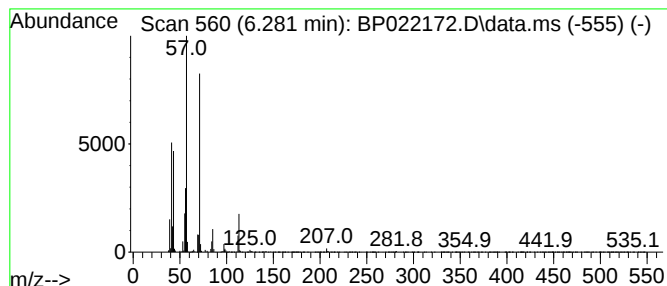
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 25

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 6.281 | 2.60 ng/ul | 119038 | 1,4-Dichlorobenzene-d4 | 7.699 |

| Hit# | of                                 | 5 | Tentative ID | MW       | MolForm      | CAS# | Qual |
|------|------------------------------------|---|--------------|----------|--------------|------|------|
| 1    | Octane, 2,6-dimethyl-              |   | 142          | C10H22   | 002051-30-1  | 91   |      |
| 2    | Nonane, 3-methyl-                  |   | 142          | C10H22   | 005911-04-6  | 86   |      |
| 3    | Octane, 3,6-dimethyl-              |   | 142          | C10H22   | 015869-94-0  | 72   |      |
| 4    | Dodecane, 2,6,11-trimethyl-        |   | 212          | C15H32   | 031295-56-4  | 72   |      |
| 5    | Oxalic acid, isobutyl pentyl ester |   | 216          | C11H20O4 | 1000309-37-0 | 64   |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

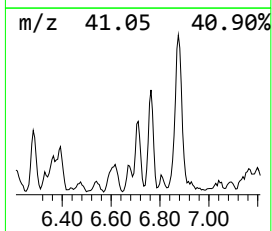
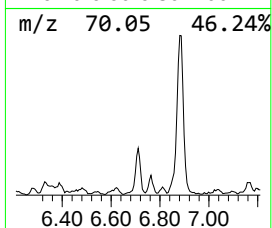
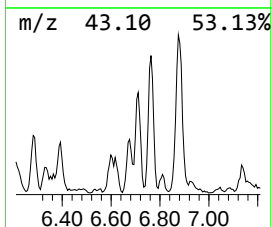
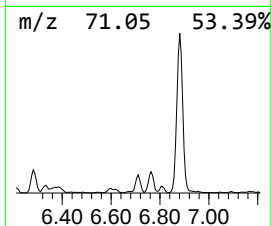
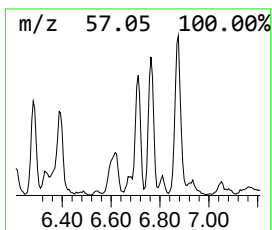
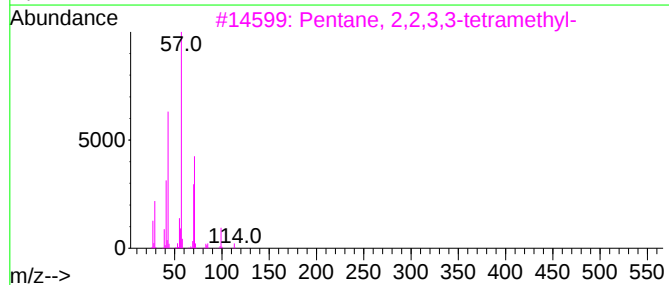
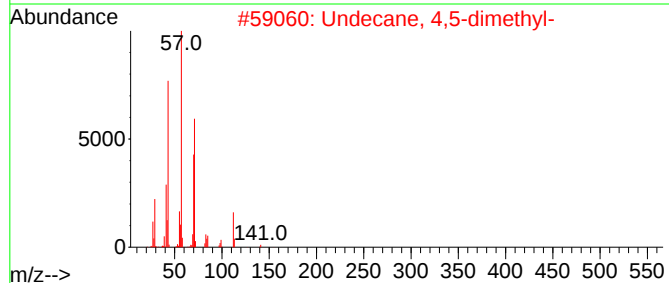
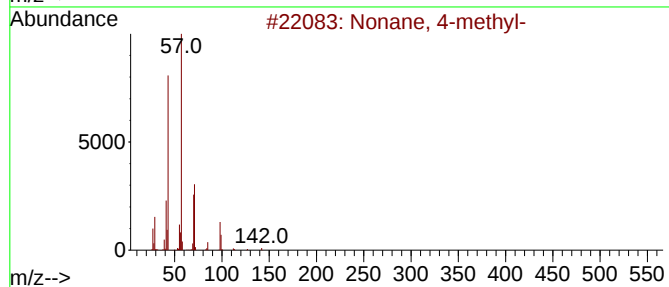
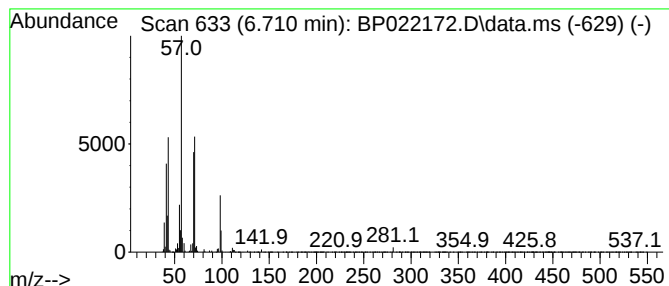
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Quant Title : SVOA CALIBRATION

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

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Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 17

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 6.710 | 3.56 ng/ul | 162568 | 1,4-Dichlorobenzene-d4 | 7.699 |

| Hit# | of 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------|-----|---------|-------------|------|
| 1    |      | Nonane, 4-methyl-             | 142 | C10H22  | 017301-94-9 | 64   |
| 2    |      | Undecane, 4,5-dimethyl-       | 184 | C13H28  | 017312-79-7 | 53   |
| 3    |      | Pentane, 2,2,3,3-tetramethyl- | 128 | C9H20   | 007154-79-2 | 50   |
| 4    |      | Nonane, 2-methyl-             | 142 | C10H22  | 000871-83-0 | 47   |
| 5    |      | Heptane, 4-propyl-            | 142 | C10H22  | 003178-29-8 | 45   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
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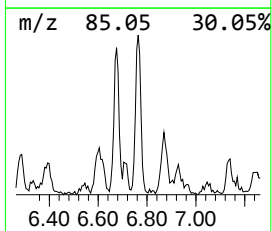
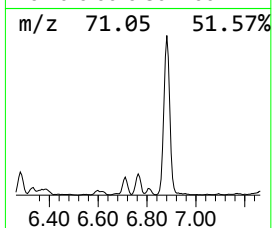
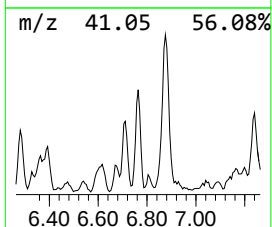
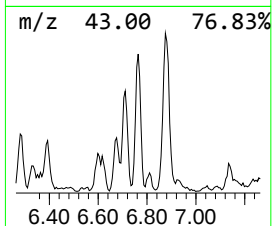
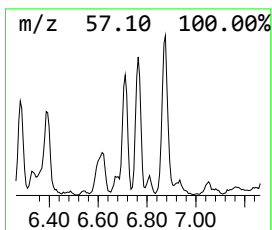
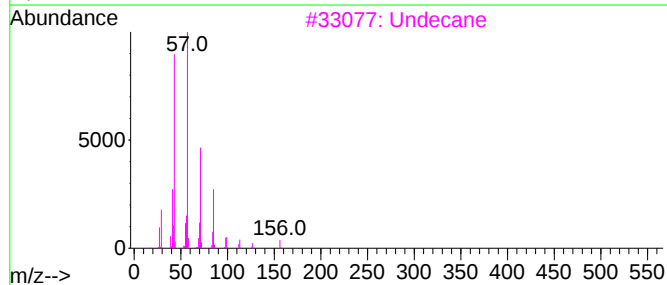
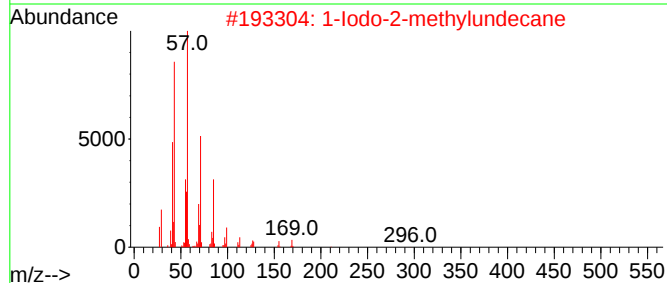
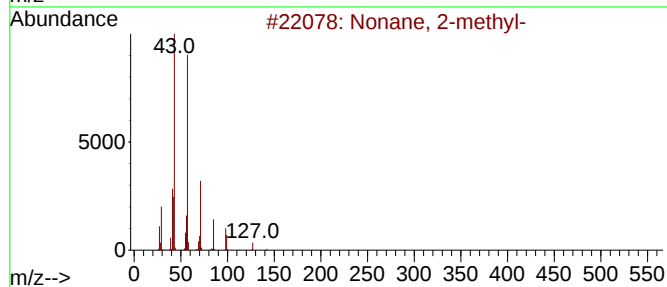
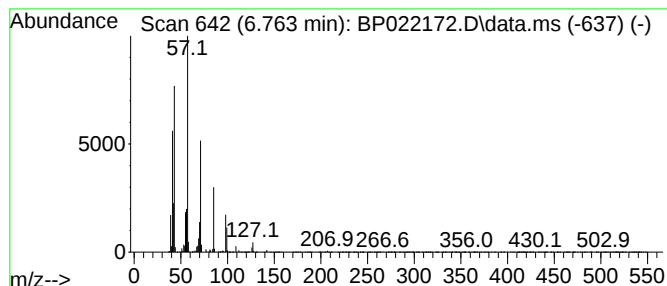
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 6.763 | 4.02 ng/ul | 183808 | 1,4-Dichlorobenzene-d4 | 7.699 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Nonane, 2-methyl-       | 142 | C10H22  | 000871-83-0 | 91   |
| 2    |    |   | 1-Iodo-2-methylundecane | 296 | C12H25I | 073105-67-6 | 81   |
| 3    |    |   | Undecane                | 156 | C11H24  | 001120-21-4 | 64   |
| 4    |    |   | Hexadecane              | 226 | C16H34  | 000544-76-3 | 59   |
| 5    |    |   | Undecane, 2,6-dimethyl- | 184 | C13H28  | 017301-23-4 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

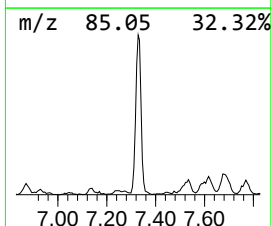
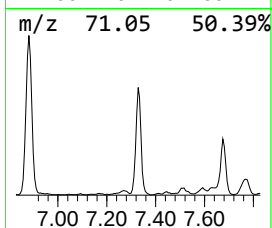
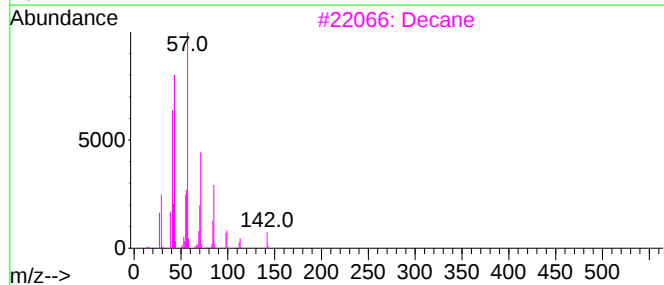
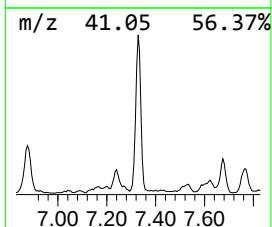
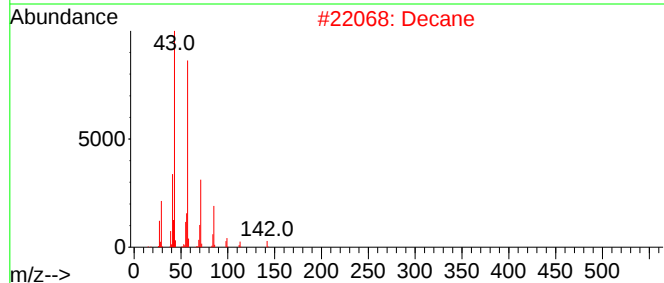
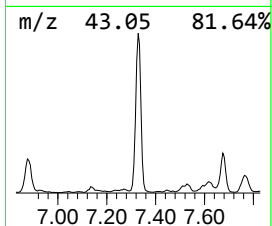
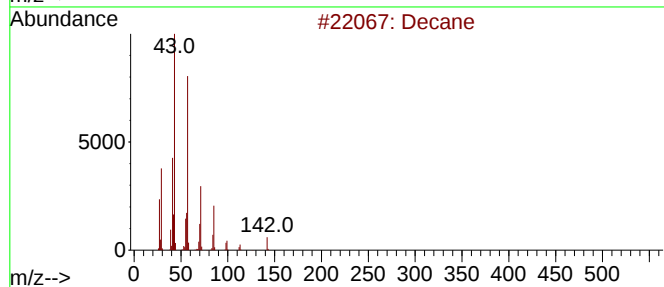
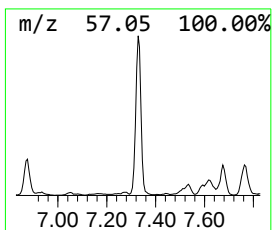
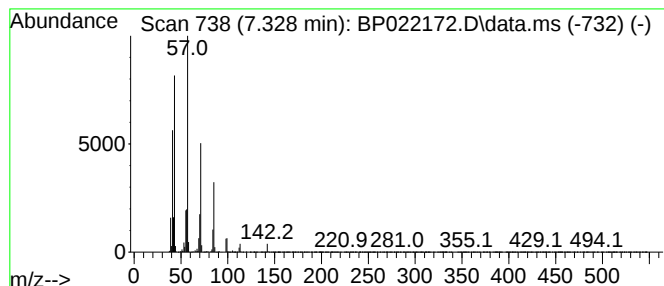
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 2

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 7.328 | 21.07 ng/ul | 963486 | 1,4-Dichlorobenzene-d4 | 7.699 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Decane                              | 142 | C10H22   | 000124-18-5  | 95   |
| 2    |    |   | Decane                              | 142 | C10H22   | 000124-18-5  | 94   |
| 3    |    |   | Decane                              | 142 | C10H22   | 000124-18-5  | 91   |
| 4    |    |   | Hexadecane                          | 226 | C16H34   | 000544-76-3  | 86   |
| 5    |    |   | Carbonic acid, prop-1-en-2-yl tr... | 284 | C17H32O3 | 1000382-90-8 | 78   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

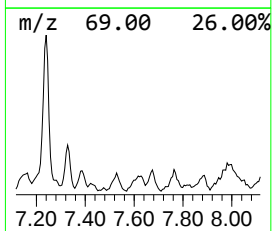
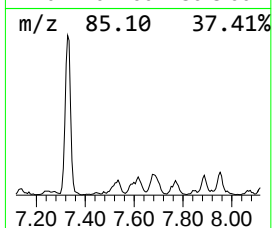
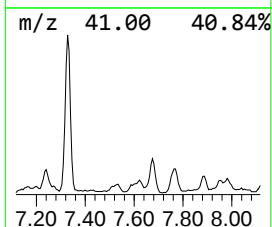
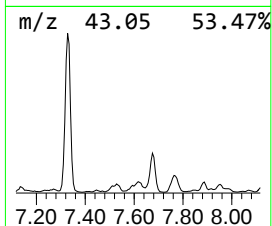
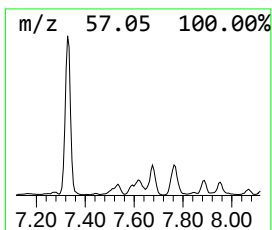
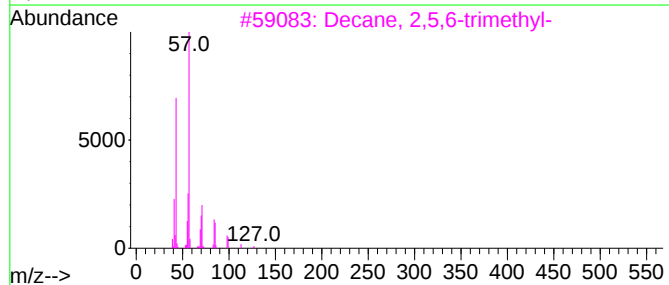
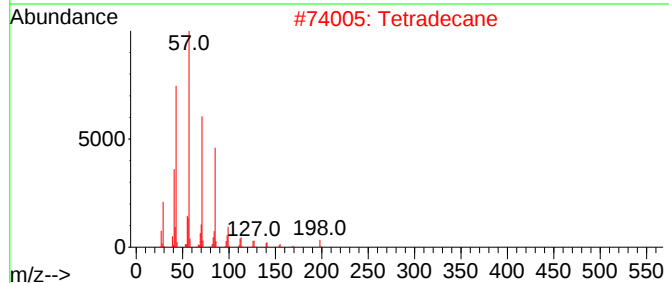
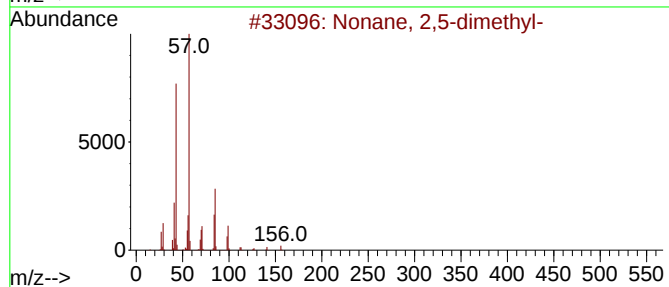
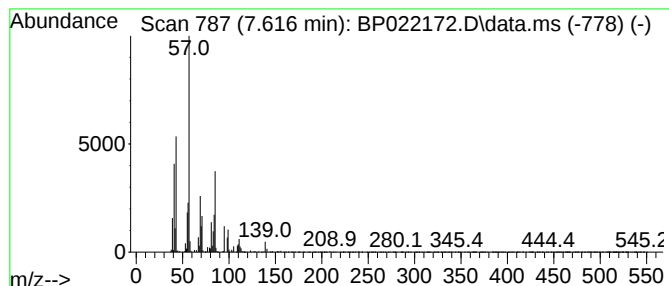
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Quant Title : SVOA CALIBRATION

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

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Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 18

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 7.616 | 3.54 ng/ul | 161676 | 1,4-Dichlorobenzene-d4 | 7.699 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-----------------------------------|-----|----------|--------------|------|
| 1    |    |   | Nonane, 2,5-dimethyl-             | 156 | C11H24   | 017302-27-1  | 50   |
| 2    |    |   | Tetradecane                       | 198 | C14H30   | 000629-59-4  | 47   |
| 3    |    |   | Decane, 2,5,6-trimethyl-          | 184 | C13H28   | 062108-23-0  | 47   |
| 4    |    |   | Oxalic acid, isobutyl hexyl ester | 230 | C12H22O4 | 1000309-37-1 | 47   |
| 5    |    |   | Pentadecane                       | 212 | C15H32   | 000629-62-9  | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

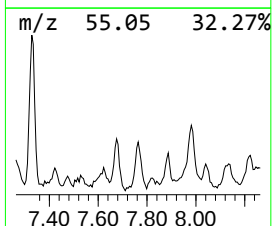
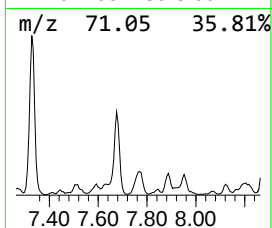
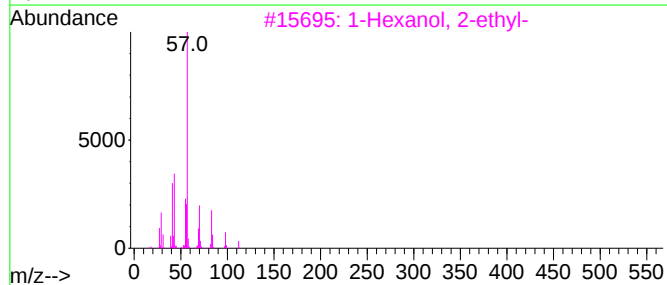
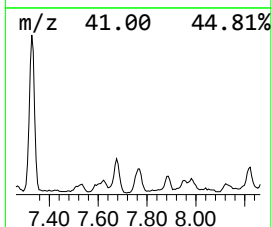
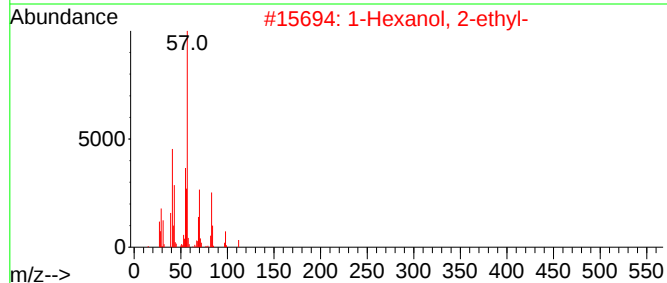
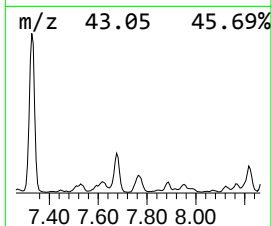
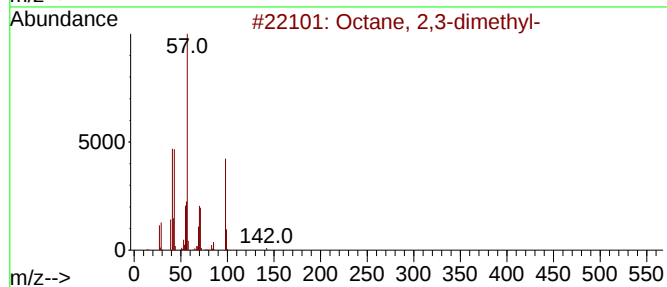
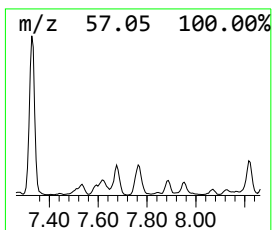
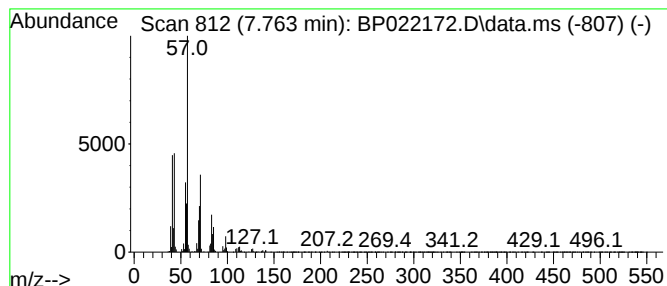
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 13

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 7.763 | 4.16 ng/ul | 190227 | 1,4-Dichlorobenzene-d4 | 7.699 |

| Hit# | of                               | 5 | Tentative ID | MW  | MolForm  | CAS#         | Qual |
|------|----------------------------------|---|--------------|-----|----------|--------------|------|
| 1    | Octane, 2,3-dimethyl-            |   |              | 142 | C10H22   | 007146-60-3  | 59   |
| 2    | 1-Hexanol, 2-ethyl-              |   |              | 130 | C8H18O   | 000104-76-7  | 59   |
| 3    | 1-Hexanol, 2-ethyl-              |   |              | 130 | C8H18O   | 000104-76-7  | 53   |
| 4    | Undecane, 2,6-dimethyl-          |   |              | 184 | C13H28   | 017301-23-4  | 53   |
| 5    | Carbonic acid, octyl vinyl ester |   |              | 200 | C11H20O3 | 1000383-25-5 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

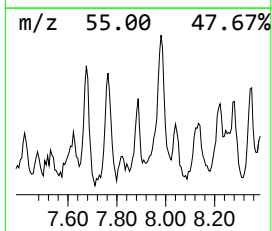
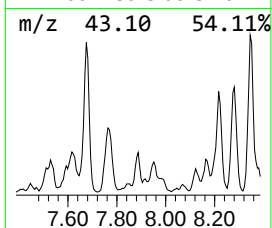
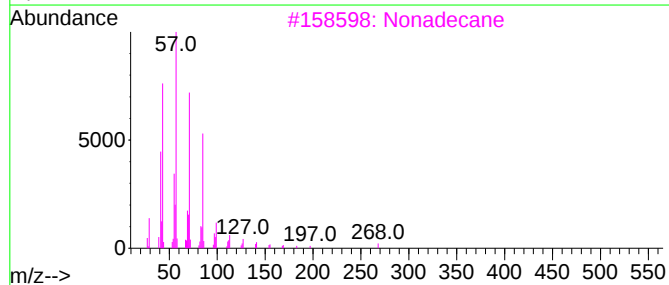
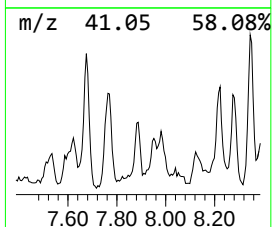
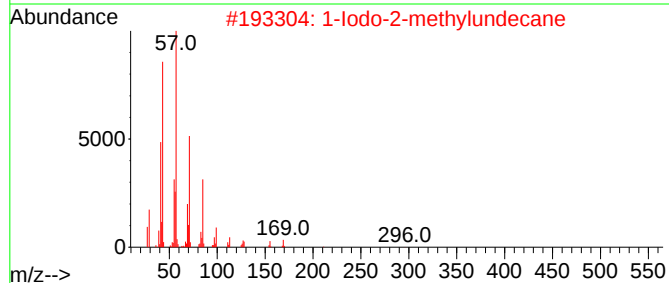
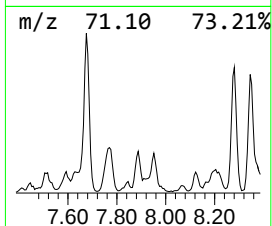
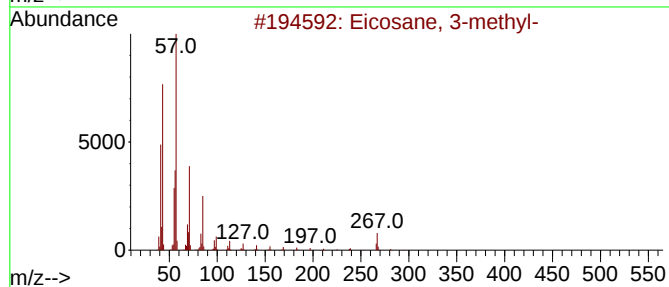
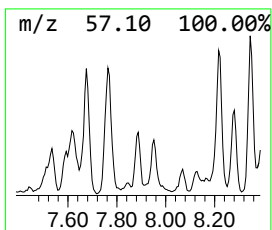
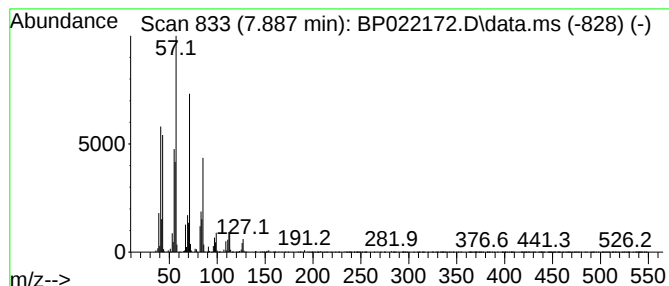
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 35

| R.T.  | EstConc    | Area  | Relative to ISTD       | R.T.  |
|-------|------------|-------|------------------------|-------|
| 7.887 | 2.04 ng/ul | 93340 | 1,4-Dichlorobenzene-d4 | 7.699 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | Eicosane, 3-methyl-                 | 296 | C21H44       | 006418-46-8 | 83      |      |      |
| 2    | 1-Iodo-2-methylundecane             | 296 | C12H25I      | 073105-67-6 | 74      |      |      |
| 3    | Nonadecane                          | 268 | C19H40       | 000629-92-5 | 64      |      |      |
| 4    | Nonane, 3-methyl-5-propyl-          | 184 | C13H28       | 031081-18-2 | 64      |      |      |
| 5    | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44       | 054833-48-6 | 64      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

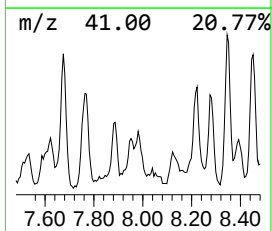
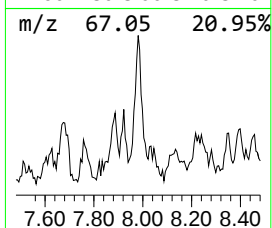
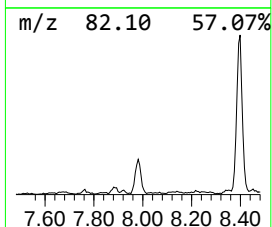
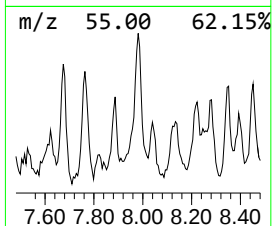
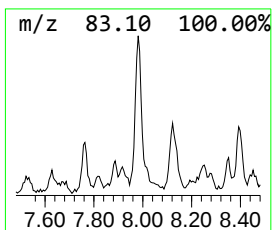
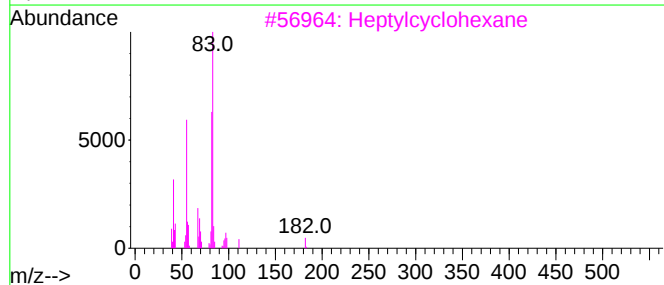
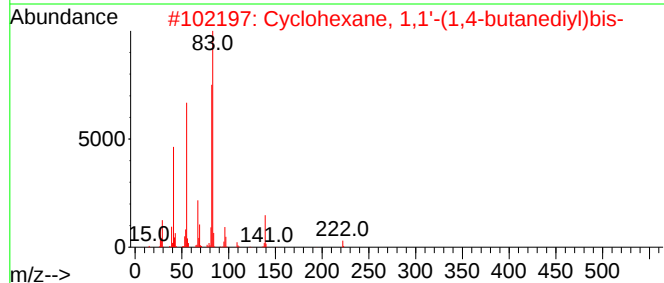
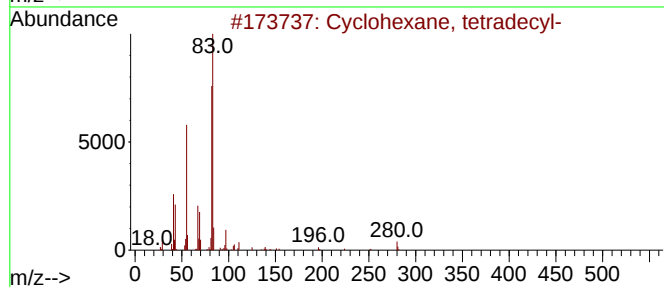
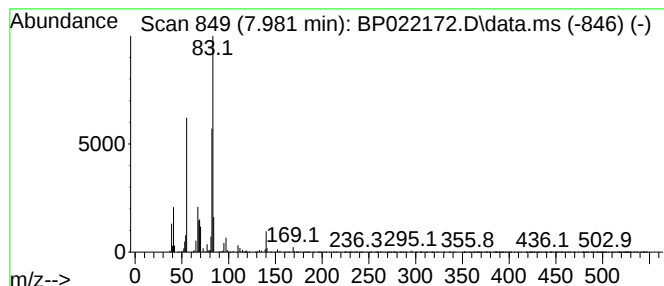
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Quant Title : SVOA CALIBRATION

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

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Peak Number 9 (DEL) Alkane: Cyclic7.981 Concentration Rank 28

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 7.981 | 2.45 ng/ul | 111964 | 1,4-Dichlorobenzene-d4 | 7.699 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, tetradecyl-            | 280 | C20H40  | 001795-18-2 | 72   |
| 2    |    |   | Cyclohexane, 1,1'-(1,4-butanediyl)- | 222 | C16H30  | 006165-44-2 | 72   |
| 3    |    |   | Heptylcyclohexane                   | 182 | C13H26  | 005617-41-4 | 72   |
| 4    |    |   | Cyclohexane, butyl-                 | 140 | C10H20  | 001678-93-9 | 72   |
| 5    |    |   | Cyclohexane, 1,1'-(1,4-butanediyl)- | 222 | C16H30  | 006165-44-2 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

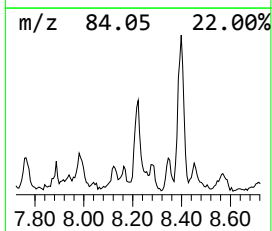
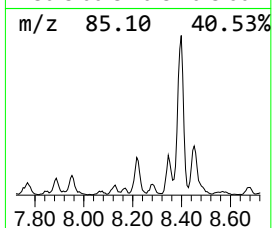
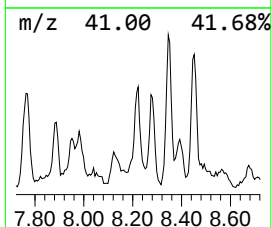
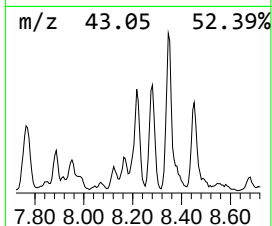
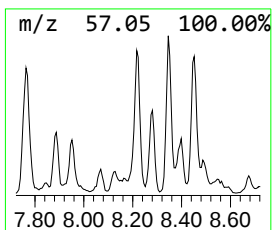
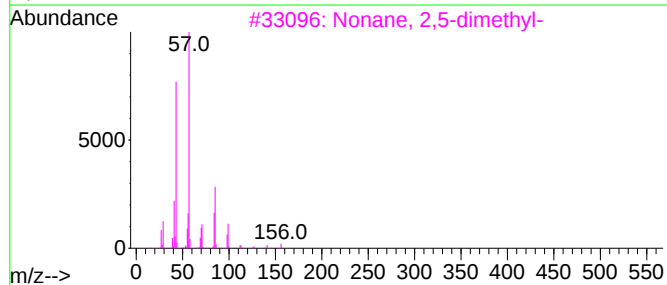
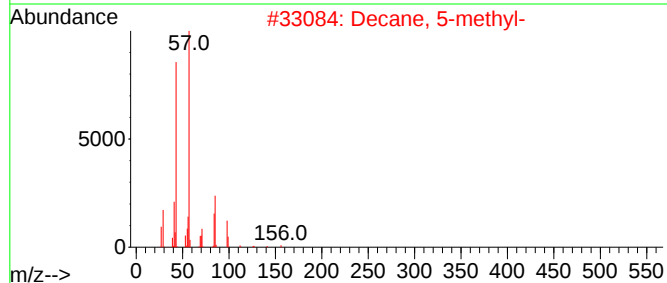
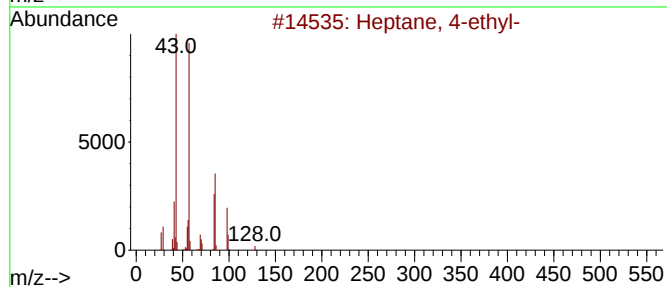
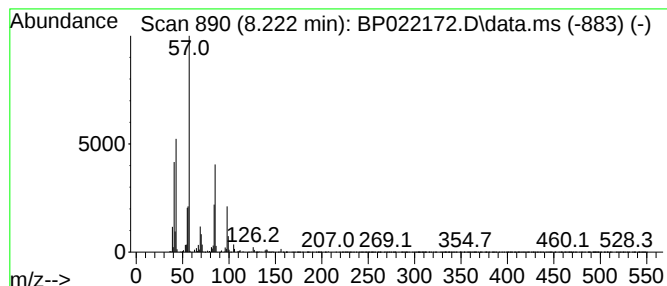
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TIC Integration Parameters: LSCINT.P

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Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.222 | 4.09 ng/ul | 186983 | 1,4-Dichlorobenzene-d4 | 7.699 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Heptane, 4-ethyl-                   | 128 | C9H20   | 002216-32-2 | 78   |
| 2    |      | Decane, 5-methyl-                   | 156 | C11H24  | 013151-35-4 | 64   |
| 3    |      | Nonane, 2,5-dimethyl-               | 156 | C11H24  | 017302-27-1 | 59   |
| 4    |      | Hexane, 2,2,3,3-tetramethyl-        | 142 | C10H22  | 013475-81-5 | 59   |
| 5    |      | Butanoic acid, 3-oxo-, 1,1-dimet... | 158 | C8H14O3 | 001694-31-1 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

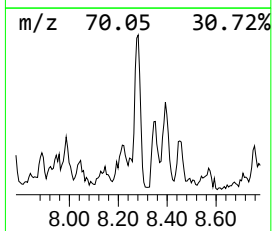
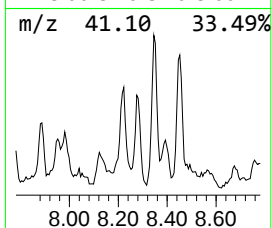
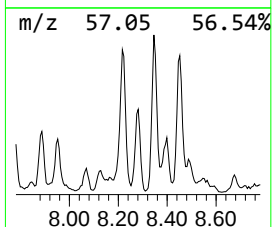
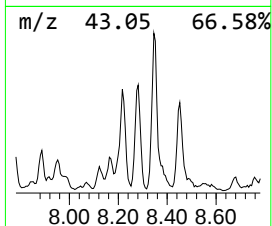
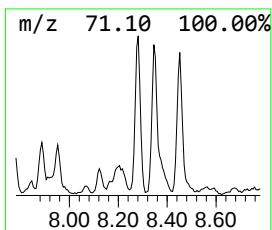
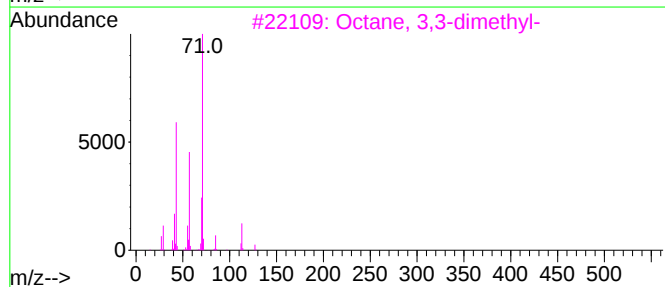
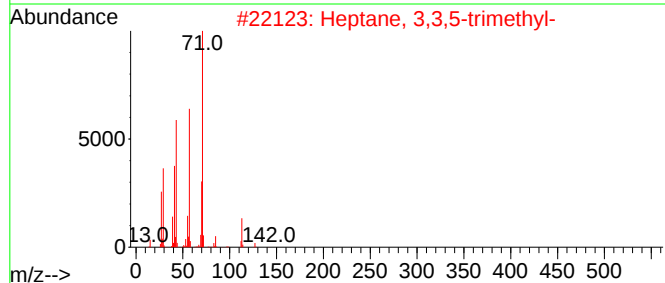
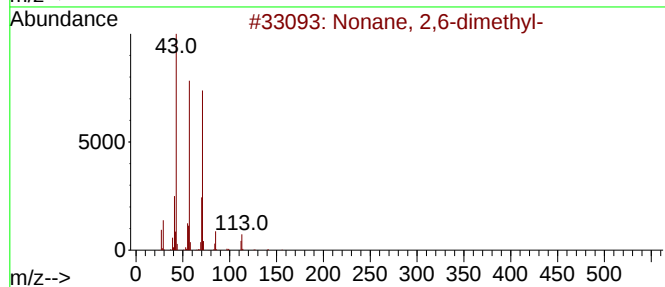
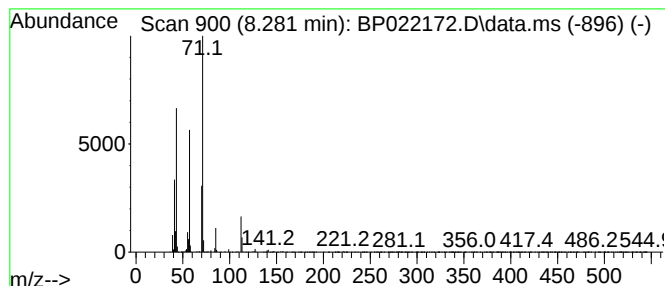
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 19

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.281 | 3.48 ng/ul | 158904 | 1,4-Dichlorobenzene-d4 | 7.699 |

| Hit# | of 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|------|------------------------------|-----|---------|-------------|------|
| 1    |      | Nonane, 2,6-dimethyl-        | 156 | C11H24  | 017302-28-2 | 72   |
| 2    |      | Heptane, 3,3,5-trimethyl-    | 142 | C10H22  | 007154-80-5 | 72   |
| 3    |      | Octane, 3,3-dimethyl-        | 142 | C10H22  | 004110-44-5 | 72   |
| 4    |      | Hexane, 3,3,4,4-tetramethyl- | 142 | C10H22  | 005171-84-6 | 64   |
| 5    |      | Heptane, 2,5,5-trimethyl-    | 142 | C10H22  | 001189-99-7 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

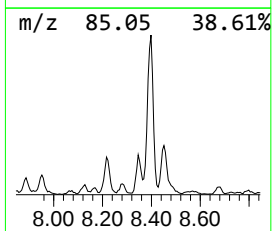
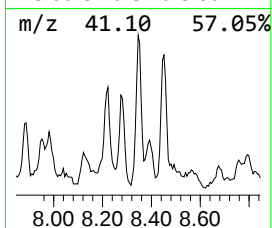
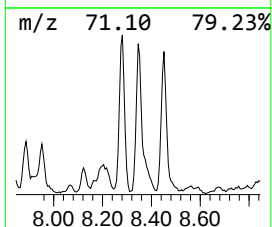
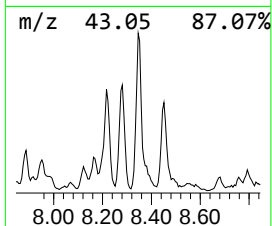
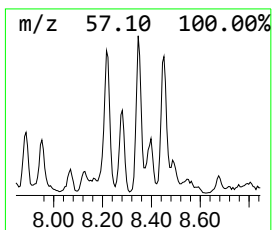
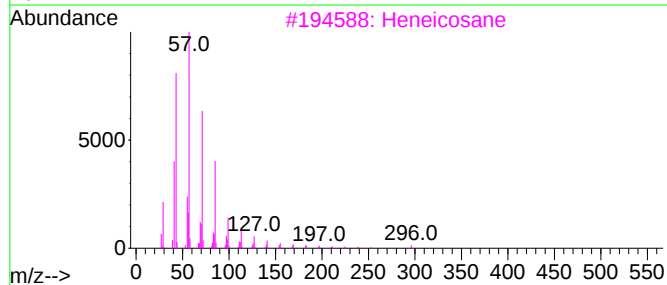
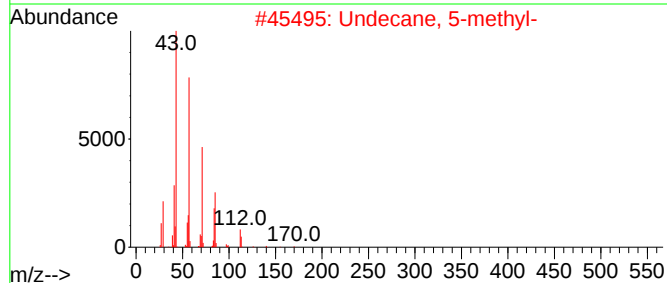
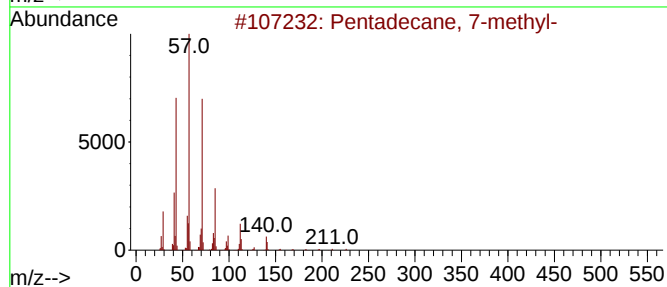
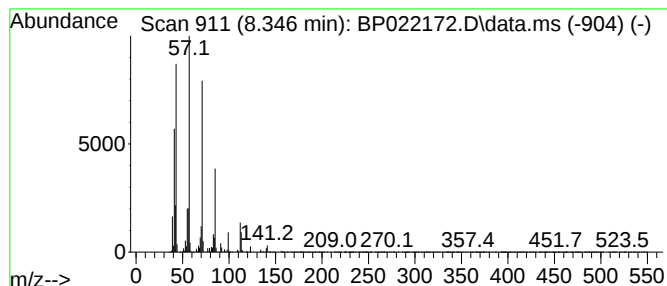
TIC Library : C:\DATABASE\NIST0.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.346 | 6.25 ng/ul | 285793 | 1,4-Dichlorobenzene-d4 | 7.699 |

| Hit# | of 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------|-----|---------|-------------|------|
| 1    |      | Pentadecane, 7-methyl-  | 226 | C16H34  | 006165-40-8 | 83   |
| 2    |      | Undecane, 5-methyl-     | 170 | C12H26  | 001632-70-8 | 80   |
| 3    |      | Heneicosane             | 296 | C21H44  | 000629-94-7 | 74   |
| 4    |      | Tetratetracontane       | 619 | C44H90  | 007098-22-8 | 72   |
| 5    |      | 1-Iodo-2-methylundecane | 296 | C12H25I | 073105-67-6 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

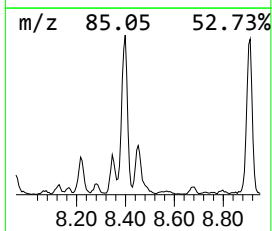
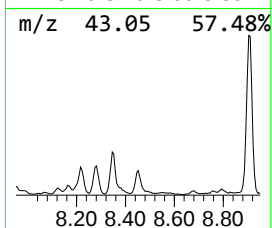
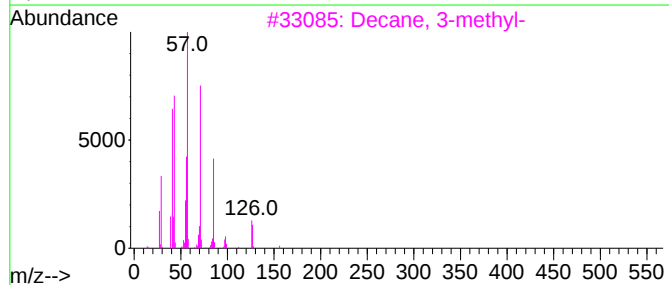
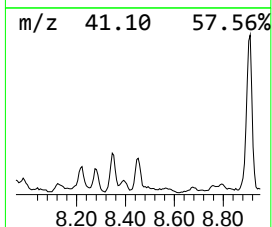
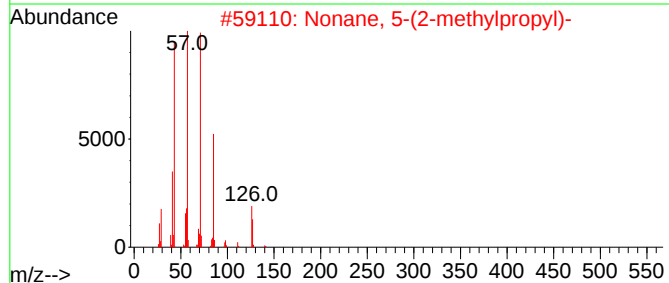
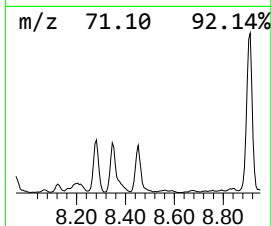
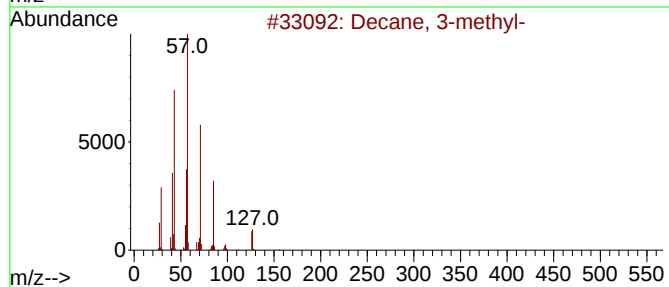
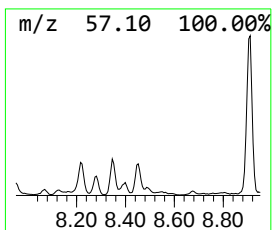
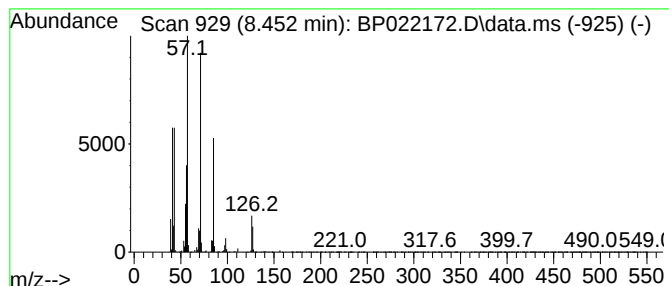
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 16

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.452 | 3.80 ng/ul | 173718 | 1,4-Dichlorobenzene-d4 | 7.699 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 3-methyl-           | 156 | C11H24  | 013151-34-3 | 91   |
| 2    |    |   | Nonane, 5-(2-methylpropyl)- | 184 | C13H28  | 062185-53-9 | 64   |
| 3    |    |   | Decane, 3-methyl-           | 156 | C11H24  | 013151-34-3 | 64   |
| 4    |    |   | Heptadecane, 8-methyl-      | 254 | C18H38  | 013287-23-5 | 59   |
| 5    |    |   | Octane, 2,6-dimethyl-       | 142 | C10H22  | 002051-30-1 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

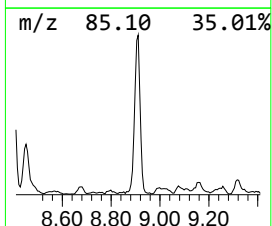
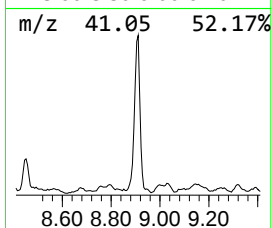
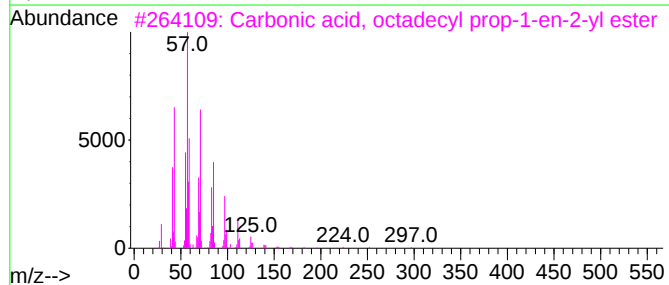
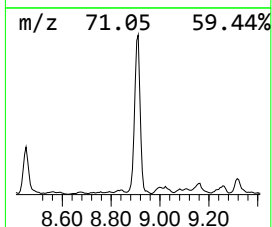
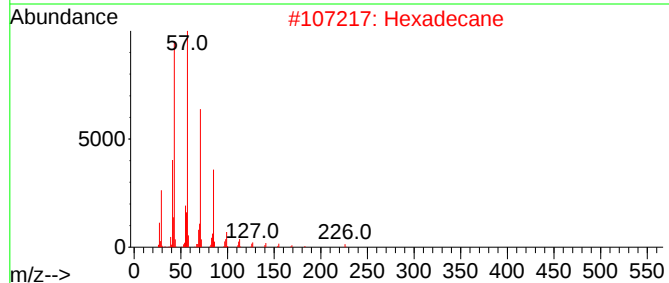
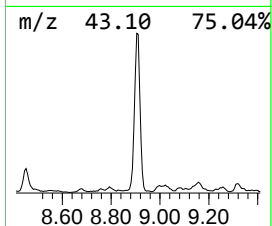
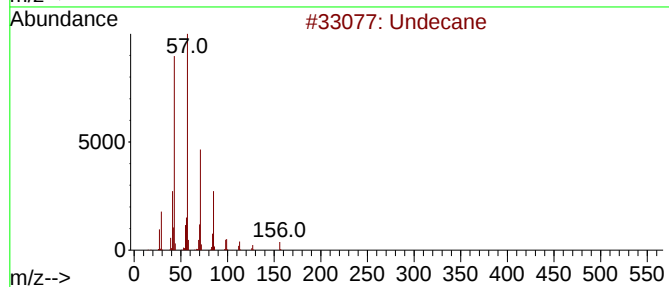
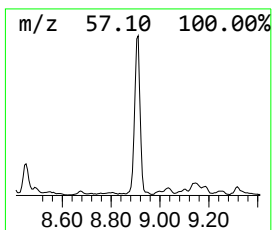
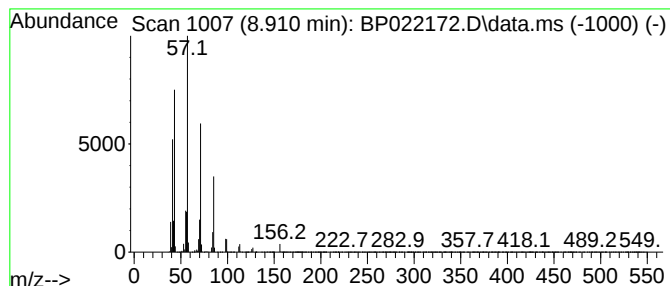
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Quant Title : SVOA CALIBRATION

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

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Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 1

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 8.910 | 22.79 ng/ul | 1042160 | 1,4-Dichlorobenzene-d4 | 7.699 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Undecane                            | 156 | C11H24   | 001120-21-4  | 90   |
| 2    |    |   | Hexadecane                          | 226 | C16H34   | 000544-76-3  | 90   |
| 3    |    |   | Carbonic acid, octadecyl prop-1-... | 354 | C22H42O3 | 1000383-11-5 | 90   |
| 4    |    |   | Undecane                            | 156 | C11H24   | 001120-21-4  | 81   |
| 5    |    |   | Tetradecane                         | 198 | C14H30   | 000629-59-4  | 78   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

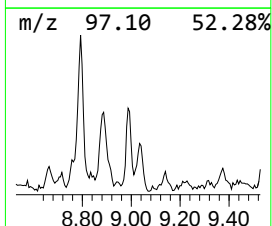
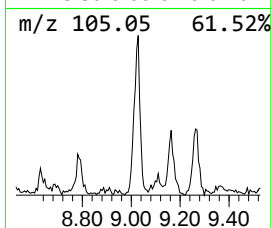
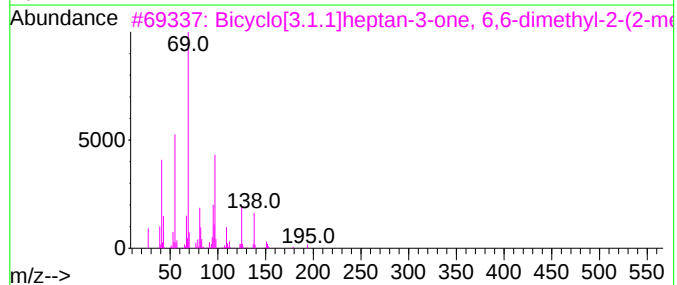
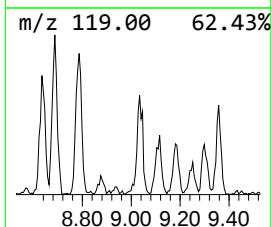
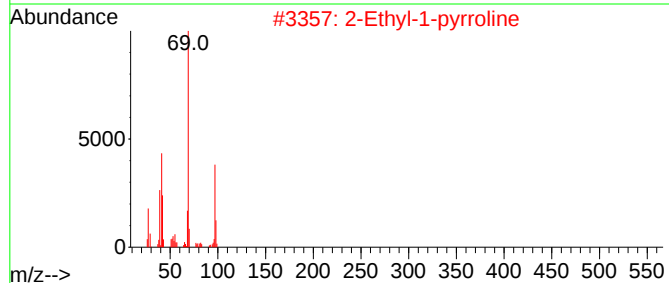
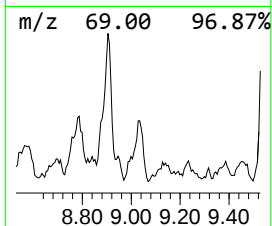
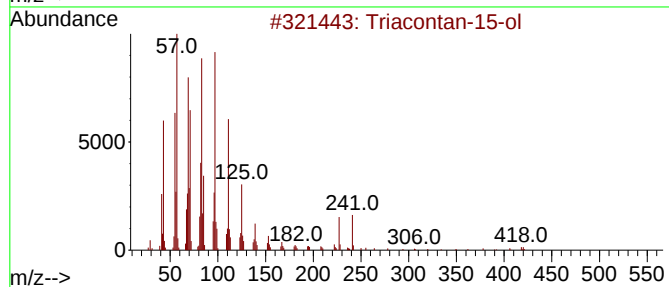
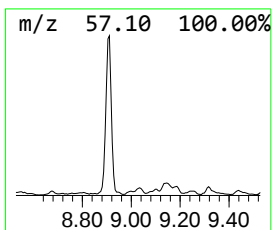
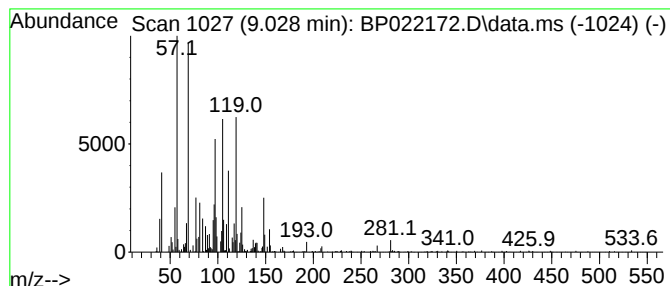
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Quant Title : SVOA CALIBRATION

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

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Peak Number 15 Triacontan-15-ol Concentration Rank 29

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 9.028 | 2.42 ng/ul | 110809 | 1,4-Dichlorobenzene-d4 | 7.699 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Triacontan-15-ol                    | 438 | C30H62O    | 031849-26-0  | 58   |
| 2    |    |   | 2-Ethyl-1-pyrroline                 | 97  | C6H11N     | 1000422-81-4 | 38   |
| 3    |    |   | Bicyclo[3.1.1]heptan-3-one, 6,6-... | 194 | C13H22O    | 1000163-97-9 | 38   |
| 4    |    |   | n-Pentadecanol                      | 228 | C15H32O    | 000629-76-5  | 35   |
| 5    |    |   | Benzeneacetic acid, 4-chloro-, 3... | 238 | C13H15ClO2 | 1000406-99-7 | 30   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

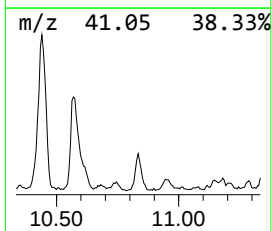
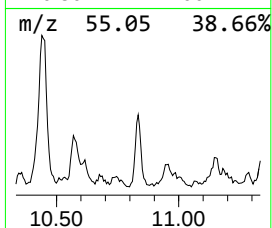
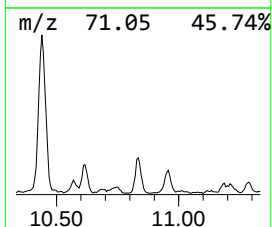
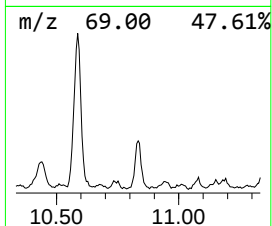
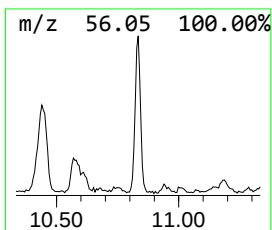
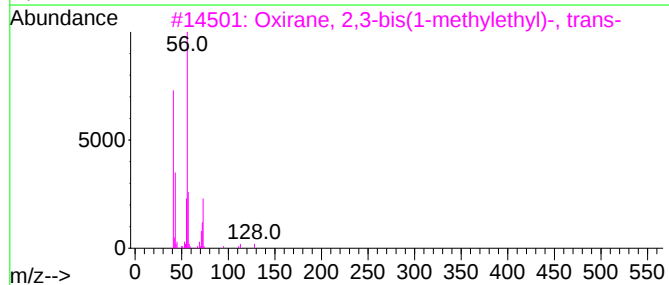
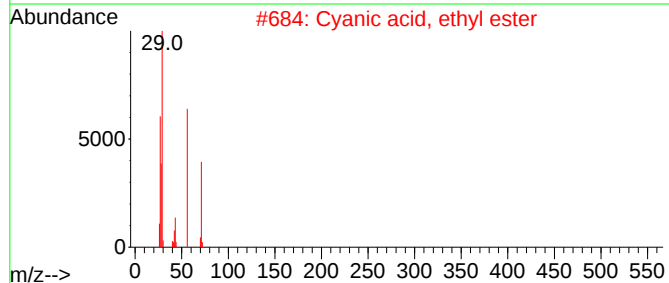
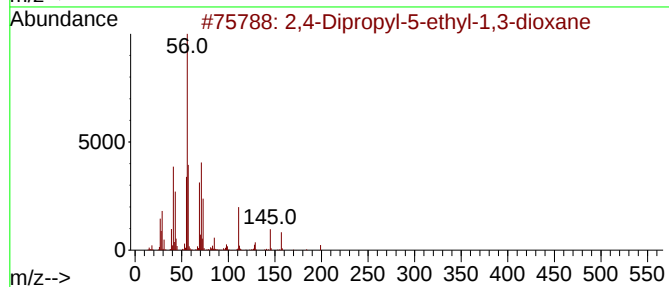
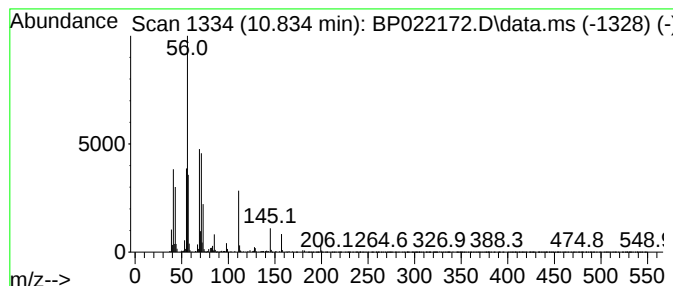
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TIC Integration Parameters: LSCINT.P

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Peak Number 16 2,4-Dipropyl-5-ethyl-1,3-di... Concentration Rank 32

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 10.834 | 2.32 ng/ul | 230419 | Naphthalene-d8   | 10.457 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|------|-------------------------------------|-----|----------|-------------|------|
| 1    |      | 2,4-Dipropyl-5-ethyl-1,3-dioxane    | 200 | C12H24O2 | 006413-83-8 | 86   |
| 2    |      | Cyanoic acid, ethyl ester           | 71  | C3H5NO   | 000627-48-5 | 52   |
| 3    |      | Oxirane, 2,3-bis(1-methylethyl)-... | 128 | C8H16O   | 054644-32-5 | 38   |
| 4    |      | 1-Pentene, 2,4-dimethyl-            | 98  | C7H14    | 002213-32-3 | 35   |
| 5    |      | 1-Hexene, 2-methyl-                 | 98  | C7H14    | 006094-02-6 | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

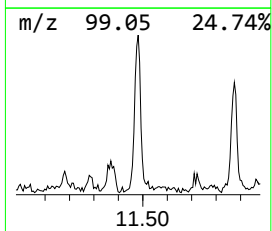
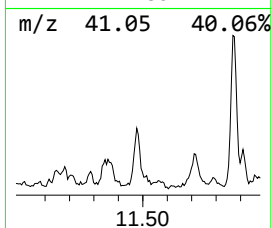
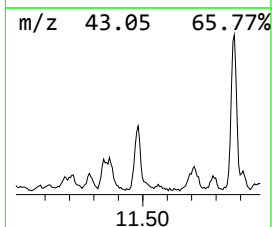
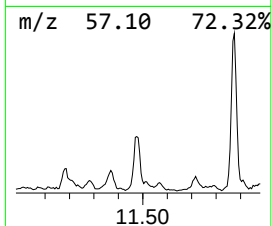
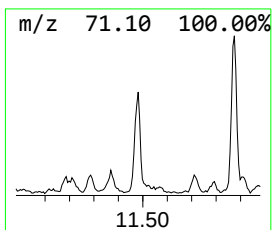
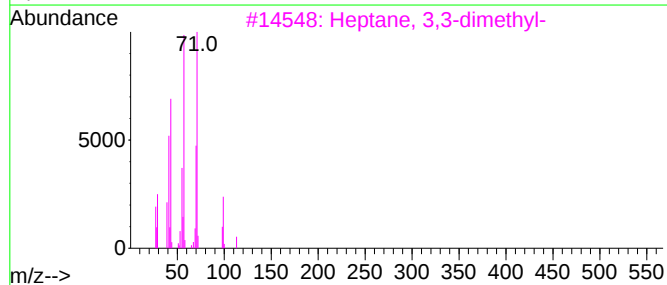
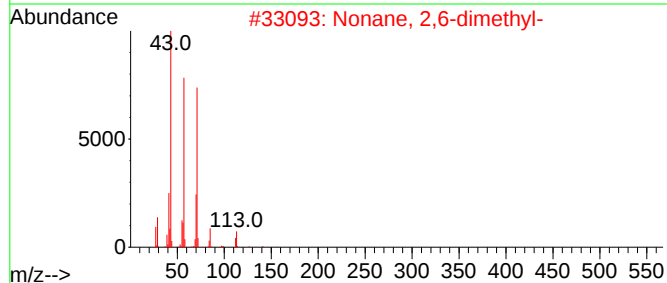
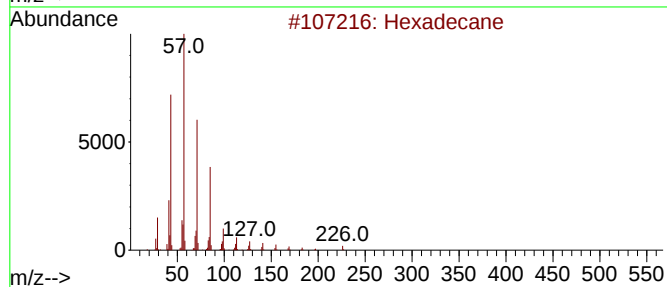
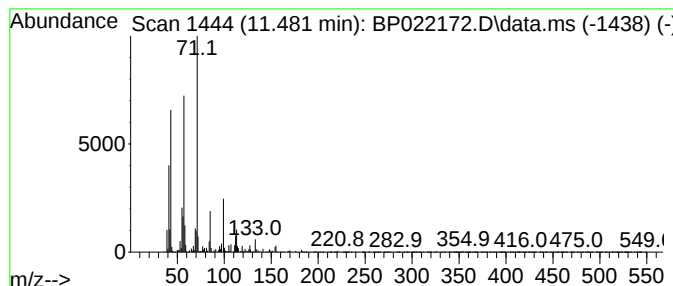
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 33

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 11.481 | 2.26 ng/ul | 224301 | Naphthalene-d8   | 10.457 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm   | CAS#         | Qual |
|------|----|---|------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Hexadecane                         | 226 | C16H34    | 000544-76-3  | 59   |
| 2    |    |   | Nonane, 2,6-dimethyl-              | 156 | C11H24    | 017302-28-2  | 53   |
| 3    |    |   | Heptane, 3,3-dimethyl-             | 128 | C9H20     | 004032-86-4  | 53   |
| 4    |    |   | Octane, 3,6-dimethyl-              | 142 | C10H22    | 015869-94-0  | 53   |
| 5    |    |   | Sulfurous acid, nonyl pentyl ester | 278 | C14H30O3S | 1000309-14-2 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

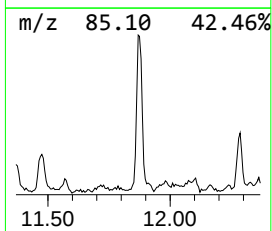
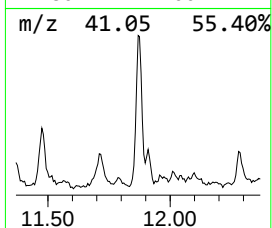
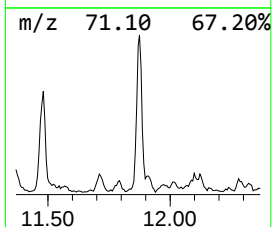
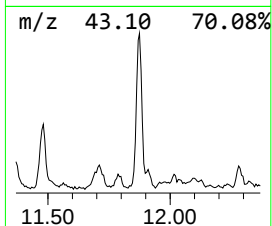
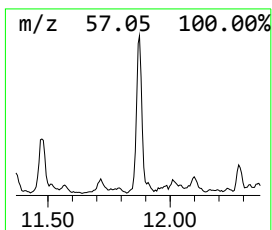
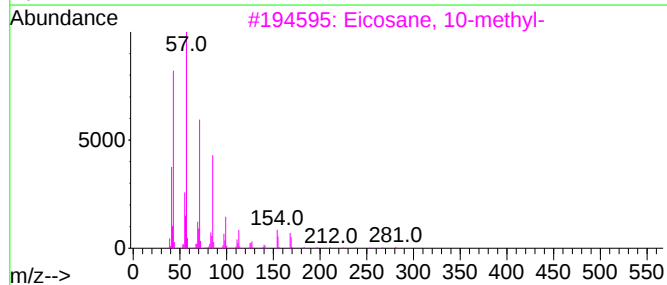
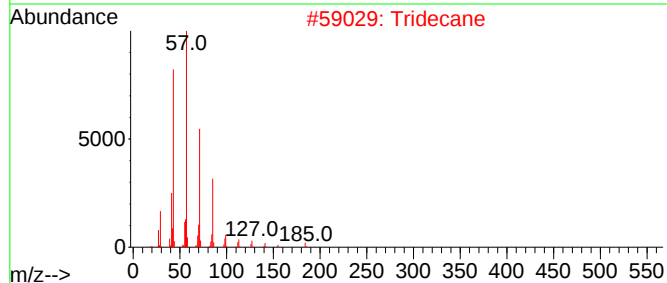
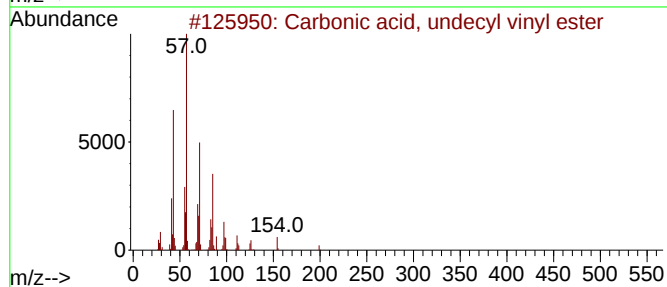
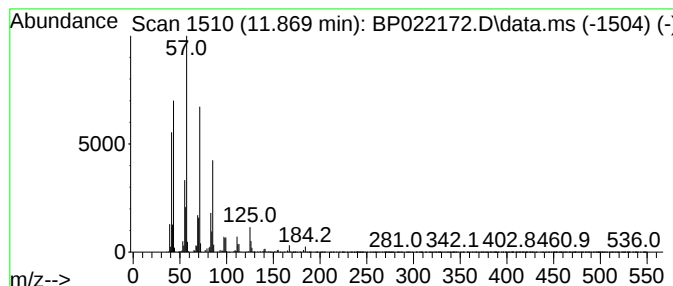
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Carbonic acid, undecyl vinyl... Concentration Rank 12

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 11.869 | 4.28 ng/ul | 424657 | Naphthalene-d8   | 10.457 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm  | CAS#         | Qual |
|------|----|---|------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Carbonic acid, undecyl vinyl ester | 242 | C14H26O3 | 1000382-54-9 | 86   |
| 2    |    |   | Tridecane                          | 184 | C13H28   | 000629-50-5  | 83   |
| 3    |    |   | Eicosane, 10-methyl-               | 296 | C21H44   | 054833-23-7  | 81   |
| 4    |    |   | Heneicosane                        | 296 | C21H44   | 000629-94-7  | 81   |
| 5    |    |   | Tetradecane                        | 198 | C14H30   | 000629-59-4  | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

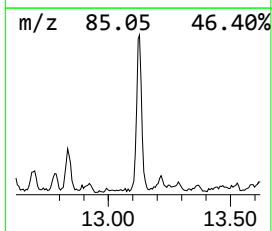
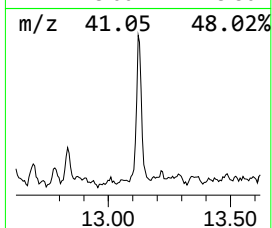
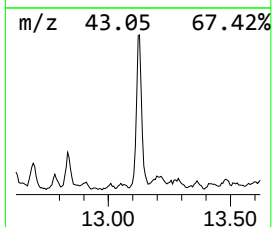
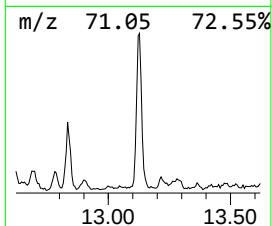
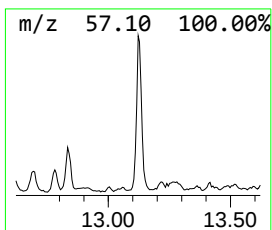
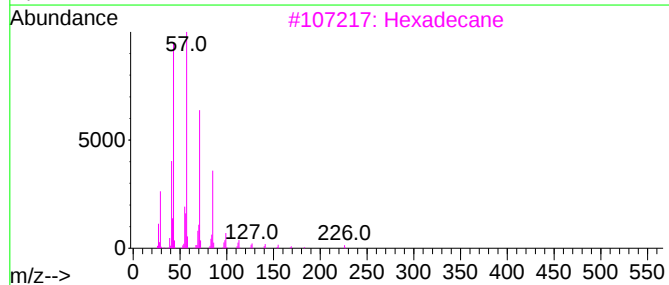
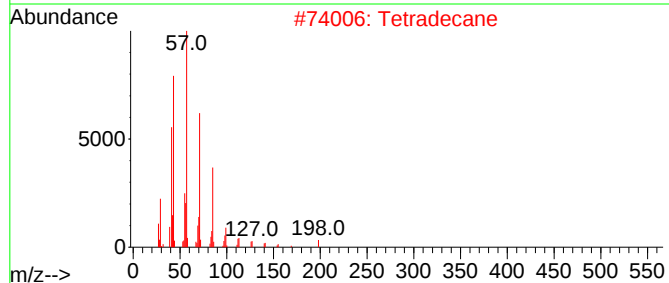
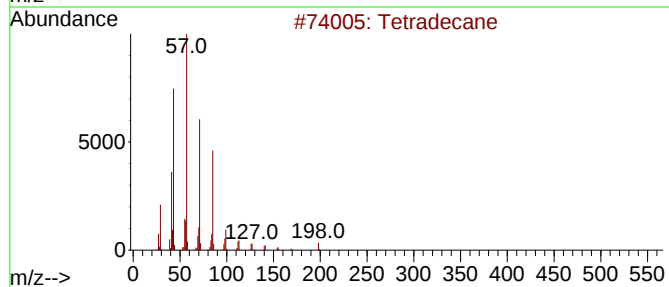
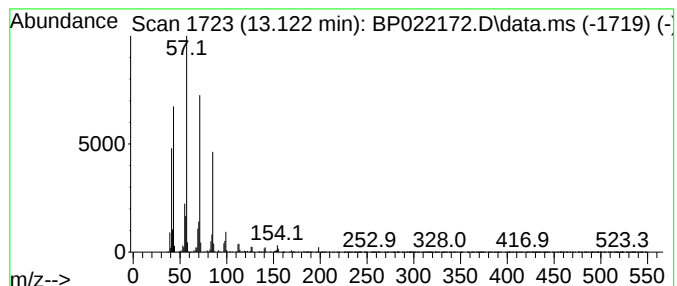
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.122 | 4.67 ng/ul | 372123 | Acenaphthene-d10 | 14.304 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Tetradecane                         | 198 | C14H30   | 000629-59-4  | 97   |
| 2    |    |   | Tetradecane                         | 198 | C14H30   | 000629-59-4  | 94   |
| 3    |    |   | Hexadecane                          | 226 | C16H34   | 000544-76-3  | 91   |
| 4    |    |   | Octadecane, 1-chloro-               | 288 | C18H37Cl | 003386-33-2  | 86   |
| 5    |    |   | Carbonic acid, octadecyl prop-1-... | 354 | C22H42O3 | 1000383-11-5 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

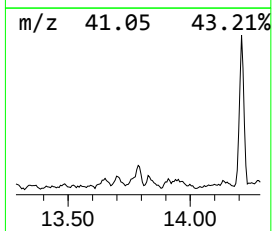
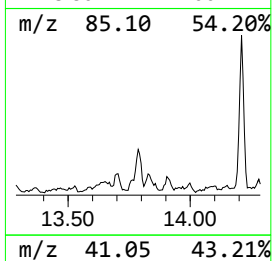
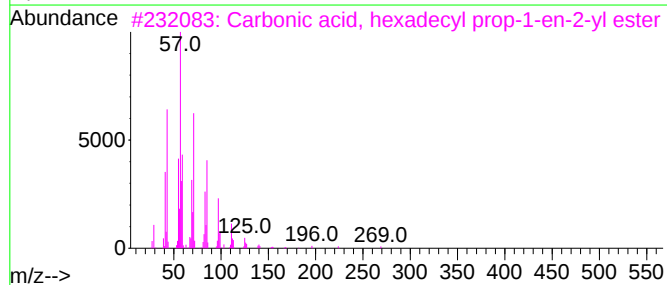
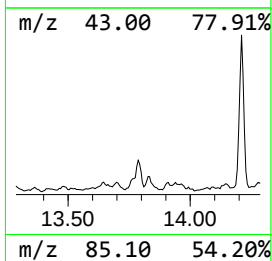
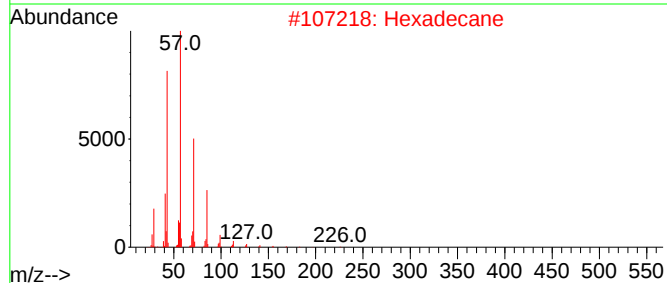
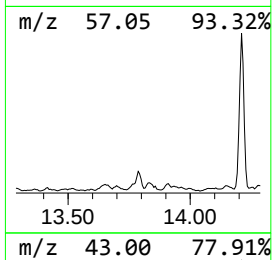
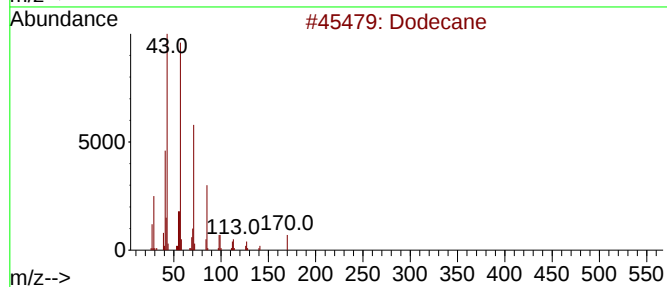
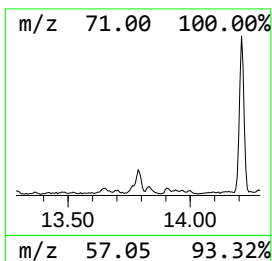
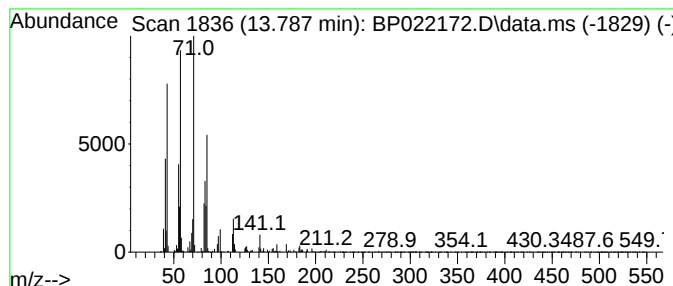
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Quant Title : SVOA CALIBRATION

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

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Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 34

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 13.787    | 2.23 ng/ul                          | 177829 | Acenaphthene-d10 | 14.304       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | Dodecane                            | 170    | C12H26           | 000112-40-3  | 59   |
| 2         | Hexadecane                          | 226    | C16H34           | 000544-76-3  | 52   |
| 3         | Carbonic acid, hexadecyl prop-1-... | 326    | C20H38O3         | 1000382-90-3 | 52   |
| 4         | Undecane, 4,6-dimethyl-             | 184    | C13H28           | 017312-82-2  | 49   |
| 5         | Heptane, 2,4,6-trimethyl-           | 142    | C10H22           | 002613-61-8  | 41   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

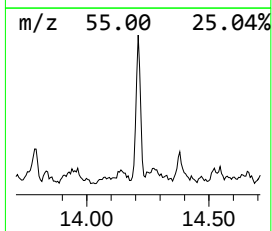
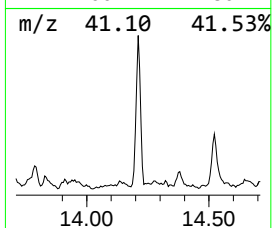
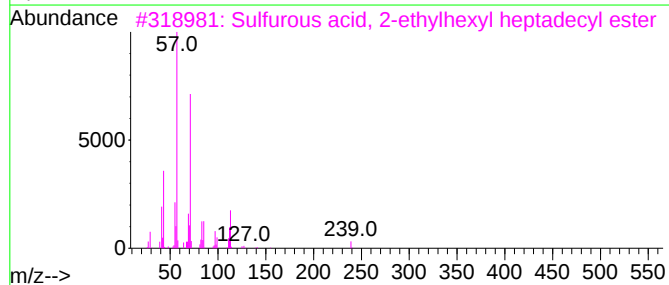
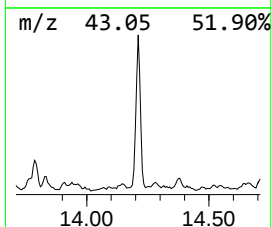
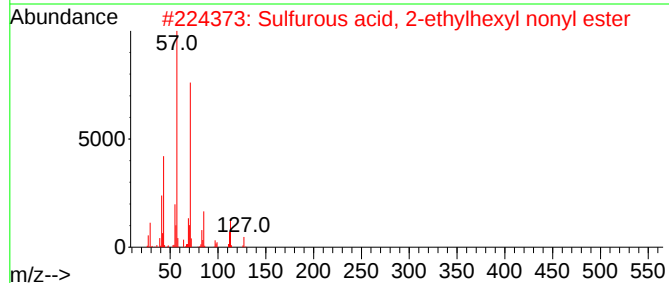
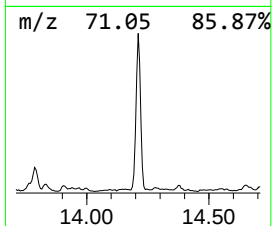
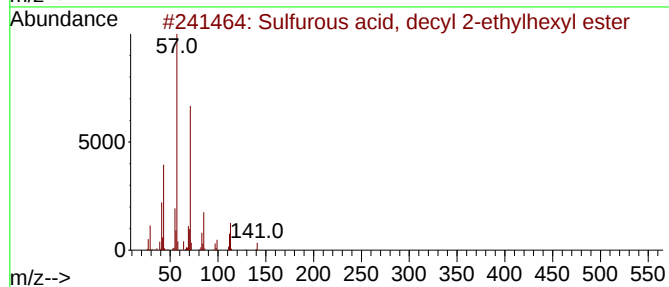
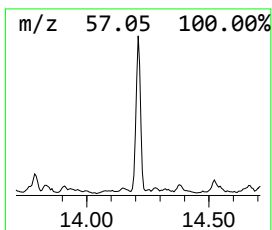
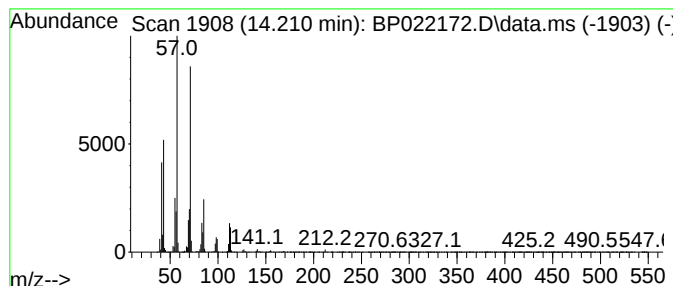
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 21 Sulfurous acid, decyl 2-eth... Concentration Rank 4

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 14.210 | 9.59 ng/ul | 764283 | Acenaphthene-d10 | 14.304 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Sulfurous acid, decyl 2-ethylhex... | 334 | C18H38O3S | 1000309-19-3 | 72   |
| 2    |    |   | Sulfurous acid, 2-ethylhexyl non... | 320 | C17H36O3S | 1010309-19-2 | 64   |
| 3    |    |   | Sulfurous acid, 2-ethylhexyl hep... | 432 | C25H52O3S | 1000309-20-0 | 64   |
| 4    |    |   | Dodecane, 2,6,10-trimethyl-         | 212 | C15H32    | 003891-98-3  | 64   |
| 5    |    |   | Heptane, 3-ethyl-5-methyl-          | 142 | C10H22    | 052896-90-9  | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

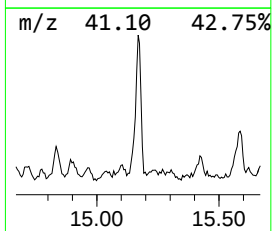
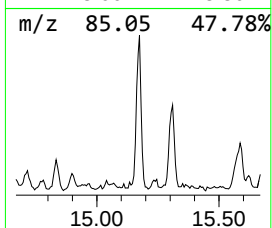
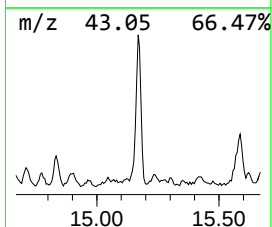
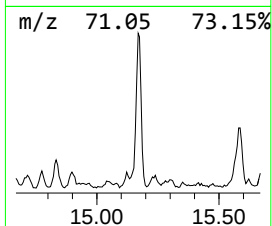
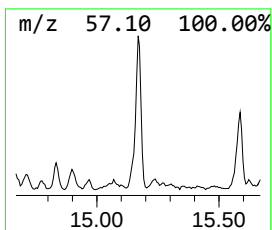
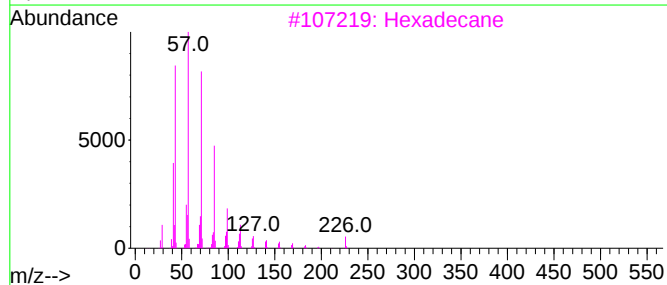
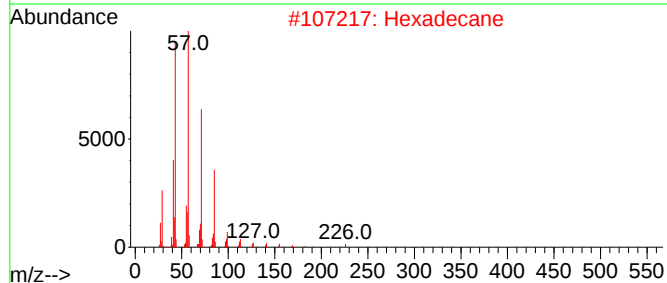
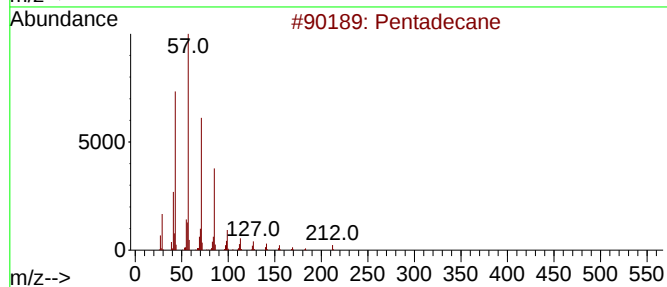
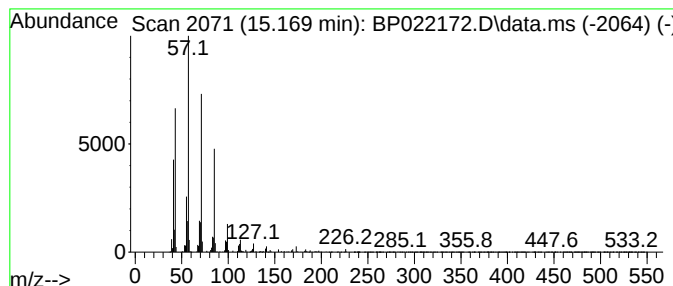
Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.      | EstConc      | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|--------------|--------|------------------|---------|-------------|------|
| 15.169    | 5.81 ng/ul   | 463386 | Acenaphthene-d10 |         | 14.304      |      |
| Hit# of 5 | Tentative ID |        | MW               | MolForm | CAS#        | Qual |
| 1         | Pentadecane  |        | 212              | C15H32  | 000629-62-9 | 93   |
| 2         | Hexadecane   |        | 226              | C16H34  | 000544-76-3 | 93   |
| 3         | Hexadecane   |        | 226              | C16H34  | 000544-76-3 | 91   |
| 4         | Tetradecane  |        | 198              | C14H30  | 000629-59-4 | 91   |
| 5         | Nonadecane   |        | 268              | C19H40  | 000629-92-5 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

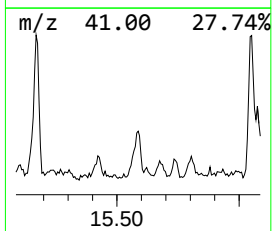
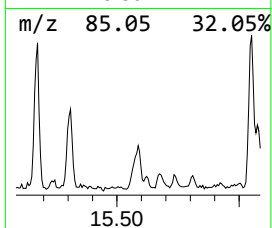
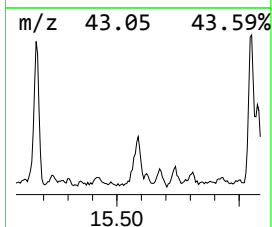
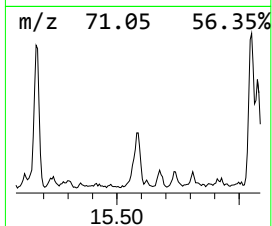
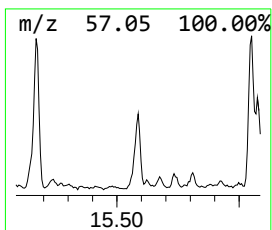
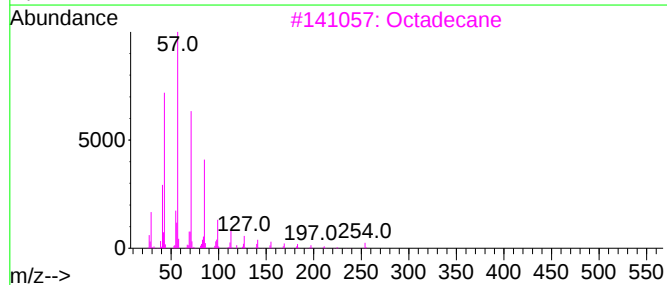
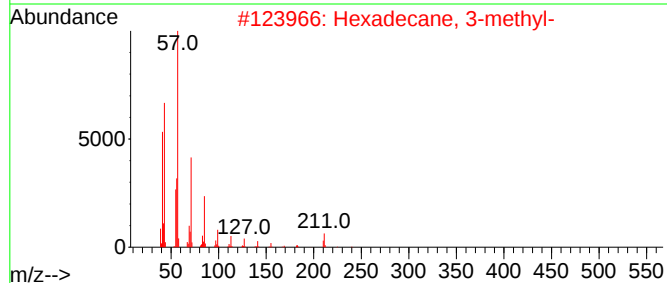
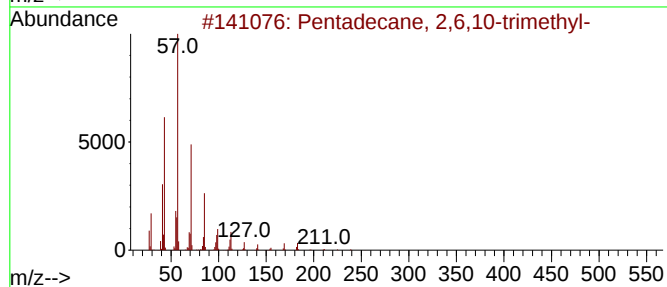
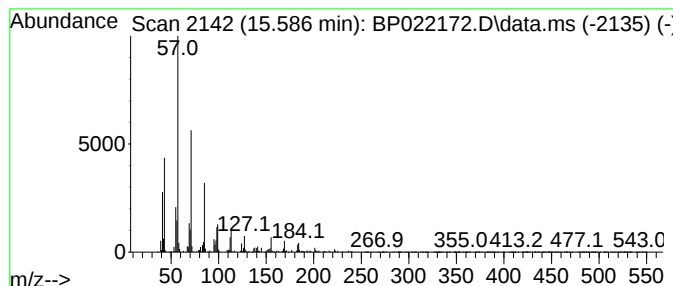
Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 30

| R.T.      | EstConc                        | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|--------------------------------|--------|------------------|---------|-------------|------|
| 15.586    | 2.41 ng/ul                     | 192510 | Acenaphthene-d10 |         | 14.304      |      |
| Hit# of 5 | Tentative ID                   |        | MW               | MolForm | CAS#        | Qual |
| 1         | Pentadecane, 2,6,10-trimethyl- |        | 254              | C18H38  | 003892-00-0 | 90   |
| 2         | Hexadecane, 3-methyl-          |        | 240              | C17H36  | 006418-43-5 | 81   |
| 3         | Octadecane                     |        | 254              | C18H38  | 000593-45-3 | 80   |
| 4         | Hentriacontane                 |        | 437              | C31H64  | 000630-04-6 | 80   |
| 5         | Decane, 2,6,8-trimethyl-       |        | 184              | C13H28  | 062108-26-3 | 76   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

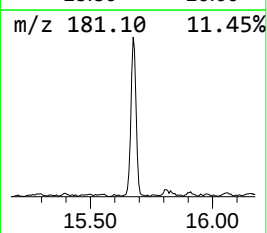
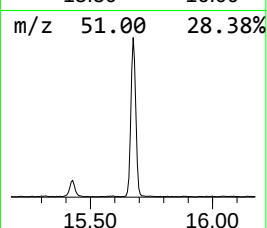
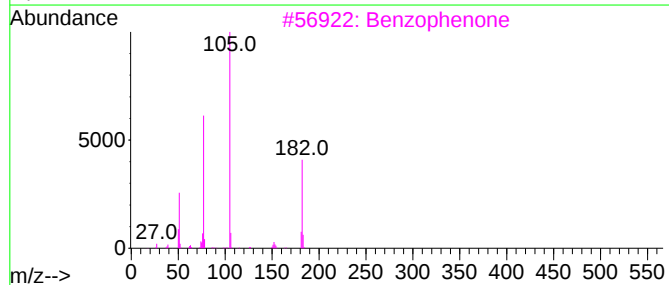
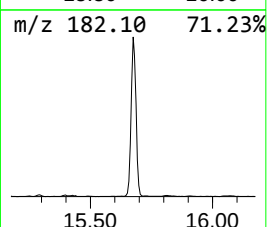
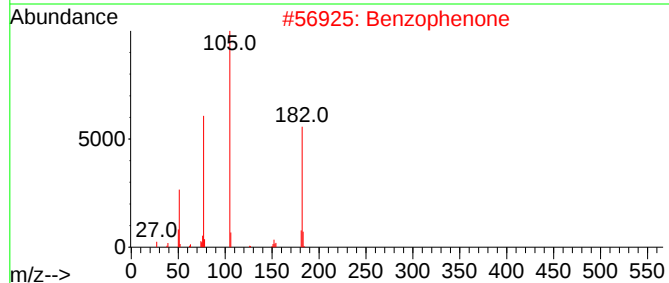
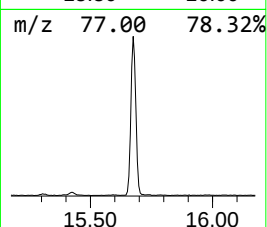
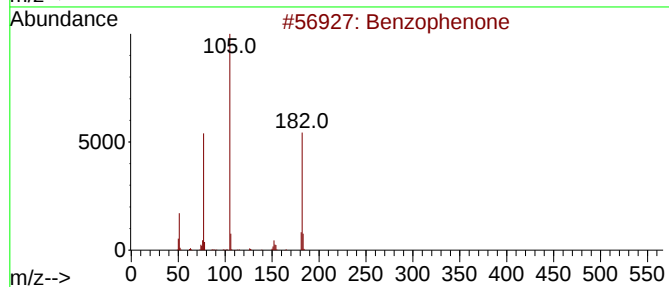
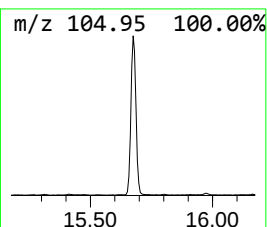
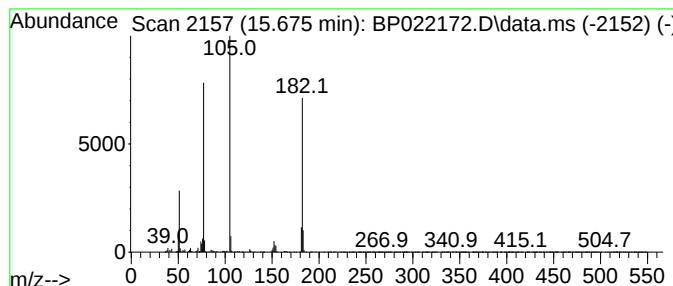
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Quant Title : SVOA CALIBRATION

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

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Peak Number 24 Benzophenone Concentration Rank 3

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 15.675 | 16.47 ng/ul | 1312990 | Acenaphthene-d10 | 14.304 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Benzophenone | 182 | C13H10O | 000119-61-9 | 97   |
| 2    |    |   | Benzophenone | 182 | C13H10O | 000119-61-9 | 96   |
| 3    |    |   | Benzophenone | 182 | C13H10O | 000119-61-9 | 94   |
| 4    |    |   | Benzophenone | 182 | C13H10O | 000119-61-9 | 89   |
| 5    |    |   | Benzophenone | 182 | C13H10O | 000119-61-9 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

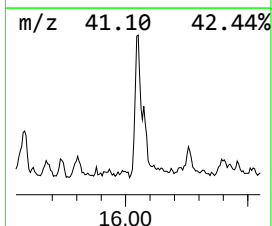
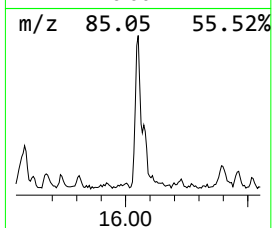
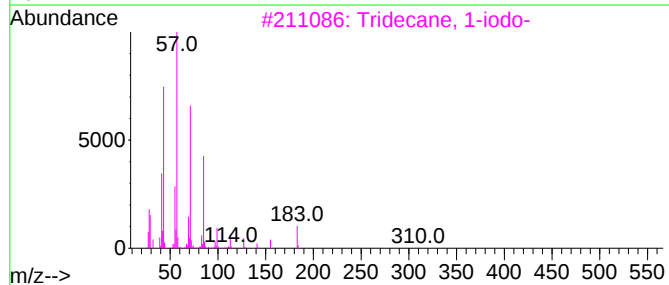
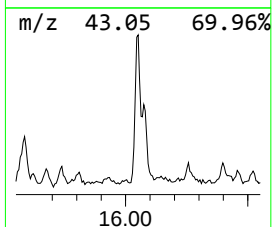
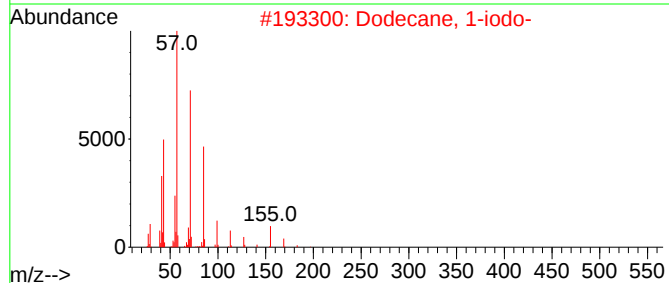
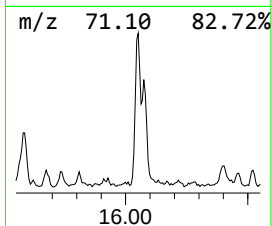
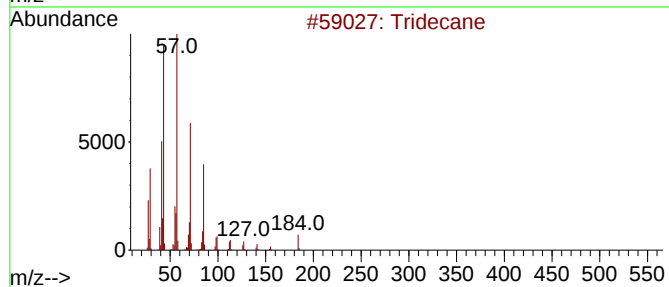
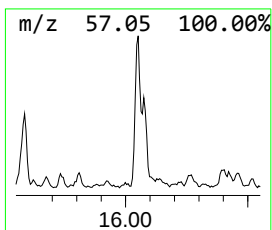
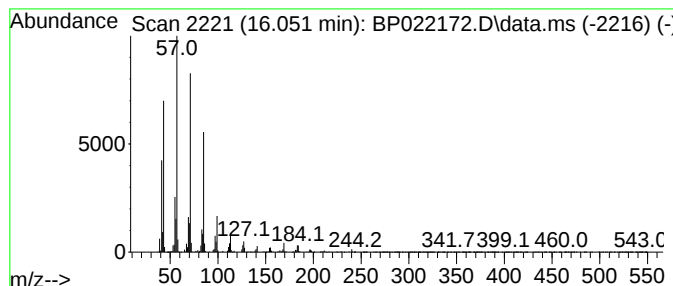
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TIC Integration Parameters: LSCINT.P

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Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.051 | 4.69 ng/ul | 477618 | Phenanthrene-d10 | 17.092 |

| Hit# | of 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|------|--------------------------------|-----|---------|-------------|------|
| 1    |      | Tridecane                      | 184 | C13H28  | 000629-50-5 | 91   |
| 2    |      | Dodecane, 1-iodo-              | 296 | C12H25I | 004292-19-7 | 90   |
| 3    |      | Tridecane, 1-iodo-             | 310 | C13H27I | 035599-77-0 | 90   |
| 4    |      | Dodecane, 2-methyl-6-propyl-   | 226 | C16H34  | 055045-08-4 | 87   |
| 5    |      | Tetradecane, 2,6,10-trimethyl- | 240 | C17H36  | 014905-56-7 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

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

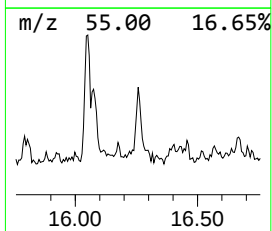
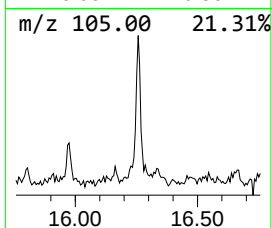
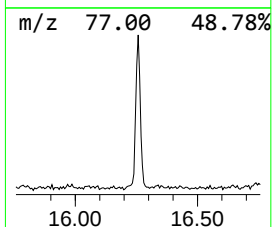
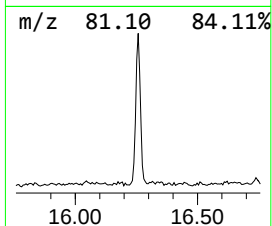
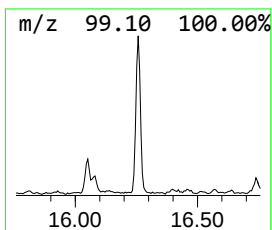
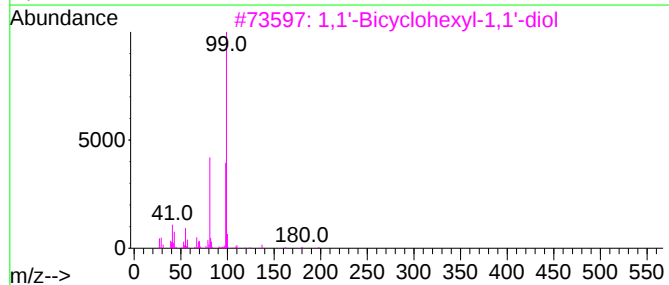
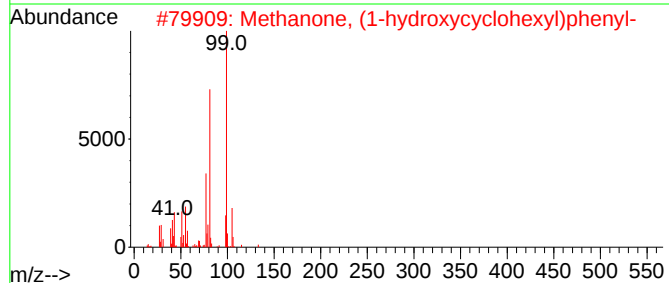
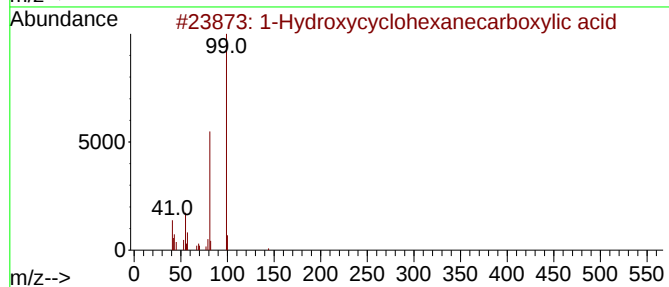
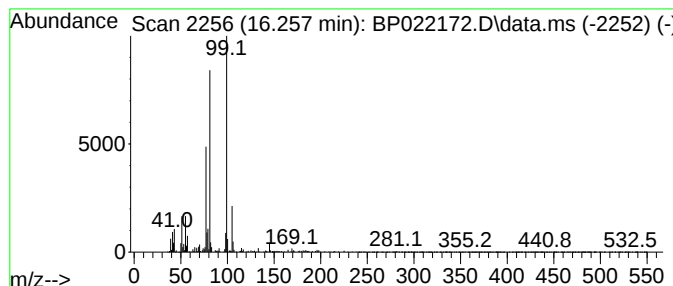
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Peak Number 26 unknown-01 Concentration Rank 26

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.257 | 2.56 ng/ul | 260625 | Phenanthrene-d10 | 17.092 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 1-Hydroxycyclohexanecarboxylic acid | 144 | C7H12O3  | 001123-28-0 | 27   |
| 2    |    |   | Methanone, (1-hydroxycyclohexyl)... | 204 | C13H16O2 | 000947-19-3 | 27   |
| 3    |    |   | 1,1'-Bicyclohexyl-1,1'-diol         | 198 | C12H22O2 | 002888-11-1 | 25   |
| 4    |    |   | 1,1'-Bicyclohexyl-1,1'-diol         | 198 | C12H22O2 | 002888-11-1 | 22   |
| 5    |    |   | Hexanethioic acid, S-heptyl ester   | 230 | C13H26OS | 002432-80-6 | 12   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

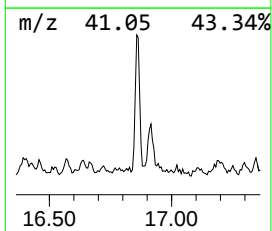
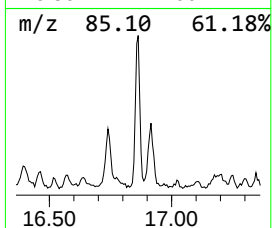
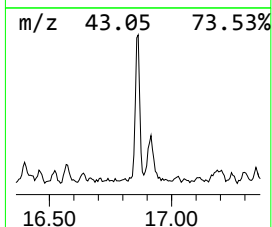
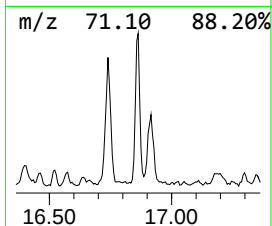
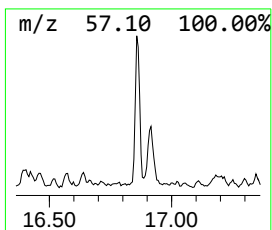
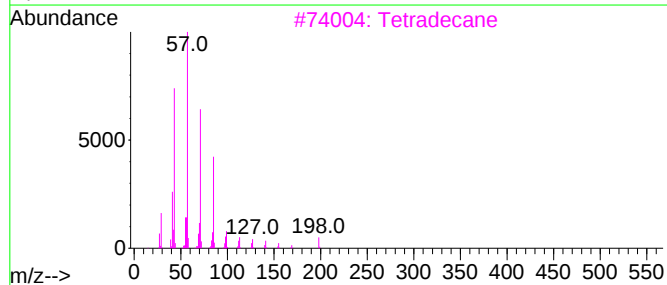
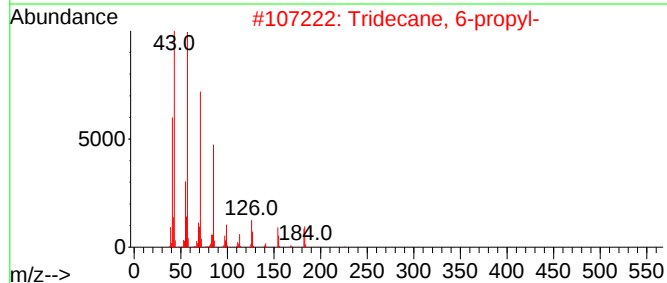
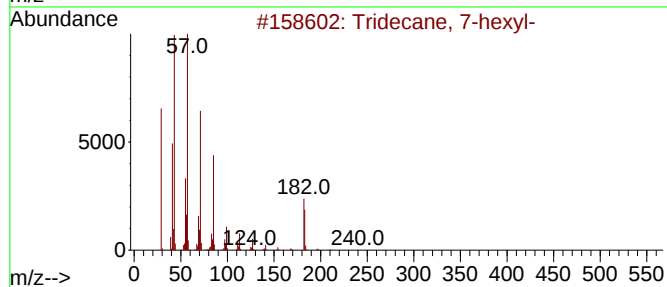
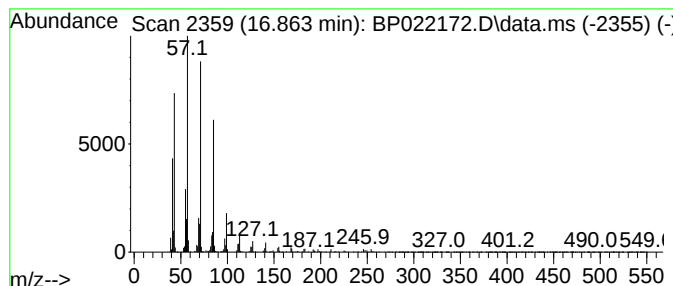
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 21

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.863 | 3.38 ng/ul | 344453 | Phenanthrene-d10 | 17.092 |

| Hit# | of 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|------|------------------------|-----|---------|-------------|------|
| 1    |      | Tridecane, 7-hexyl-    | 268 | C19H40  | 007225-66-3 | 90   |
| 2    |      | Tridecane, 6-propyl-   | 226 | C16H34  | 055045-10-8 | 87   |
| 3    |      | Tetradecane            | 198 | C14H30  | 000629-59-4 | 87   |
| 4    |      | Heptacosane            | 380 | C27H56  | 000593-49-7 | 80   |
| 5    |      | Heptadecane, 2-methyl- | 254 | C18H38  | 001560-89-0 | 74   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

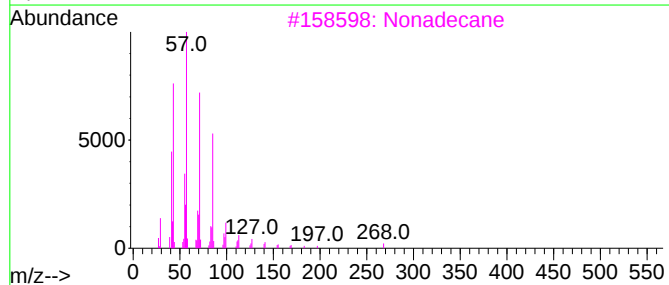
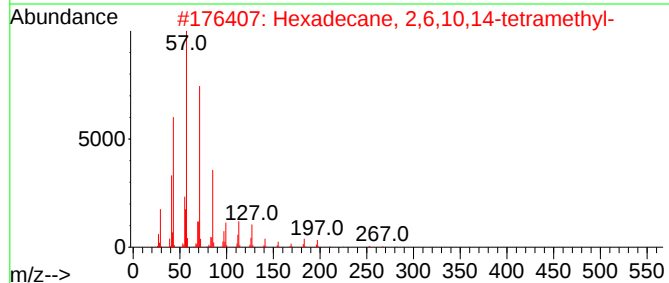
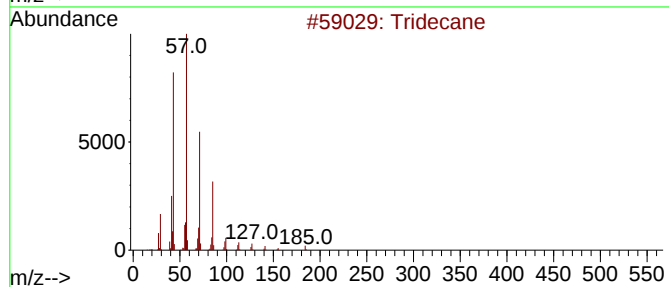
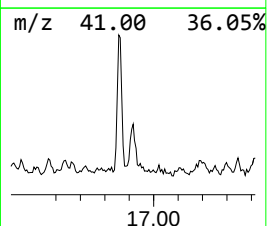
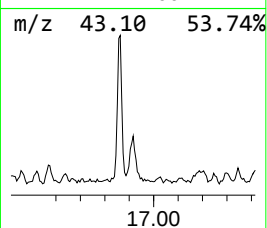
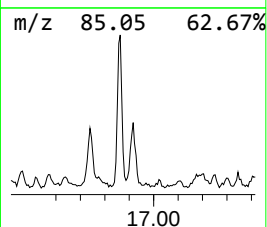
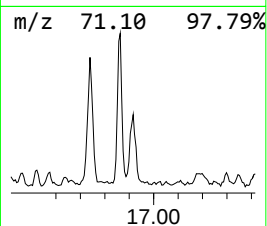
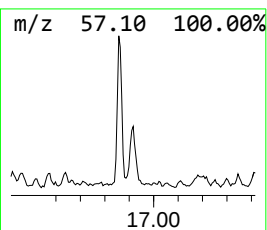
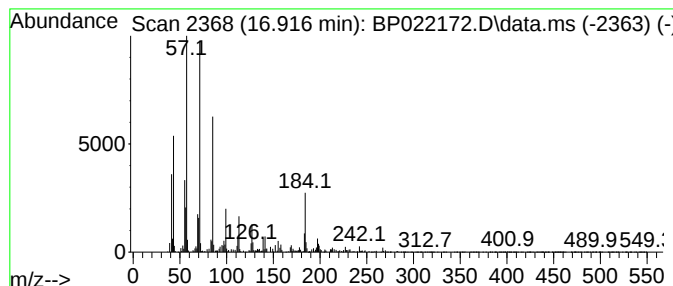
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Quant Title : SVOA CALIBRATION

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

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Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 27

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.916 | 2.46 ng/ul | 250963 | Phenanthrene-d10 | 17.092 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Tridecane                           | 184 | C13H28  | 000629-50-5 | 90   |
| 2    |      | Hexadecane, 2,6,10,14-tetramethyl-  | 282 | C20H42  | 000638-36-8 | 89   |
| 3    |      | Nonadecane                          | 268 | C19H40  | 000629-92-5 | 87   |
| 4    |      | Tridecane                           | 184 | C13H28  | 000629-50-5 | 86   |
| 5    |      | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40  | 001921-70-6 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

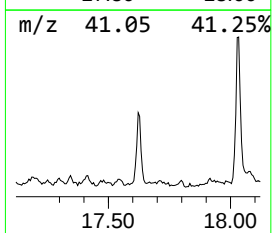
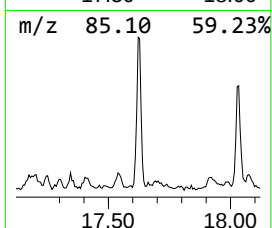
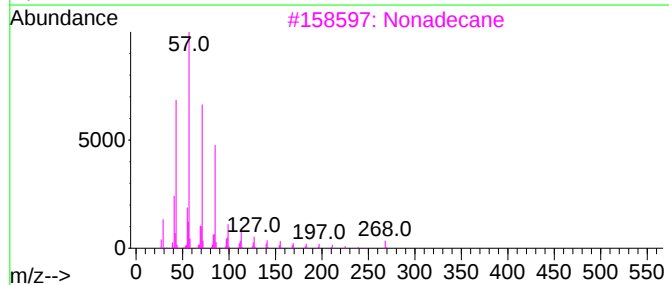
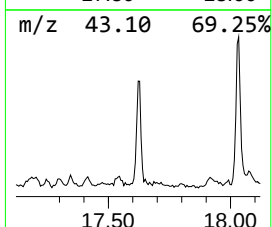
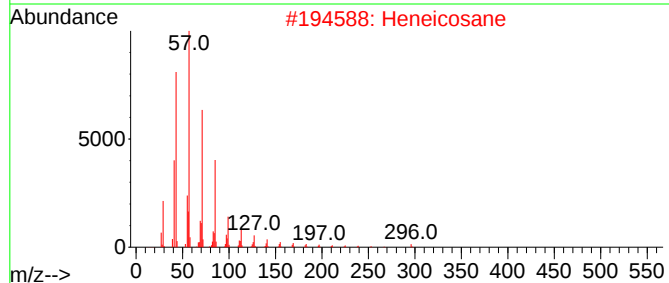
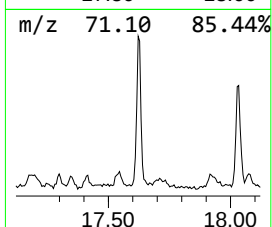
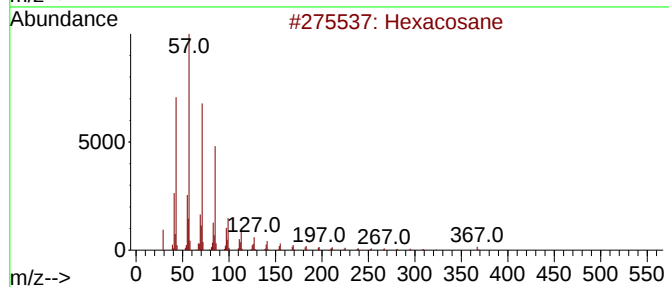
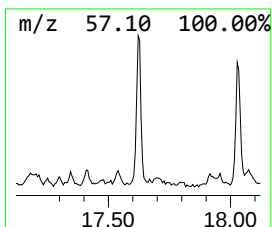
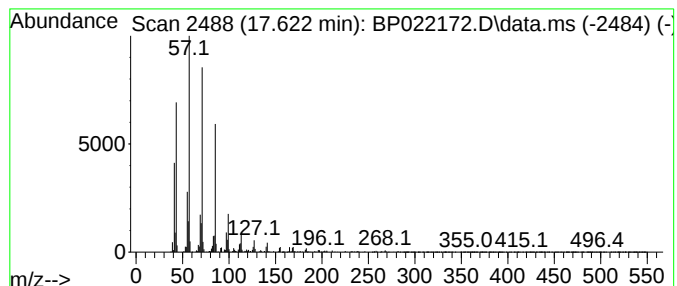
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 22

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.622 | 3.31 ng/ul | 336979 | Phenanthrene-d10 | 17.092 |

| Hit# | of 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|------|---------------------|-----|---------|-------------|------|
| 1    |      | Hexacosane          | 366 | C26H54  | 000630-01-3 | 91   |
| 2    |      | Heneicosane         | 296 | C21H44  | 000629-94-7 | 91   |
| 3    |      | Nonadecane          | 268 | C19H40  | 000629-92-5 | 90   |
| 4    |      | Tetradecane         | 198 | C14H30  | 000629-59-4 | 87   |
| 5    |      | Dodecane, 4-methyl- | 184 | C13H28  | 006117-97-1 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

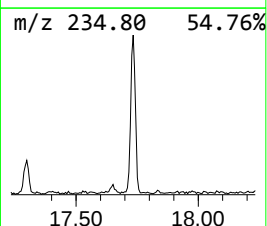
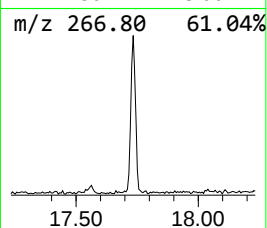
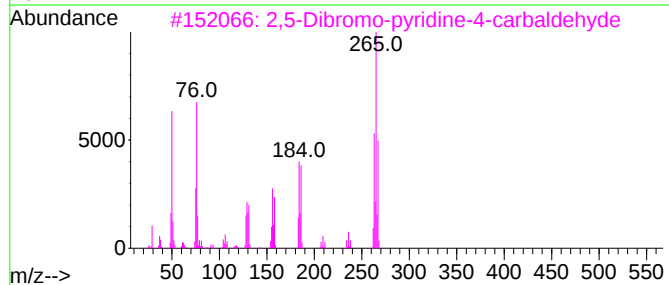
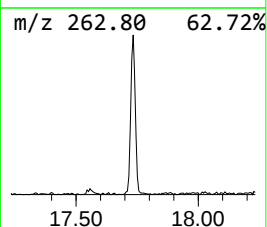
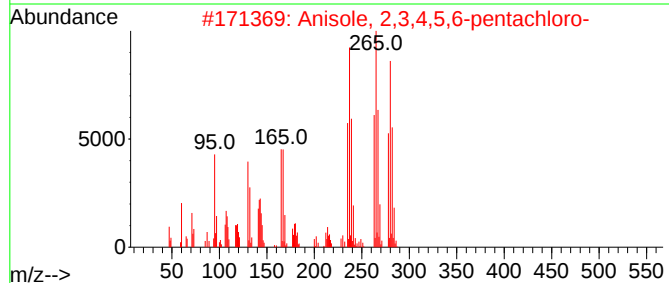
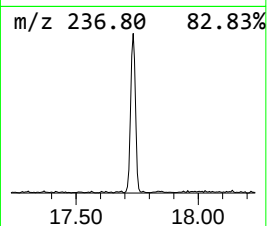
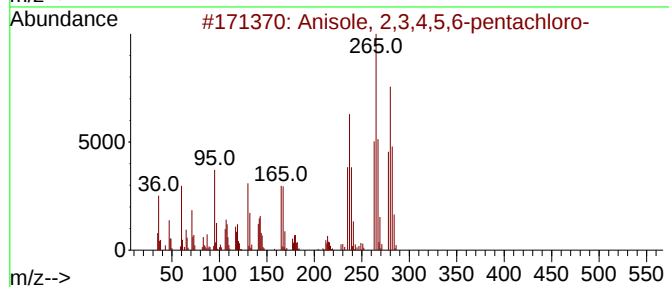
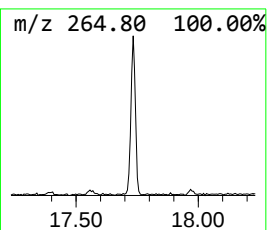
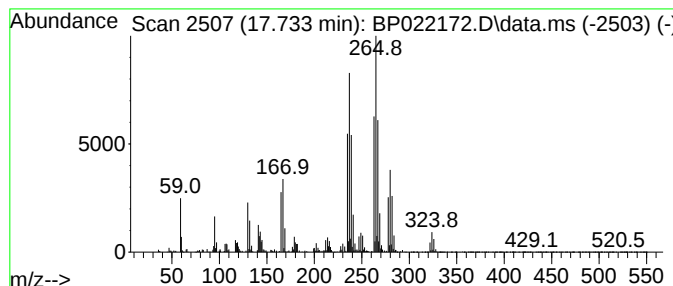
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 30 Anisole, 2,3,4,5,6-pentachl... Concentration Rank 23

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.733 | 3.13 ng/ul | 319261 | Phenanthrene-d10 | 17.092 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Anisole, 2,3,4,5,6-pentachloro-     | 278 | C7H3Cl5O    | 001825-21-4  | 55   |
| 2    |    |   | Anisole, 2,3,4,5,6-pentachloro-     | 278 | C7H3Cl5O    | 001825-21-4  | 46   |
| 3    |    |   | 2,5-Dibromo-pyridine-4-carbaldehyde | 263 | C6H3Br2NO   | 959244-28-1  | 42   |
| 4    |    |   | 2,3,4,5,6-Pentachloroaniline        | 263 | C6H2Cl5N    | 000527-20-8  | 42   |
| 5    |    |   | 1H-1,2,3-benzotriazole, 7-bromo-... | 265 | C6H2BrCl2N3 | 1000396-93-1 | 30   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

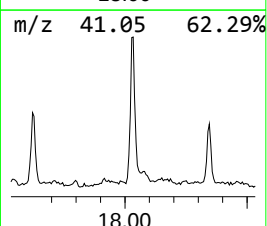
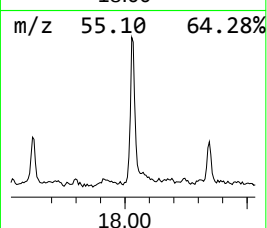
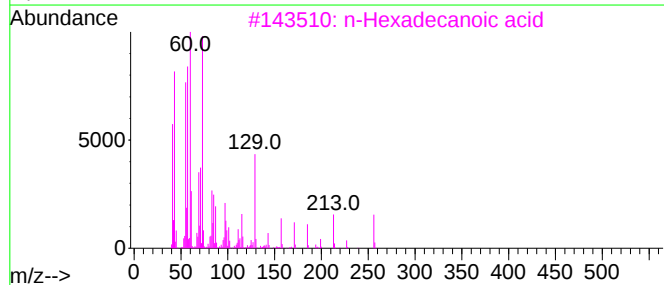
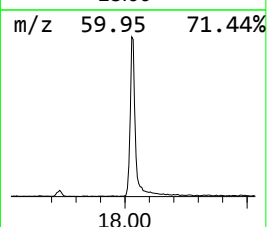
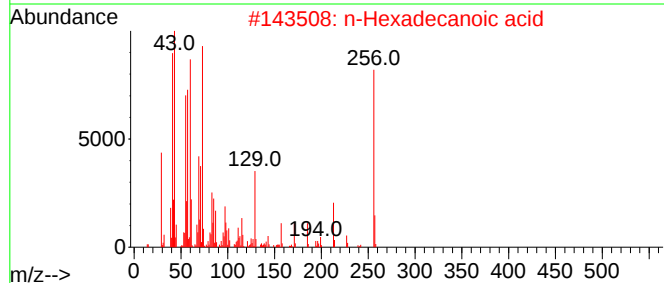
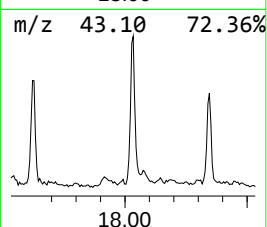
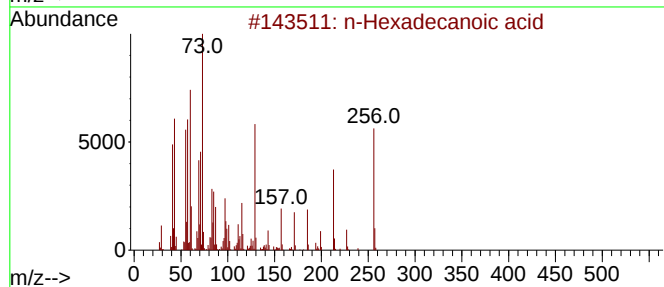
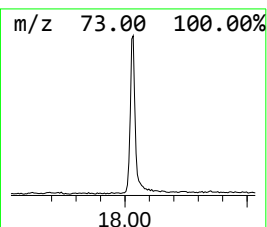
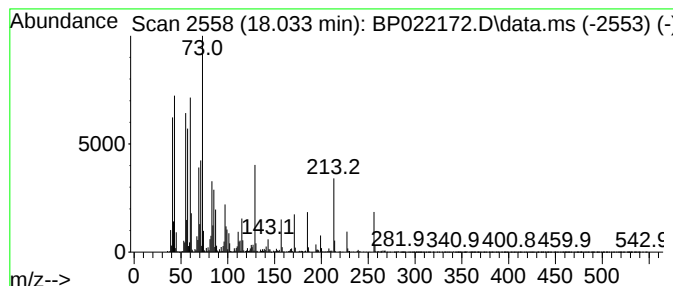
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Quant Title : SVOA CALIBRATION

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

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Peak Number 31 n-Hexadecanoic acid Concentration Rank 5

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.033 | 9.20 ng/ul | 937543 | Phenanthrene-d10 | 17.092 |

| Hit# | of                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|---------------------|-----|--------------|-------------|---------|------|------|
| 1    | n-Hexadecanoic acid | 256 | C16H32O2     | 000057-10-3 | 98      |      |      |
| 2    | n-Hexadecanoic acid | 256 | C16H32O2     | 000057-10-3 | 98      |      |      |
| 3    | n-Hexadecanoic acid | 256 | C16H32O2     | 000057-10-3 | 95      |      |      |
| 4    | n-Hexadecanoic acid | 256 | C16H32O2     | 000057-10-3 | 94      |      |      |
| 5    | n-Hexadecanoic acid | 256 | C16H32O2     | 000057-10-3 | 94      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

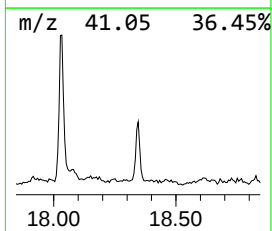
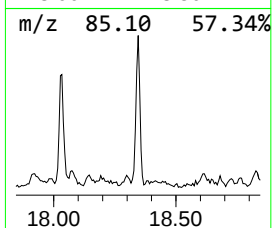
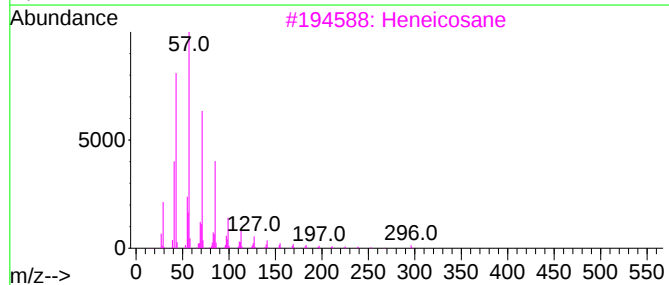
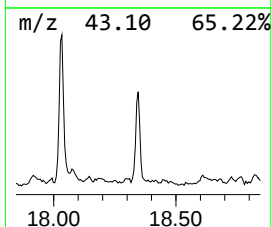
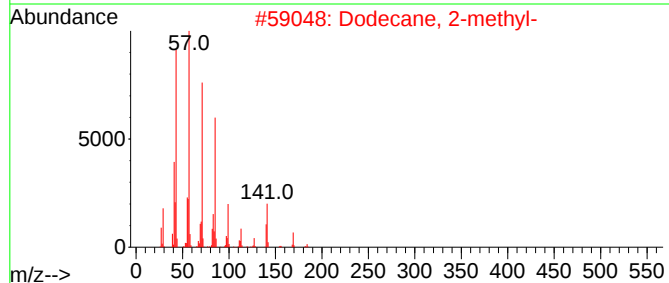
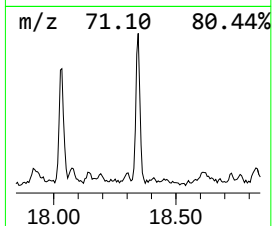
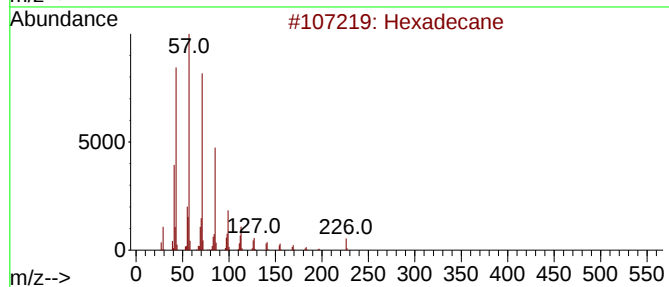
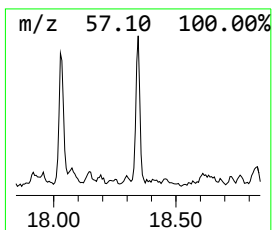
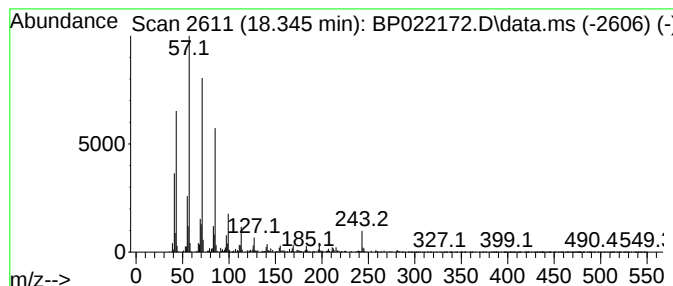
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Quant Title : SVOA CALIBRATION

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

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Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 20

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.345 | 3.40 ng/ul | 346202 | Phenanthrene-d10 | 17.092 |

| Hit# | of 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|------|---------------------|-----|---------|-------------|------|
| 1    |      | Hexadecane          | 226 | C16H34  | 000544-76-3 | 91   |
| 2    |      | Dodecane, 2-methyl- | 184 | C13H28  | 001560-97-0 | 91   |
| 3    |      | Heneicosane         | 296 | C21H44  | 000629-94-7 | 87   |
| 4    |      | Octadecane          | 254 | C18H38  | 000593-45-3 | 87   |
| 5    |      | Docosane            | 310 | C22H46  | 000629-97-0 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

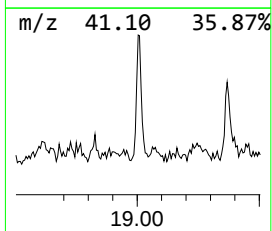
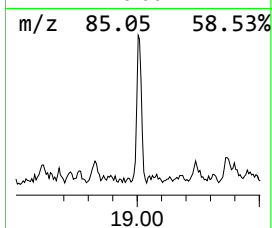
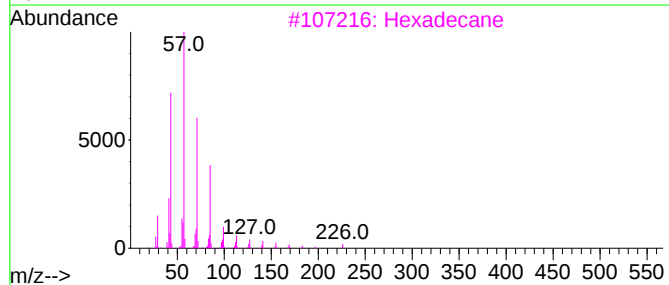
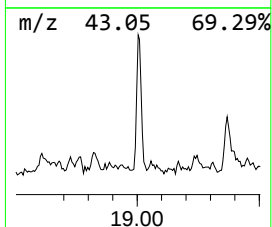
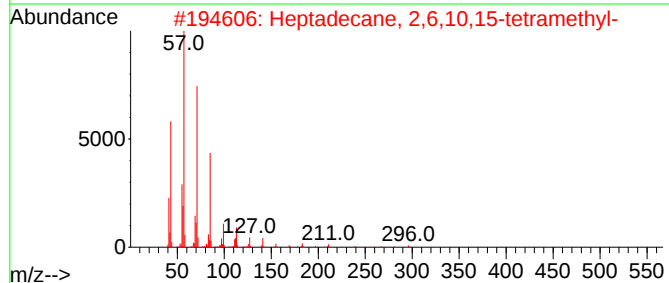
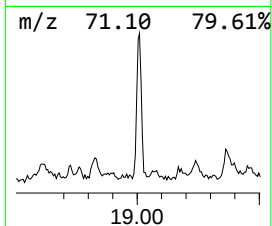
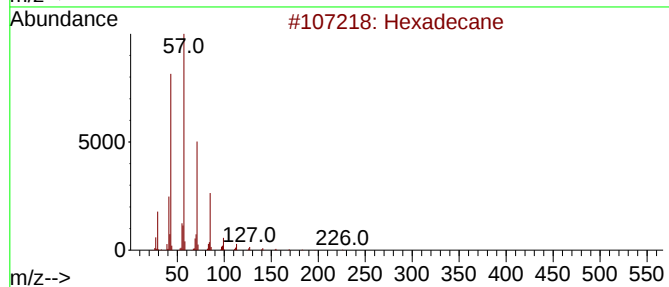
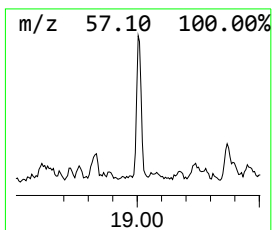
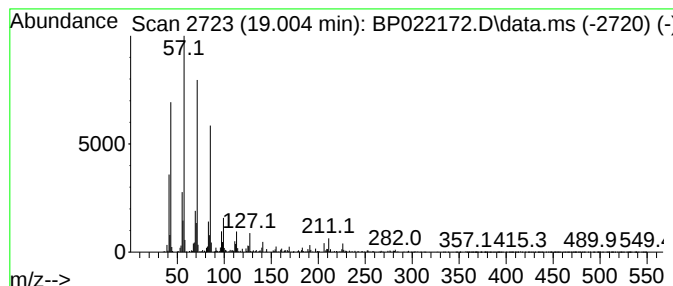
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 31

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 19.004 | 2.40 ng/ul | 244336 | Phenanthrene-d10 | 17.092 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Hexadecane                          | 226 | C16H34  | 000544-76-3 | 95   |
| 2    |      | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 93   |
| 3    |      | Hexadecane                          | 226 | C16H34  | 000544-76-3 | 93   |
| 4    |      | Hexadecane                          | 226 | C16H34  | 000544-76-3 | 91   |
| 5    |      | Hexacosane                          | 366 | C26H54  | 000630-01-3 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

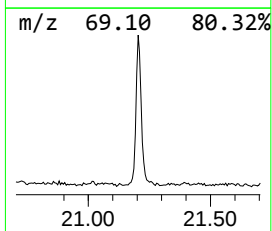
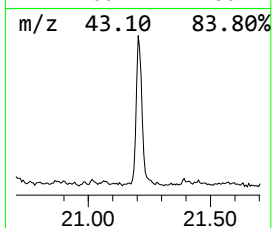
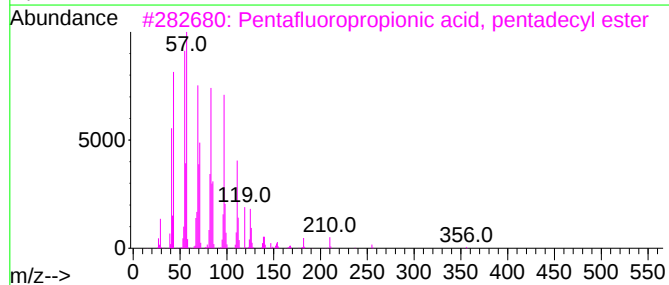
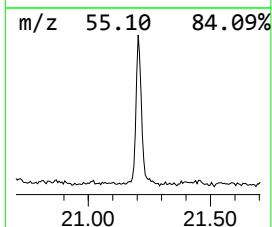
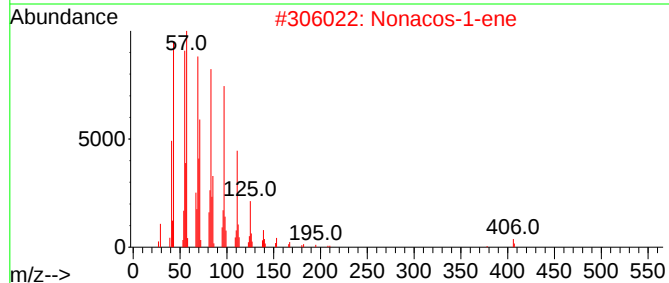
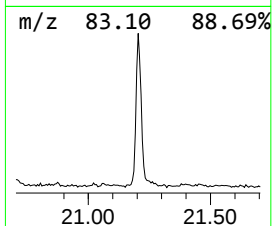
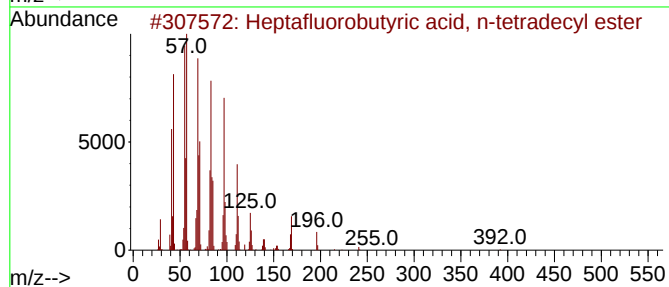
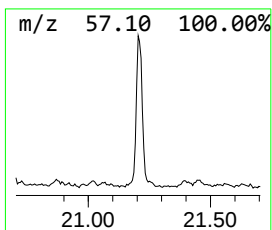
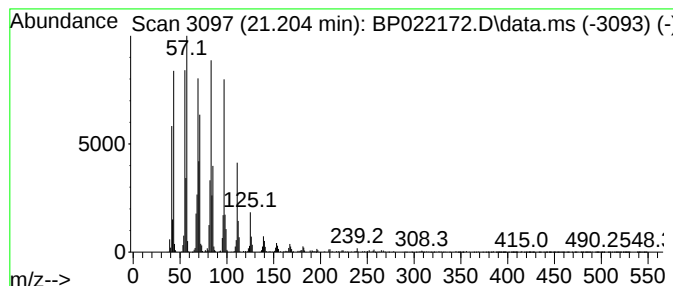
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 34 Heptafluorobutyric acid, n-... Concentration Rank 6

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 21.204 | 7.23 ng/ul | 926949 | Chrysene-d12     | 21.533 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|------------|-------------|------|
| 1    |    |   | Heptafluorobutyric acid, n-tetra... | 410 | C18H29F7O2 | 007365-36-8 | 94   |
| 2    |    |   | Nonacos-1-ene                       | 406 | C29H58     | 018835-35-3 | 94   |
| 3    |    |   | Pentafluoropropionic acid, penta... | 374 | C18H31F5O2 | 959092-08-1 | 94   |
| 4    |    |   | Heptadecyl heptafluorobutyrate      | 452 | C21H35F7O2 | 959085-66-6 | 94   |
| 5    |    |   | 5-Octadecene, (E)-                  | 252 | C18H36     | 007206-21-5 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

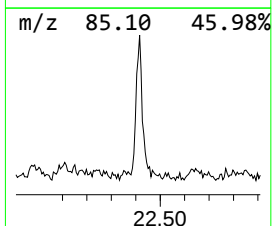
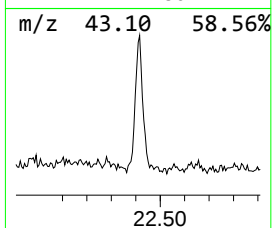
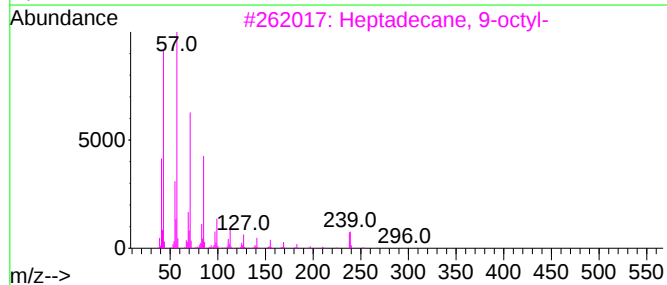
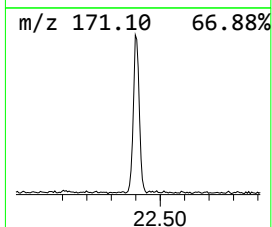
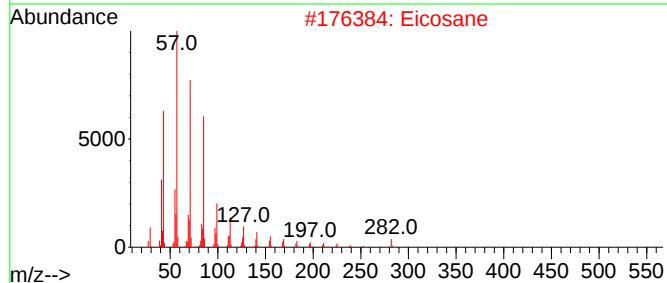
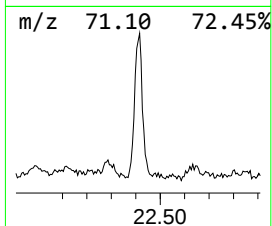
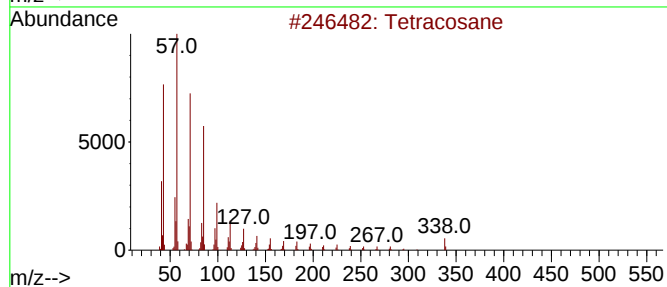
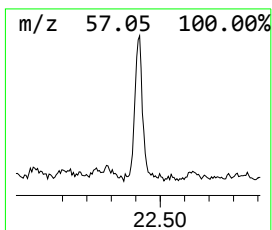
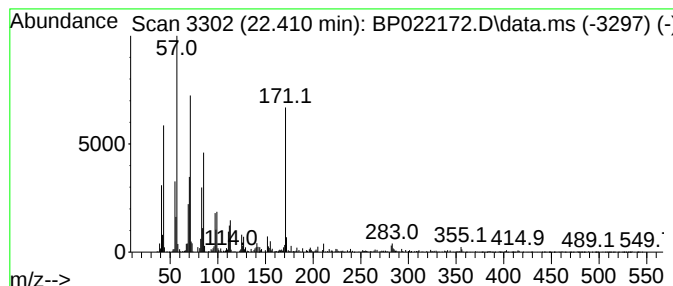
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 24

| R.T.      | EstConc                    | Area   | Relative to ISTD |          | R.T.        |      |
|-----------|----------------------------|--------|------------------|----------|-------------|------|
| 22.410    | 2.84 ng/ul                 | 363769 | Chrysene-d12     |          | 21.533      |      |
| Hit# of 5 | Tentative ID               |        | MW               | MolForm  | CAS#        | Qual |
| 1         | Tetracosane                |        | 338              | C24H50   | 000646-31-1 | 91   |
| 2         | Eicosane                   |        | 282              | C20H42   | 000112-95-8 | 91   |
| 3         | Heptadecane, 9-octyl-      |        | 352              | C25H52   | 007225-64-1 | 78   |
| 4         | Disulfide, di-tert-dodecyl |        | 402              | C24H50S2 | 027458-90-8 | 70   |
| 5         | Eicosane                   |        | 282              | C20H42   | 000112-95-8 | 70   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100224\  
Data File : BP022172.D  
Acq On : 02 Oct 2024 22:14  
Operator : RC/JU  
Sample : P4016-09  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCVW0

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION

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

|                    |        |         |       |          | --Internal Standard-- |        |         |      |  |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|--|
| TIC Top Hit name   | RT     | EstConc | Units | Response | #                     | RT     | Resp    | Conc |  |
| (DEL) Alkane: S... | 5.752  | 5.7     | ng/ul | 258221   | 1                     | 7.699  | 914383  | 20.0 |  |
| (DEL) Alkane: S... | 6.281  | 2.6     | ng/ul | 119038   | 1                     | 7.699  | 914383  | 20.0 |  |
| (DEL) Alkane: S... | 6.710  | 3.6     | ng/ul | 162568   | 1                     | 7.699  | 914383  | 20.0 |  |
| (DEL) Alkane: S... | 6.763  | 4.0     | ng/ul | 183808   | 1                     | 7.699  | 914383  | 20.0 |  |
| (DEL) Alkane: S... | 7.328  | 21.1    | ng/ul | 963486   | 1                     | 7.699  | 914383  | 20.0 |  |
| (DEL) Alkane: S... | 7.616  | 3.5     | ng/ul | 161676   | 1                     | 7.699  | 914383  | 20.0 |  |
| (DEL) Alkane: S... | 7.763  | 4.2     | ng/ul | 190227   | 1                     | 7.699  | 914383  | 20.0 |  |
| (DEL) Alkane: S... | 7.887  | 2.0     | ng/ul | 93340    | 1                     | 7.699  | 914383  | 20.0 |  |
| (DEL) Alkane: C... | 7.981  | 2.5     | ng/ul | 111964   | 1                     | 7.699  | 914383  | 20.0 |  |
| (DEL) Alkane: S... | 8.222  | 4.1     | ng/ul | 186983   | 1                     | 7.699  | 914383  | 20.0 |  |
| (DEL) Alkane: S... | 8.281  | 3.5     | ng/ul | 158904   | 1                     | 7.699  | 914383  | 20.0 |  |
| (DEL) Alkane: S... | 8.346  | 6.3     | ng/ul | 285793   | 1                     | 7.699  | 914383  | 20.0 |  |
| (DEL) Alkane: S... | 8.452  | 3.8     | ng/ul | 173718   | 1                     | 7.699  | 914383  | 20.0 |  |
| (DEL) Alkane: S... | 8.910  | 22.8    | ng/ul | 1042160  | 1                     | 7.699  | 914383  | 20.0 |  |
| Triacontan-15-ol   | 9.028  | 2.4     | ng/ul | 110809   | 1                     | 7.699  | 914383  | 20.0 |  |
| 2,4-Dipropyl-5-... | 10.834 | 2.3     | ng/ul | 230419   | 2                     | 10.457 | 1986010 | 20.0 |  |
| (DEL) Alkane: S... | 11.481 | 2.3     | ng/ul | 224301   | 2                     | 10.457 | 1986010 | 20.0 |  |
| Carbonic acid, ... | 11.869 | 4.3     | ng/ul | 424657   | 2                     | 10.457 | 1986010 | 20.0 |  |
| (DEL) Alkane: S... | 13.122 | 4.7     | ng/ul | 372123   | 3                     | 14.304 | 1594550 | 20.0 |  |
| (DEL) Alkane: S... | 13.787 | 2.2     | ng/ul | 177829   | 3                     | 14.304 | 1594550 | 20.0 |  |
| Sulfurous acid,... | 14.210 | 9.6     | ng/ul | 764283   | 3                     | 14.304 | 1594550 | 20.0 |  |
| (DEL) Alkane: S... | 15.169 | 5.8     | ng/ul | 463386   | 3                     | 14.304 | 1594550 | 20.0 |  |
| (DEL) Alkane: S... | 15.586 | 2.4     | ng/ul | 192510   | 3                     | 14.304 | 1594550 | 20.0 |  |
| Benzophenone       | 15.675 | 16.5    | ng/ul | 1312990  | 3                     | 14.304 | 1594550 | 20.0 |  |
| (DEL) Alkane: S... | 16.051 | 4.7     | ng/ul | 477618   | 4                     | 17.092 | 2038210 | 20.0 |  |
| unknown-01         | 16.257 | 2.6     | ng/ul | 260625   | 4                     | 17.092 | 2038210 | 20.0 |  |
| (DEL) Alkane: S... | 16.863 | 3.4     | ng/ul | 344453   | 4                     | 17.092 | 2038210 | 20.0 |  |
| (DEL) Alkane: S... | 16.916 | 2.5     | ng/ul | 250963   | 4                     | 17.092 | 2038210 | 20.0 |  |
| (DEL) Alkane: S... | 17.622 | 3.3     | ng/ul | 336979   | 4                     | 17.092 | 2038210 | 20.0 |  |
| Anisole, 2,3,4,... | 17.733 | 3.1     | ng/ul | 319261   | 4                     | 17.092 | 2038210 | 20.0 |  |
| n-Hexadecanoic ... | 18.033 | 9.2     | ng/ul | 937543   | 4                     | 17.092 | 2038210 | 20.0 |  |
| (DEL) Alkane: S... | 18.345 | 3.4     | ng/ul | 346202   | 4                     | 17.092 | 2038210 | 20.0 |  |
| (DEL) Alkane: S... | 19.004 | 2.4     | ng/ul | 244336   | 4                     | 17.092 | 2038210 | 20.0 |  |
| Heptafluorobuty... | 21.204 | 7.2     | ng/ul | 926949   | 5                     | 21.533 | 2564490 | 20.0 |  |
| (DEL) Alkane: S... | 22.410 | 2.8     | ng/ul | 363769   | 5                     | 21.533 | 2564490 | 20.0 |  |