

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
 Data File : BP011263.D  
 Acq On : 04 Aug 2022 14:20  
 Operator : CG/JU  
 Sample : N3956-08  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EXF05

Integration Parameters: LSCINT.P

Integrator: RTE

Smoother: OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
 Title : SVOA CALIBRATION

Signal : TIC: BP011263.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         | 3.228       | 36            | 41          | 53           | rVB      | 1555753        | 2170050       | 0.81%           | 0.151%        |
| 2         | 3.558       | 92            | 97          | 113          | rVB      | 7348197        | 9760603       | 3.64%           | 0.677%        |
| 3         | 3.834       | 137           | 144         | 149          | rBV      | 12299420       | 18979004      | 7.07%           | 1.317%        |
| 4         | 4.081       | 174           | 186         | 190          | rVB      | 20237410       | 53954578      | 20.11%          | 3.745%        |
| 5         | 4.446       | 244           | 248         | 251          | rVV      | 1648864        | 2285370       | 0.85%           | 0.159%        |
| 6         | 4.481       | 251           | 254         | 261          | rVB      | 1448422        | 1909871       | 0.71%           | 0.133%        |
| 7         | 4.934       | 325           | 331         | 336          | rVV      | 9712814        | 14000383      | 5.22%           | 0.972%        |
| 8         | 5.005       | 336           | 343         | 349          | rVV2     | 869648         | 1482089       | 0.55%           | 0.103%        |
| 9         | 5.217       | 373           | 379         | 385          | rVV      | 1865186        | 2767769       | 1.03%           | 0.192%        |
| 10        | 5.469       | 417           | 422         | 426          | rVV      | 934801         | 1357699       | 0.51%           | 0.094%        |
| 11        | 5.787       | 460           | 476         | 481          | rVV      | 16500820       | 33079354      | 12.33%          | 2.296%        |
| 12        | 6.287       | 553           | 561         | 567          | rVV2     | 6054280        | 10677777      | 3.98%           | 0.741%        |
| 13        | 6.605       | 611           | 615         | 623          | rVB      | 1594454        | 2337894       | 0.87%           | 0.162%        |
| 14        | 6.852       | 646           | 657         | 669          | rVB3     | 5700872        | 13281373      | 4.95%           | 0.922%        |
| 15        | 6.958       | 669           | 675         | 682          | rVB      | 1019983        | 1544409       | 0.58%           | 0.107%        |
| 16        | 7.040       | 682           | 689         | 695          | rBV      | 6389829        | 9318061       | 3.47%           | 0.647%        |
| 17        | 7.187       | 702           | 714         | 720          | rBV      | 14588058       | 29590531      | 11.03%          | 2.054%        |
| 18        | 7.605       | 774           | 785         | 792          | rBV4     | 1678796        | 3587594       | 1.34%           | 0.249%        |
| 19        | 7.681       | 792           | 798         | 809          | rVB      | 2852371        | 4959915       | 1.85%           | 0.344%        |
| 20        | 7.940       | 829           | 842         | 847          | rBV      | 28246092       | 68792089      | 25.64%          | 4.774%        |
| 21        | 8.011       | 847           | 854         | 862          | rVV      | 6729157        | 10712867      | 3.99%           | 0.744%        |
| 22        | 8.210       | 879           | 888         | 895          | rBV2     | 3662237        | 6654956       | 2.48%           | 0.462%        |
| 23        | 8.305       | 895           | 904         | 908          | rBV      | 856719         | 1509450       | 0.56%           | 0.105%        |
| 24        | 8.381       | 908           | 917         | 922          | rVB2     | 1400481        | 2539727       | 0.95%           | 0.176%        |
| 25        | 8.434       | 922           | 926         | 931          | rVB      | 1059230        | 1665091       | 0.62%           | 0.116%        |
| 26        | 8.493       | 931           | 936         | 941          | rVB      | 1974940        | 2865373       | 1.07%           | 0.199%        |
| 27        | 8.587       | 947           | 952         | 957          | rVB      | 1852980        | 2753698       | 1.03%           | 0.191%        |
| 28        | 8.716       | 957           | 974         | 979          | rBV      | 18001978       | 54959545      | 20.48%          | 3.814%        |
| 29        | 8.769       | 979           | 983         | 989          | rVB2     | 1511282        | 2515311       | 0.94%           | 0.175%        |
| 30        | 8.963       | 1010          | 1016        | 1020         | rBV3     | 1100423        | 2629017       | 0.98%           | 0.182%        |
| 31        | 9.099       | 1033          | 1039        | 1047         | rVV      | 8828327        | 14087720      | 5.25%           | 0.978%        |
| 32        | 9.299       | 1064          | 1073        | 1085         | rVB3     | 682771         | 1641623       | 0.61%           | 0.114%        |
| 33        | 9.428       | 1085          | 1095        | 1099         | rBV      | 17962192       | 31375146      | 11.69%          | 2.178%        |
| 34        | 9.481       | 1099          | 1104        | 1112         | rVV3     | 2853790        | 7914961       | 2.95%           | 0.549%        |
| 35        | 9.557       | 1112          | 1117        | 1123         | rVB      | 2449442        | 4103898       | 1.53%           | 0.285%        |
| 36        | 9.681       | 1130          | 1138        | 1148         | rBV      | 1878496        | 4639214       | 1.73%           | 0.322%        |

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Integrator: RTE

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

Peak separation: 5

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 Title : SVOA CALIBRATION

|    |        |      |      |      |      |          |           |         |         |
|----|--------|------|------|------|------|----------|-----------|---------|---------|
| 37 | 9.822  | 1157 | 1162 | 1171 | rVB2 | 1279660  | 2312025   | 0.86%   | 0.160%  |
| 38 | 10.105 | 1171 | 1210 | 1214 | rBV7 | 27830982 | 268311747 | 100.00% | 18.622% |
| 39 | 10.428 | 1256 | 1265 | 1277 | rVB2 | 12232759 | 24039552  | 8.96%   | 1.668%  |
| 40 | 10.622 | 1285 | 1298 | 1301 | rBV  | 12009845 | 24094537  | 8.98%   | 1.672%  |
| 41 | 10.657 | 1301 | 1304 | 1306 | rVV  | 5010526  | 7433335   | 2.77%   | 0.516%  |
| 42 | 10.740 | 1306 | 1318 | 1336 | rVB  | 20747033 | 65654750  | 24.47%  | 4.557%  |
| 43 | 10.957 | 1341 | 1355 | 1357 | rBV2 | 20402331 | 55825310  | 20.81%  | 3.874%  |
| 44 | 10.999 | 1357 | 1362 | 1367 | rVB2 | 30794245 | 66222479  | 24.68%  | 4.596%  |
| 45 | 11.299 | 1406 | 1413 | 1420 | rBV  | 2807656  | 5163158   | 1.92%   | 0.358%  |
| 46 | 11.375 | 1420 | 1426 | 1432 | rVB2 | 3761870  | 6464158   | 2.41%   | 0.449%  |
| 47 | 11.457 | 1432 | 1440 | 1446 | rBV4 | 2183543  | 5257559   | 1.96%   | 0.365%  |
| 48 | 11.651 | 1463 | 1473 | 1480 | rBV9 | 593855   | 2058850   | 0.77%   | 0.143%  |
| 49 | 12.187 | 1546 | 1564 | 1569 | rVB2 | 34605953 | 106062010 | 39.53%  | 7.361%  |
| 50 | 12.328 | 1575 | 1588 | 1594 | rBV  | 21012939 | 48883062  | 18.22%  | 3.393%  |
| 51 | 12.393 | 1594 | 1599 | 1605 | rBV  | 2810103  | 4757489   | 1.77%   | 0.330%  |
| 52 | 12.487 | 1605 | 1615 | 1625 | rVB5 | 3589260  | 10480950  | 3.91%   | 0.727%  |
| 53 | 12.616 | 1626 | 1637 | 1648 | rBV  | 22069486 | 52682944  | 19.63%  | 3.656%  |
| 54 | 12.728 | 1648 | 1656 | 1660 | rBV2 | 11198765 | 23788150  | 8.87%   | 1.651%  |
| 55 | 12.769 | 1660 | 1663 | 1664 | rBV2 | 1247083  | 1420592   | 0.53%   | 0.099%  |
| 56 | 12.998 | 1696 | 1702 | 1707 | rBV  | 3668930  | 6043545   | 2.25%   | 0.419%  |
| 57 | 13.051 | 1707 | 1711 | 1717 | rVB2 | 1953149  | 3226628   | 1.20%   | 0.224%  |
| 58 | 13.116 | 1717 | 1722 | 1728 | rBV2 | 3203005  | 4884001   | 1.82%   | 0.339%  |
| 59 | 13.234 | 1735 | 1742 | 1747 | rBV  | 3877790  | 6495435   | 2.42%   | 0.451%  |
| 60 | 13.410 | 1769 | 1772 | 1778 | rVB  | 1040335  | 1503515   | 0.56%   | 0.104%  |
| 61 | 13.604 | 1799 | 1805 | 1812 | rBV4 | 1341974  | 2937245   | 1.09%   | 0.204%  |
| 62 | 13.722 | 1812 | 1825 | 1840 | rVV3 | 16914396 | 33041123  | 12.31%  | 2.293%  |
| 63 | 13.869 | 1840 | 1850 | 1859 | rVV  | 5191824  | 7845812   | 2.92%   | 0.545%  |
| 64 | 13.969 | 1860 | 1867 | 1875 | rVV  | 2975555  | 5502472   | 2.05%   | 0.382%  |
| 65 | 14.051 | 1875 | 1881 | 1889 | rVB  | 6710010  | 10535815  | 3.93%   | 0.731%  |
| 66 | 14.198 | 1900 | 1906 | 1912 | rBV2 | 1133951  | 1853587   | 0.69%   | 0.129%  |
| 67 | 14.281 | 1914 | 1920 | 1927 | rVV  | 2453857  | 3731152   | 1.39%   | 0.259%  |
| 68 | 14.345 | 1927 | 1931 | 1939 | rVV  | 1550242  | 2451339   | 0.91%   | 0.170%  |
| 69 | 14.428 | 1939 | 1945 | 1950 | rBV  | 2051919  | 2908071   | 1.08%   | 0.202%  |
| 70 | 14.481 | 1950 | 1954 | 1958 | rVV  | 3007647  | 3766762   | 1.40%   | 0.261%  |
| 71 | 14.528 | 1958 | 1962 | 1970 | rVB5 | 812287   | 1474912   | 0.55%   | 0.102%  |
| 72 | 14.610 | 1970 | 1976 | 1980 | rBV  | 3487096  | 4534901   | 1.69%   | 0.315%  |
| 73 | 14.940 | 2027 | 2032 | 2039 | rVB4 | 741622   | 1380939   | 0.51%   | 0.096%  |
| 74 | 15.087 | 2051 | 2057 | 2062 | rBV  | 3795206  | 5337374   | 1.99%   | 0.370%  |
| 75 | 15.192 | 2067 | 2075 | 2083 | rVB  | 4032038  | 6482133   | 2.42%   | 0.450%  |
| 76 | 15.281 | 2084 | 2090 | 2094 | rBV  | 3824453  | 5506895   | 2.05%   | 0.382%  |

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 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EXF05

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-BP080222.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |          |       |        |
|-----|--------|------|------|------|------|---------|----------|-------|--------|
| 77  | 15.328 | 2094 | 2098 | 2102 | rVV3 | 2399155 | 3653548  | 1.36% | 0.254% |
| 78  | 15.369 | 2102 | 2105 | 2109 | rVV2 | 1651073 | 2179234  | 0.81% | 0.151% |
| 79  | 15.416 | 2109 | 2113 | 2117 | rVB  | 1228670 | 1584293  | 0.59% | 0.110% |
| 80  | 15.510 | 2124 | 2129 | 2135 | rVB  | 4138042 | 5973552  | 2.23% | 0.415% |
| 81  | 15.598 | 2139 | 2144 | 2147 | rBV2 | 4333230 | 6691426  | 2.49% | 0.464% |
| 82  | 15.640 | 2147 | 2151 | 2154 | rVB  | 2555152 | 3730958  | 1.39% | 0.259% |
| 83  | 15.851 | 2181 | 2187 | 2192 | rBV3 | 831353  | 2078897  | 0.77% | 0.144% |
| 84  | 15.904 | 2192 | 2196 | 2203 | rVB  | 1624006 | 2370799  | 0.88% | 0.165% |
| 85  | 16.016 | 2209 | 2215 | 2218 | rBV  | 1359749 | 2048175  | 0.76% | 0.142% |
| 86  | 16.104 | 2226 | 2230 | 2232 | rBV  | 1979158 | 2751969  | 1.03% | 0.191% |
| 87  | 16.192 | 2241 | 2245 | 2249 | rBV  | 1244442 | 1617773  | 0.60% | 0.112% |
| 88  | 16.710 | 2326 | 2333 | 2340 | rVB2 | 672631  | 1429756  | 0.53% | 0.099% |
| 89  | 16.810 | 2340 | 2350 | 2352 | rBV  | 5885863 | 11304760 | 4.21% | 0.785% |
| 90  | 16.839 | 2352 | 2355 | 2360 | rVB  | 8890551 | 11462714 | 4.27% | 0.796% |
| 91  | 17.045 | 2384 | 2390 | 2397 | rVB  | 2497682 | 4243690  | 1.58% | 0.295% |
| 92  | 17.151 | 2402 | 2408 | 2412 | rBV2 | 3642856 | 5256092  | 1.96% | 0.365% |
| 93  | 17.722 | 2501 | 2505 | 2513 | rVB  | 1446140 | 2326286  | 0.87% | 0.161% |
| 94  | 17.845 | 2519 | 2526 | 2531 | rVB2 | 3490960 | 6402259  | 2.39% | 0.444% |
| 95  | 17.898 | 2531 | 2535 | 2538 | rBV  | 2013813 | 2557774  | 0.95% | 0.178% |
| 96  | 19.410 | 2787 | 2792 | 2798 | rVB  | 4373752 | 5833586  | 2.17% | 0.405% |
| 97  | 19.475 | 2799 | 2803 | 2808 | rVB  | 2027532 | 2369151  | 0.88% | 0.164% |
| 98  | 21.175 | 3087 | 3092 | 3096 | rBV  | 2539332 | 3228780  | 1.20% | 0.224% |
| 99  | 23.339 | 3454 | 3460 | 3468 | rVB2 | 3034610 | 5669160  | 2.11% | 0.393% |
| 100 | 23.486 | 3480 | 3485 | 3497 | rVB2 | 1708242 | 3344736  | 1.25% | 0.232% |

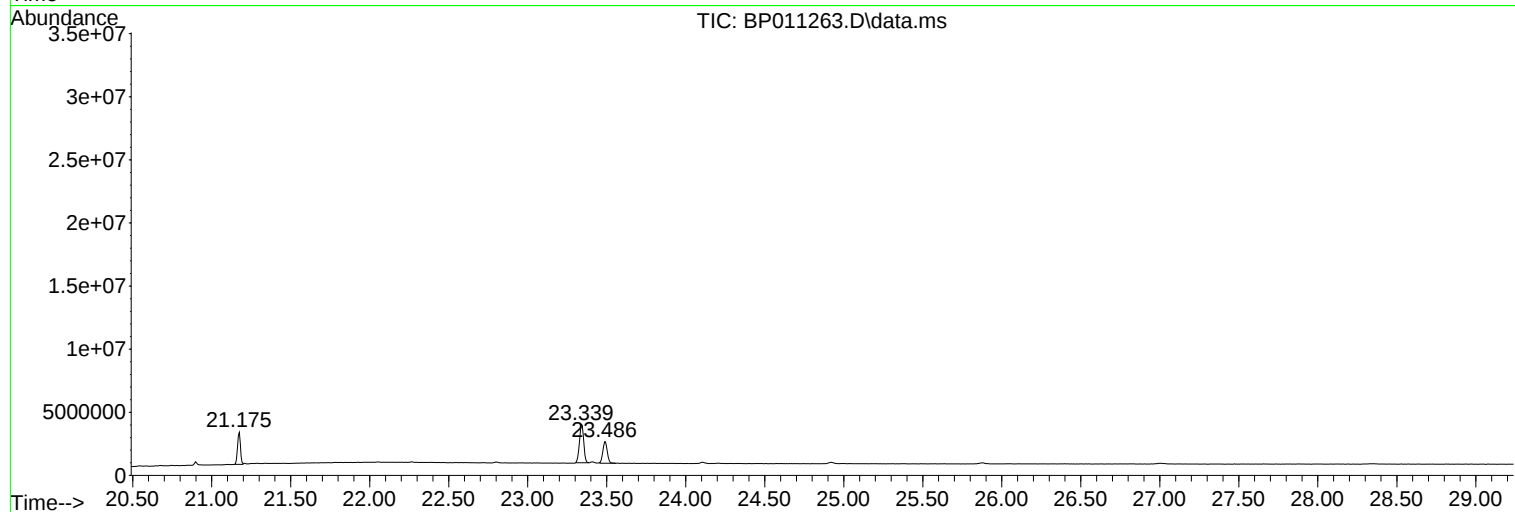
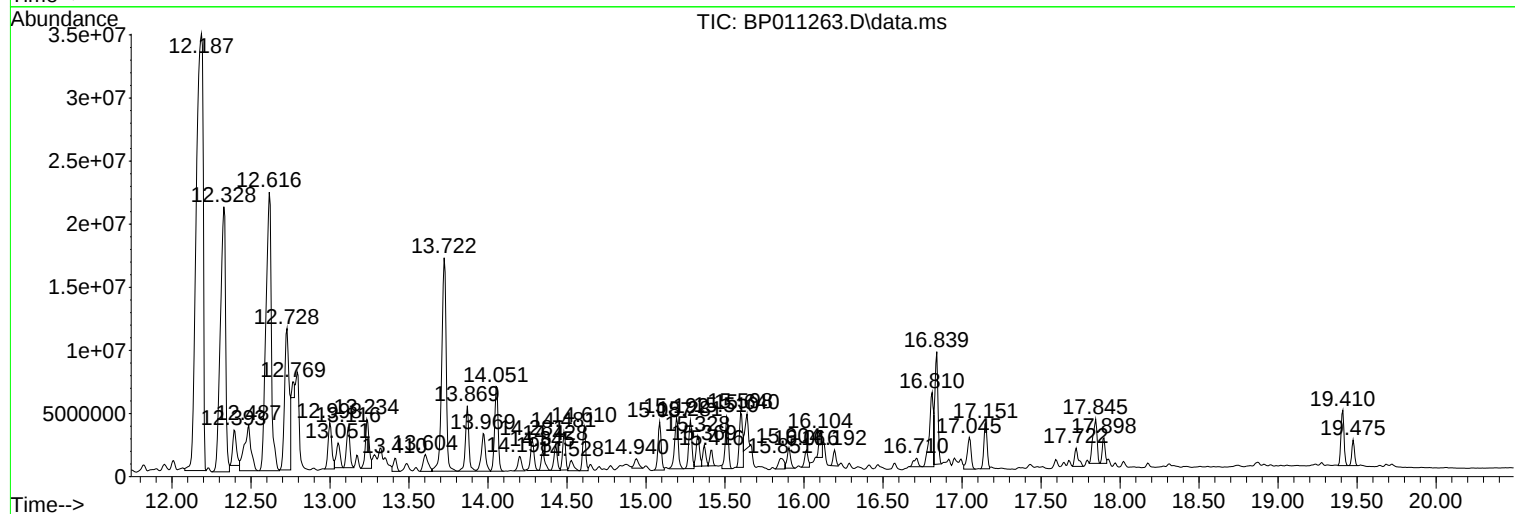
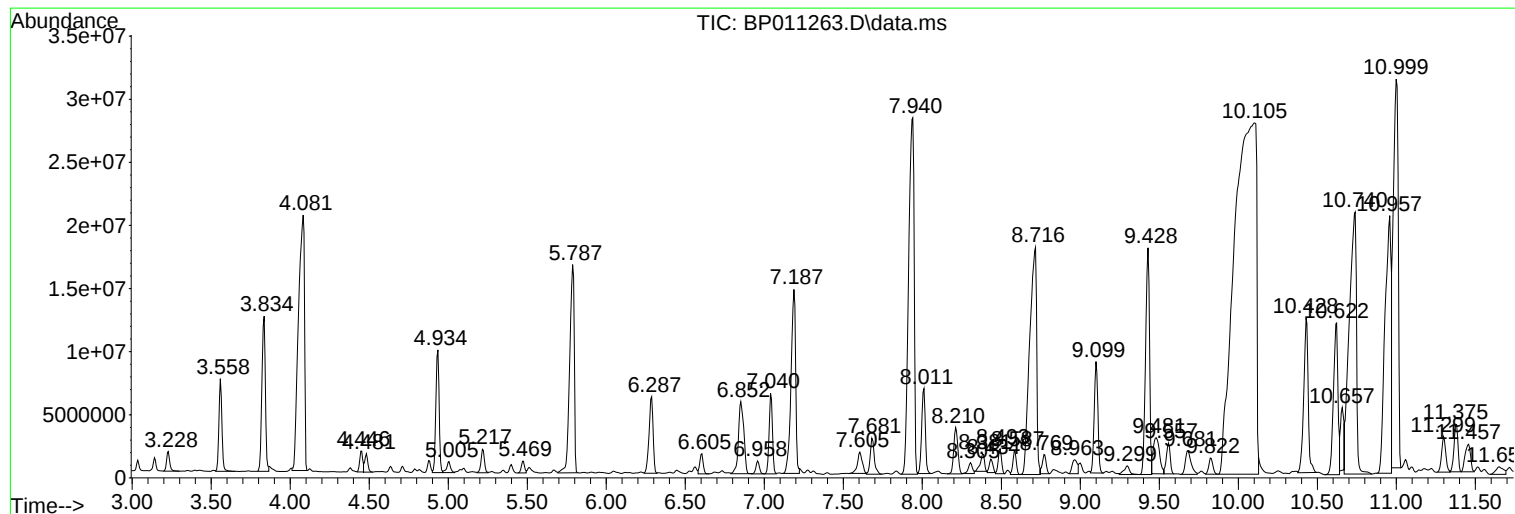
Sum of corrected areas: 1440869691

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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

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



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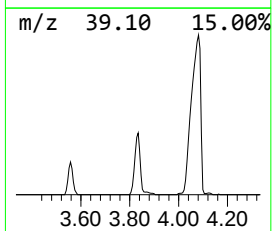
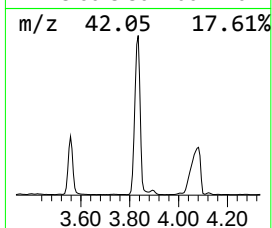
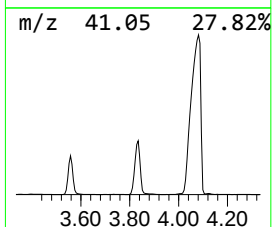
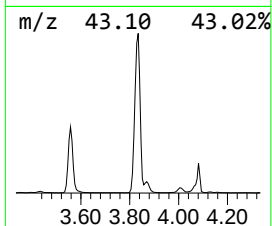
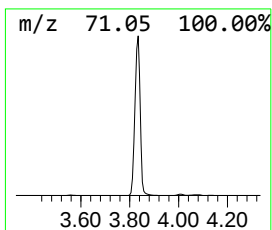
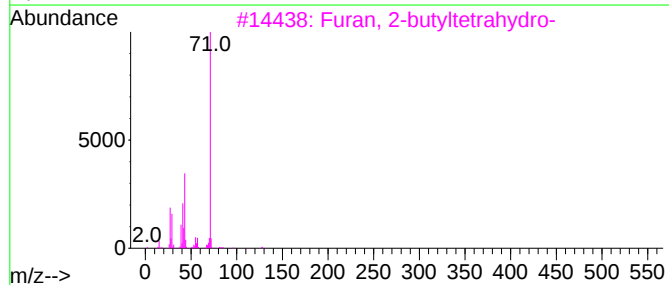
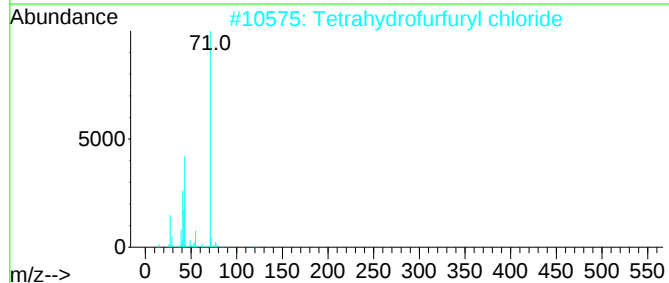
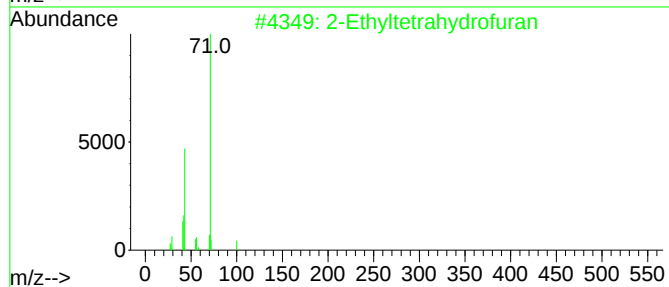
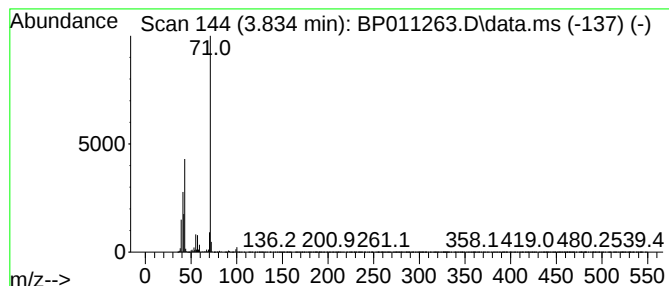
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 2 2-Ethyltetrahydrofuran Concentration Rank 7

| R.T.  | EstConc      | Area     | Relative to ISTD       | R.T.  |
|-------|--------------|----------|------------------------|-------|
| 3.834 | 105.80 ng/ul | 18979000 | 1,4-Dichlorobenzene-d4 | 7.605 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#         | Qual |
|------|----|---|------------------------------|-----|---------|--------------|------|
| 1    |    |   | 2-Ethyltetrahydrofuran       | 100 | C6H12O  | 1010431-64-0 | 78   |
| 2    |    |   | Tetrahydrofurfuryl chloride  | 120 | C5H9ClO | 003003-84-7  | 64   |
| 3    |    |   | Furan, 2-butyltetrahydro-    | 128 | C8H16O  | 001004-29-1  | 56   |
| 4    |    |   | 2-Furanmethanol, tetrahydro- | 102 | C5H10O2 | 000097-99-4  | 56   |
| 5    |    |   | Oxirane, 3-hydroxypropyl-    | 102 | C5H10O2 | 021915-56-0  | 56   |



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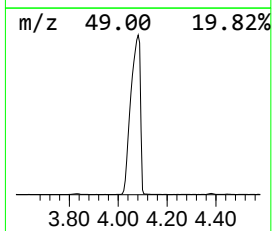
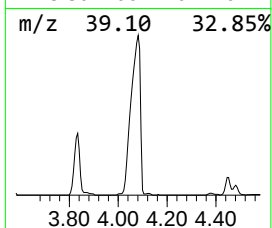
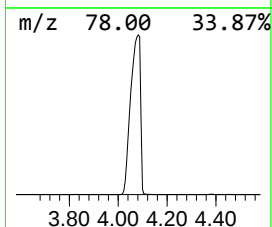
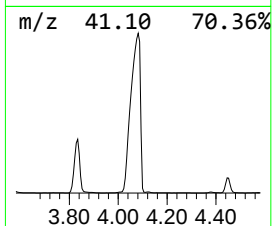
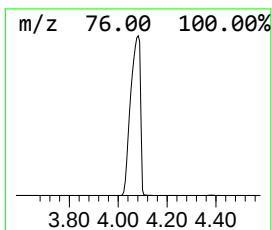
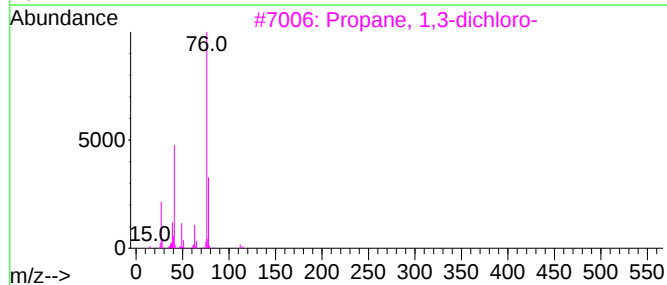
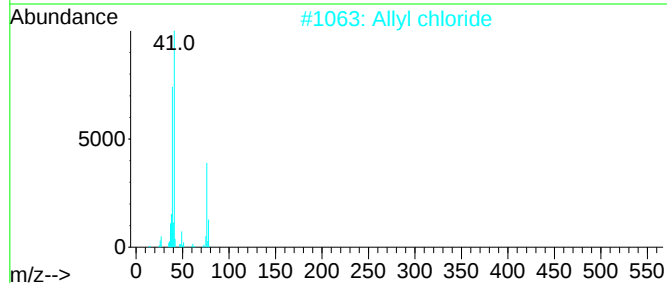
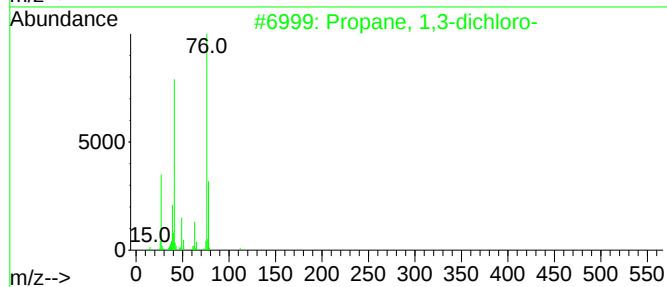
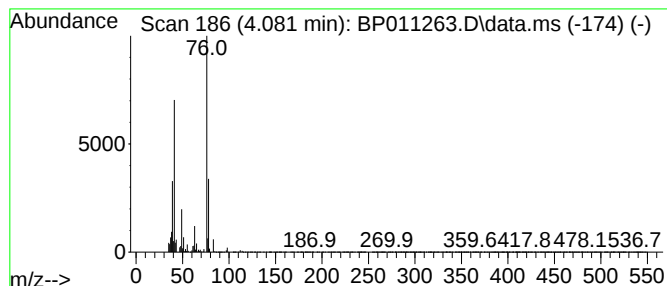
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 Propane, 1,3-dichloro- Concentration Rank 3

| R.T.  | EstConc      | Area     | Relative to ISTD       | R.T.  |
|-------|--------------|----------|------------------------|-------|
| 4.081 | 300.78 ng/ul | 53954600 | 1,4-Dichlorobenzene-d4 | 7.605 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Propane, 1,3-dichloro- | 112 | C3H6Cl2 | 000142-28-9 | 91   |
| 2    |    |   | Allyl chloride         | 76  | C3H5Cl  | 000107-05-1 | 72   |
| 3    |    |   | Propane, 1,3-dichloro- | 112 | C3H6Cl2 | 000142-28-9 | 64   |
| 4    |    |   | Propane, 1,3-dichloro- | 112 | C3H6Cl2 | 000142-28-9 | 64   |
| 5    |    |   | 1-Propene, 1-chloro-   | 76  | C3H5Cl  | 000590-21-6 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

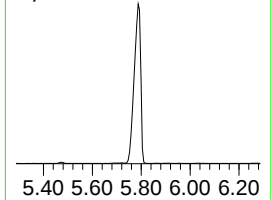
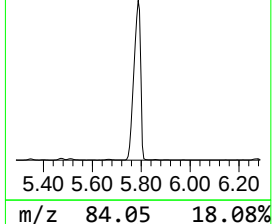
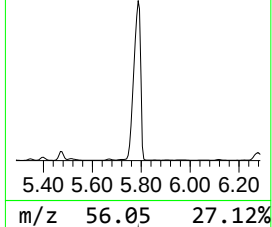
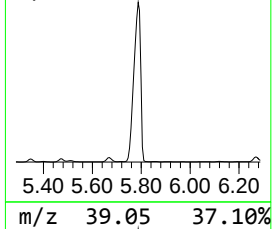
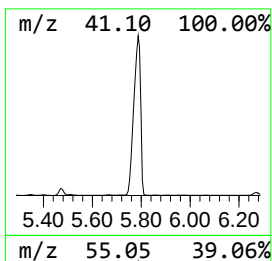
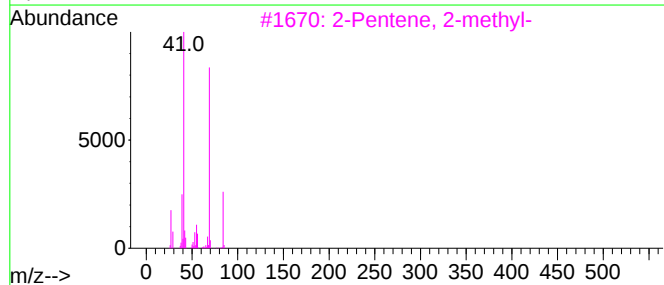
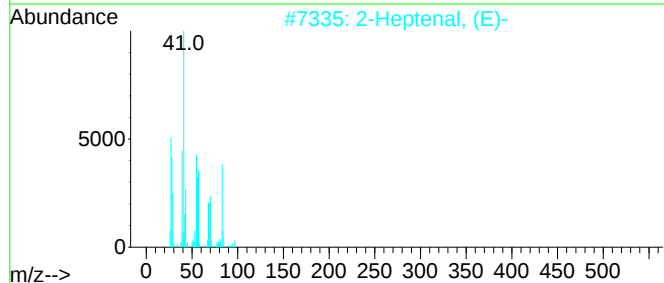
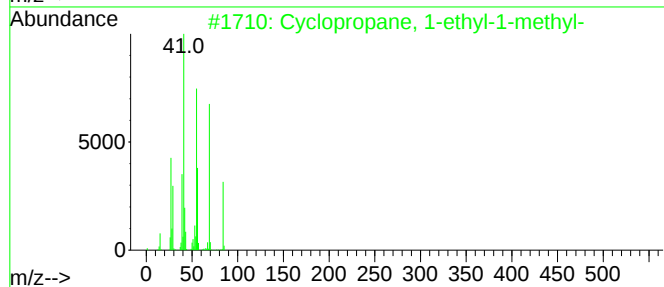
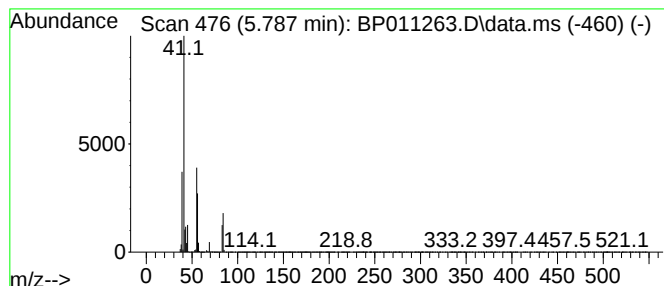
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Quant Title : SVOA CALIBRATION

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

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Peak Number 10 (DEL) Alkane: Cyclic5.787 Concentration Rank 5

| R.T.  | EstConc      | Area     | Relative to ISTD       | R.T.  |
|-------|--------------|----------|------------------------|-------|
| 5.787 | 184.41 ng/ul | 33079400 | 1,4-Dichlorobenzene-d4 | 7.605 |

| Hit# | of 5 | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|------|---------------------------------|-----|---------|-------------|------|
| 1    |      | Cyclopropane, 1-ethyl-1-methyl- | 84  | C6H12   | 053778-43-1 | 47   |
| 2    |      | 2-Heptenal, (E)-                | 112 | C7H12O  | 018829-55-5 | 28   |
| 3    |      | 2-Pentene, 2-methyl-            | 84  | C6H12   | 000625-27-4 | 23   |
| 4    |      | 1-Hexene                        | 84  | C6H12   | 000592-41-6 | 12   |
| 5    |      | 1-Propene, 3-(ethenyloxy)-      | 84  | C5H8O   | 003917-15-5 | 9    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

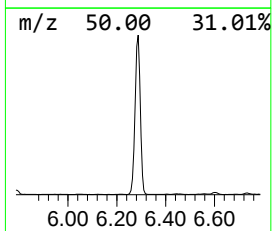
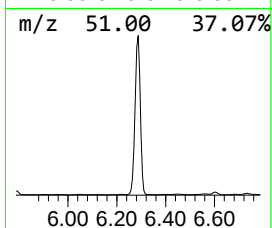
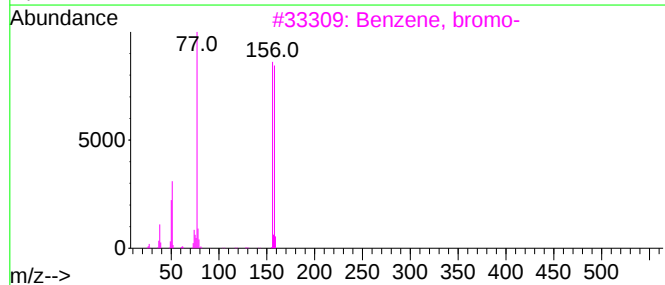
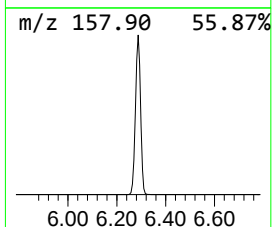
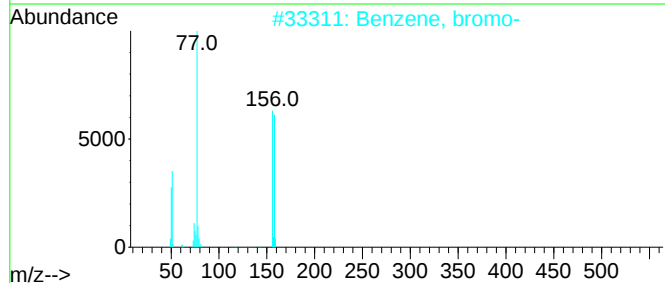
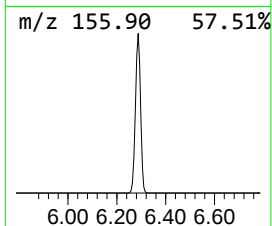
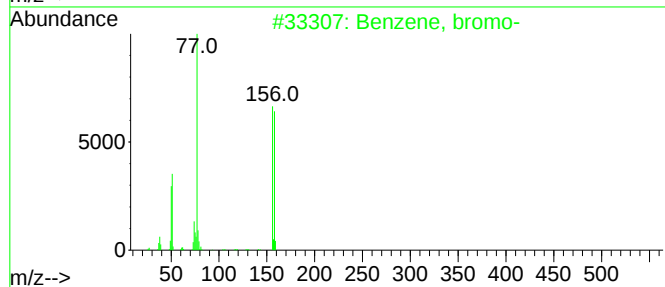
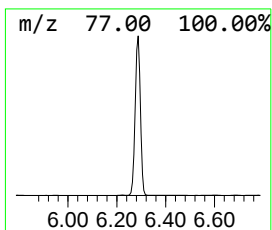
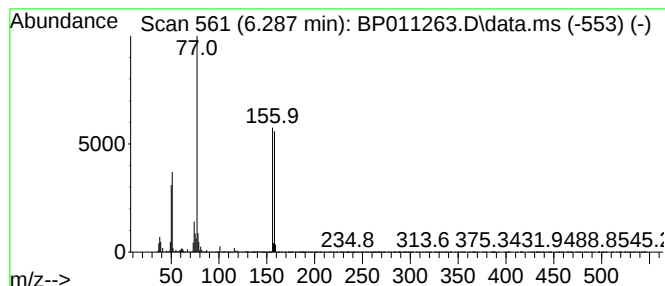
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Quant Title : SVOA CALIBRATION

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

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Peak Number 11 Benzene, bromo- Concentration Rank 11

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 6.287 | 59.53 ng/ul | 10677800 | 1,4-Dichlorobenzene-d4 | 7.605 |

| Hit# | of | 5 | Tentative ID    | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, bromo- | 156 | C6H5Br  | 000108-86-1 | 97   |
| 2    |    |   | Benzene, bromo- | 156 | C6H5Br  | 000108-86-1 | 96   |
| 3    |    |   | Benzene, bromo- | 156 | C6H5Br  | 000108-86-1 | 94   |
| 4    |    |   | Benzene, bromo- | 156 | C6H5Br  | 000108-86-1 | 90   |
| 5    |    |   | Benzene, bromo- | 156 | C6H5Br  | 000108-86-1 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

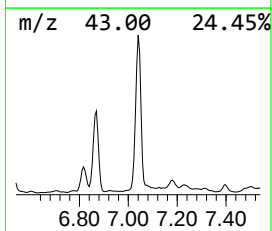
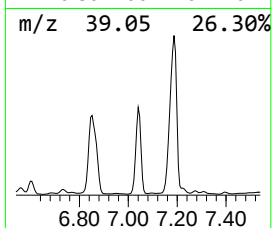
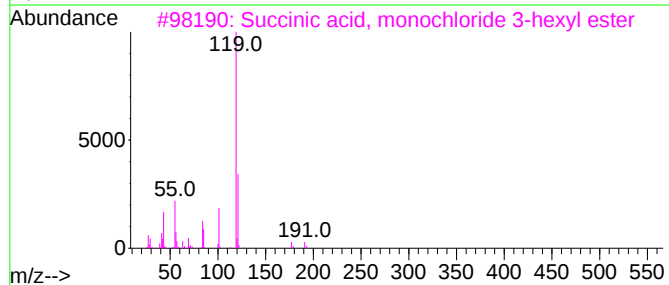
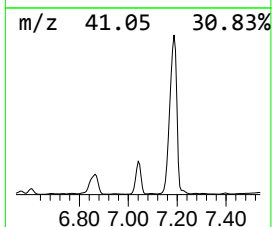
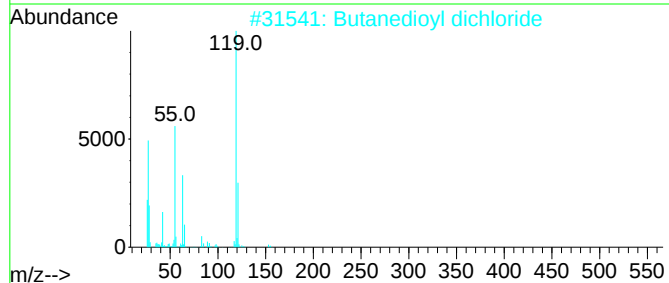
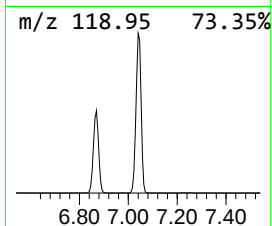
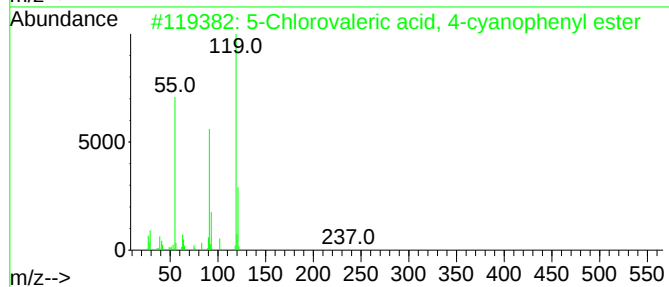
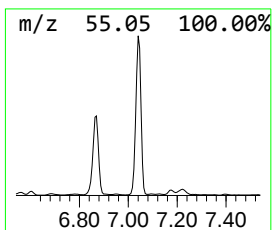
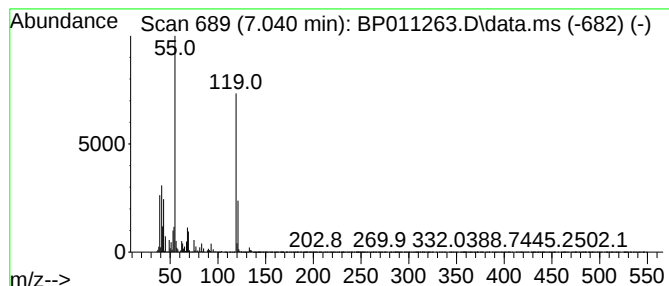
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 13 unknown-01 Concentration Rank 18

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 7.040 | 51.95 ng/ul | 9318060 | 1,4-Dichlorobenzene-d4 | 7.605 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | 5-Chlorovaleric acid, 4-cyanophe... | 237 | C12H12ClNO2 | 1000307-97-8 | 40   |
| 2    |    |   | Butanedioyl dichloride              | 154 | C4H4Cl2O2   | 000543-20-4  | 33   |
| 3    |    |   | Succinic acid, monochloride 3-he... | 220 | C10H17ClO3  | 1010349-46-1 | 33   |
| 4    |    |   | 2,5-Pyrrolidinedione, 1-[(2-meth... | 233 | C12H11NO4   | 110920-14-4  | 32   |
| 5    |    |   | Butanedioyl dichloride              | 154 | C4H4Cl2O2   | 000543-20-4  | 28   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

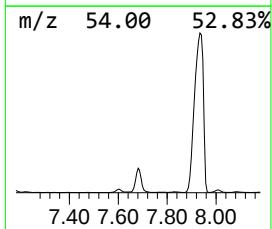
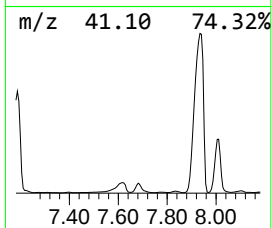
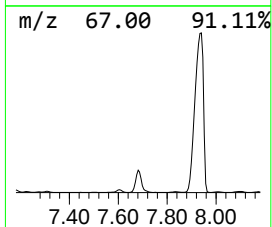
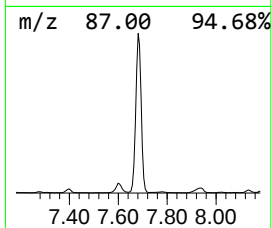
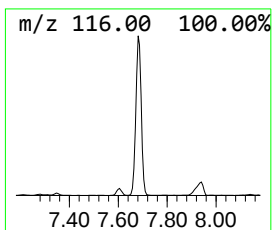
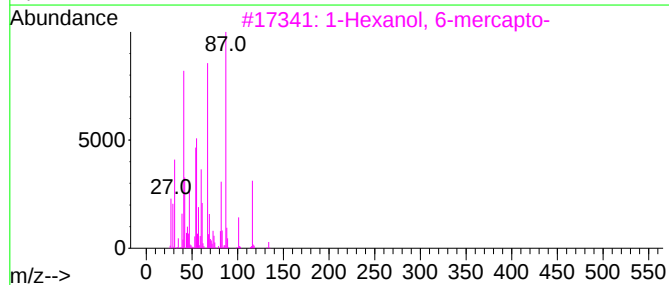
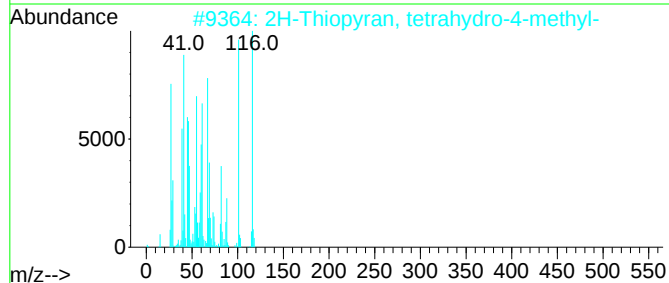
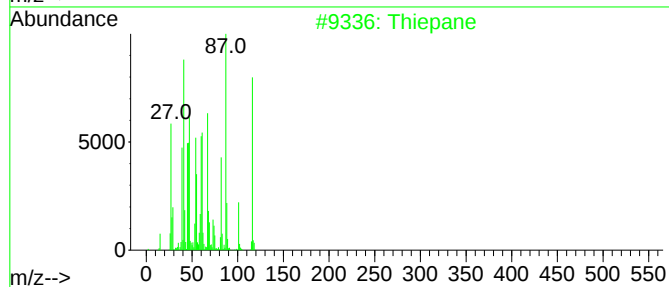
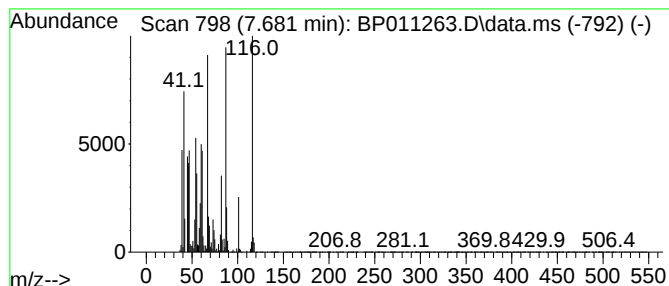
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Thiepane Concentration Rank 30

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 7.681 | 27.65 ng/ul | 4959920 | 1,4-Dichlorobenzene-d4 | 7.605 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Thiepane                           | 116 | C6H12S  | 004753-80-4 | 68   |
| 2    |    |   | 2H-Thiopyran, tetrahydro-4-methyl- | 116 | C6H12S  | 005161-17-1 | 64   |
| 3    |    |   | 1-Hexanol, 6-mercapto-             | 134 | C6H14OS | 001633-78-9 | 47   |
| 4    |    |   | 1-Propene, 1-(ethylthio)-2-methyl- | 116 | C6H12S  | 027482-14-0 | 38   |
| 5    |    |   | 2H-Thiopyran-3(4H)-one, dihydro-   | 116 | C5H8OS  | 019090-03-0 | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

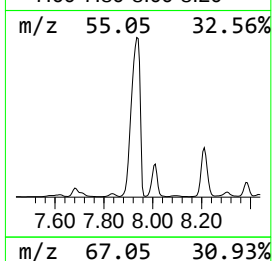
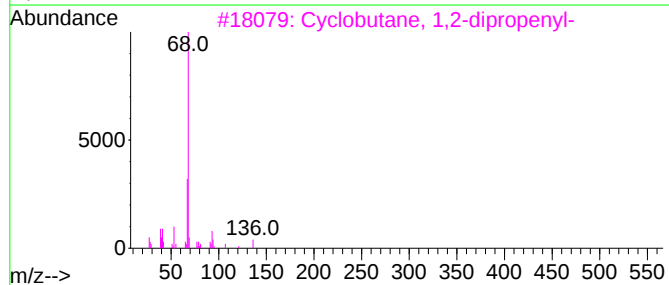
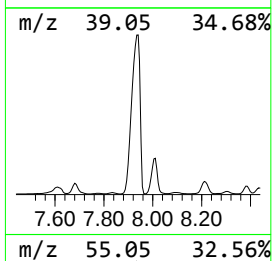
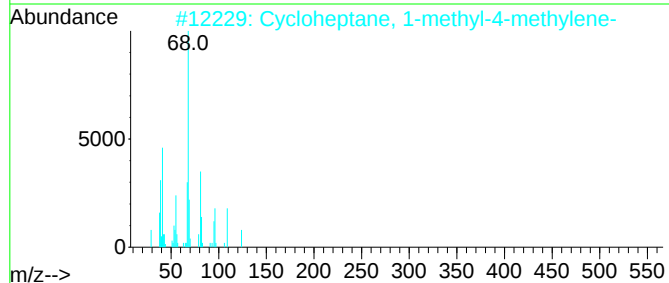
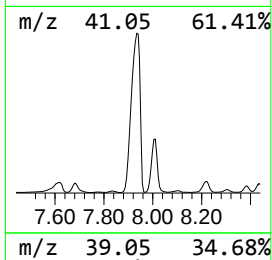
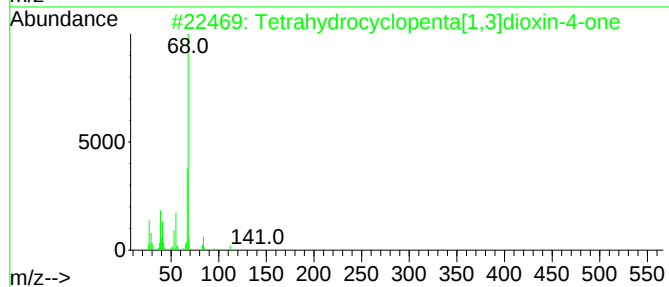
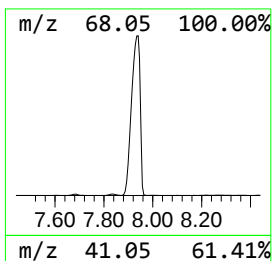
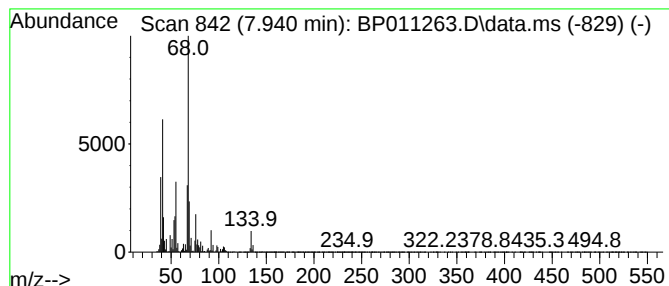
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Quant Title : SVOA CALIBRATION

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

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Peak Number 15 unknown-02 Concentration Rank 1

| R.T.  | EstConc      | Area     | Relative to ISTD       | R.T.  |
|-------|--------------|----------|------------------------|-------|
| 7.940 | 383.50 ng/ul | 68792100 | 1,4-Dichlorobenzene-d4 | 7.605 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Tetrahydrocyclopenta[1,3]dioxin-... | 142 | C7H10O3 | 124898-85-7  | 33   |
| 2    |    |   | Cycloheptane, 1-methyl-4-methylene- | 124 | C9H16   | 023799-25-9  | 33   |
| 3    |    |   | Cyclobutane, 1,2-dipropenyl-        | 136 | C10H16  | 022769-00-2  | 33   |
| 4    |    |   | Cyclopentanemethanol                | 100 | C6H12O  | 003637-61-4  | 33   |
| 5    |    |   | 2-Methyl-6-methylene-7-octen-4-ol   | 154 | C10H18O | 1000424-70-9 | 33   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

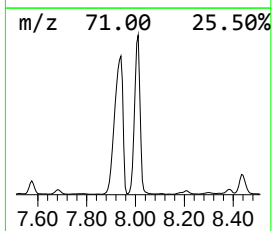
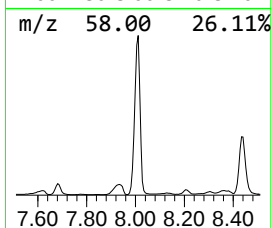
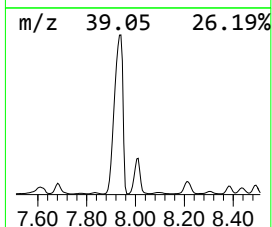
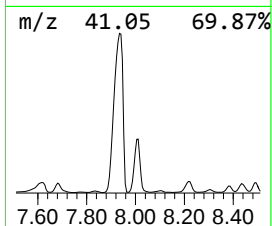
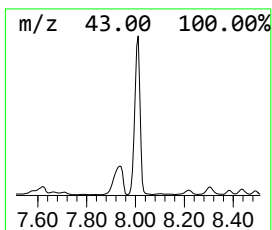
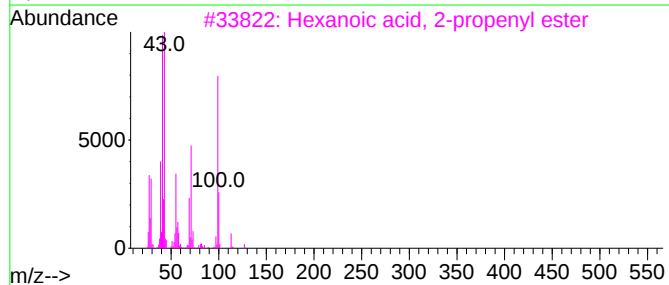
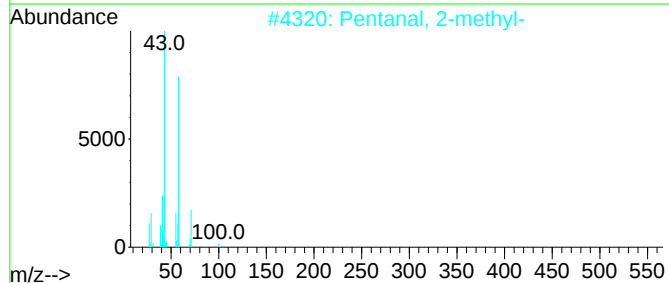
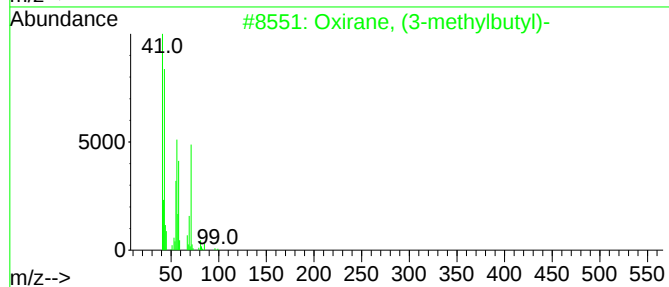
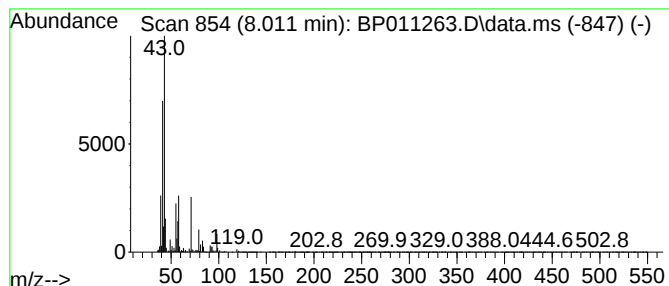
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Quant Title : SVOA CALIBRATION

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

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Peak Number 16 unknown-03 Concentration Rank 10

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 8.011 | 59.72 ng/ul | 10712900 | 1,4-Dichlorobenzene-d4 | 7.605 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Oxirane, (3-methylbutyl)-           | 114 | C7H14O  | 053229-41-7 | 36   |
| 2    |    |   | Pentanal, 2-methyl-                 | 100 | C6H12O  | 000123-15-9 | 25   |
| 3    |    |   | Hexanoic acid, 2-propenyl ester     | 156 | C9H16O2 | 000123-68-2 | 25   |
| 4    |    |   | 2-Hexanone, 5-methyl-               | 114 | C7H14O  | 000110-12-3 | 23   |
| 5    |    |   | Propanoic acid, 2-methyl-, 1-met... | 130 | C7H14O2 | 000617-50-5 | 17   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

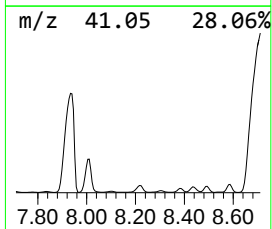
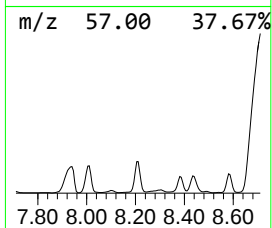
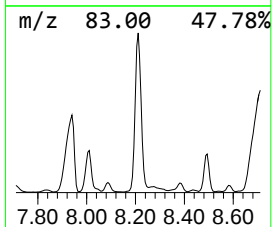
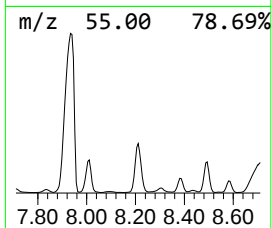
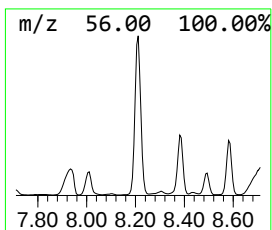
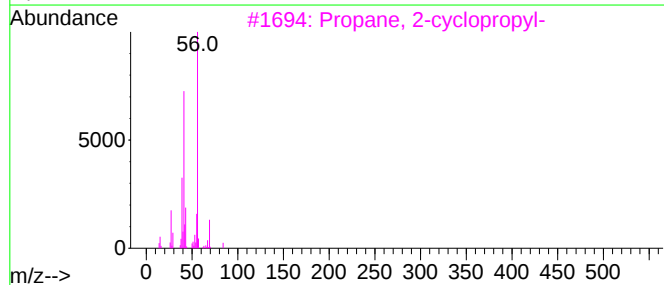
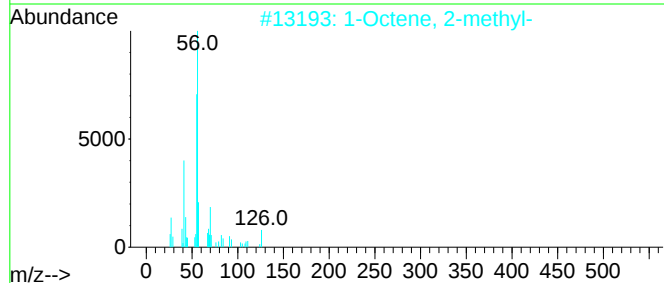
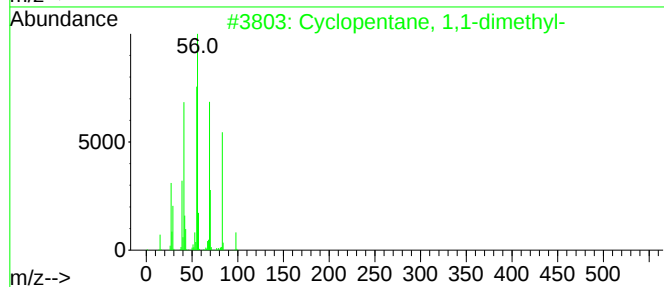
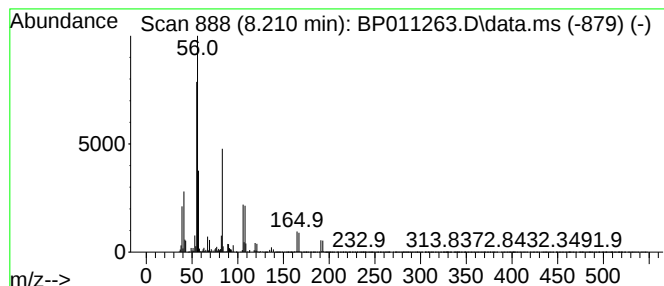
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 17 (DEL) Alkane: Cyclic8.210 Concentration Rank 22

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 8.210 | 37.10 ng/ul | 6654960 | 1,4-Dichlorobenzene-d4 | 7.605 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Cyclopentane, 1,1-dimethyl-         | 98  | C7H14    | 001638-26-2 | 33   |
| 2    |    |   | 1-Octene, 2-methyl-                 | 126 | C9H18    | 004588-18-5 | 17   |
| 3    |    |   | Propane, 2-cyclopropyl-             | 84  | C6H12    | 003638-35-5 | 16   |
| 4    |    |   | Cyclobutyl bromide                  | 134 | C4H7Br   | 004399-47-7 | 14   |
| 5    |    |   | Cyclopentanecarboxylic acid, 2-a... | 129 | C6H11NO2 | 040482-05-1 | 12   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

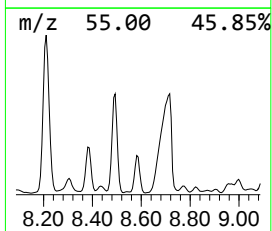
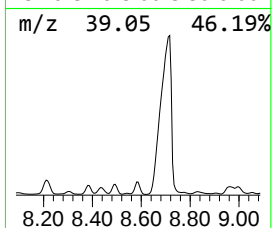
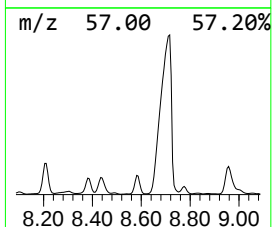
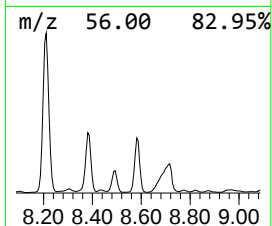
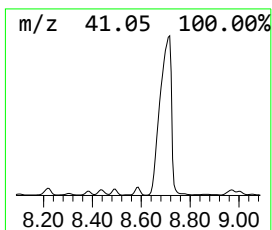
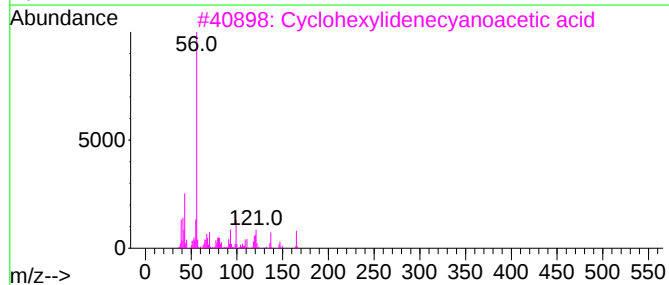
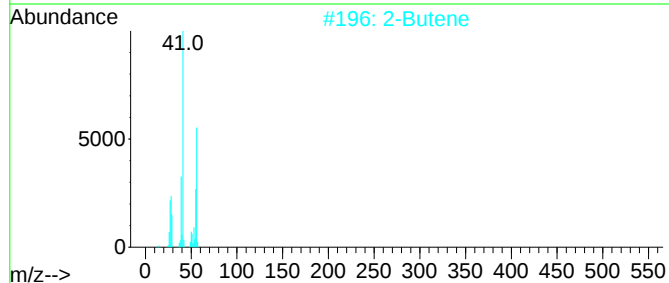
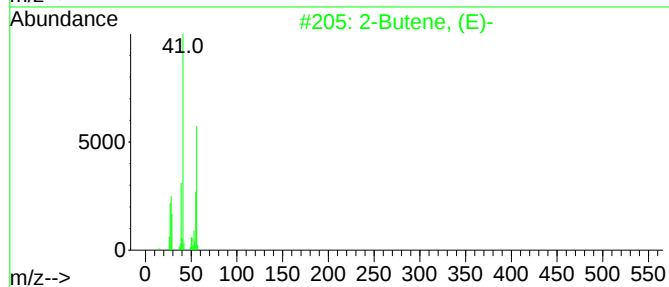
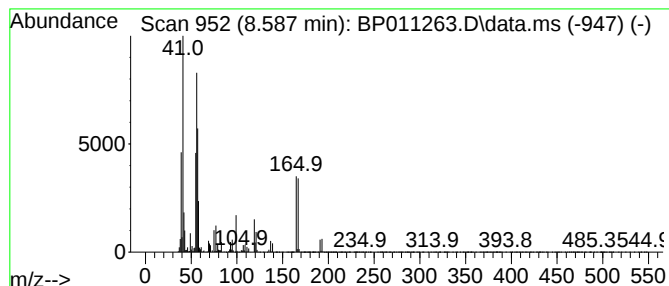
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 42

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 8.587 | 15.35 ng/ul | 2753700 | 1,4-Dichlorobenzene-d4 | 7.605 |

| Hit# | of                                  | 5 | Tentative ID | MW  | MolForm  | CAS#         | Qual |
|------|-------------------------------------|---|--------------|-----|----------|--------------|------|
| 1    | 2-Butene, (E)-                      |   |              | 56  | C4H8     | 000624-64-6  | 37   |
| 2    | 2-Butene                            |   |              | 56  | C4H8     | 000107-01-7  | 37   |
| 3    | Cyclohexylidenecyanoacetic acid     |   |              | 165 | C9H11NO2 | 037107-50-9  | 32   |
| 4    | 1-Heptene                           |   |              | 98  | C7H14    | 000592-76-7  | 25   |
| 5    | 3-[N-Aziridylmethyl]-2-norbornanone |   |              | 165 | C10H15NO | 1000257-03-5 | 25   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

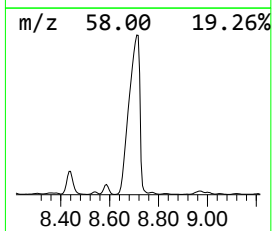
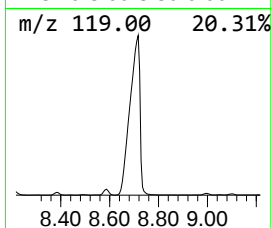
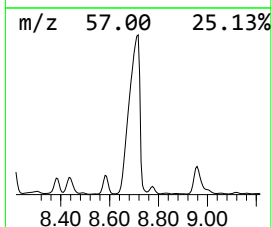
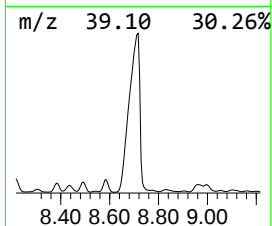
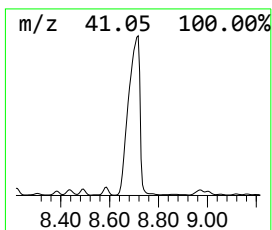
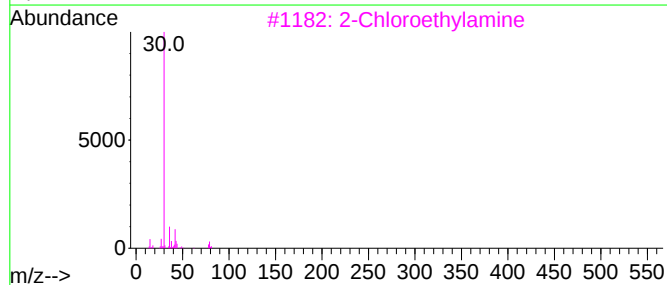
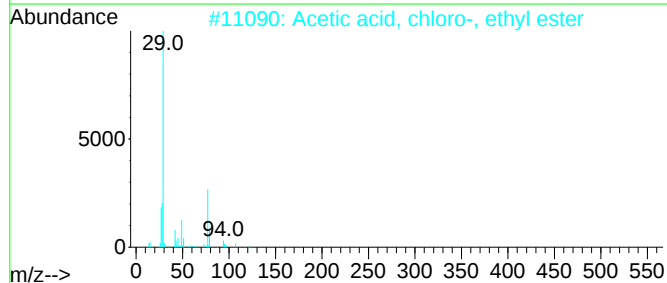
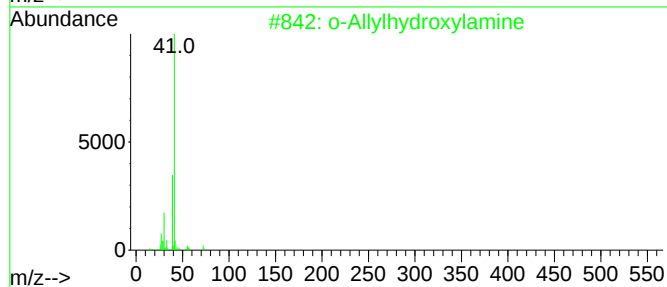
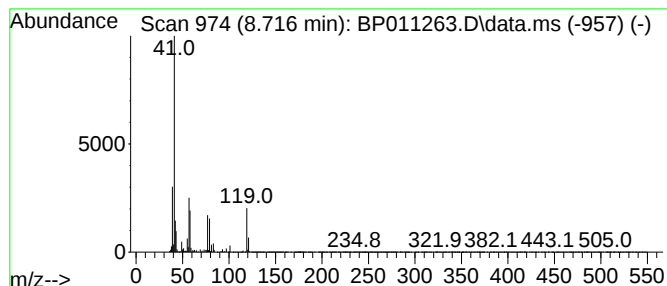
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 unknown-04 Concentration Rank 2

| R.T.  | EstConc      | Area     | Relative to ISTD       | R.T.  |
|-------|--------------|----------|------------------------|-------|
| 8.716 | 306.39 ng/ul | 54959500 | 1,4-Dichlorobenzene-d4 | 7.605 |

| Hit# | of 5 | Tentative ID                      | MW  | MolForm  | CAS#         | Qual |
|------|------|-----------------------------------|-----|----------|--------------|------|
| 1    |      | o-Allylhydroxylamine              | 73  | C3H7NO   | 006542-54-7  | 7    |
| 2    |      | Acetic acid, chloro-, ethyl ester | 122 | C4H7ClO2 | 000105-39-5  | 7    |
| 3    |      | 2-Chloroethylamine                | 79  | C2H6ClN  | 000689-98-5  | 4    |
| 4    |      | 1-Allyl-cyclohexane-1,2-diol      | 156 | C9H16O2  | 1000143-28-3 | 4    |
| 5    |      | Chloroacetic acid allyl ester     | 134 | C5H7ClO2 | 002916-14-5  | 4    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

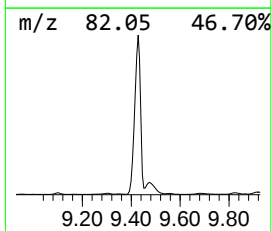
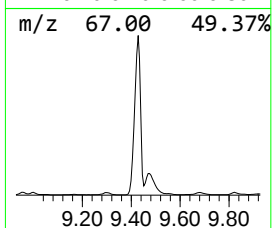
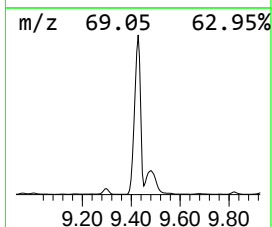
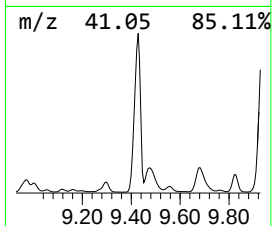
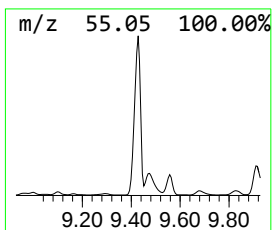
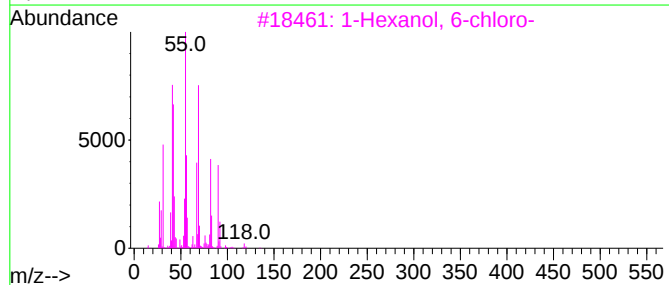
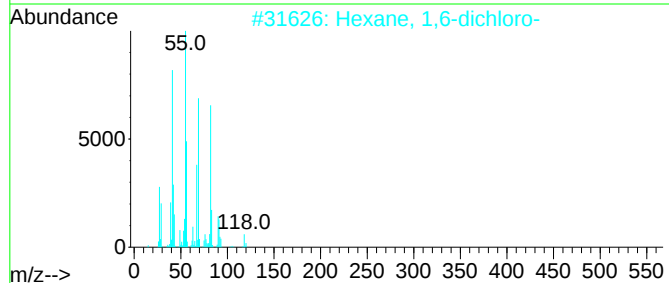
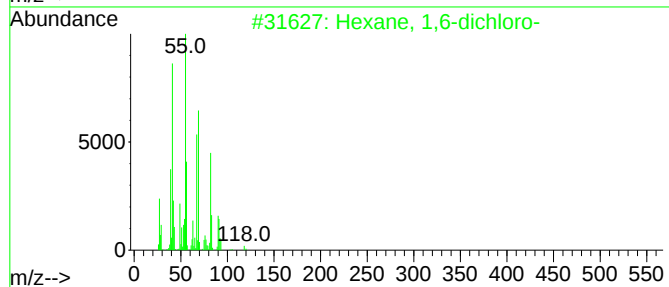
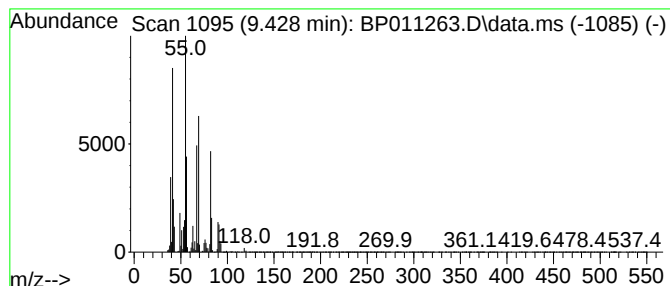
Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 24 Hexane, 1,6-dichloro- Concentration Rank 32

| R.T.      | EstConc                        | Area     | Relative to ISTD | R.T.        |      |
|-----------|--------------------------------|----------|------------------|-------------|------|
| 9.428     | 26.10 ng/ul                    | 31375100 | Naphthalene-d8   | 10.410      |      |
| Hit# of 5 | Tentative ID                   | MW       | MolForm          | CAS#        | Qual |
| 1         | Hexane, 1,6-dichloro-          | 154      | C6H12Cl2         | 002163-00-0 | 91   |
| 2         | Hexane, 1,6-dichloro-          | 154      | C6H12Cl2         | 002163-00-0 | 80   |
| 3         | 1-Hexanol, 6-chloro-           | 136      | C6H13ClO         | 002009-83-8 | 45   |
| 4         | 1-Hexene, 6-chloro-            | 118      | C6H11Cl          | 000928-89-2 | 45   |
| 5         | Cyclohexane, (1-methylpropyl)- | 140      | C10H20           | 007058-01-7 | 40   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

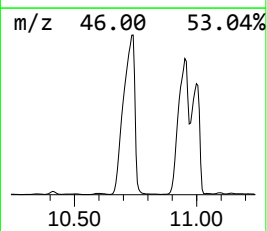
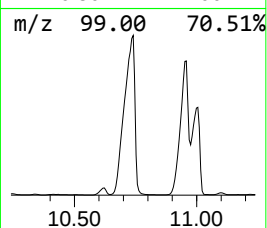
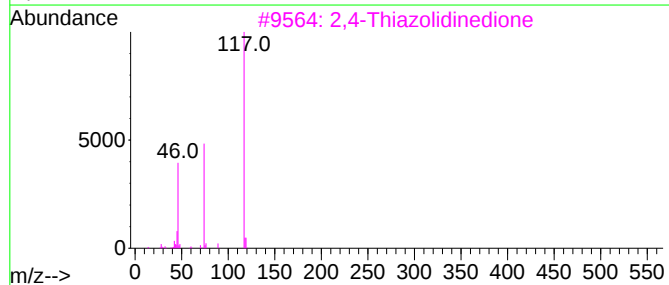
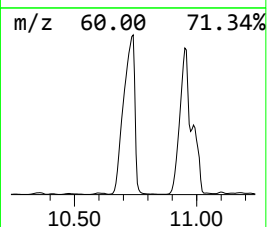
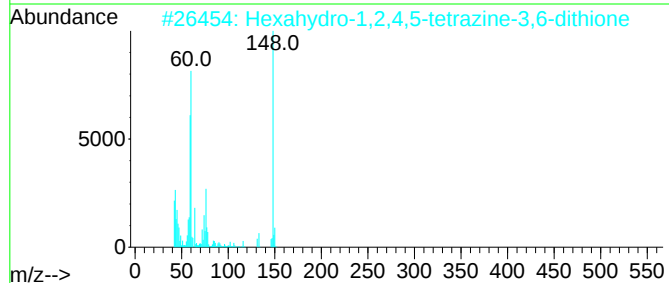
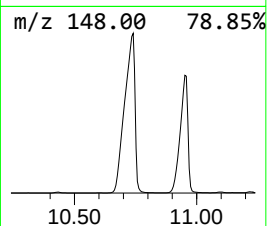
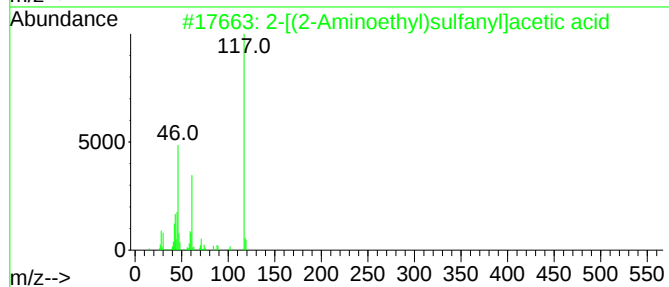
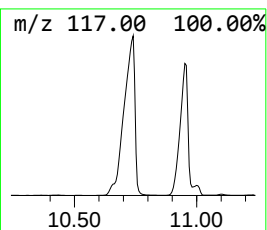
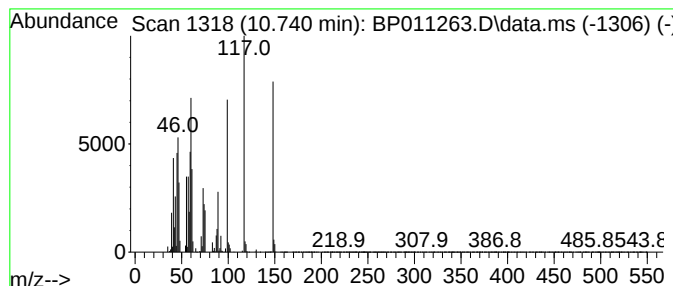
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TIC Integration Parameters: LSCINT.P

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Peak Number 28 unknown-05 Concentration Rank 15

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 10.740 | 54.62 ng/ul | 65654800 | Naphthalene-d8   | 10.410 |

| Hit# | of 5                                | Tentative ID | MW       | MolForm     | CAS# | Qual |
|------|-------------------------------------|--------------|----------|-------------|------|------|
| 1    | 2-[(2-Aminoethyl)sulfanyl]acetic... | 135          | C4H9NO2S | 003335-52-2 | 27   |      |
| 2    | Hexahydro-1,2,4,5-tetrazine-3,6-... | 148          | C2H4N4S2 | 036239-33-5 | 27   |      |
| 3    | 2,4-Thiazolidinedione               | 117          | C3H3NO2S | 002295-31-0 | 14   |      |
| 4    | Ethylvinyl sulfide                  | 88           | C4H8S    | 000627-50-9 | 11   |      |
| 5    | Ethylvinyl sulfide                  | 88           | C4H8S    | 000627-50-9 | 11   |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

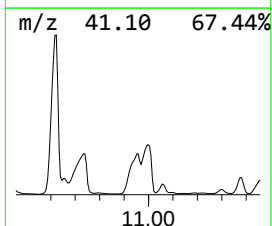
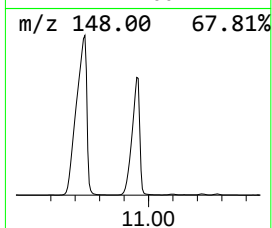
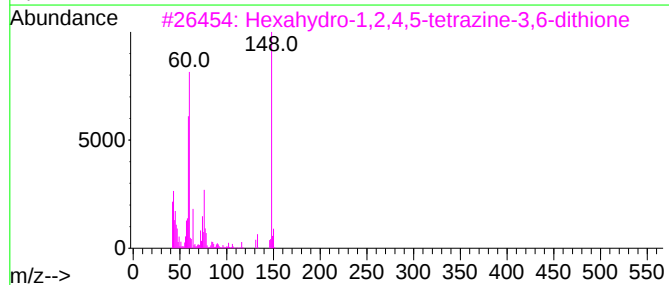
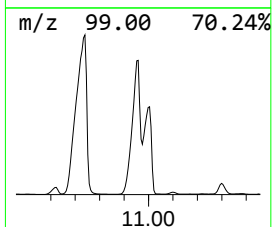
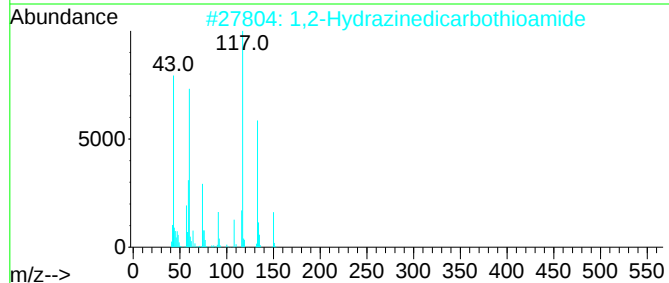
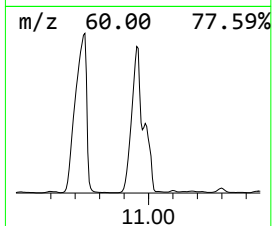
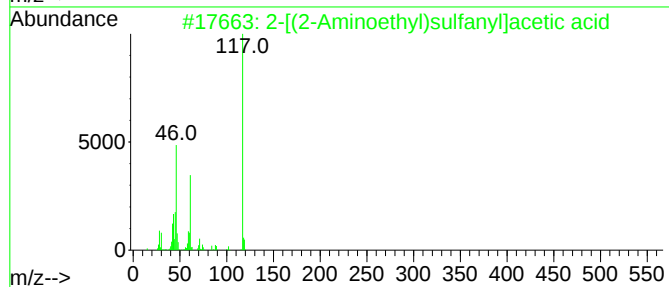
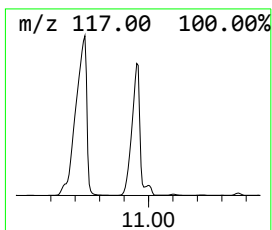
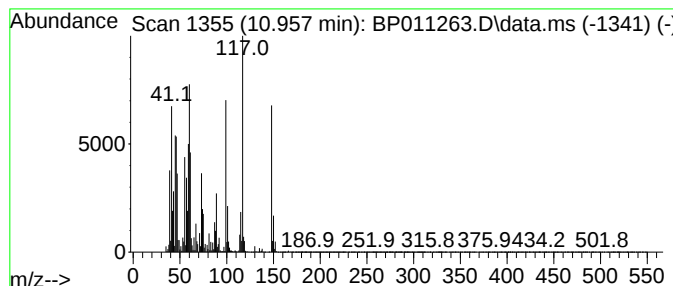
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 29 unknown-06 Concentration Rank 19

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 10.957 | 46.44 ng/ul | 55825300 | Naphthalene-d8   | 10.410 |

| Hit# | of 5              | Tentative ID                      | MW  | MolForm  | CAS#        | Qual |
|------|-------------------|-----------------------------------|-----|----------|-------------|------|
| 1    | 2                 | [(2-Aminoethyl)sulfanyl]acetic... | 135 | C4H9NO2S | 003335-52-2 | 35   |
| 2    | 1,2               | Hydrazinedicarbothioamide         | 150 | C2H6N4S2 | 000142-46-1 | 35   |
| 3    | Hexahydro-1,2,4,5 | tetrazine-3,6-...                 | 148 | C2H4N4S2 | 036239-33-5 | 30   |
| 4    | Thiazole, 4,5     | dihydro-2-methyl-                 | 101 | C4H7NS   | 002346-00-1 | 22   |
| 5    | 2-Thio-2,4        | oxazolidinedione                  | 117 | C3H3NO2S | 002346-24-9 | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

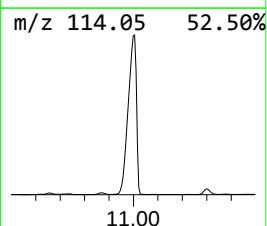
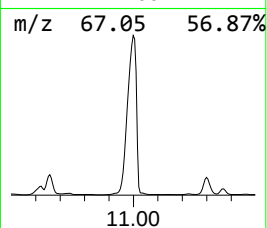
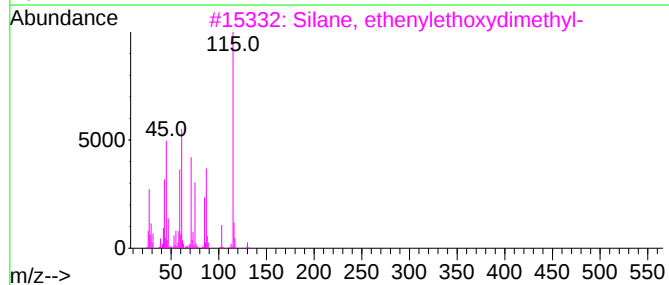
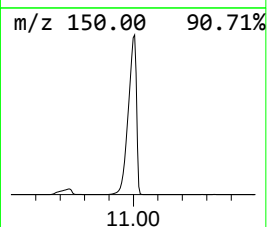
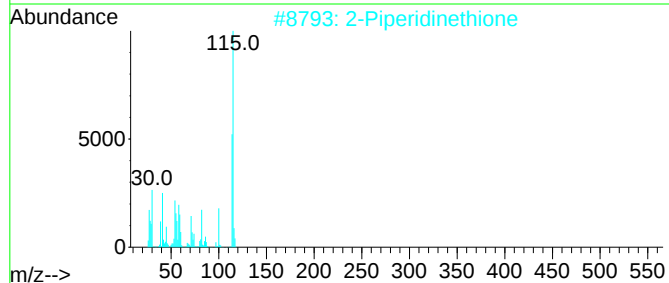
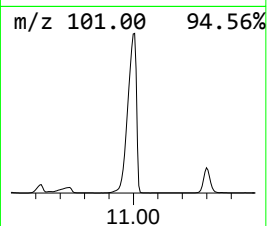
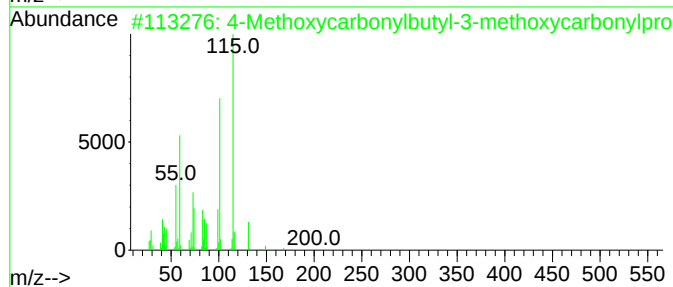
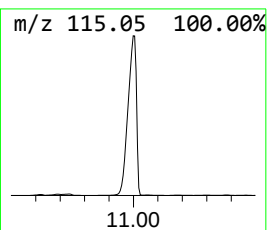
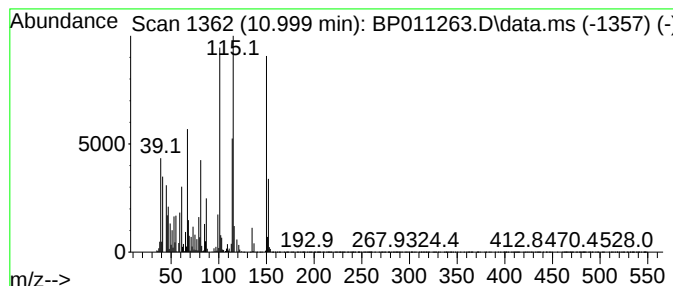
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Quant Title : SVOA CALIBRATION

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

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Peak Number 30 unknown-07 Concentration Rank 14

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 10.999 | 55.09 ng/ul | 66222500 | Naphthalene-d8   | 10.410 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 4-Methoxycarbonylbutyl-3-methoxy... | 232 | C11H20O5 | 017361-53-4 | 16   |
| 2    |    |   | 2-Piperidinethione                  | 115 | C5H9NS   | 013070-01-4 | 14   |
| 3    |    |   | Silane, ethenylethoxydimethyl-      | 130 | C6H14OSi | 005356-83-2 | 14   |
| 4    |    |   | 1,3,4-Thiadiazol-2-amine            | 101 | C2H3N3S  | 004005-51-0 | 11   |
| 5    |    |   | 3H-1,2,4-Triazole-3-thione, 1,2-... | 101 | C2H3N3S  | 003179-31-5 | 11   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

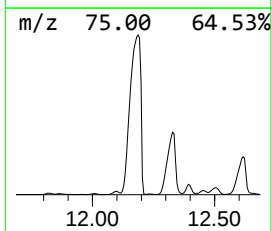
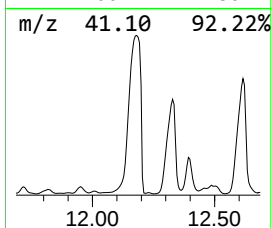
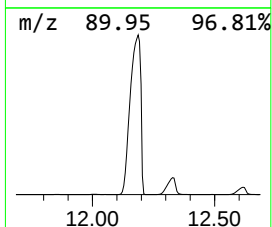
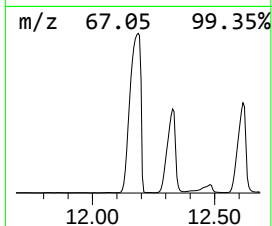
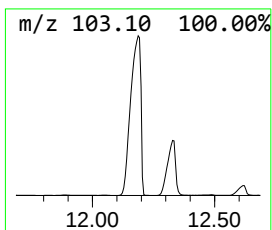
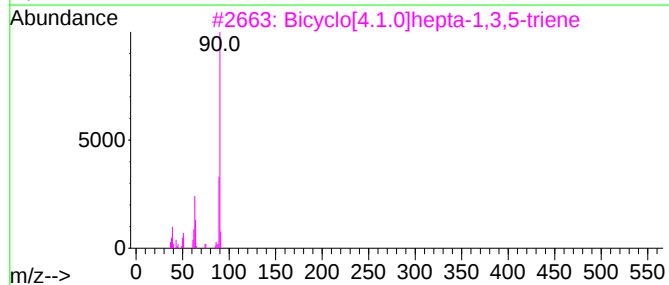
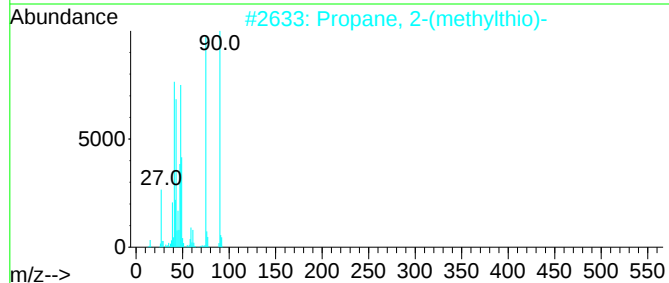
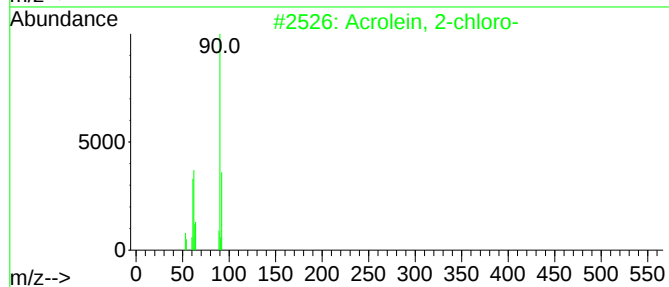
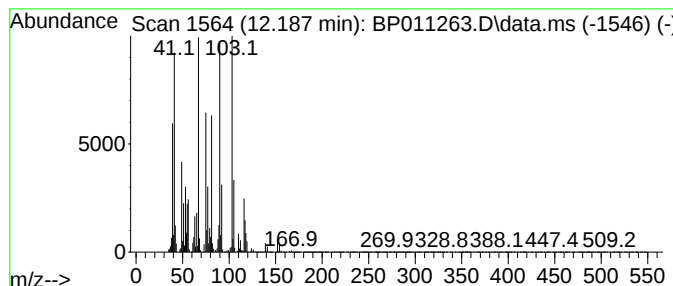
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 34 unknown-08 Concentration Rank 8

| R.T.   | EstConc     | Area      | Relative to ISTD | R.T.   |
|--------|-------------|-----------|------------------|--------|
| 12.187 | 88.24 ng/ul | 106062000 | Naphthalene-d8   | 10.410 |

| Hit# | of 5 | Tentative ID                     | MW  | MolForm   | CAS#        | Qual |
|------|------|----------------------------------|-----|-----------|-------------|------|
| 1    |      | Acrolein, 2-chloro-              | 90  | C3H3ClO   | 000683-51-2 | 43   |
| 2    |      | Propane, 2-(methylthio)-         | 90  | C4H10S    | 001551-21-9 | 32   |
| 3    |      | Bicyclo[4.1.0]hepta-1,3,5-triene | 90  | C7H6      | 004646-69-9 | 30   |
| 4    |      | Thiourea, methyl-                | 90  | C2H6N2S   | 000598-52-7 | 27   |
| 5    |      | 2,2-Bis(chloromethyl)-1-propanol | 156 | C5H10Cl2O | 005355-54-4 | 25   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

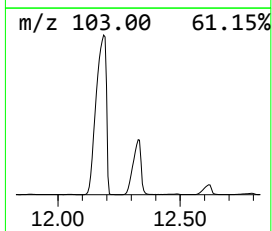
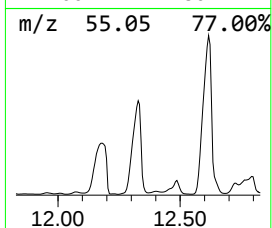
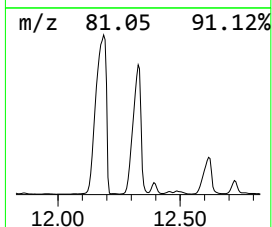
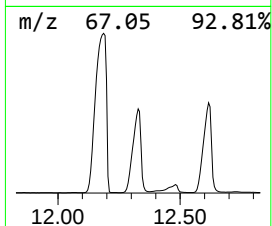
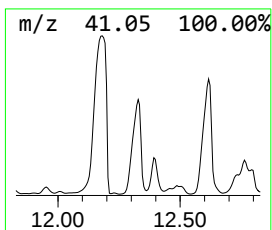
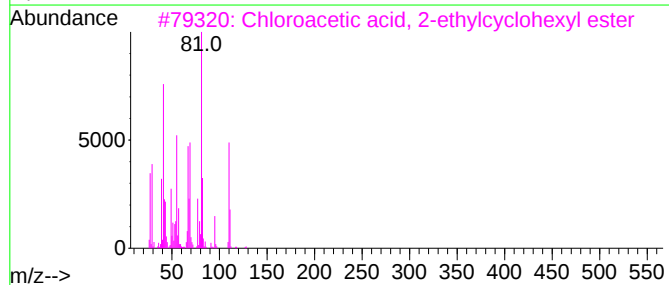
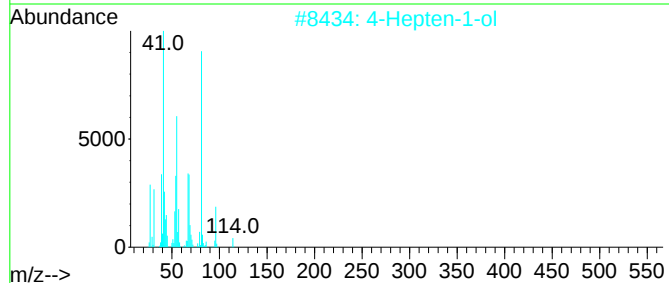
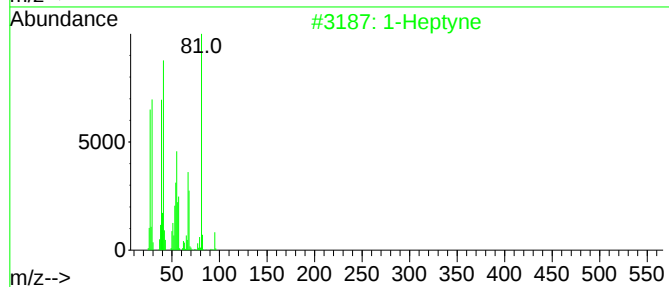
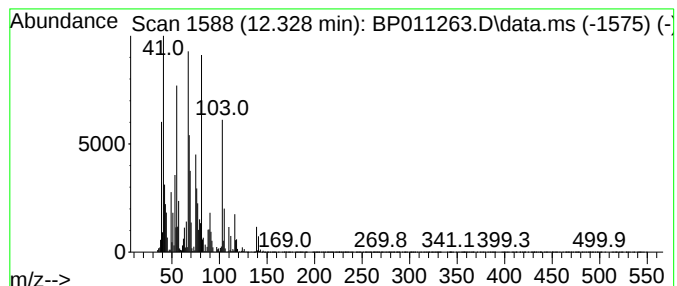
Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 35 unknown-09 Concentration Rank 21

| R.T.      | EstConc                                    | Area     | Relative to ISTD | R.T.         |      |
|-----------|--------------------------------------------|----------|------------------|--------------|------|
| 12.328    | 40.67 ng/ul                                | 48883100 | Naphthalene-d8   | 10.410       |      |
| Hit# of 5 | Tentative ID                               | MW       | MolForm          | CAS#         | Qual |
| 1         | 1-Heptyne                                  | 96       | C7H12            | 000628-71-7  | 27   |
| 2         | 4-Hepten-1-ol                              | 114      | C7H14O           | 020851-55-2  | 27   |
| 3         | Chloroacetic acid, 2-ethylcyclohexyl ester | 204      | C10H17ClO2       | 1000279-89-1 | 25   |
| 4         | 3-Octyne                                   | 110      | C8H14            | 015232-76-5  | 22   |
| 5         | 5-Hexenal, 4-methylene-                    | 110      | C7H10O           | 017844-21-2  | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

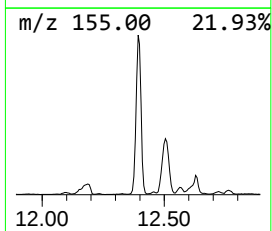
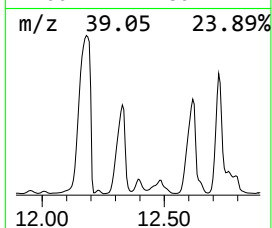
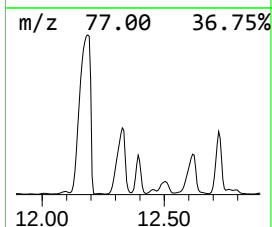
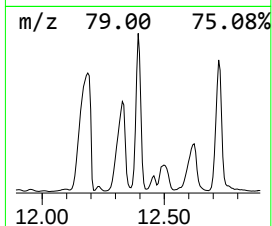
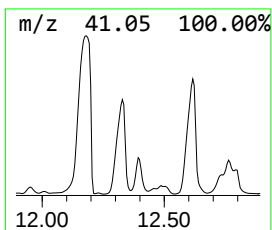
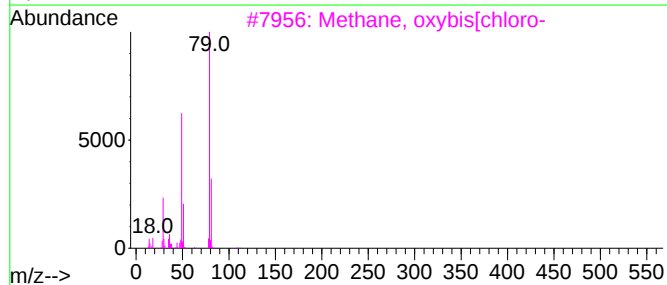
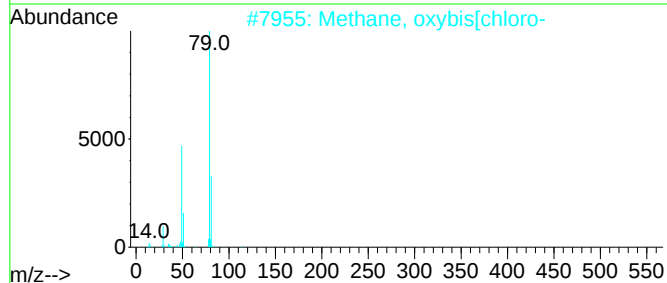
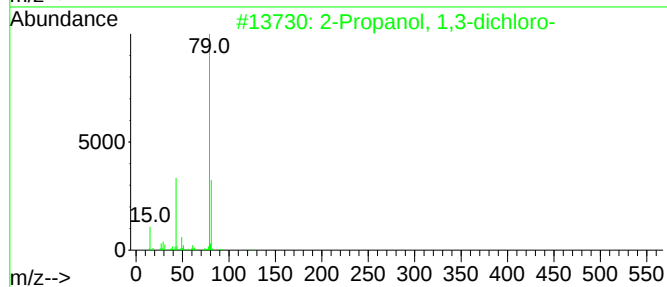
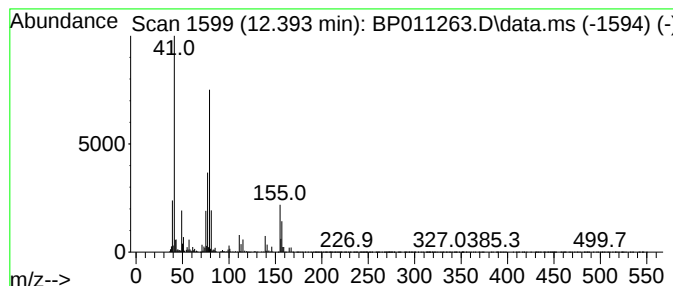
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Quant Title : SVOA CALIBRATION

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

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Peak Number 36 unknown-10 Concentration Rank 33

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 12.393 | 25.50 ng/ul | 4757490 | Acenaphthene-d10 | 14.281 |

| Hit# | of                            | 5 | Tentative ID | MW  | MolForm  | CAS#        | Qual |
|------|-------------------------------|---|--------------|-----|----------|-------------|------|
| 1    | 2-Propanol, 1,3-dichloro-     |   |              | 128 | C3H6Cl2O | 000096-23-1 | 28   |
| 2    | Methane, oxybis[chloro-       |   |              | 114 | C2H4Cl2O | 000542-88-1 | 17   |
| 3    | Methane, oxybis[chloro-       |   |              | 114 | C2H4Cl2O | 000542-88-1 | 12   |
| 4    | Chloroacetic acid allyl ester |   |              | 134 | C5H7ClO2 | 002916-14-5 | 12   |
| 5    | Chloroacetic acid allyl ester |   |              | 134 | C5H7ClO2 | 002916-14-5 | 12   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

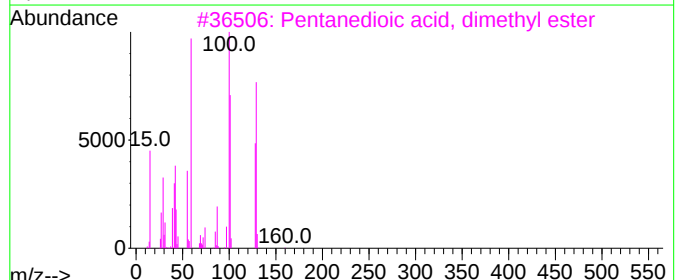
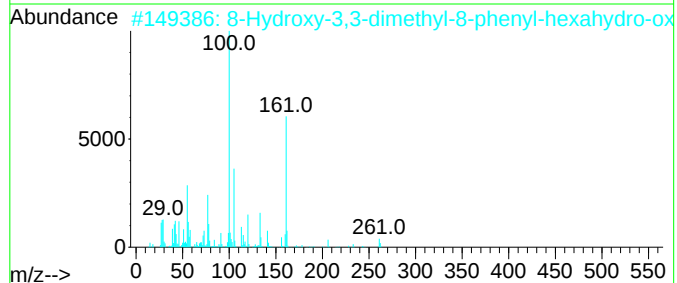
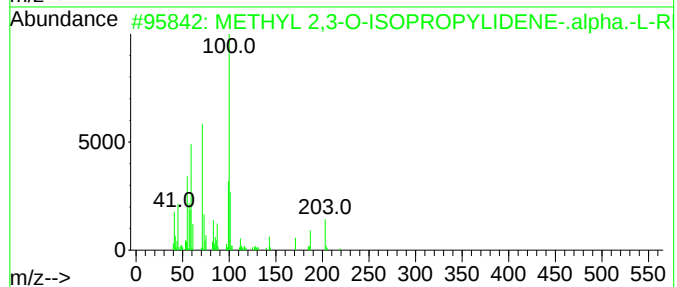
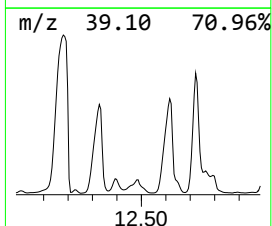
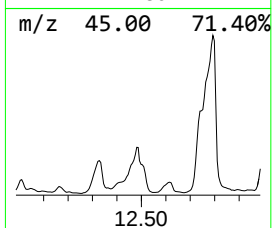
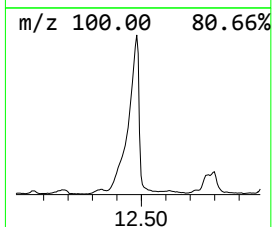
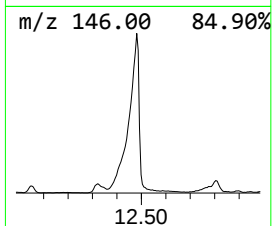
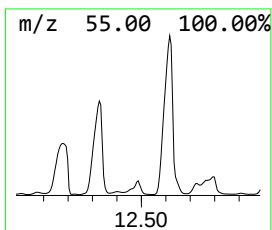
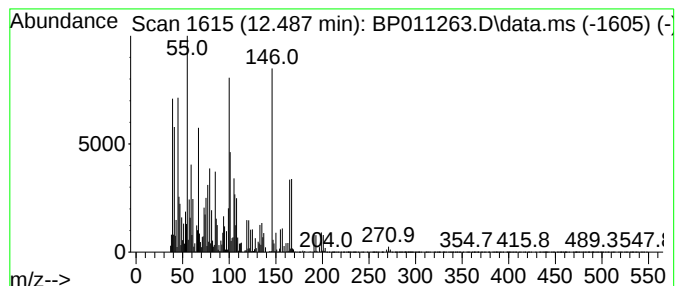
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Quant Title : SVOA CALIBRATION

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

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Peak Number 37 unknown-11 Concentration Rank 13

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 12.487 | 56.18 ng/ul | 10481000 | Acenaphthene-d10 | 14.281 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | METHYL 2,3-O-ISOPROPYLIDENE-.alp... | 218 | C10H18O5     | 1000428-66-6 | 22      |      |      |
| 2    | 8-Hydroxy-3,3-dimethyl-8-phenyl-... | 261 | C15H19NO3    | 1000287-81-1 | 14      |      |      |
| 3    | Pentanedioic acid, dimethyl ester   | 160 | C7H12O4      | 001119-40-0  | 14      |      |      |
| 4    | N,N-Diethyl-2-phenylacetamide       | 191 | C12H17NO     | 1000411-34-2 | 11      |      |      |
| 5    | Cyclobutanecarboxamide, N-(2-phe... | 273 | C18H27NO     | 1000451-60-5 | 11      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

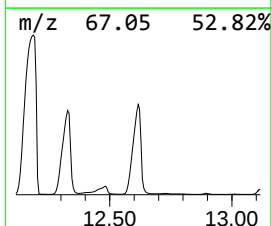
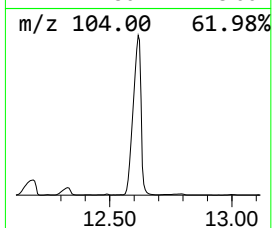
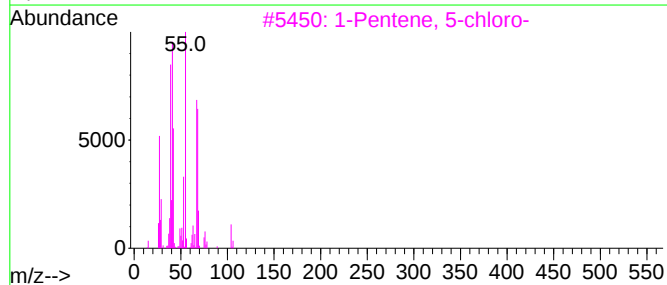
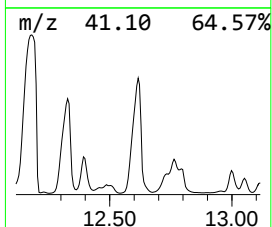
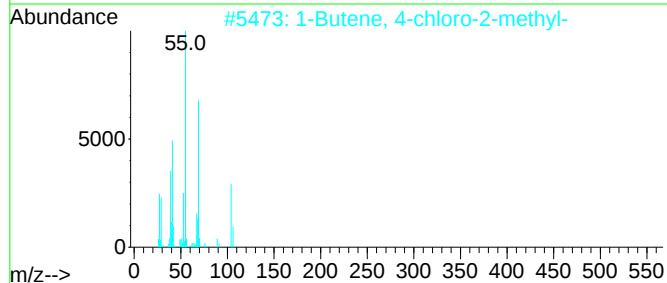
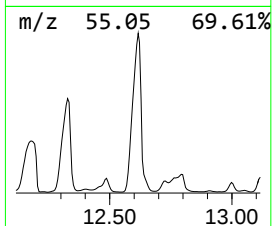
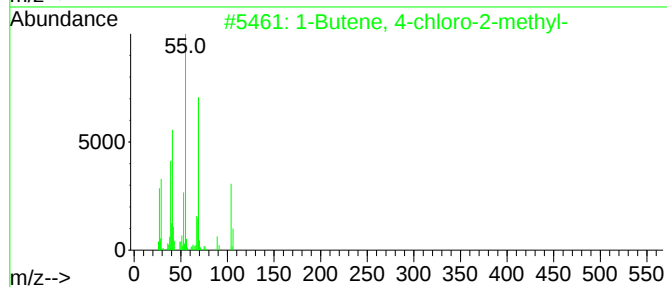
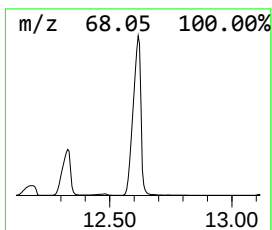
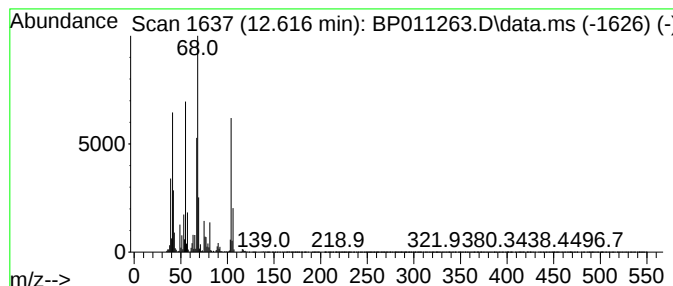
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 38 unknown-12 Concentration Rank 4

| R.T.   | EstConc      | Area     | Relative to ISTD | R.T.   |
|--------|--------------|----------|------------------|--------|
| 12.616 | 282.39 ng/ul | 52682900 | Acenaphthene-d10 | 14.281 |

| Hit# | of 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|------|------------------------------|-----|---------|-------------|------|
| 1    |      | 1-Butene, 4-chloro-2-methyl- | 104 | C5H9Cl  | 010523-96-3 | 43   |
| 2    |      | 1-Butene, 4-chloro-2-methyl- | 104 | C5H9Cl  | 010523-96-3 | 43   |
| 3    |      | 1-Pentene, 5-chloro-         | 104 | C5H9Cl  | 000928-50-7 | 40   |
| 4    |      | 1,6-Octadiene                | 110 | C8H14   | 003710-41-6 | 38   |
| 5    |      | 1,6-Octadiene, (E)-          | 110 | C8H14   | 019036-81-8 | 32   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

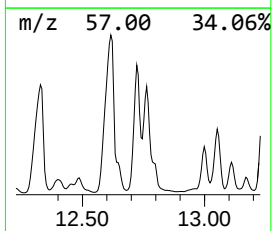
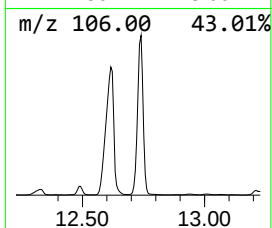
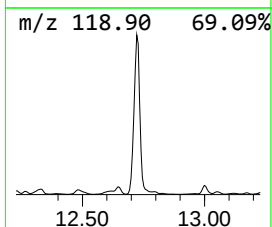
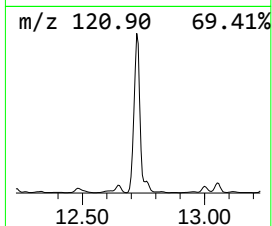
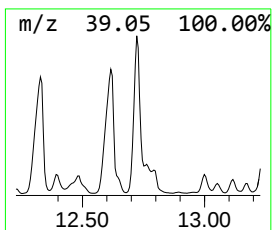
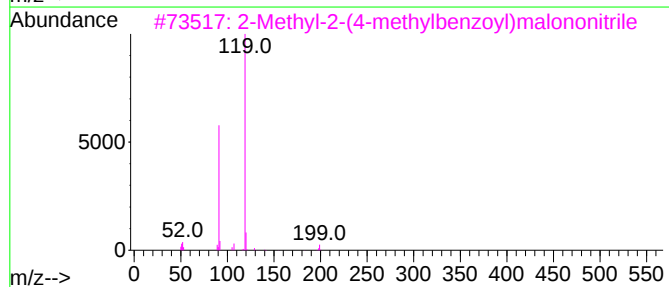
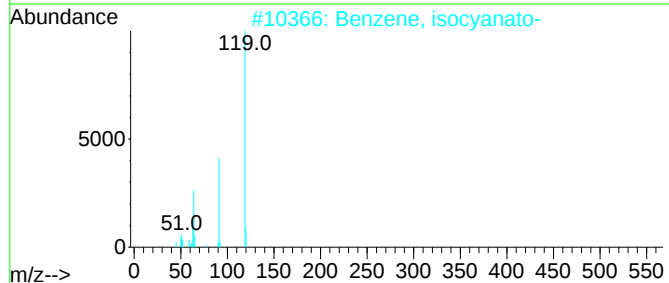
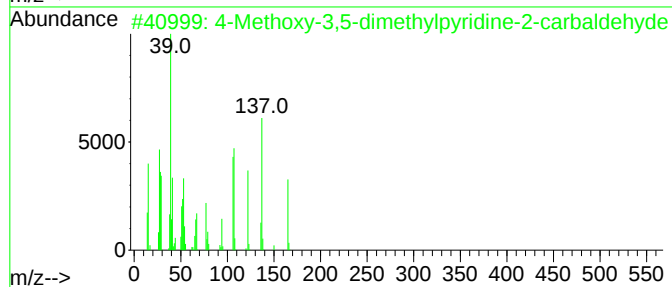
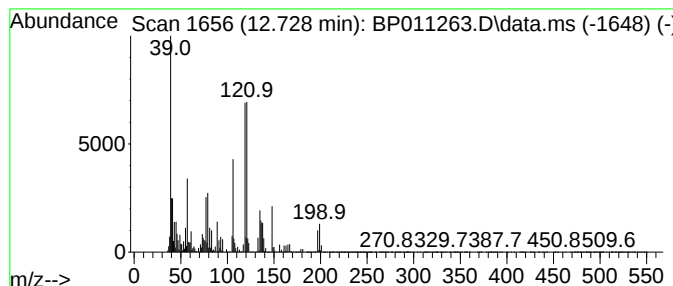
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 39 unknown-13 Concentration Rank 6

| R.T.   | EstConc      | Area     | Relative to ISTD | R.T.   |
|--------|--------------|----------|------------------|--------|
| 12.728 | 127.51 ng/ul | 23788200 | Acenaphthene-d10 | 14.281 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|------|-------------------------------------|-----|-----------|--------------|------|
| 1    |      | 4-Methoxy-3,5-dimethylpyridine-2... | 165 | C9H11NO2  | 1000486-86-4 | 36   |
| 2    |      | Benzene, isocyanato-                | 119 | C7H5NO    | 000103-71-9  | 10   |
| 3    |      | 2-Methyl-2-(4-methylbenzoyl)malo... | 198 | C12H10N2O | 1000326-92-9 | 9    |
| 4    |      | 3(2H)-Benzofuranone, 6-methyl-      | 148 | C9H8O2    | 020895-41-4  | 9    |
| 5    |      | Ethanone, 2-bromo-1-(4-methylphe... | 212 | C9H9BrO   | 000619-41-0  | 9    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

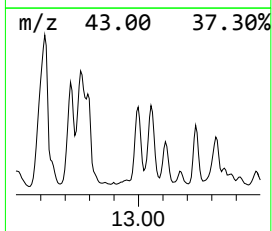
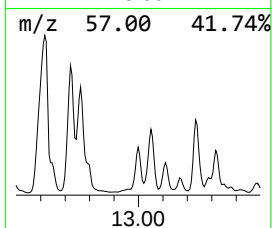
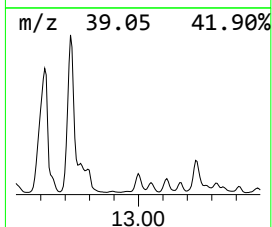
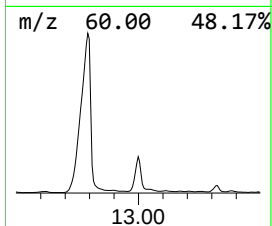
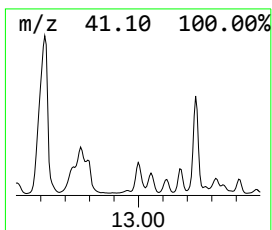
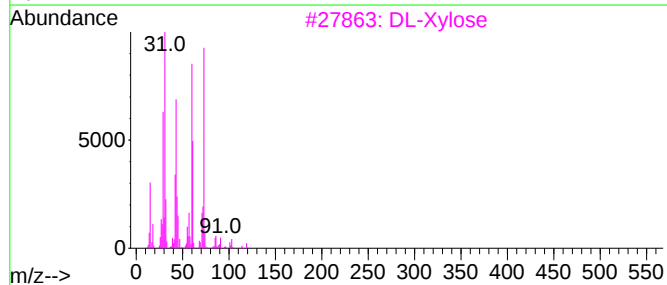
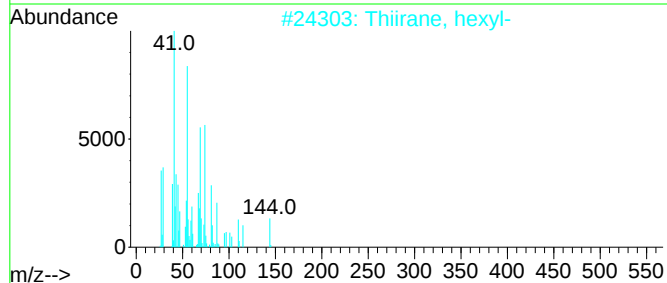
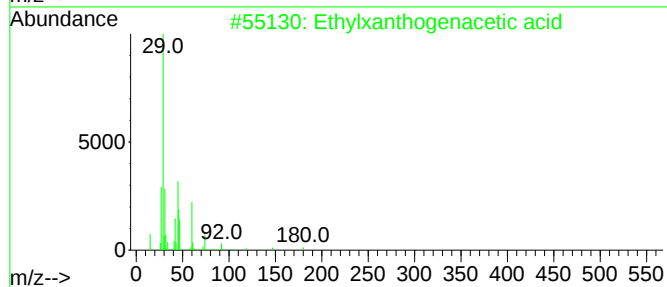
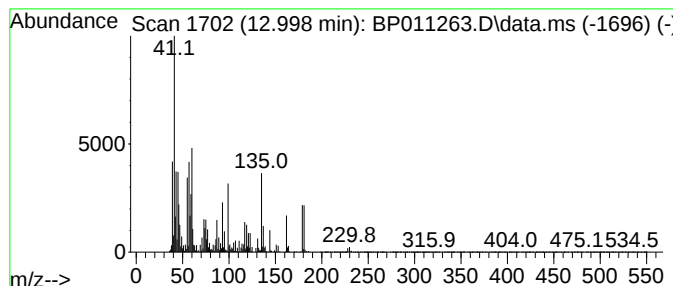
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TIC Integration Parameters: LSCINT.P

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Peak Number 41 unknown-14 Concentration Rank 26

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 12.999 | 32.40 ng/ul | 6043550 | Acenaphthene-d10 | 14.281 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Ethylxanthogenacetic acid           | 180 | C5H8O3S2 | 025554-84-1 | 11   |
| 2    |    |   | Thiirane, hexyl-                    | 144 | C8H16S   | 005633-78-3 | 10   |
| 3    |    |   | DL-Xylose                           | 150 | C5H10O5  | 041247-05-6 | 9    |
| 4    |    |   | Thiirane, methyl-                   | 74  | C3H6S    | 001072-43-1 | 9    |
| 5    |    |   | Pseudourea, 2-(2-aminoethyl)-2-t... | 119 | C3H9N3S  | 000151-16-6 | 9    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

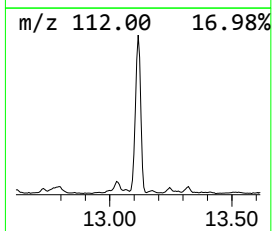
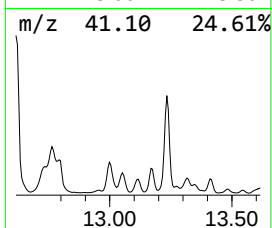
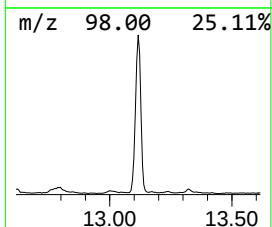
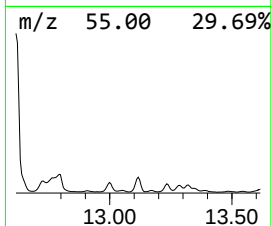
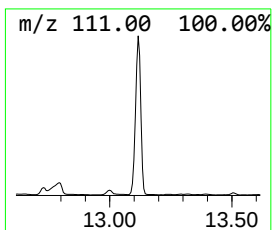
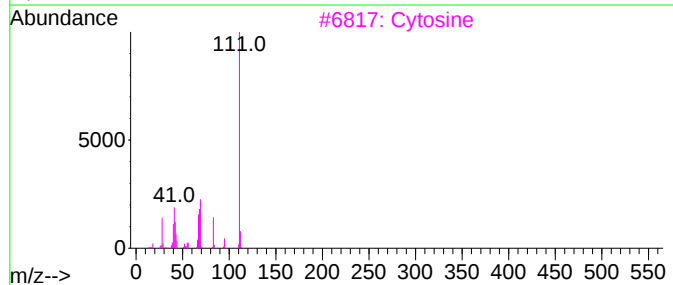
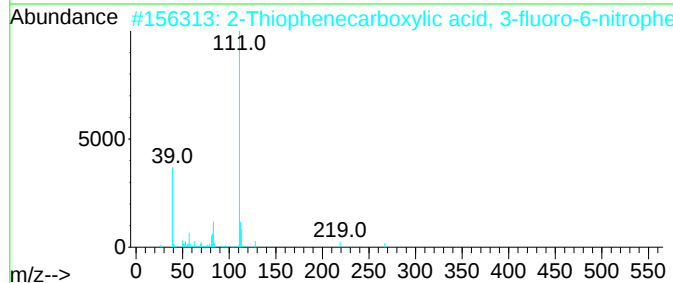
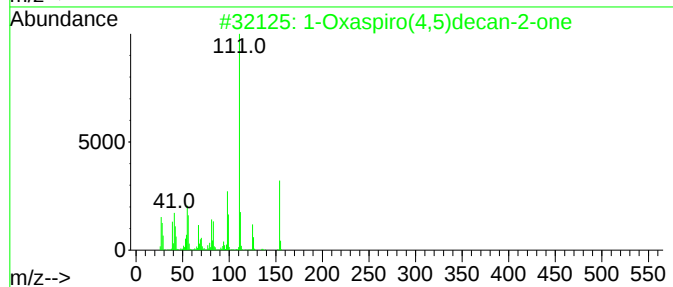
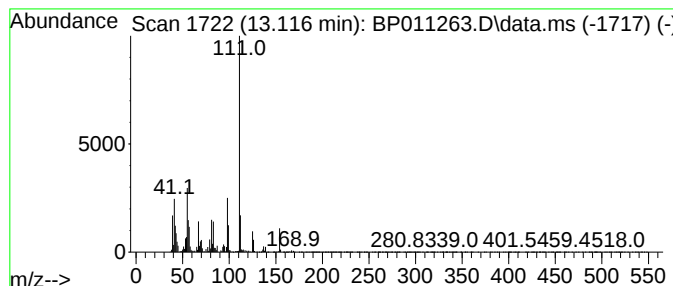
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Quant Title : SVOA CALIBRATION

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

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Peak Number 43 1-Oxaspiro(4,5)decan-2-one Concentration Rank 31

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 13.116 | 26.18 ng/ul | 4884000 | Acenaphthene-d10 | 14.281 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 1-Oxaspiro(4,5)decan-2-one          | 154 | C9H14O2    | 000699-61-6  | 91   |
| 2    |    |   | 2-Thiophenecarboxylic acid, 3-fl... | 267 | C11H6FN04S | 1000293-63-5 | 47   |
| 3    |    |   | Cytosine                            | 111 | C4H5N3O    | 000071-30-7  | 43   |
| 4    |    |   | 2-Aminopyrimidine-1-oxide           | 111 | C4H5N3O    | 035034-15-2  | 43   |
| 5    |    |   | Isocytosine                         | 111 | C4H5N3O    | 000108-53-2  | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

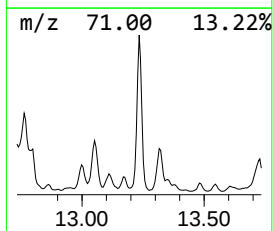
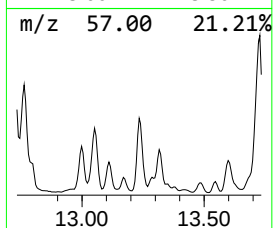
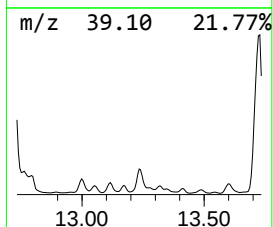
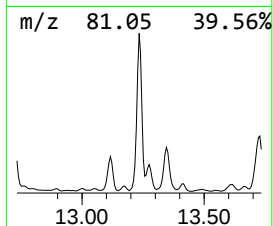
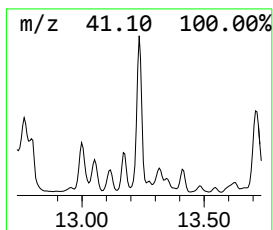
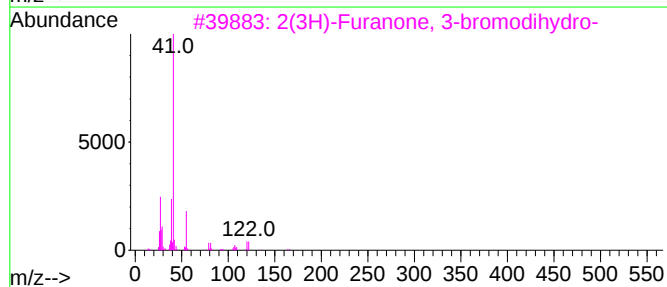
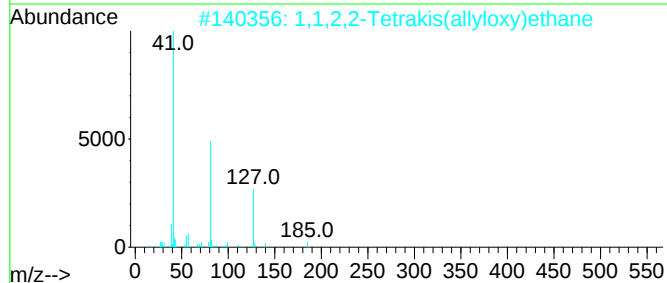
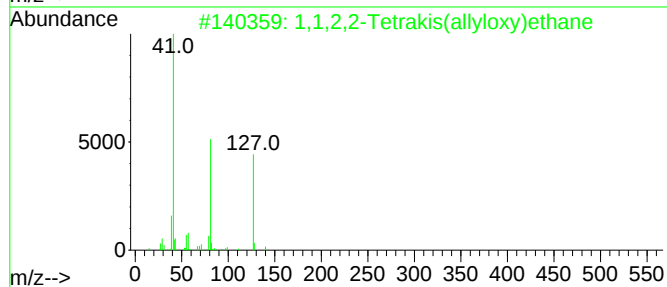
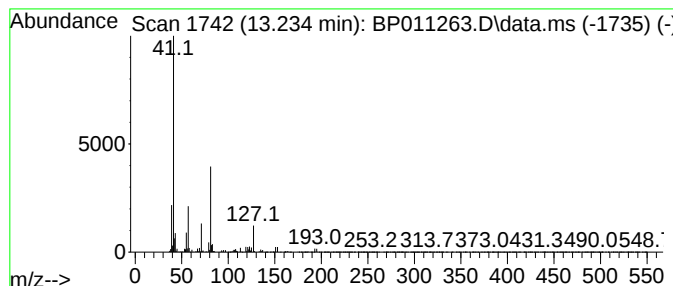
Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 44 unknown-15 Concentration Rank 24

| R.T.      | EstConc                          | Area    | Relative to ISTD | R.T.         |      |
|-----------|----------------------------------|---------|------------------|--------------|------|
| 13.234    | 34.82 ng/ul                      | 6495440 | Acenaphthene-d10 | 14.281       |      |
| Hit# of 5 | Tentative ID                     | MW      | MolForm          | CAS#         | Qual |
| 1         | 1,1,2,2-Tetrakis(allyloxy)ethane | 254     | C14H22O4         | 016646-44-9  | 40   |
| 2         | 1,1,2,2-Tetrakis(allyloxy)ethane | 254     | C14H22O4         | 016646-44-9  | 40   |
| 3         | 2(3H)-Furanone, 3-bromodihydro-  | 164     | C4H5BrO2         | 005061-21-2  | 9    |
| 4         | Octanal diallyl acetal           | 226     | C14H26O2         | 1010430-98-7 | 9    |
| 5         | 2-Azacyclooctanone               | 127     | C7H13NO          | 000673-66-5  | 7    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

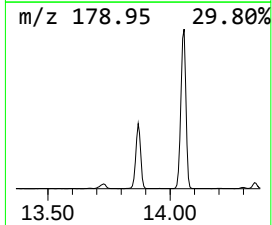
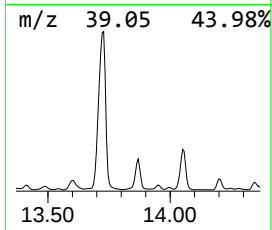
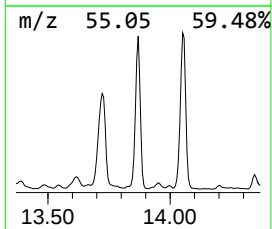
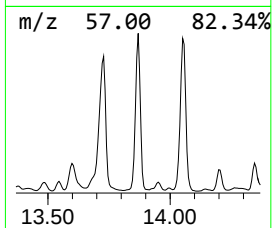
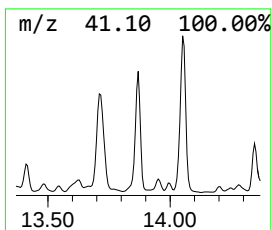
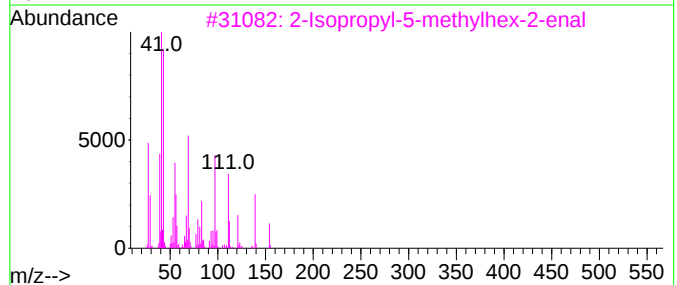
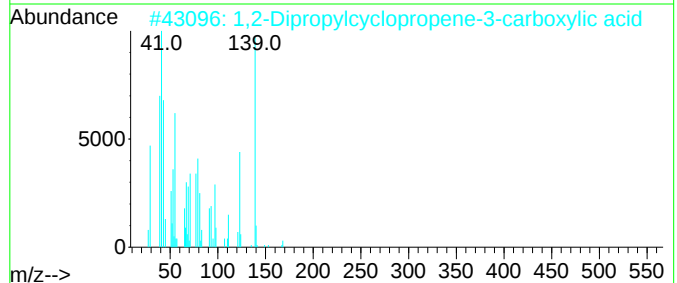
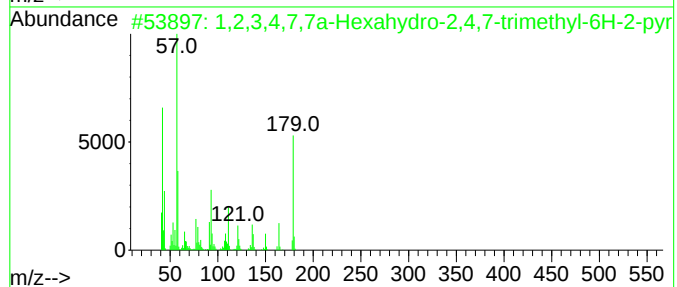
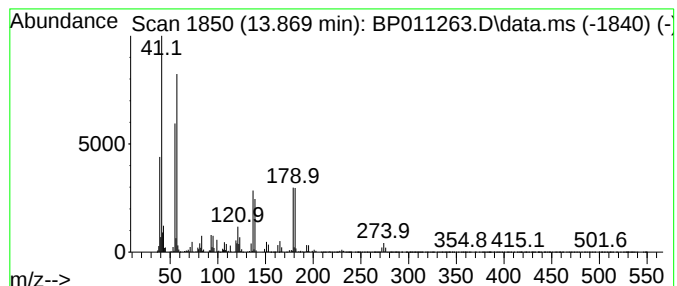
Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 47 unknown-16 Concentration Rank 20

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 13.869    | 42.06 ng/ul                         | 7845810 | Acenaphthene-d10 | 14.281      |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | 1,2,3,4,7,7a-Hexahydro-2,4,7-tri... | 179     | C11H17NO         | 006878-83-7 | 25   |
| 2         | 1,2-Dipropylcyclopropene-3-carbo... | 168     | C10H16O2         | 001605-38-5 | 12   |
| 3         | 2-Isopropyl-5-methylhex-2-enal      | 154     | C10H18O          | 035158-25-9 | 12   |
| 4         | Butane, 2,3-dibromo-                | 214     | C4H8Br2          | 005408-86-6 | 12   |
| 5         | Heptane, 1-bromo-                   | 178     | C7H15Br          | 000629-04-9 | 10   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

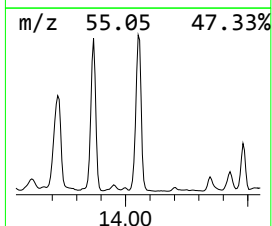
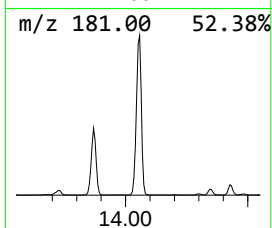
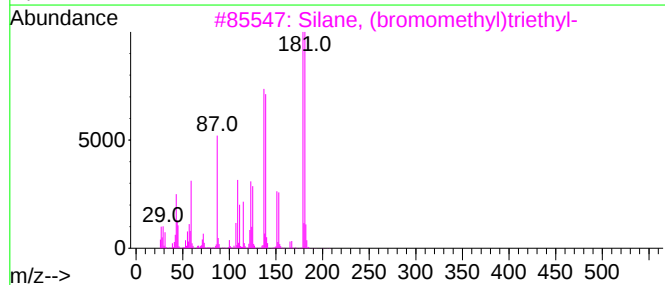
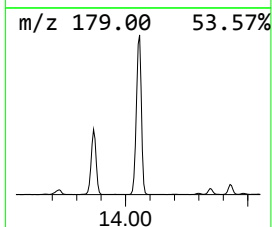
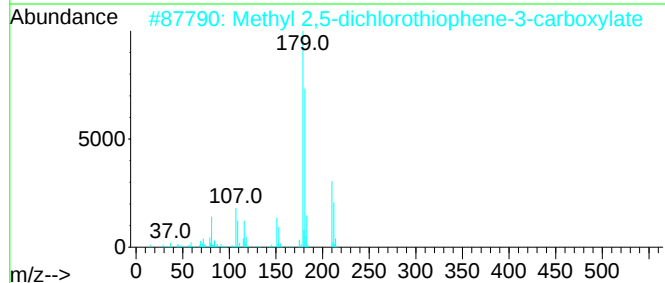
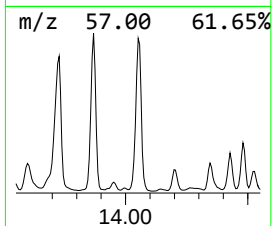
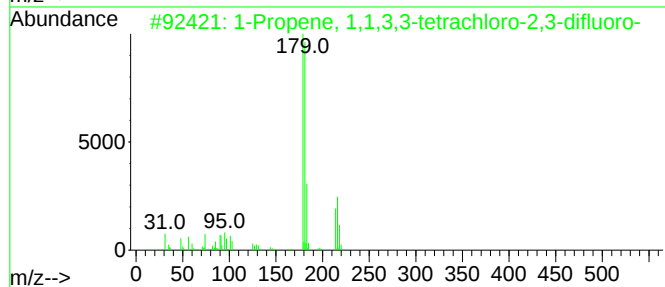
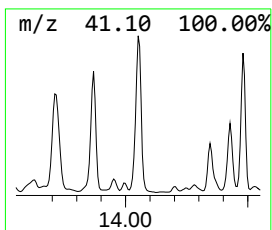
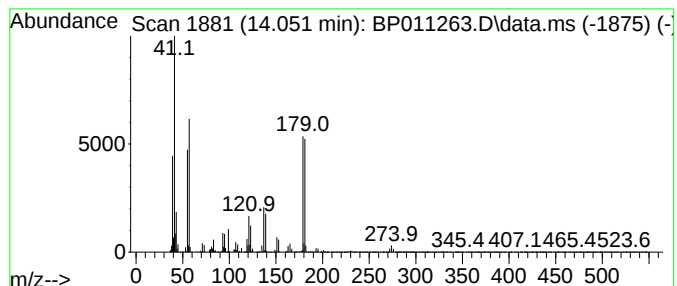
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 48 unknown-17 Concentration Rank 12

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 14.051 | 56.47 ng/ul | 10535800 | Acenaphthene-d10 | 14.281 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1-Propene, 1,1,3,3-tetrachloro-2... | 214 | C3Cl4F2      | 000815-16-7 | 27      |      |      |
| 2    | Methyl 2,5-dichlorothiophene-3-c... | 210 | C6H4Cl2O2S   | 145129-54-0 | 16      |      |      |
| 3    | Silane, (bromomethyl)triethyl-      | 208 | C7H17BrSi    | 001112-53-4 | 16      |      |      |
| 4    | 2,5-Dichloro-4-fluoroaniline        | 179 | C6H4Cl2FN    | 002729-37-5 | 14      |      |      |
| 5    | 2,4-Dichloro-3-fluoroaniline        | 179 | C6H4Cl2FN    | 000443-93-6 | 14      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

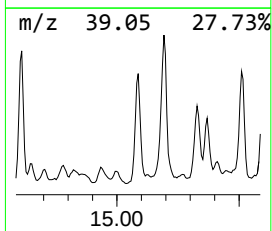
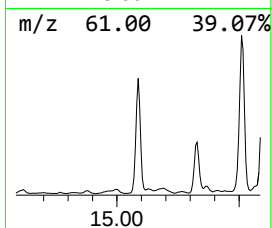
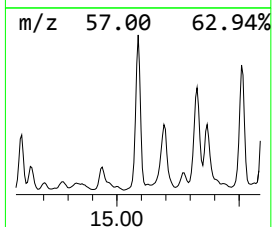
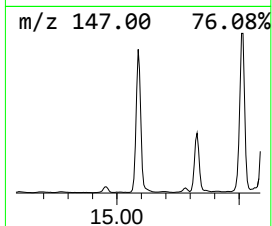
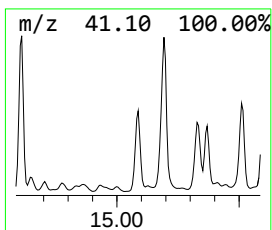
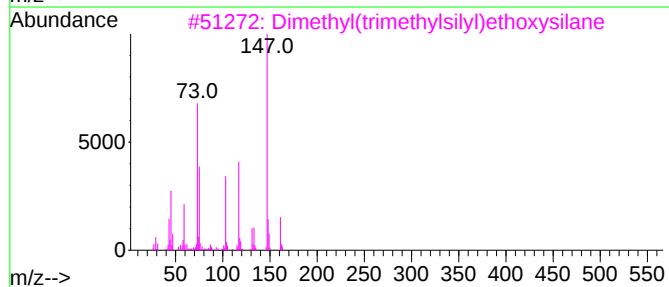
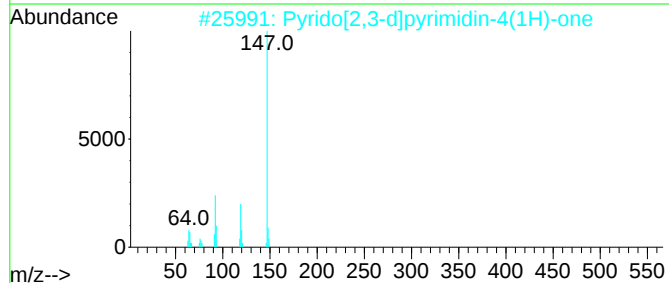
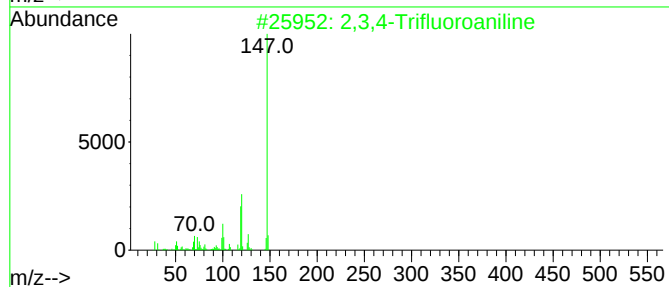
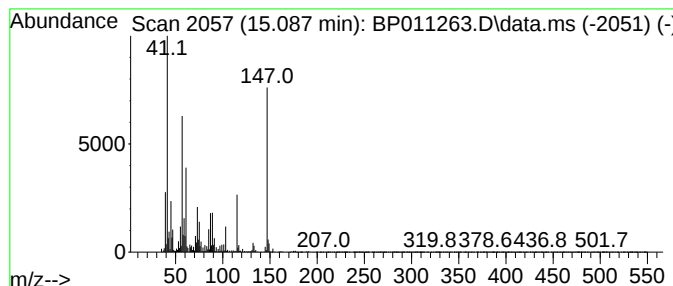
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ClientSampleId :  
EXF05

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 55 unknown-18 Concentration Rank 29

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|--------------|------|
| 15.087    | 28.61 ng/ul                         | 5337370 | Acenaphthene-d10 | 14.281       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#         | Qual |
| 1         | 2,3,4-Trifluoroaniline              | 147     | C6H4F3N          | 003862-73-5  | 35   |
| 2         | Pyrido[2,3-d]pyrimidin-4(1H)-one    | 147     | C7H5N3O          | 024410-19-3  | 32   |
| 3         | Dimethyl(trimethylsilyl)ethoxysi... | 176     | C7H20OSi2        | 018297-47-7  | 32   |
| 4         | (3,5,6-Trifluoropyridin-2-yl)ace... | 191     | C7H4F3NO2        | 1000444-85-1 | 25   |
| 5         | 6-Ethyl-3-pentamethyldisilyloxyd... | 316     | C17H40OSi2       | 1000216-89-7 | 25   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

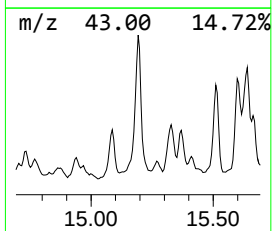
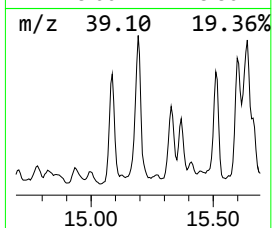
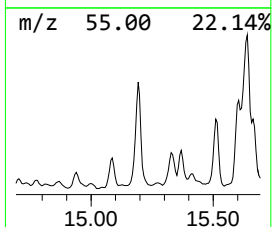
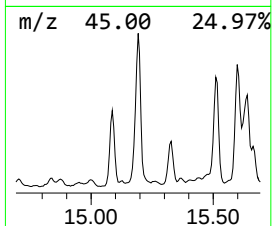
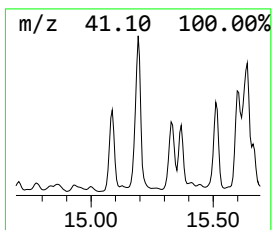
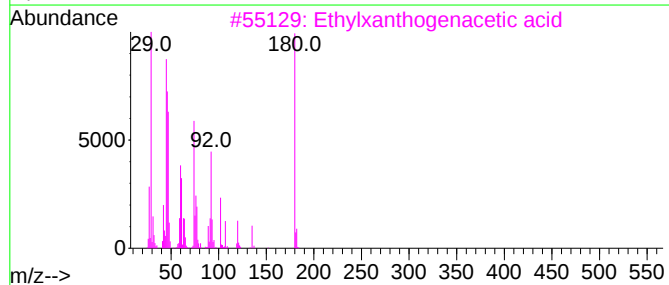
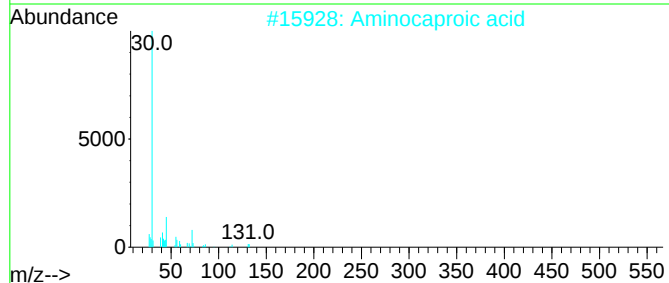
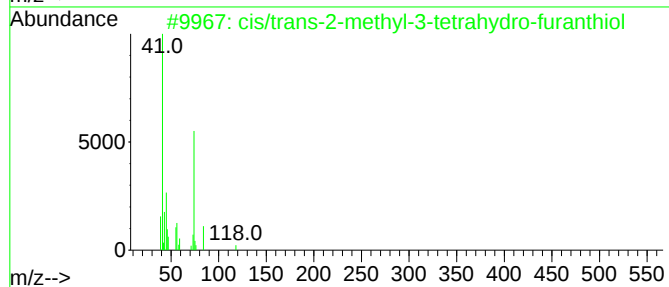
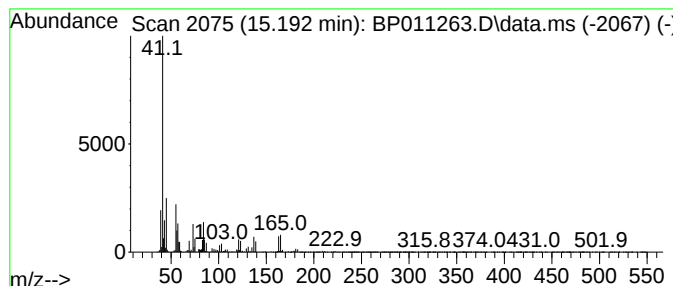
Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 56 unknown-19 Concentration Rank 25

| R.T.      | EstConc                             | Area    | Relative to ISTD |          | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|----------|--------------|------|
| 15.193    | 34.75 ng/ul                         | 6482130 | Acenaphthene-d10 |          | 14.281       |      |
| Hit# of 5 | Tentative ID                        |         | MW               | MolForm  | CAS#         | Qual |
| 1         | cis/trans-2-methyl-3-tetrahydro-... |         | 118              | C5H10O5  | 1000357-16-8 | 17   |
| 2         | Aminocaproic acid                   |         | 131              | C6H13NO2 | 000060-32-2  | 9    |
| 3         | Ethylxanthogenacetic acid           |         | 180              | C5H8O3S2 | 025554-84-1  | 9    |
| 4         | Hexane, 1,5-dibromo-                |         | 242              | C6H12Br2 | 000627-96-3  | 9    |
| 5         | Ethylene glycol diallyl ether       |         | 142              | C8H14O2  | 007529-27-3  | 9    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

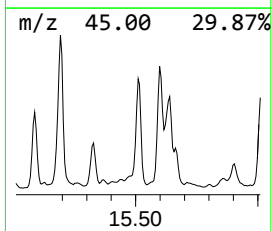
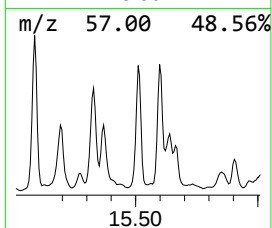
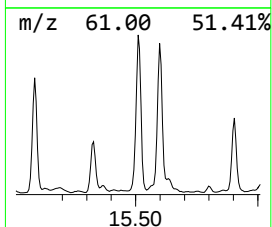
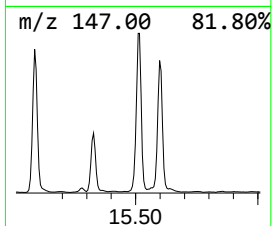
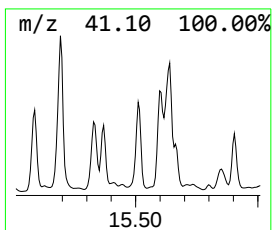
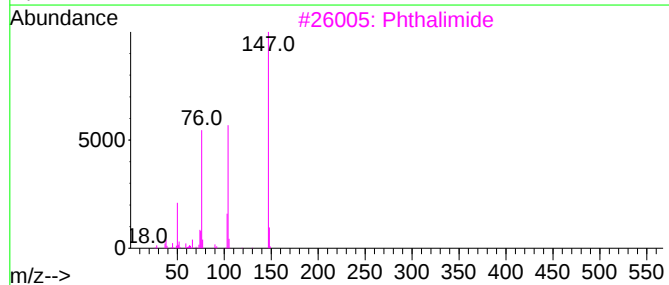
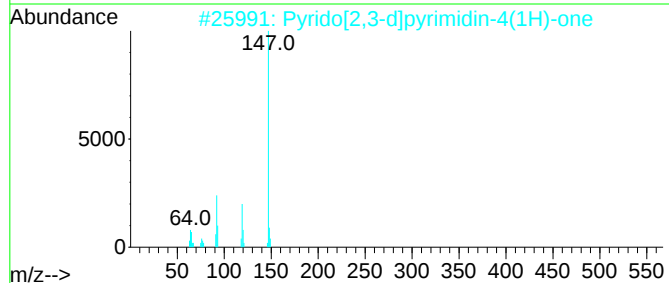
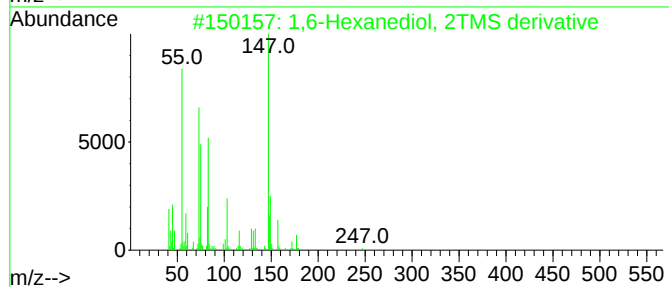
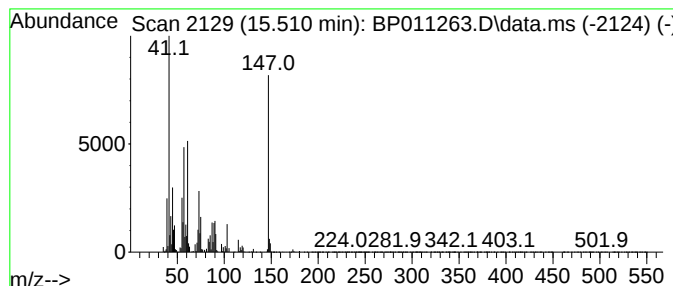
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 57 unknown-20 Concentration Rank 27

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 15.510 | 32.02 ng/ul | 5973550 | Acenaphthene-d10 | 14.281 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-------------|-------------|------|
| 1    |    |   | 1,6-Hexanediol, 2TMS derivative     | 262 | C12H30O2Si2 | 006222-22-6 | 43   |
| 2    |    |   | Pyrido[2,3-d]pyrimidin-4(1H)-one    | 147 | C7H5N3O     | 024410-19-3 | 38   |
| 3    |    |   | Phthalimide                         | 147 | C8H5NO2     | 000085-41-6 | 35   |
| 4    |    |   | (3-Methoxyphenyl)acetonitrile       | 147 | C9H9NO      | 019924-43-7 | 35   |
| 5    |    |   | Pyrimidine, 2-fluoro-5-chloro-4-... | 147 | C4H3ClFN3   | 000155-10-2 | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

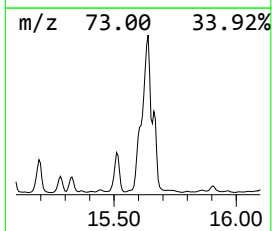
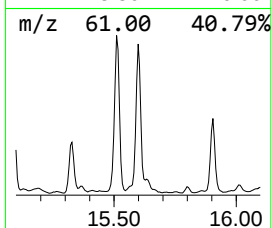
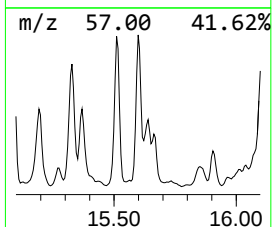
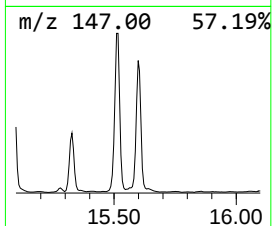
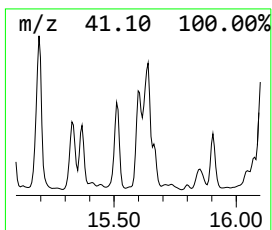
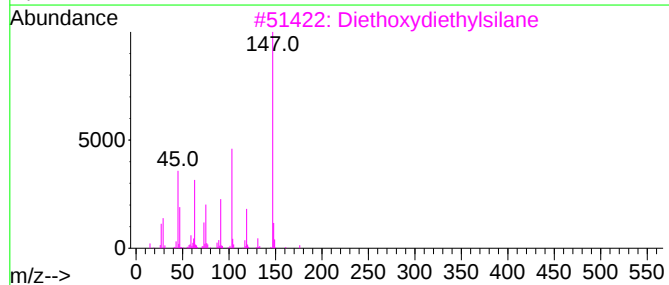
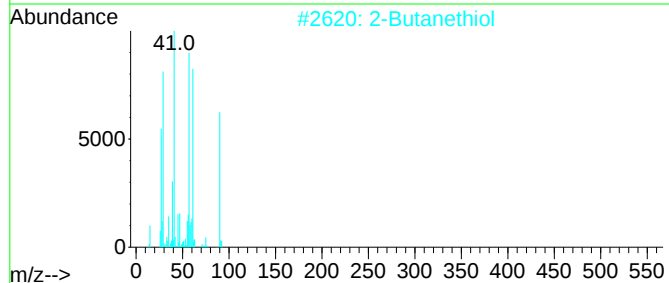
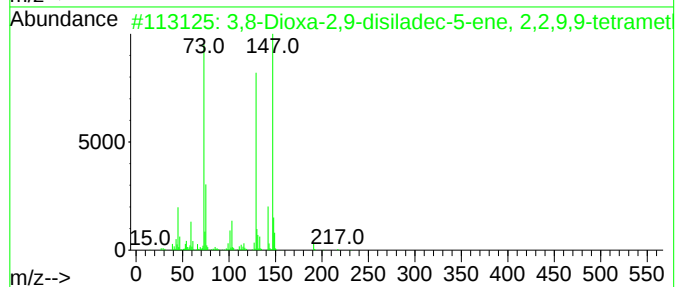
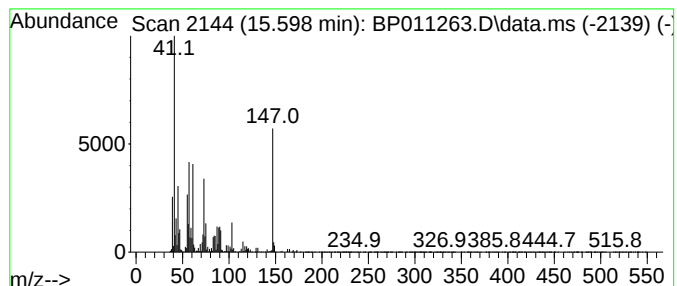
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 58 unknown-21 Concentration Rank 23

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 15.598 | 35.87 ng/ul | 6691430 | Acenaphthene-d10 | 14.281 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | 3,8-Dioxa-2,9-disiladec-5-ene, 2... | 232 | C10H24O2Si2 | 1000473-41-0 | 32   |
| 2    |    |   | 2-Butanethiol                       | 90  | C4H10S      | 000513-53-1  | 27   |
| 3    |    |   | Diethoxydiethylsilane               | 176 | C8H20O2Si   | 005021-93-2  | 16   |
| 4    |    |   | (3-Methoxyphenyl)acetonitrile       | 147 | C9H9NO      | 019924-43-7  | 14   |
| 5    |    |   | Phthalimide                         | 147 | C8H5NO2     | 000085-41-6  | 14   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

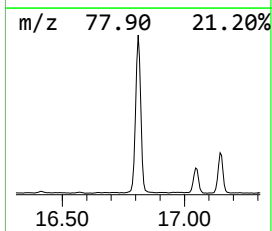
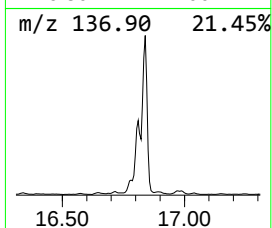
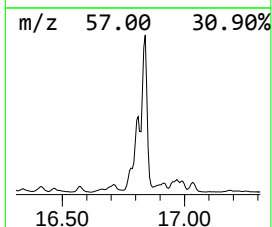
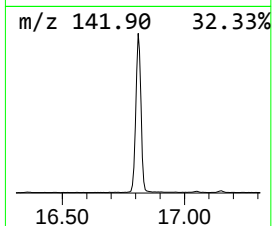
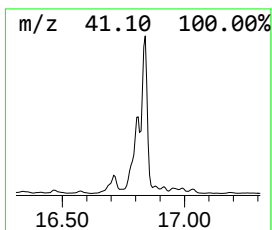
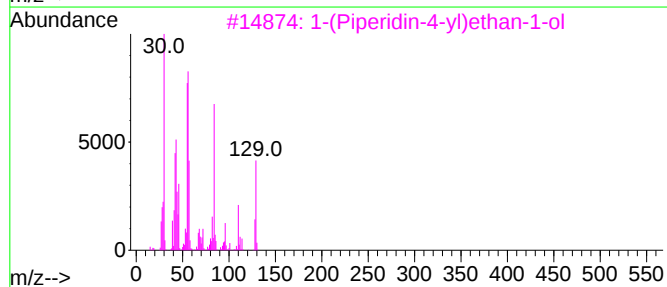
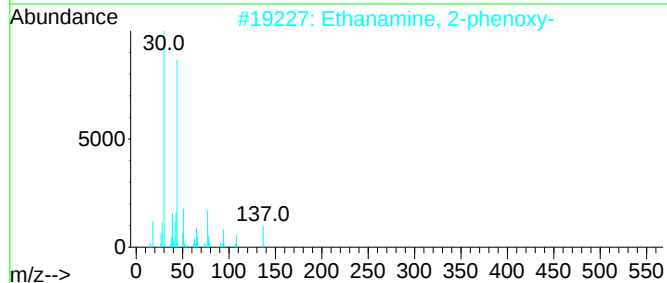
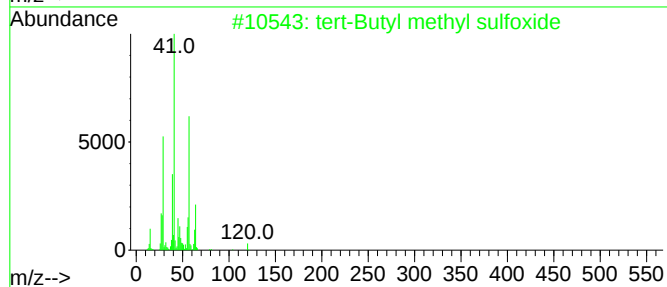
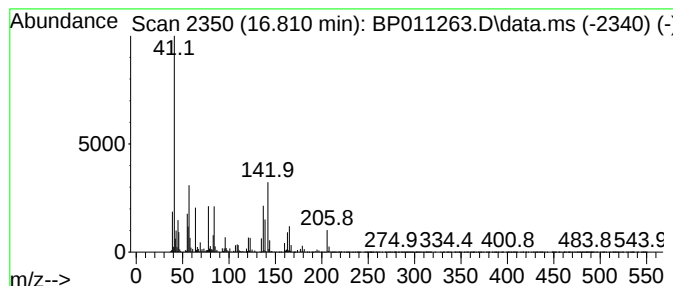
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 66 unknown-22 Concentration Rank 17

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 16.810 | 53.28 ng/ul | 11304800 | Phenanthrene-d10 | 17.045 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | tert-Butyl methyl sulfoxide        | 120 | C5H12OS | 014094-11-2 | 7    |
| 2    |    |   | Ethanamine, 2-phenoxy-             | 137 | C8H11NO | 001758-46-9 | 3    |
| 3    |    |   | 1-(Piperidin-4-yl)ethan-1-ol       | 129 | C7H15NO | 006457-48-3 | 2    |
| 4    |    |   | Cyclopropaneethanol, 2,2-difluoro- | 122 | C5H8F2O | 117284-59-0 | 2    |
| 5    |    |   | Phenylethanamine                   | 137 | C8H11NO | 007568-93-6 | 2    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

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

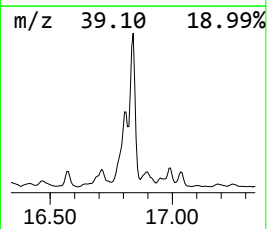
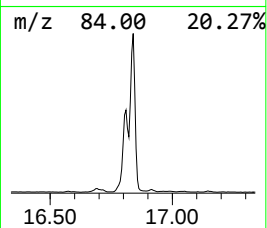
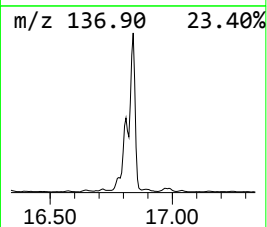
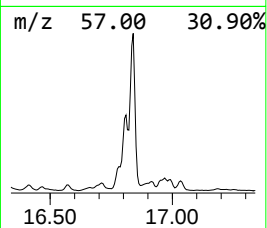
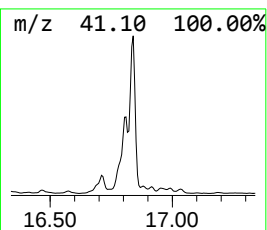
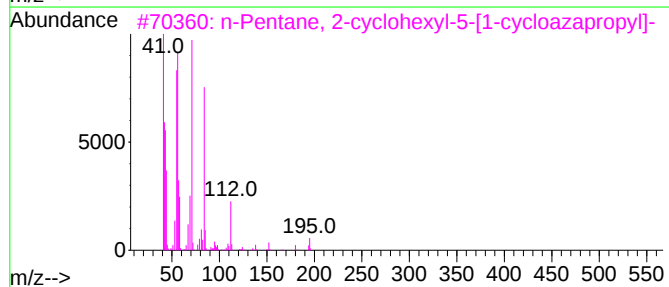
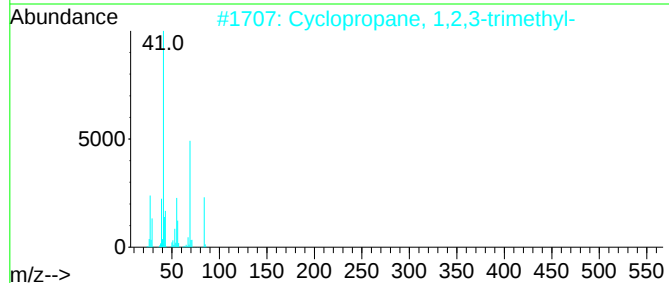
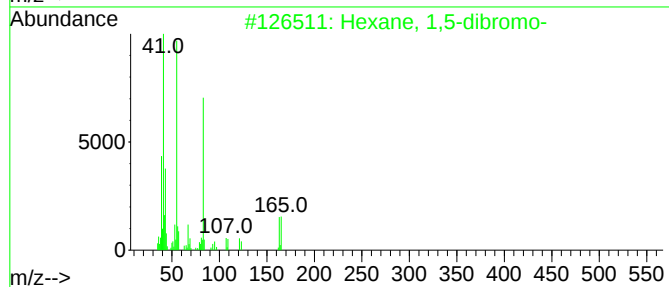
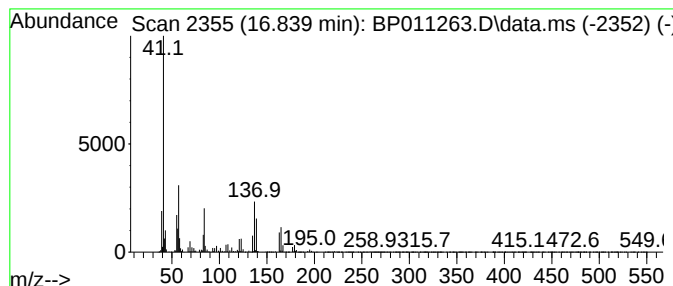
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Peak Number 67 unknown-23

Concentration Rank 16

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 16.839 | 54.02 ng/ul | 11462700 | Phenanthrene-d10 | 17.045 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Hexane, 1,5-dibromo-                | 242 | C6H12Br2 | 000627-96-3 | 13   |
| 2    |    |   | Cyclopropane, 1,2,3-trimethyl-      | 84  | C6H12    | 042984-19-0 | 9    |
| 3    |    |   | n-Pentane, 2-cyclohexyl-5-[1-cyc... | 195 | C13H25N  | 038914-53-3 | 9    |
| 4    |    |   | 1-Butene, 2,3-dimethyl-             | 84  | C6H12    | 000563-78-0 | 9    |
| 5    |    |   | Pentane, 3-methylene-               | 84  | C6H12    | 000760-21-4 | 9    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
Data File : BP011263.D  
Acq On : 04 Aug 2022 14:20  
Operator : CG/JU  
Sample : N3956-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EXF05

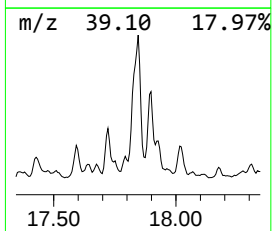
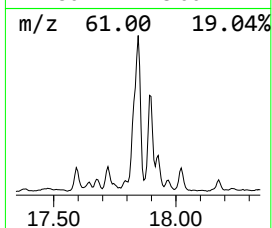
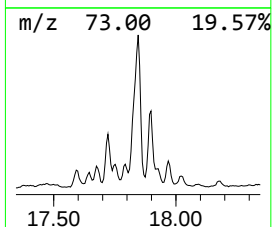
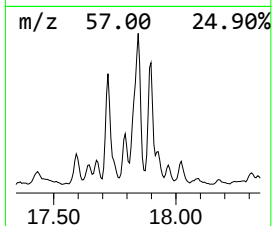
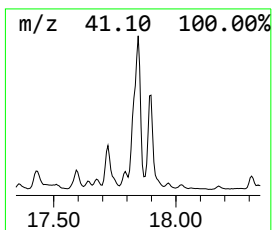
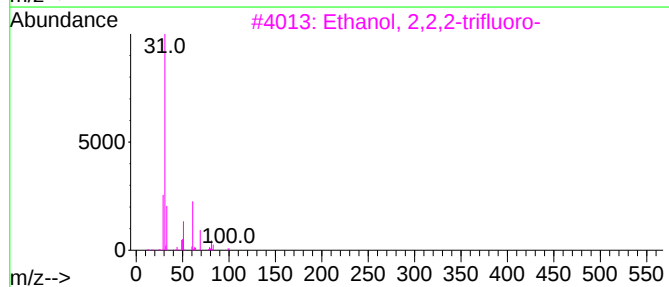
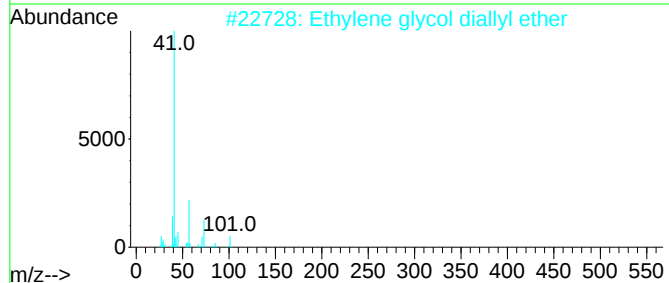
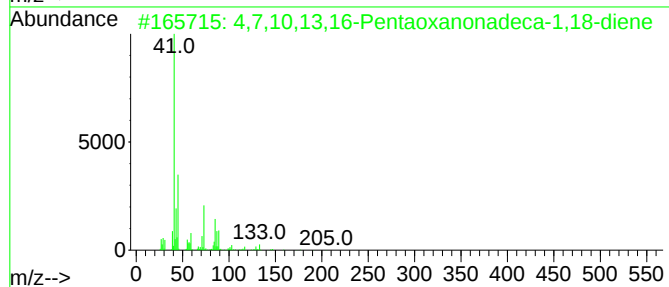
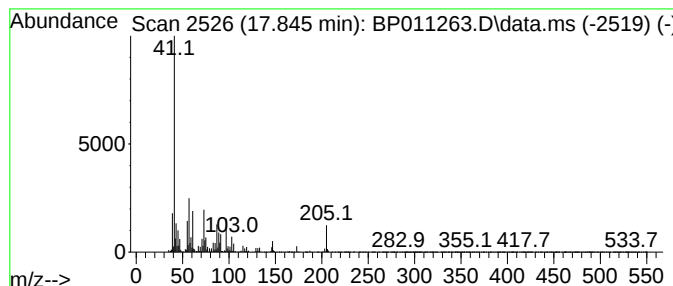
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 69 unknown-24 Concentration Rank 28

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 17.845 | 30.17 ng/ul | 6402260 | Phenanthrene-d10 | 17.045 |

| Hit# | of                                 | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 4,7,10,13,16-Pentaoxonadeca-1,...  | 274 | C14H26O5     | 058185-54-9 | 27      |      |      |
| 2    | Ethylene glycol diallyl ether      | 142 | C8H14O2      | 007529-27-3 | 10      |      |      |
| 3    | Ethanol, 2,2,2-trifluoro-          | 100 | C2H3F3O      | 000075-89-8 | 9       |      |      |
| 4    | Isobutylamine                      | 73  | C4H11N       | 000078-81-9 | 9       |      |      |
| 5    | Acetic acid, hydroxy-, ethyl ester | 104 | C4H8O3       | 000623-50-7 | 9       |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080422\  
 Data File : BP011263.D  
 Acq On : 04 Aug 2022 14:20  
 Operator : CG/JU  
 Sample : N3956-08  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EXF05

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
 Quant Title : SVOA CALIBRATION

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

| TIC Top Hit name   | RT     | EstConc | Units | Response  | --Internal Standard-- |        |          |      |
|--------------------|--------|---------|-------|-----------|-----------------------|--------|----------|------|
|                    |        |         |       |           | #                     | RT     | Resp     | Conc |
| 2-Ethyltetrahyd... | 3.834  | 105.8   | ng/ul | 18979000  | 1                     | 7.605  | 3587590  | 20.0 |
| Propane, 1,3-di... | 4.081  | 300.8   | ng/ul | 53954600  | 1                     | 7.605  | 3587590  | 20.0 |
| (DEL) Alkane: C... | 5.787  | 184.4   | ng/ul | 33079400  | 1                     | 7.605  | 3587590  | 20.0 |
| Benzene, bromo-    | 6.287  | 59.5    | ng/ul | 10677800  | 1                     | 7.605  | 3587590  | 20.0 |
| unknown-01         | 7.040  | 52.0    | ng/ul | 9318060   | 1                     | 7.605  | 3587590  | 20.0 |
| Thiepane           | 7.681  | 27.6    | ng/ul | 4959920   | 1                     | 7.605  | 3587590  | 20.0 |
| unknown-02         | 7.940  | 383.5   | ng/ul | 68792100  | 1                     | 7.605  | 3587590  | 20.0 |
| unknown-03         | 8.011  | 59.7    | ng/ul | 10712900  | 1                     | 7.605  | 3587590  | 20.0 |
| (DEL) Alkane: C... | 8.210  | 37.1    | ng/ul | 6654960   | 1                     | 7.605  | 3587590  | 20.0 |
| (DEL) Alkane: S... | 8.587  | 15.4    | ng/ul | 2753700   | 1                     | 7.605  | 3587590  | 20.0 |
| unknown-04         | 8.716  | 306.4   | ng/ul | 54959500  | 1                     | 7.605  | 3587590  | 20.0 |
| Hexane, 1,6-dic... | 9.428  | 26.1    | ng/ul | 31375100  | 2                     | 10.410 | 24039600 | 20.0 |
| unknown-05         | 10.740 | 54.6    | ng/ul | 65654800  | 2                     | 10.410 | 24039600 | 20.0 |
| unknown-06         | 10.957 | 46.4    | ng/ul | 55825300  | 2                     | 10.410 | 24039600 | 20.0 |
| unknown-07         | 10.999 | 55.1    | ng/ul | 66222500  | 2                     | 10.410 | 24039600 | 20.0 |
| unknown-08         | 12.187 | 88.2    | ng/ul | 106062000 | 2                     | 10.410 | 24039600 | 20.0 |
| unknown-09         | 12.328 | 40.7    | ng/ul | 48883100  | 2                     | 10.410 | 24039600 | 20.0 |
| unknown-10         | 12.393 | 25.5    | ng/ul | 4757490   | 3                     | 14.281 | 3731150  | 20.0 |
| unknown-11         | 12.487 | 56.2    | ng/ul | 10481000  | 3                     | 14.281 | 3731150  | 20.0 |
| unknown-12         | 12.616 | 282.4   | ng/ul | 52682900  | 3                     | 14.281 | 3731150  | 20.0 |
| unknown-13         | 12.728 | 127.5   | ng/ul | 23788200  | 3                     | 14.281 | 3731150  | 20.0 |
| unknown-14         | 12.999 | 32.4    | ng/ul | 6043550   | 3                     | 14.281 | 3731150  | 20.0 |
| 1-Oxaspiro(4,5)... | 13.116 | 26.2    | ng/ul | 4884000   | 3                     | 14.281 | 3731150  | 20.0 |
| unknown-15         | 13.234 | 34.8    | ng/ul | 6495440   | 3                     | 14.281 | 3731150  | 20.0 |
| unknown-16         | 13.869 | 42.1    | ng/ul | 7845810   | 3                     | 14.281 | 3731150  | 20.0 |
| unknown-17         | 14.051 | 56.5    | ng/ul | 10535800  | 3                     | 14.281 | 3731150  | 20.0 |
| unknown-18         | 15.087 | 28.6    | ng/ul | 5337370   | 3                     | 14.281 | 3731150  | 20.0 |
| unknown-19         | 15.193 | 34.8    | ng/ul | 6482130   | 3                     | 14.281 | 3731150  | 20.0 |
| unknown-20         | 15.510 | 32.0    | ng/ul | 5973550   | 3                     | 14.281 | 3731150  | 20.0 |
| unknown-21         | 15.598 | 35.9    | ng/ul | 6691430   | 3                     | 14.281 | 3731150  | 20.0 |
| unknown-22         | 16.810 | 53.3    | ng/ul | 11304800  | 4                     | 17.045 | 4243690  | 20.0 |
| unknown-23         | 16.839 | 54.0    | ng/ul | 11462700  | 4                     | 17.045 | 4243690  | 20.0 |
| unknown-24         | 17.845 | 30.2    | ng/ul | 6402260   | 4                     | 17.045 | 4243690  | 20.0 |