

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092823\  
 Data File : BP017427.D  
 Acq On : 28 Sep 2023 23:10  
 Operator : MA/JU  
 Sample : 04417-39  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EZYE6

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-BP092023.MA.M  
 Title : SVOA CALIBRATION

Signal : TIC: BP017427.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         | 3.299       | 32            | 36          | 40           | rVB      | 40930          | 45652         | 1.12%           | 0.100%        |
| 2         | 3.740       | 108           | 111         | 119          | rBV      | 141094         | 208896        | 5.13%           | 0.458%        |
| 3         | 3.846       | 127           | 129         | 134          | rVB2     | 15823          | 19730         | 0.48%           | 0.043%        |
| 4         | 3.928       | 138           | 143         | 148          | rBV      | 70947          | 98086         | 2.41%           | 0.215%        |
| 5         | 4.046       | 160           | 163         | 170          | rVB      | 22146          | 28922         | 0.71%           | 0.063%        |
| 6         | 4.193       | 183           | 188         | 194          | rBV5     | 30697          | 50115         | 1.23%           | 0.110%        |
| 7         | 4.275       | 197           | 202         | 209          | rVB3     | 25819          | 43126         | 1.06%           | 0.095%        |
| 8         | 4.434       | 225           | 229         | 233          | rVB2     | 23232          | 33580         | 0.82%           | 0.074%        |
| 9         | 4.987       | 318           | 323         | 327          | rBV      | 39361          | 58824         | 1.44%           | 0.129%        |
| 10        | 5.028       | 327           | 330         | 338          | rVV2     | 23748          | 50161         | 1.23%           | 0.110%        |
| 11        | 5.534       | 410           | 416         | 418          | rBV2     | 29823          | 49374         | 1.21%           | 0.108%        |
| 12        | 5.758       | 448           | 454         | 458          | rVB2     | 13809          | 23541         | 0.58%           | 0.052%        |
| 13        | 5.863       | 468           | 472         | 476          | rBV2     | 20863          | 28364         | 0.70%           | 0.062%        |
| 14        | 5.911       | 476           | 480         | 484          | rVB      | 16514          | 26148         | 0.64%           | 0.057%        |
| 15        | 6.358       | 550           | 556         | 560          | rBV      | 24526          | 40134         | 0.99%           | 0.088%        |
| 16        | 6.393       | 560           | 562         | 572          | rVV10    | 9542           | 21223         | 0.52%           | 0.047%        |
| 17        | 6.516       | 576           | 583         | 589          | rBV5     | 19157          | 33899         | 0.83%           | 0.074%        |
| 18        | 6.799       | 626           | 631         | 636          | rVB2     | 17095          | 27105         | 0.67%           | 0.059%        |
| 19        | 7.116       | 678           | 685         | 698          | rBV      | 706399         | 1139166       | 27.97%          | 2.499%        |
| 20        | 7.299       | 707           | 716         | 725          | rBV      | 540750         | 851539        | 20.90%          | 1.868%        |
| 21        | 7.481       | 737           | 747         | 756          | rBV      | 741284         | 1212181       | 29.76%          | 2.660%        |
| 22        | 7.575       | 758           | 763         | 767          | rBV3     | 13906          | 20820         | 0.51%           | 0.046%        |
| 23        | 7.846       | 805           | 809         | 813          | rVB4     | 13876          | 20557         | 0.50%           | 0.045%        |
| 24        | 7.963       | 819           | 829         | 837          | rBV      | 472995         | 751670        | 18.45%          | 1.649%        |
| 25        | 8.469       | 909           | 915         | 921          | rVV8     | 9072           | 18981         | 0.47%           | 0.042%        |
| 26        | 8.546       | 922           | 928         | 939          | rVB6     | 17502          | 43085         | 1.06%           | 0.095%        |
| 27        | 8.681       | 942           | 951         | 961          | rBV      | 829499         | 1368165       | 33.59%          | 3.002%        |
| 28        | 8.752       | 961           | 963         | 968          | rVV6     | 13671          | 23912         | 0.59%           | 0.052%        |
| 29        | 8.840       | 968           | 978         | 983          | rVV2     | 60578          | 147120        | 3.61%           | 0.323%        |
| 30        | 9.122       | 1021          | 1026        | 1029         | rBV      | 82421          | 118979        | 2.92%           | 0.261%        |
| 31        | 9.169       | 1029          | 1034        | 1040         | rVB      | 662908         | 1015956       | 24.94%          | 2.229%        |
| 32        | 9.281       | 1049          | 1053        | 1058         | rVB8     | 12408          | 20808         | 0.51%           | 0.046%        |
| 33        | 9.357       | 1062          | 1066        | 1073         | rVB4     | 14964          | 27648         | 0.68%           | 0.061%        |
| 34        | 9.646       | 1108          | 1115        | 1124         | rVB10    | 21080          | 49403         | 1.21%           | 0.108%        |
| 35        | 9.887       | 1149          | 1156        | 1165         | rVV      | 510692         | 849331        | 20.85%          | 1.864%        |
| 36        | 10.052      | 1175          | 1184        | 1187         | rBV8     | 15210          | 31285         | 0.77%           | 0.069%        |

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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-BP092023.MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |       |         |         |        |        |
|----|--------|------|------|------|-------|---------|---------|--------|--------|
| 37 | 10.310 | 1223 | 1228 | 1234 | rVB9  | 10850   | 20011   | 0.49%  | 0.044% |
| 38 | 10.416 | 1234 | 1246 | 1256 | rBV   | 848652  | 1457876 | 35.79% | 3.199% |
| 39 | 10.504 | 1257 | 1261 | 1265 | rVV4  | 18193   | 32936   | 0.81%  | 0.072% |
| 40 | 10.569 | 1269 | 1272 | 1276 | rVV6  | 13609   | 21014   | 0.52%  | 0.046% |
| 41 | 10.616 | 1276 | 1280 | 1286 | rVV5  | 16780   | 29650   | 0.73%  | 0.065% |
| 42 | 10.687 | 1286 | 1292 | 1297 | rVV   | 69611   | 113573  | 2.79%  | 0.249% |
| 43 | 10.757 | 1297 | 1304 | 1306 | rVV5  | 20608   | 45189   | 1.11%  | 0.099% |
| 44 | 10.799 | 1306 | 1311 | 1317 | rVV   | 608653  | 1032989 | 25.36% | 2.266% |
| 45 | 10.857 | 1317 | 1321 | 1329 | rVB2  | 80767   | 142085  | 3.49%  | 0.312% |
| 46 | 10.969 | 1330 | 1340 | 1349 | rBV   | 619458  | 1048821 | 25.75% | 2.301% |
| 47 | 11.204 | 1376 | 1380 | 1389 | rVB3  | 32603   | 60883   | 1.49%  | 0.134% |
| 48 | 11.404 | 1403 | 1414 | 1415 | rBV10 | 15715   | 35222   | 0.86%  | 0.077% |
| 49 | 11.581 | 1437 | 1444 | 1449 | rVB9  | 21800   | 54976   | 1.35%  | 0.121% |
| 50 | 11.640 | 1449 | 1454 | 1458 | rBV7  | 21176   | 45712   | 1.12%  | 0.100% |
| 51 | 11.734 | 1465 | 1470 | 1474 | rBV   | 46653   | 63530   | 1.56%  | 0.139% |
| 52 | 11.787 | 1474 | 1479 | 1484 | rBV4  | 31547   | 59184   | 1.45%  | 0.130% |
| 53 | 11.834 | 1484 | 1487 | 1493 | rVB8  | 14991   | 24028   | 0.59%  | 0.053% |
| 54 | 11.940 | 1500 | 1505 | 1510 | rBV5  | 36448   | 64814   | 1.59%  | 0.142% |
| 55 | 12.004 | 1511 | 1516 | 1521 | rBV2  | 50345   | 88310   | 2.17%  | 0.194% |
| 56 | 12.146 | 1536 | 1540 | 1545 | rVB   | 108778  | 143203  | 3.52%  | 0.314% |
| 57 | 12.275 | 1559 | 1562 | 1564 | rBV4  | 18147   | 19458   | 0.48%  | 0.043% |
| 58 | 12.322 | 1565 | 1570 | 1579 | rVB   | 216213  | 357812  | 8.78%  | 0.785% |
| 59 | 12.469 | 1589 | 1595 | 1604 | rVB3  | 80540   | 175939  | 4.32%  | 0.386% |
| 60 | 12.551 | 1605 | 1609 | 1613 | rBV2  | 33679   | 60587   | 1.49%  | 0.133% |
| 61 | 12.687 | 1628 | 1632 | 1637 | rVB   | 52314   | 76428   | 1.88%  | 0.168% |
| 62 | 12.893 | 1664 | 1667 | 1671 | rVB5  | 16611   | 23947   | 0.59%  | 0.053% |
| 63 | 12.957 | 1675 | 1678 | 1685 | rVV5  | 31369   | 55196   | 1.36%  | 0.121% |
| 64 | 13.034 | 1686 | 1691 | 1696 | rVV5  | 55027   | 109660  | 2.69%  | 0.241% |
| 65 | 13.098 | 1698 | 1702 | 1709 | rVB2  | 54151   | 97880   | 2.40%  | 0.215% |
| 66 | 13.193 | 1714 | 1718 | 1726 | rVB5  | 34683   | 63669   | 1.56%  | 0.140% |
| 67 | 13.310 | 1735 | 1738 | 1742 | rVB6  | 17481   | 21205   | 0.52%  | 0.047% |
| 68 | 13.398 | 1742 | 1753 | 1758 | rBV3  | 154357  | 362435  | 8.90%  | 0.795% |
| 69 | 13.493 | 1762 | 1769 | 1778 | rVV   | 588046  | 989150  | 24.28% | 2.170% |
| 70 | 13.798 | 1817 | 1821 | 1823 | rBV4  | 34045   | 51103   | 1.25%  | 0.112% |
| 71 | 13.963 | 1845 | 1849 | 1853 | rBV3  | 44883   | 76444   | 1.88%  | 0.168% |
| 72 | 14.016 | 1853 | 1858 | 1861 | rVV2  | 264449  | 397028  | 9.75%  | 0.871% |
| 73 | 14.069 | 1861 | 1867 | 1874 | rVB   | 1443230 | 2011046 | 49.37% | 4.412% |
| 74 | 14.351 | 1907 | 1915 | 1922 | rBV   | 1780318 | 2954013 | 72.52% | 6.481% |
| 75 | 14.428 | 1922 | 1928 | 1933 | rVB2  | 278903  | 428840  | 10.53% | 0.941% |
| 76 | 14.481 | 1933 | 1937 | 1940 | rBV   | 167306  | 233656  | 5.74%  | 0.513% |

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Integrator: RTE

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Min Area: 0.1 % of largest Peak

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |        |
|-----|--------|------|------|------|------|---------|---------|---------|--------|
| 77  | 14.575 | 1949 | 1953 | 1958 | rBV3 | 82522   | 113187  | 2.78%   | 0.248% |
| 78  | 14.657 | 1962 | 1967 | 1972 | rBV  | 927810  | 1303813 | 32.01%  | 2.861% |
| 79  | 14.892 | 2003 | 2007 | 2013 | rBV  | 244942  | 425022  | 10.43%  | 0.933% |
| 80  | 15.051 | 2030 | 2034 | 2038 | rVB3 | 55569   | 106722  | 2.62%   | 0.234% |
| 81  | 15.228 | 2061 | 2064 | 2065 | rBV3 | 48129   | 51206   | 1.26%   | 0.112% |
| 82  | 15.257 | 2066 | 2069 | 2076 | rVB6 | 99228   | 152608  | 3.75%   | 0.335% |
| 83  | 15.322 | 2077 | 2080 | 2087 | rVB7 | 87025   | 171291  | 4.21%   | 0.376% |
| 84  | 15.398 | 2088 | 2093 | 2097 | rBV  | 247796  | 402458  | 9.88%   | 0.883% |
| 85  | 15.439 | 2097 | 2100 | 2102 | rBV  | 122060  | 128323  | 3.15%   | 0.282% |
| 86  | 15.657 | 2132 | 2137 | 2145 | rBV  | 2297519 | 3195918 | 78.46%  | 7.012% |
| 87  | 15.839 | 2165 | 2168 | 2174 | rVB2 | 136494  | 201417  | 4.94%   | 0.442% |
| 88  | 16.004 | 2193 | 2196 | 2199 | rBV2 | 75951   | 110419  | 2.71%   | 0.242% |
| 89  | 16.322 | 2245 | 2250 | 2256 | rBV  | 503040  | 897917  | 22.04%  | 1.970% |
| 90  | 16.698 | 2311 | 2314 | 2324 | rVB6 | 177345  | 316431  | 7.77%   | 0.694% |
| 91  | 17.163 | 2389 | 2393 | 2398 | rVB  | 510163  | 688117  | 16.89%  | 1.510% |
| 92  | 17.445 | 2436 | 2441 | 2445 | rBV  | 1138073 | 1654954 | 40.63%  | 3.631% |
| 93  | 17.486 | 2445 | 2448 | 2452 | rVB  | 461979  | 621264  | 15.25%  | 1.363% |
| 94  | 17.545 | 2453 | 2458 | 2465 | rBV  | 2218354 | 3363287 | 82.57%  | 7.379% |
| 95  | 17.869 | 2511 | 2513 | 2518 | rVB  | 424097  | 465570  | 11.43%  | 1.022% |
| 96  | 18.292 | 2582 | 2585 | 2590 | rBV4 | 526719  | 855357  | 21.00%  | 1.877% |
| 97  | 19.286 | 2751 | 2754 | 2757 | rBV  | 904643  | 1210438 | 29.72%  | 2.656% |
| 98  | 19.469 | 2782 | 2785 | 2788 | rVB  | 885520  | 982957  | 24.13%  | 2.157% |
| 99  | 19.798 | 2837 | 2841 | 2845 | rBV  | 2868829 | 4073391 | 100.00% | 8.937% |
| 100 | 20.204 | 2907 | 2910 | 2914 | rVB2 | 1960459 | 2337392 | 57.38%  | 5.129% |

Sum of corrected areas: 45576447



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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
Quant Title : SVOA CALIBRATION

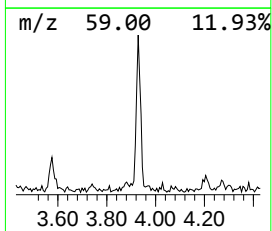
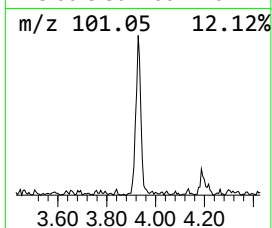
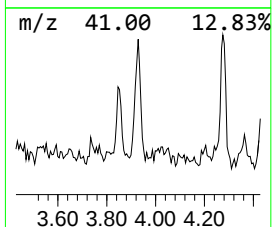
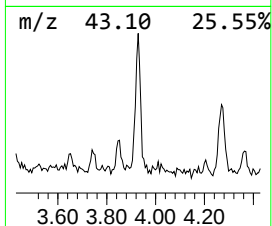
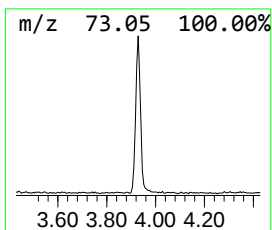
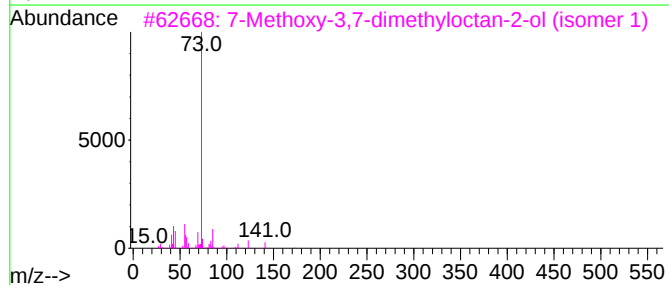
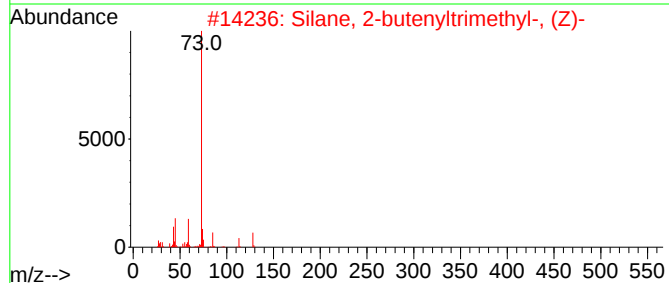
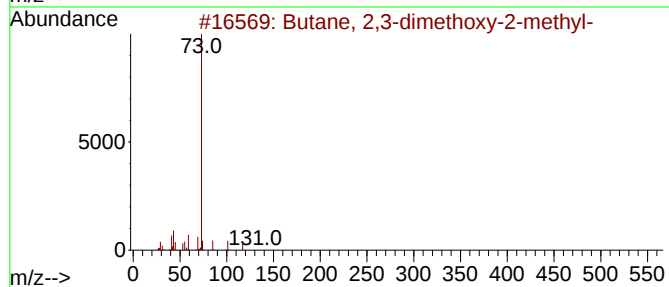
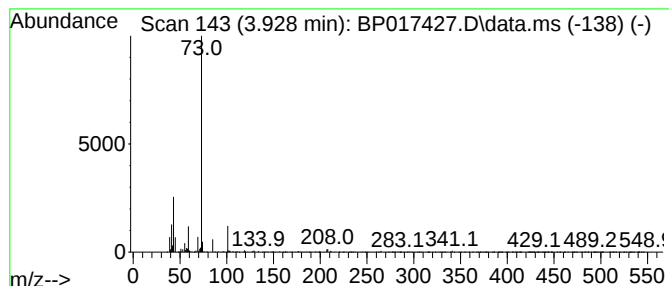
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Butane, 2,3-dimethoxy-2-met... Concentration Rank 16

| R.T.  | EstConc    | Area  | Relative to ISTD       | R.T.  |
|-------|------------|-------|------------------------|-------|
| 3.928 | 2.61 ng/ul | 98086 | 1,4-Dichlorobenzene-d4 | 7.963 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Butane, 2,3-dimethoxy-2-methyl-     | 132 | C7H16O2  | 074421-00-4  | 50   |
| 2    |    |   | Silane, 2-butenyltrimethyl-, (Z)-   | 128 | C7H16Si  | 017486-13-4  | 25   |
| 3    |    |   | 7-Methoxy-3,7-dimethyloctan-2-ol... | 188 | C11H24O2 | 1000511-09-7 | 17   |
| 4    |    |   | 1-(2-Methoxyethoxy)-2-methyl-2-p... | 162 | C8H18O3  | 1000364-24-4 | 14   |
| 5    |    |   | Silane, trimethyl(1-methyl-1-pro... | 128 | C7H16Si  | 010111-13-4  | 12   |



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Quant Title : SVOA CALIBRATION

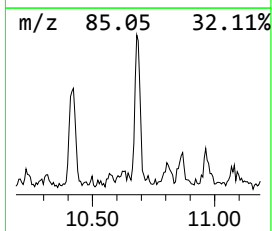
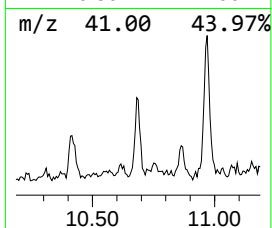
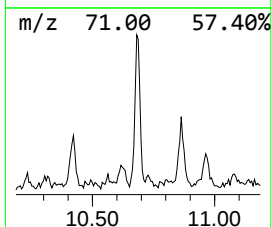
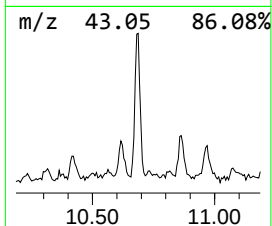
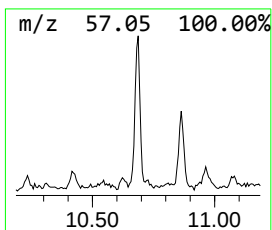
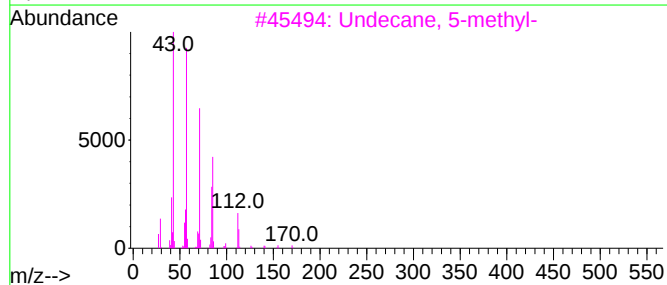
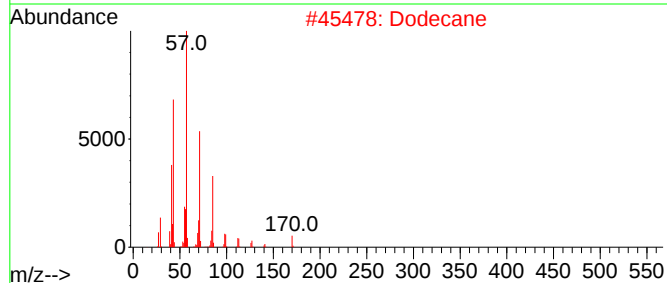
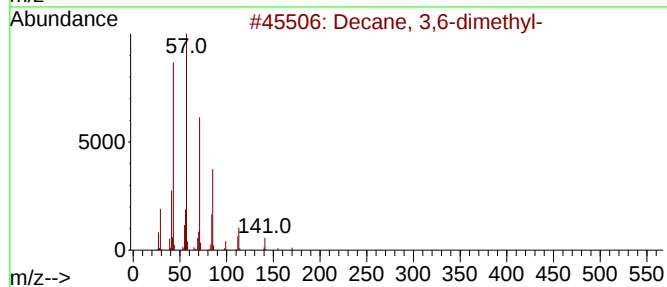
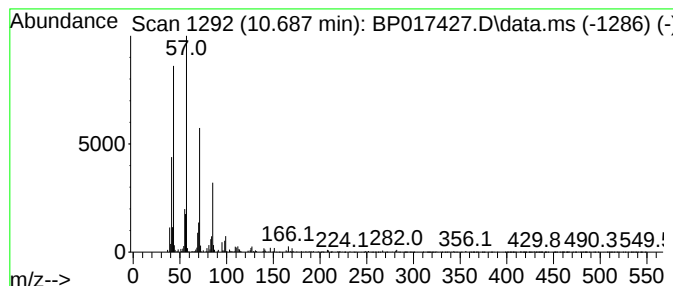
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 18

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 10.687 | 2.20 ng/ul | 113573 | Naphthalene-d8   | 10.804 |

| Hit# | of 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|------|------------------------------------|-----|---------|-------------|------|
| 1    |      | Decane, 3,6-dimethyl-              | 170 | C12H26  | 017312-53-7 | 87   |
| 2    |      | Dodecane                           | 170 | C12H26  | 000112-40-3 | 87   |
| 3    |      | Undecane, 5-methyl-                | 170 | C12H26  | 001632-70-8 | 87   |
| 4    |      | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8 | 83   |
| 5    |      | Tetradecane                        | 198 | C14H30  | 000629-59-4 | 81   |



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ClientSampleId :  
EZY6

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Quant Title : SVOA CALIBRATION

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

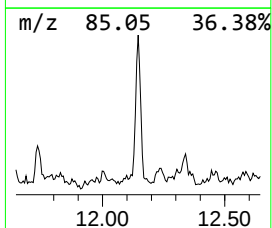
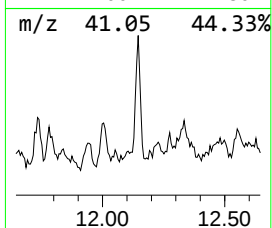
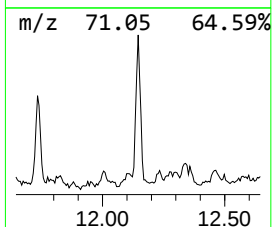
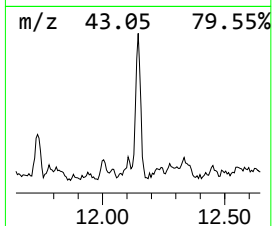
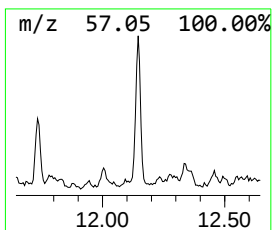
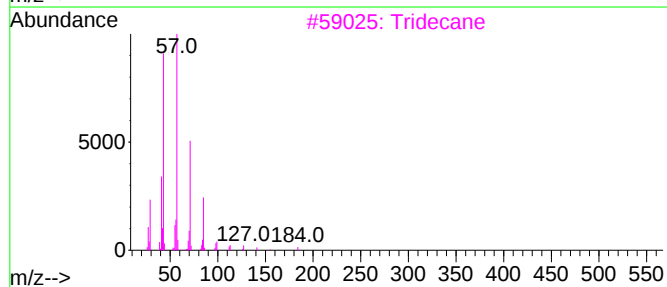
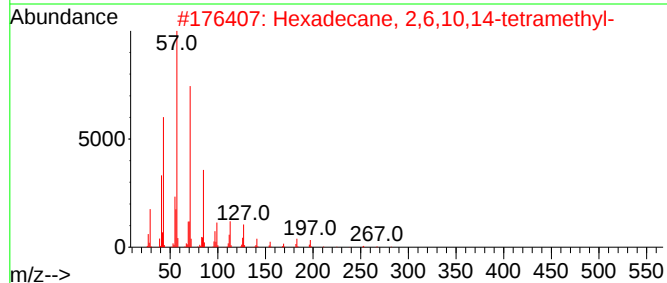
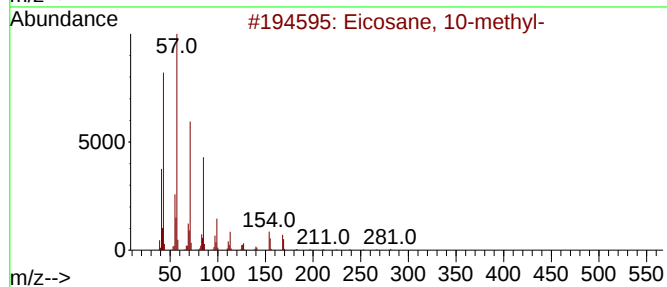
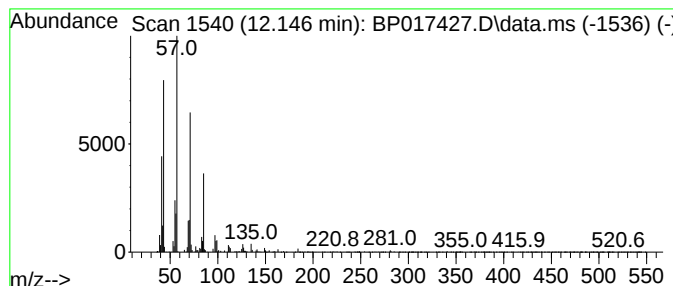
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Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 12.146 | 2.77 ng/ul | 143203 | Naphthalene-d8   | 10.804 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Eicosane, 10-methyl-                | 296 | C21H44  | 054833-23-7 | 87   |
| 2    |      | Hexadecane, 2,6,10,14-tetramethyl-  | 282 | C20H42  | 000638-36-8 | 87   |
| 3    |      | Tridecane                           | 184 | C13H28  | 000629-50-5 | 87   |
| 4    |      | Hexadecane                          | 226 | C16H34  | 000544-76-3 | 83   |
| 5    |      | Heptadecane, 2,6,10,14-tetramethyl- | 296 | C21H44  | 018344-37-1 | 81   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092823\  
Data File : BP017427.D  
Acq On : 28 Sep 2023 23:10  
Operator : MA/JU  
Sample : 04417-39  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EZY6

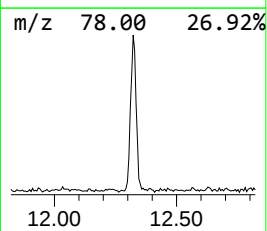
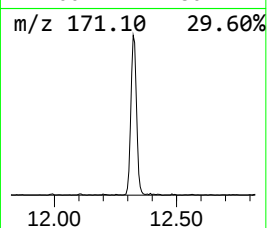
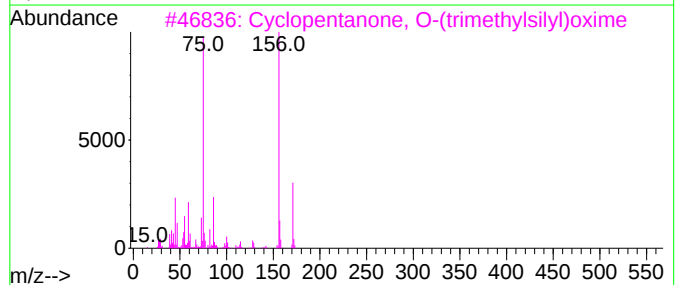
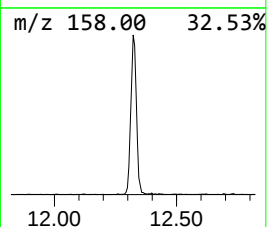
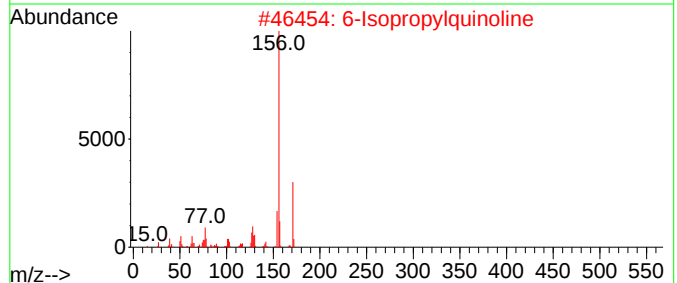
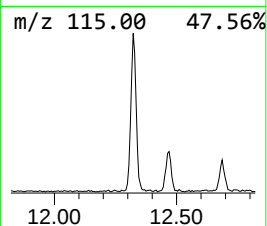
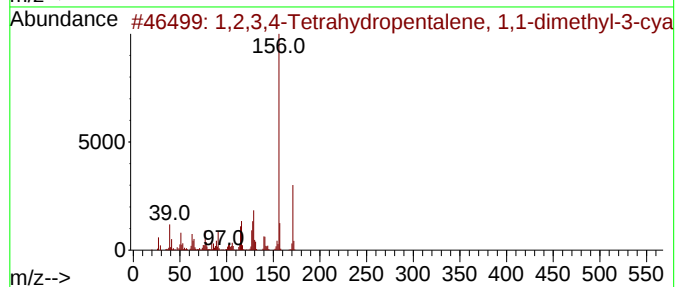
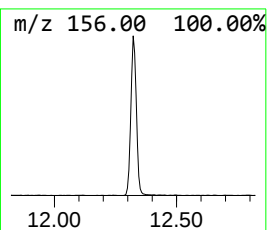
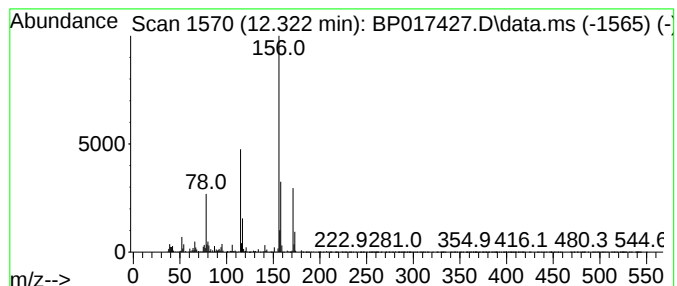
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Quant Title : SVOA CALIBRATION

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

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Peak Number 4 unknown-01 Concentration Rank 7

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 12.322 | 6.93 ng/ul | 357812 | Naphthalene-d8   | 10.804 |

| Hit# | of                                   | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|--------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 1,2,3,4-Tetrahydropentalene, 1,1...  | 171 | C12H13N      | 1000217-18-6 | 43      |      |      |
| 2    | 6-Isopropylquinoline                 | 171 | C12H13N      | 000135-79-5  | 38      |      |      |
| 3    | Cyclopentanone, O-(trimethylsilyl... | 171 | C8H17NOSi    | 026208-87-7  | 32      |      |      |
| 4    | Chloroxylenol                        | 156 | C8H9ClO      | 000088-04-0  | 32      |      |      |
| 5    | 2-Chloro-4,6-dimethyl-3-pyridina...  | 156 | C7H9ClN2     | 140413-40-7  | 32      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092823\  
Data File : BP017427.D  
Acq On : 28 Sep 2023 23:10  
Operator : MA/JU  
Sample : 04417-39  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
Quant Title : SVOA CALIBRATION

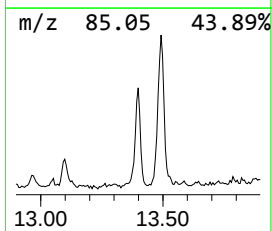
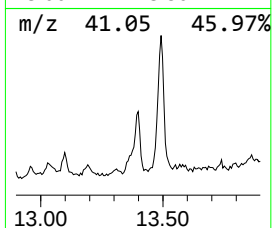
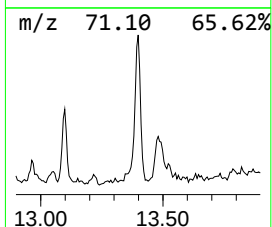
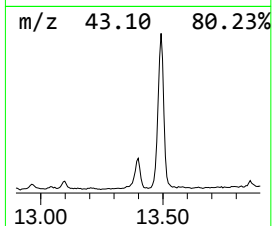
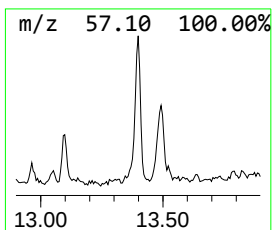
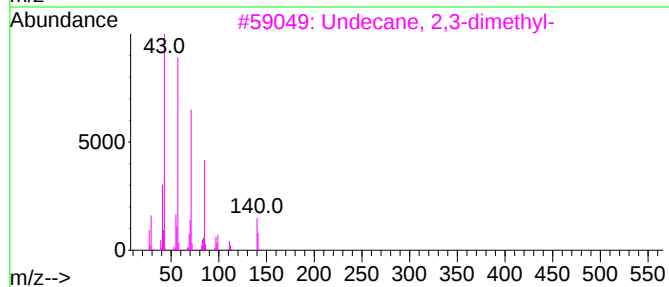
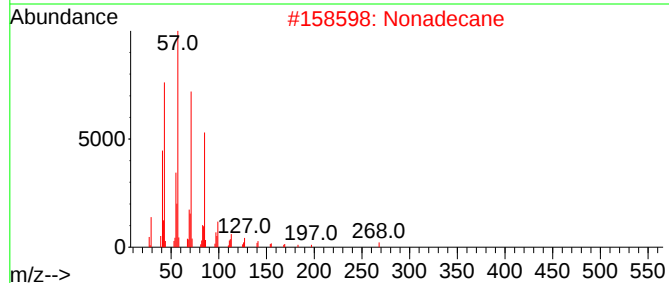
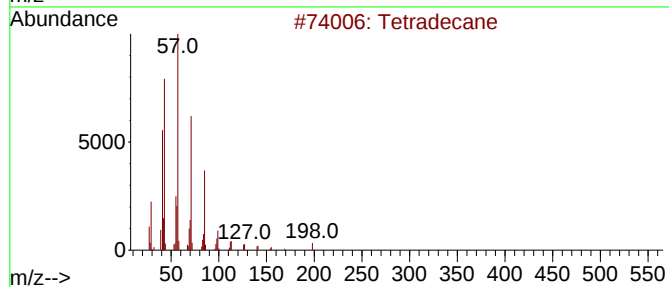
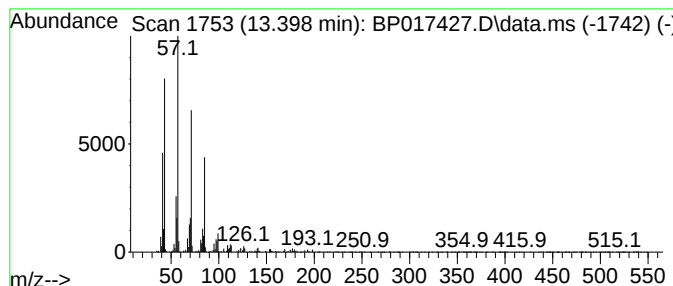
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 12

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.398 | 5.56 ng/ul | 362435 | Acenaphthene-d10 | 14.657 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm  | CAS#         | Qual |
|------|----|---|------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Tetradecane                        | 198 | C14H30   | 000629-59-4  | 93   |
| 2    |    |   | Nonadecane                         | 268 | C19H40   | 000629-92-5  | 91   |
| 3    |    |   | Undecane, 2,3-dimethyl-            | 184 | C13H28   | 017312-77-5  | 90   |
| 4    |    |   | Pentadecane                        | 212 | C15H32   | 000629-62-9  | 90   |
| 5    |    |   | Carbonic acid, eicosyl vinyl ester | 368 | C23H44O3 | 1000382-54-3 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092823\  
Data File : BP017427.D  
Acq On : 28 Sep 2023 23:10  
Operator : MA/JU  
Sample : 04417-39  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EZY6

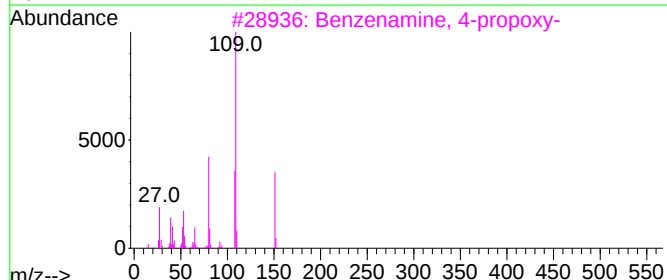
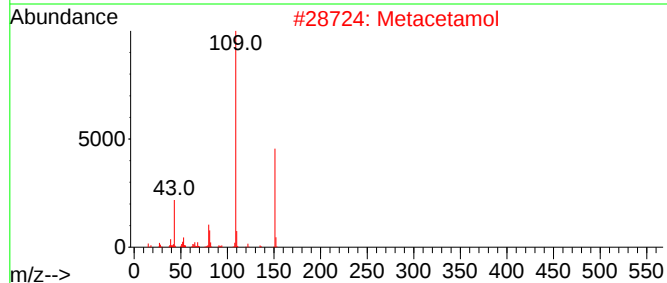
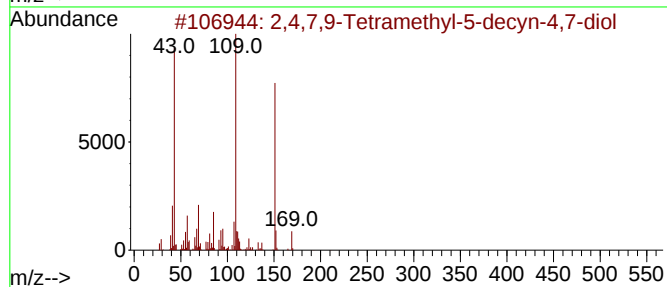
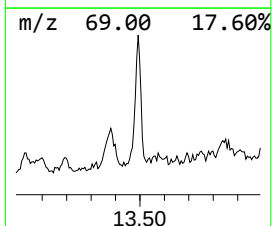
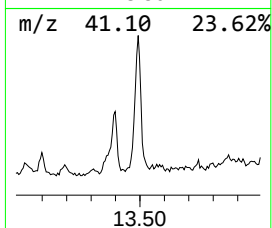
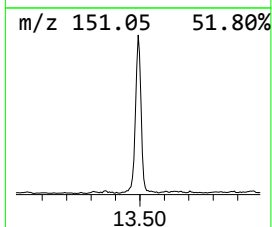
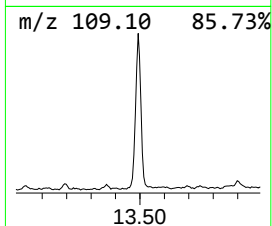
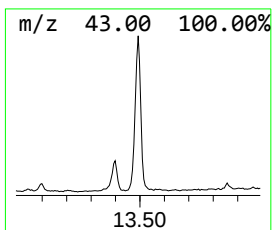
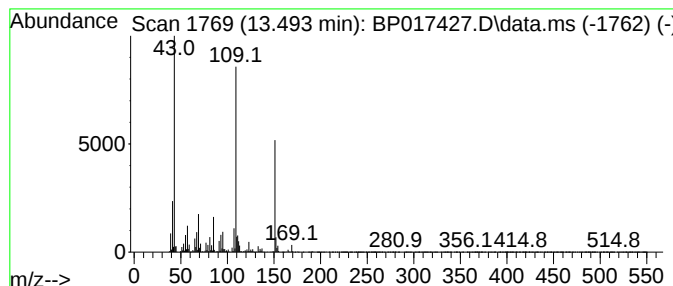
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Quant Title : SVOA CALIBRATION

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

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Peak Number 6 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 2

| R.T.      | EstConc                             | Area   | Relative to ISTD |          | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|----------|-------------|------|
| 13.493    | 15.17 ng/ul                         | 989150 | Acenaphthene-d10 |          | 14.657      |      |
| Hit# of 5 | Tentative ID                        |        | MW               | MolForm  | CAS#        | Qual |
| 1         | 2,4,7,9-Tetramethyl-5-decyn-4,7-... |        | 226              | C14H26O2 | 000126-86-3 | 91   |
| 2         | Metacetamol                         |        | 151              | C8H9NO2  | 000621-42-1 | 59   |
| 3         | Benzenamine, 4-propoxy-             |        | 151              | C9H13NO  | 004469-80-1 | 59   |
| 4         | 3-Pyridinol, 4-methyl-, acetate ... |        | 151              | C8H9NO2  | 001006-96-8 | 59   |
| 5         | 2,4,7,9-Tetramethyl-5-decyn-4,7-... |        | 226              | C14H26O2 | 000126-86-3 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092823\  
Data File : BP017427.D  
Acq On : 28 Sep 2023 23:10  
Operator : MA/JU  
Sample : 04417-39  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EZY6

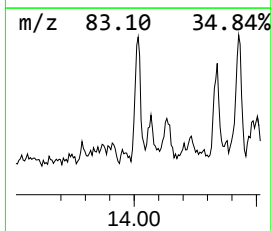
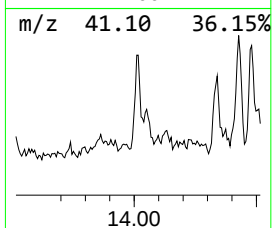
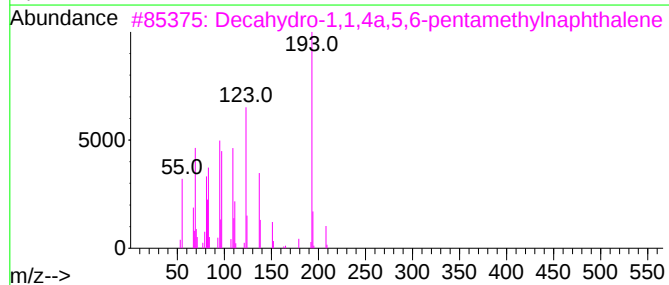
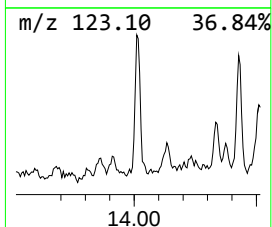
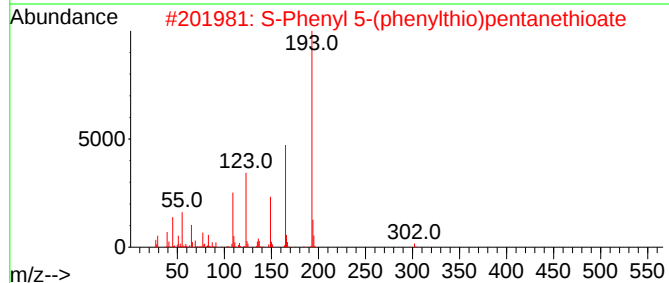
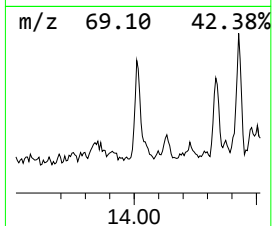
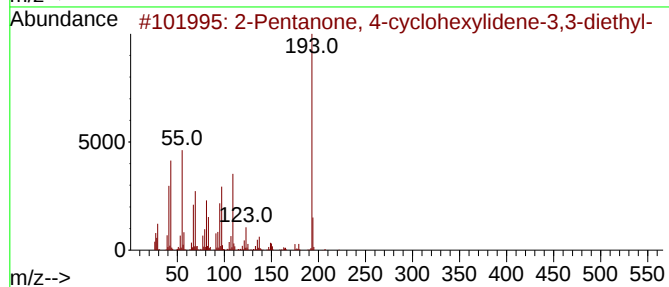
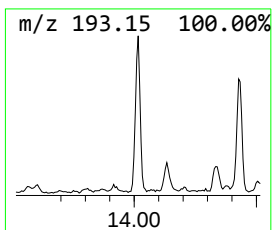
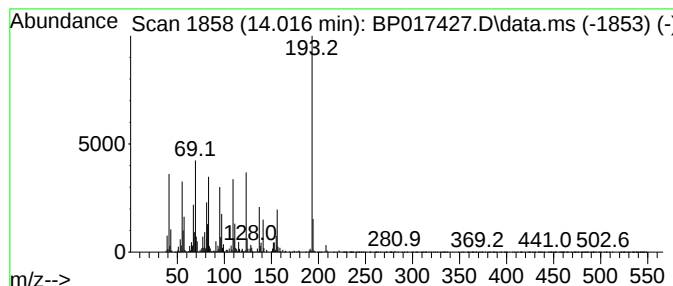
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Quant Title : SVOA CALIBRATION

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

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Peak Number 7 unknown-02 Concentration Rank 10

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 14.016 | 6.09 ng/ul | 397028 | Acenaphthene-d10 | 14.657 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 2-Pentanone, 4-cyclohexylidene-3... | 222 | C15H26O   | 313253-65-5  | 40   |
| 2    |    |   | S-Phenyl 5-(phenylthio)pentaneth... | 302 | C17H18OS2 | 1000234-40-7 | 38   |
| 3    |    |   | Decahydro-1,1,4a,5,6-pentamethyl... | 208 | C15H28    | 080655-44-3  | 37   |
| 4    |    |   | 3H-3,10a-Methano-1,2-benzodioxoc... | 226 | C13H22O3  | 095906-83-5  | 36   |
| 5    |    |   | 1H-Indene, octahydro-2,2,4,4,7,7... | 208 | C15H28    | 054832-83-6  | 36   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092823\  
Data File : BP017427.D  
Acq On : 28 Sep 2023 23:10  
Operator : MA/JU  
Sample : 04417-39  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

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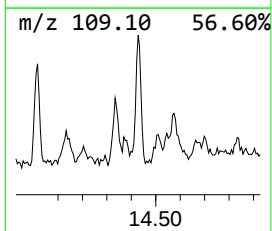
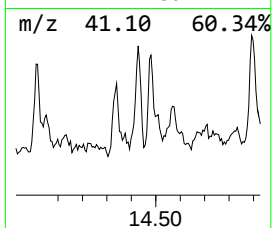
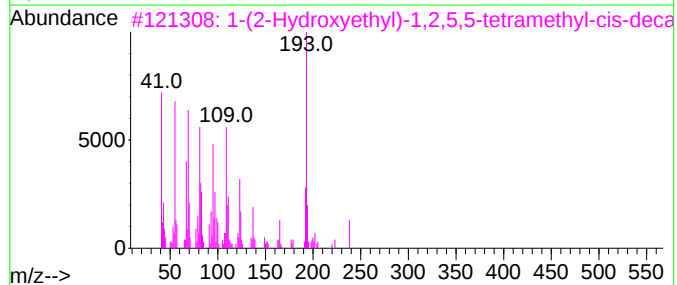
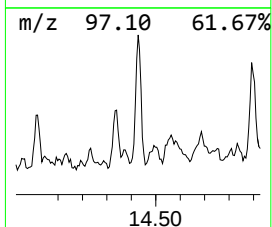
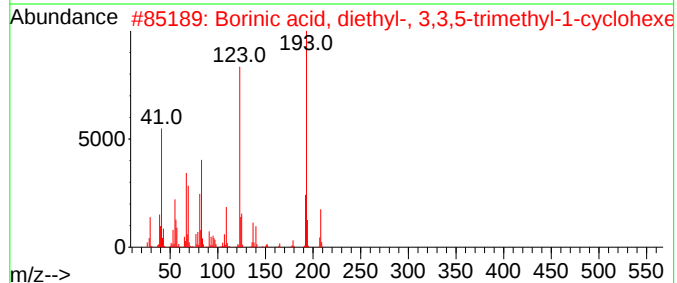
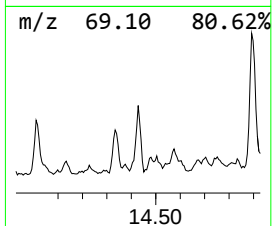
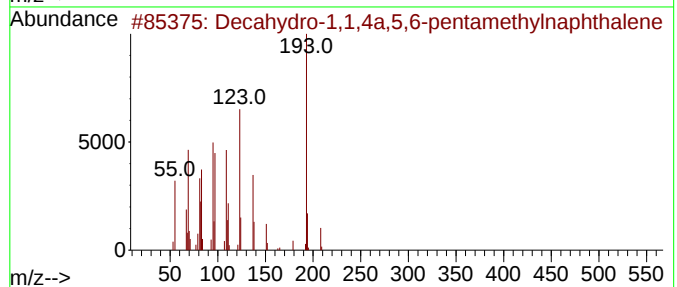
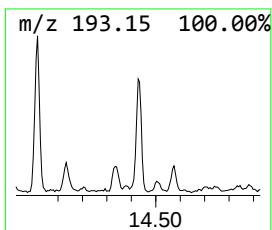
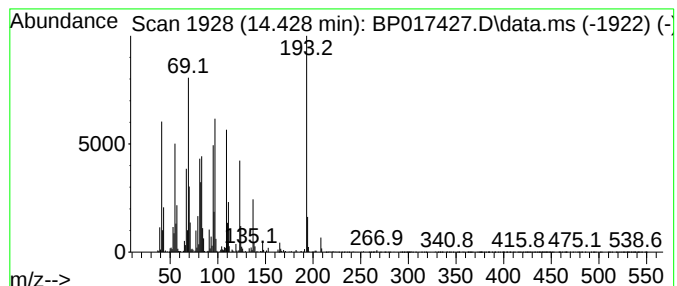
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Quant Title : SVOA CALIBRATION

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

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Peak Number 8 unknown-03 Concentration Rank 8

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 14.428 | 6.58 ng/ul | 428840 | Acenaphthene-d10 | 14.657 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Decahydro-1,1,4a,5,6-pentamethyl... | 208 | C15H28      | 080655-44-3  | 38   |
| 2    |    |   | Borinic acid, diethyl-, 3,3,5-tr... | 208 | C13H25BO    | 057387-76-5  | 35   |
| 3    |    |   | 1-(2-Hydroxyethyl)-1,2,5,5-tetra... | 238 | C16H30O     | 1010298-97-4 | 32   |
| 4    |    |   | 3-((2S,5S)-7-(Methylsulfonyl)-1-... | 272 | C12H20N2O3S | 1465868-22-7 | 27   |
| 5    |    |   | Silane, chlorodiethyl(2-methylpe... | 222 | C10H23ClOSi | 1000363-79-1 | 16   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092823\  
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Operator : MA/JU  
Sample : 04417-39  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EZY6

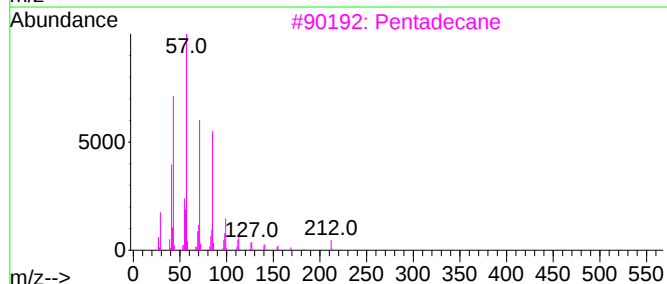
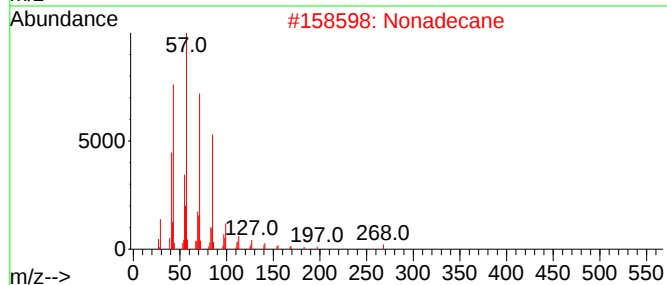
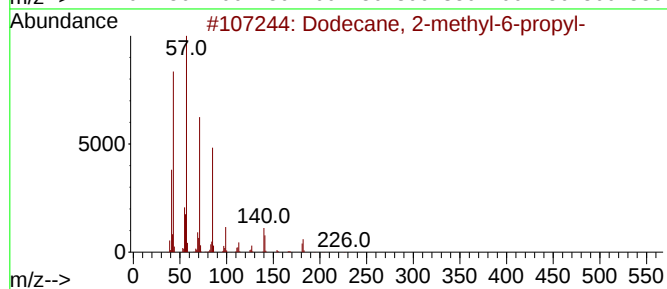
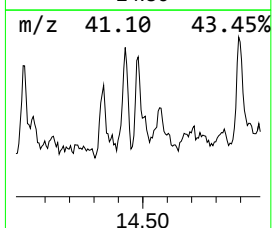
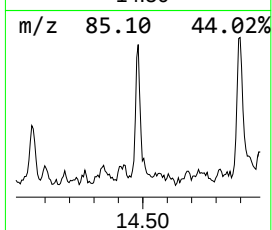
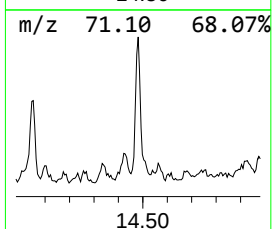
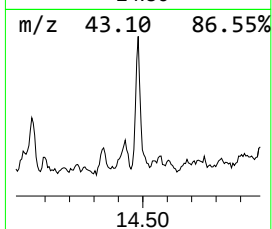
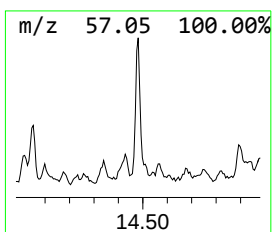
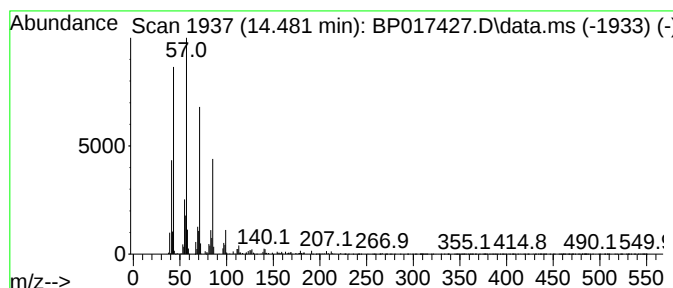
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Quant Title : SVOA CALIBRATION

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

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Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 13

| R.T.      | EstConc                      | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|------------------------------|--------|------------------|---------|-------------|------|
| 14.481    | 3.58 ng/ul                   | 233656 | Acenaphthene-d10 |         | 14.657      |      |
| Hit# of 5 | Tentative ID                 |        | MW               | MolForm | CAS#        | Qual |
| 1         | Dodecane, 2-methyl-6-propyl- |        | 226              | C16H34  | 055045-08-4 | 90   |
| 2         | Nonadecane                   |        | 268              | C19H40  | 000629-92-5 | 90   |
| 3         | Pentadecane                  |        | 212              | C15H32  | 000629-62-9 | 90   |
| 4         | Pentadecane                  |        | 212              | C15H32  | 000629-62-9 | 87   |
| 5         | Tridecane, 7-propyl-         |        | 226              | C16H34  | 055045-09-5 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092823\  
Data File : BP017427.D  
Acq On : 28 Sep 2023 23:10  
Operator : MA/JU  
Sample : 04417-39  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EZY6

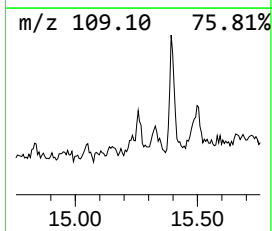
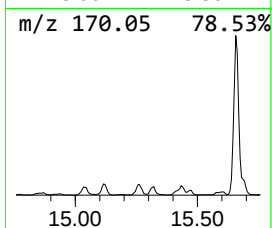
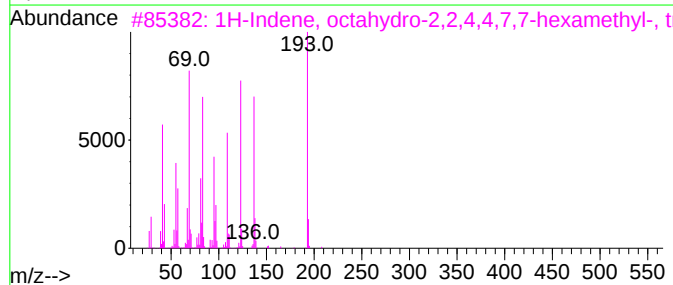
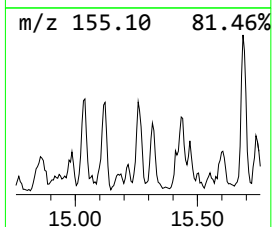
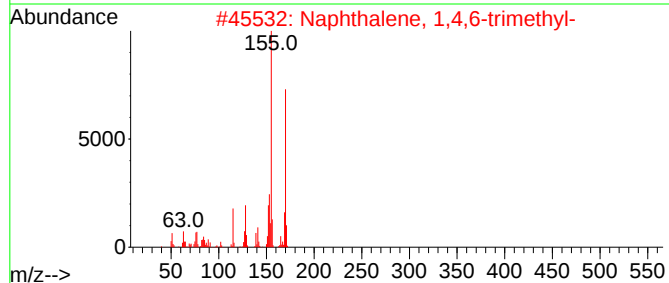
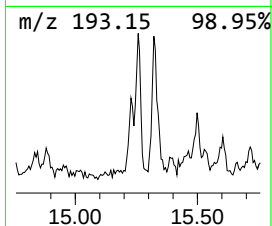
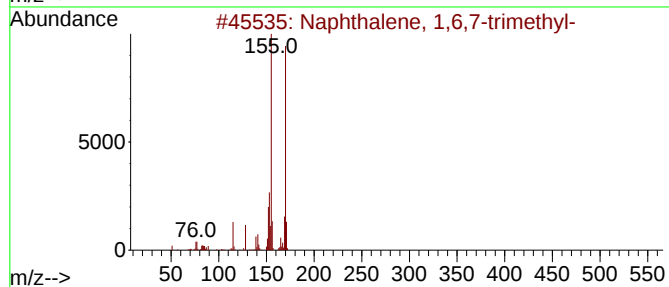
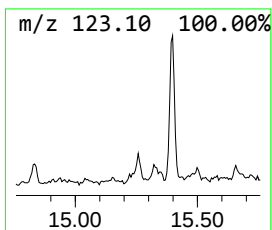
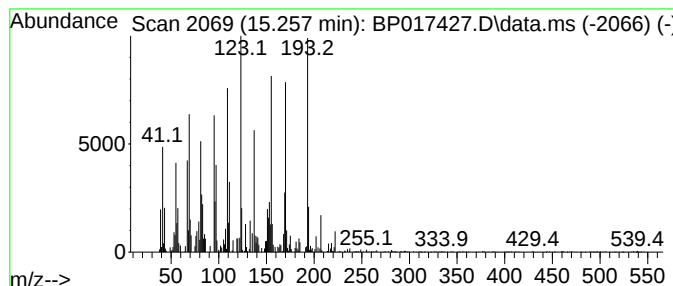
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Quant Title : SVOA CALIBRATION

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

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Peak Number 10 unknown-04 Concentration Rank 17

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 15.257 | 2.34 ng/ul | 152608 | Acenaphthene-d10 | 14.657 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,6,7-trimethyl-       | 170 | C13H14  | 002245-38-7 | 47   |
| 2    |    |   | Naphthalene, 1,4,6-trimethyl-       | 170 | C13H14  | 002131-42-2 | 43   |
| 3    |    |   | 1H-Indene, octahydro-2,2,4,4,7,7... | 208 | C15H28  | 054832-83-6 | 42   |
| 4    |    |   | Neopentylidenecyclohexane           | 152 | C11H20  | 039546-80-0 | 35   |
| 5    |    |   | Naphthalene, 1,6,7-trimethyl-       | 170 | C13H14  | 002245-38-7 | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092823\  
Data File : BP017427.D  
Acq On : 28 Sep 2023 23:10  
Operator : MA/JU  
Sample : 04417-39  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EZY6

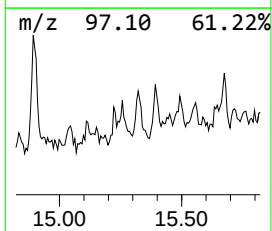
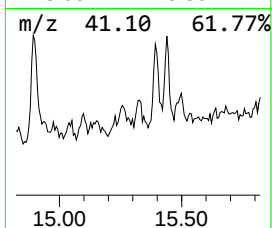
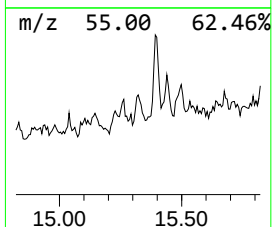
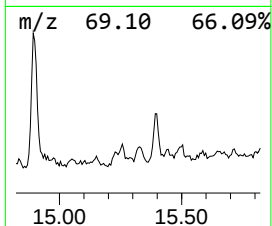
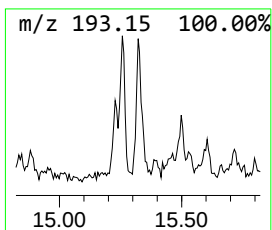
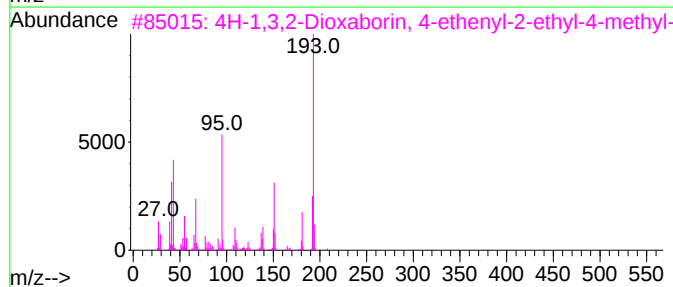
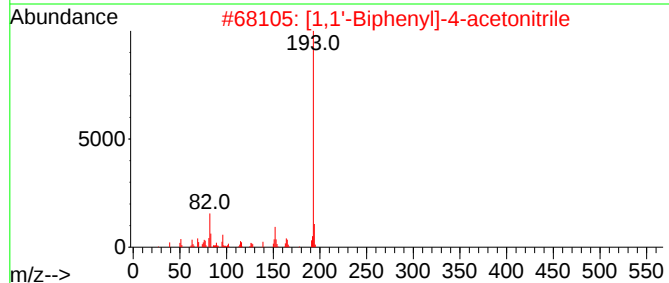
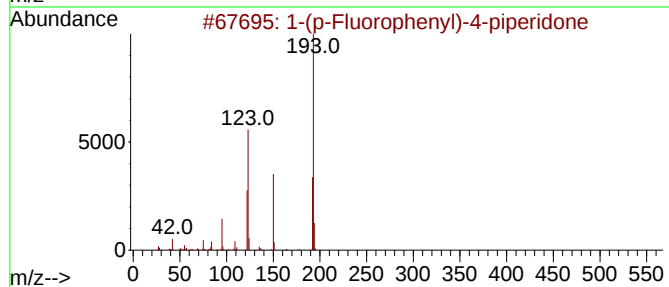
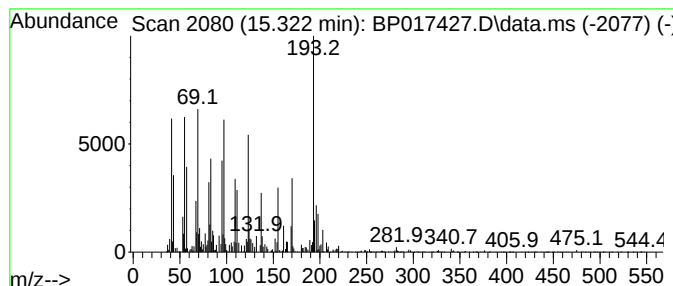
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Quant Title : SVOA CALIBRATION

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

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Peak Number 11 unknown-05 Concentration Rank 15

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 15.322    | 2.63 ng/ul                          | 171291 | Acenaphthene-d10 | 14.657       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 1-(p-Fluorophenyl)-4-piperidone     | 193    | C11H12FNO        | 1000238-56-7 | 38   |
| 2         | [1,1'-Biphenyl]-4-acetonitrile      | 193    | C14H11N          | 031603-77-7  | 35   |
| 3         | 4H-1,3,2-Dioxaborin, 4-ethenyl-2... | 208    | C12H21BO2        | 074630-04-9  | 35   |
| 4         | 2-Anthracenamine                    | 193    | C14H11N          | 000613-13-8  | 27   |
| 5         | 9-Phenanthrenamine                  | 193    | C14H11N          | 000947-73-9  | 27   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092823\  
Data File : BP017427.D  
Acq On : 28 Sep 2023 23:10  
Operator : MA/JU  
Sample : 04417-39  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EZY6

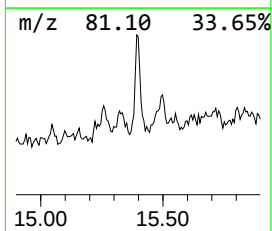
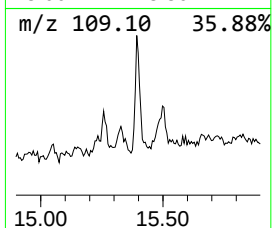
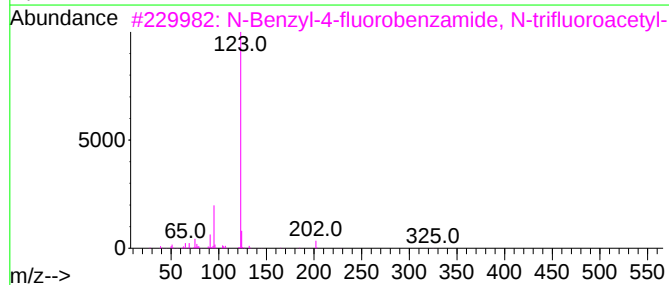
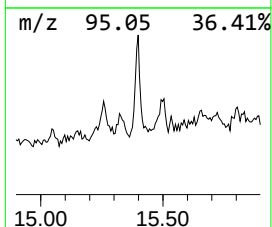
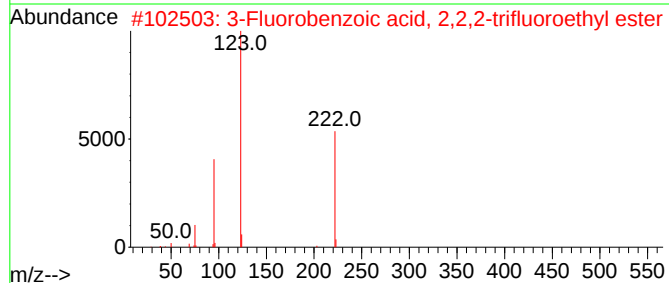
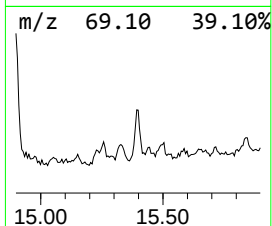
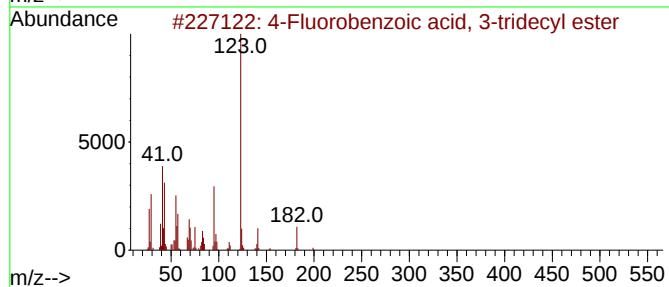
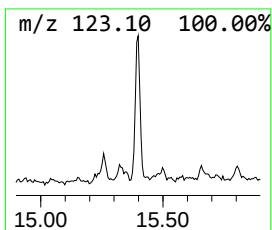
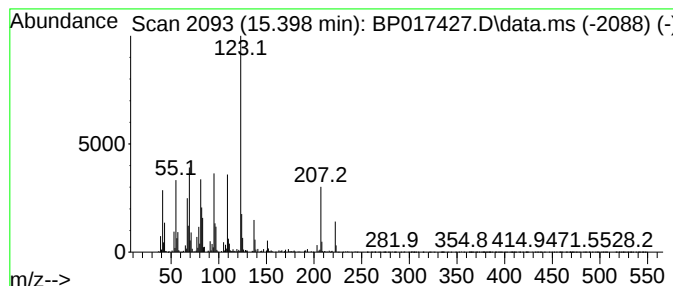
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Quant Title : SVOA CALIBRATION

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

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Peak Number 12 4-Fluorobenzoic acid, 3-tri... Concentration Rank 9

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 15.398 | 6.17 ng/ul | 402458 | Acenaphthene-d10 | 14.657 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | 4-Fluorobenzoic acid, 3-tridecyl... | 322 | C20H31F02   | 1000280-67-9 | 50   |
| 2    |    |   | 3-Fluorobenzoic acid, 2,2,2-trif... | 222 | C9H6F4O2    | 1000325-53-8 | 49   |
| 3    |    |   | N-Benzyl-4-fluorobenzamide, N-tr... | 325 | C16H11F4N02 | 1000492-36-8 | 43   |
| 4    |    |   | Bicyclo[4.1.0]heptane-7-carboxam... | 222 | C11H14N2O5  | 329912-72-3  | 43   |
| 5    |    |   | 4-Diacetylamino-3-methylphenyl a... | 249 | C13H15N04   | 1000432-47-5 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092823\  
Data File : BP017427.D  
Acq On : 28 Sep 2023 23:10  
Operator : MA/JU  
Sample : 04417-39  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EZY6

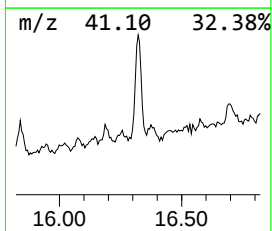
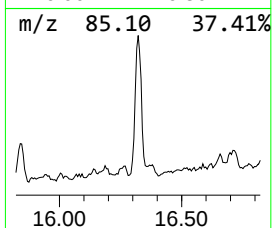
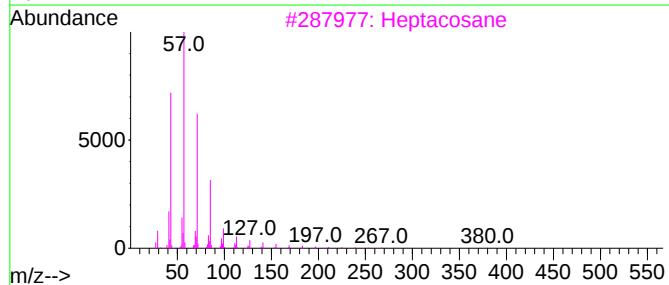
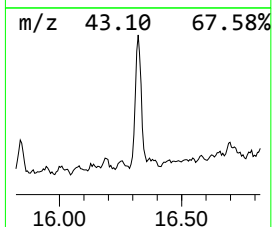
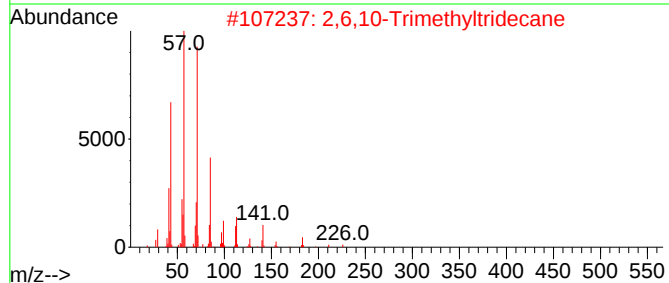
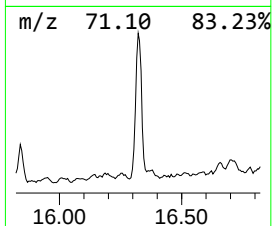
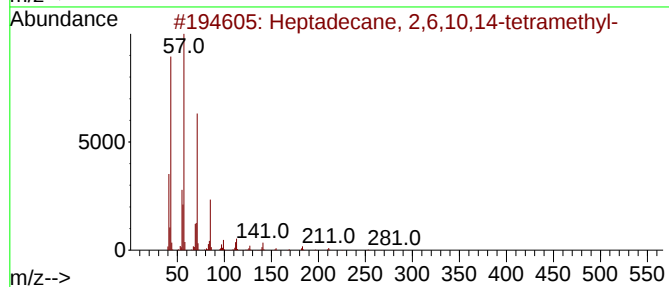
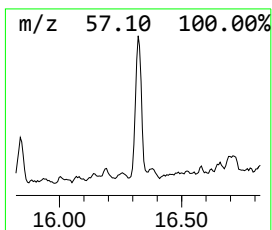
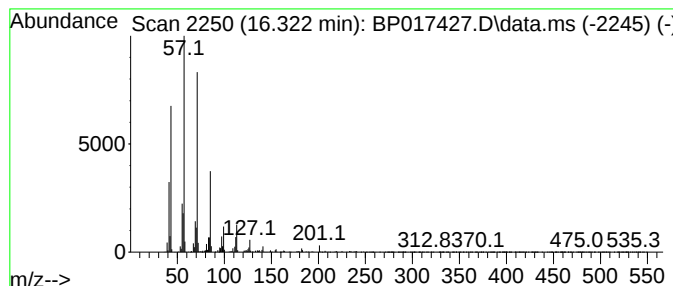
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Quant Title : SVOA CALIBRATION

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

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Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 16.322 | 10.85 ng/ul | 897917 | Phenanthrene-d10 | 17.445 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Heptadecane, 2,6,10,14-tetramethyl- | 296 | C21H44  | 018344-37-1 | 91   |
| 2    |      | 2,6,10-Trimethyltridecane           | 226 | C16H34  | 003891-99-4 | 91   |
| 3    |      | Heptacosane                         | 380 | C27H56  | 000593-49-7 | 91   |
| 4    |      | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 90   |
| 5    |      | Hexadecane, 2,6,10,14-tetramethyl-  | 282 | C20H42  | 000638-36-8 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092823\  
Data File : BP017427.D  
Acq On : 28 Sep 2023 23:10  
Operator : MA/JU  
Sample : 04417-39  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
Quant Title : SVOA CALIBRATION

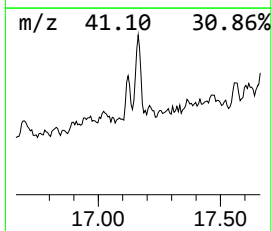
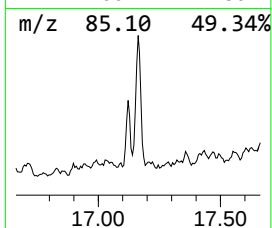
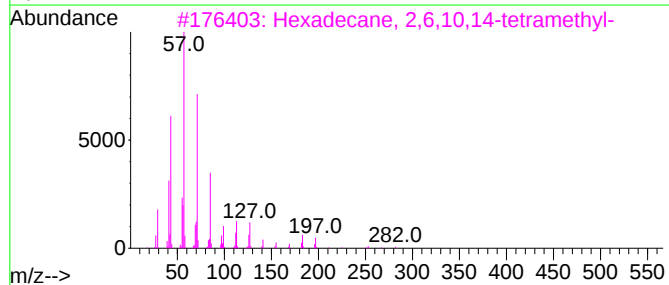
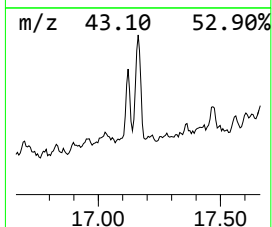
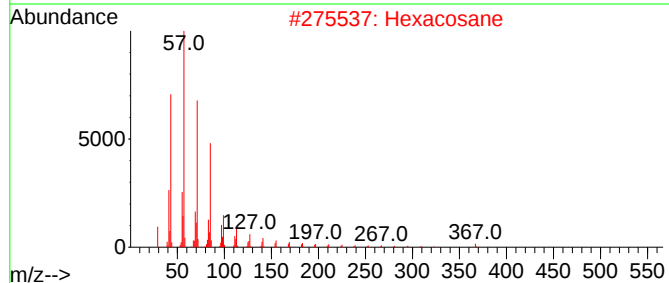
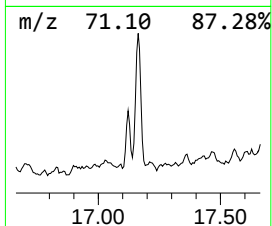
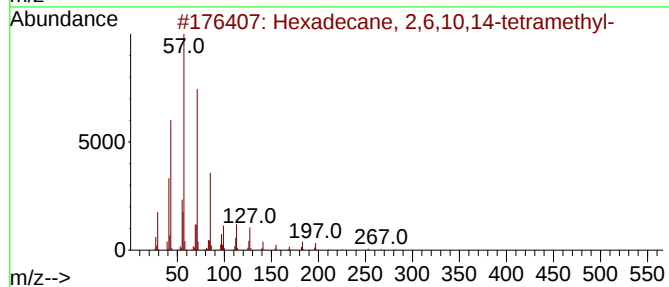
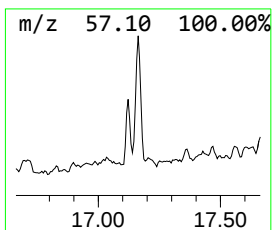
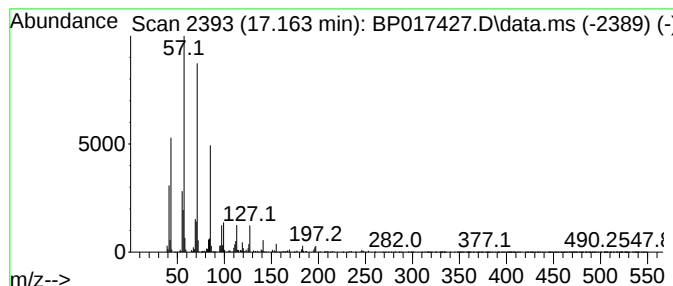
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TIC Integration Parameters: LSCINT.P

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Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.163 | 8.32 ng/ul | 688117 | Phenanthrene-d10 | 17.445 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane, 2,6,10,14-tetramethyl-  | 282 | C20H42  | 000638-36-8 | 93   |
| 2    |    |   | Hexacosane                          | 366 | C26H54  | 000630-01-3 | 91   |
| 3    |    |   | Hexadecane, 2,6,10,14-tetramethyl-  | 282 | C20H42  | 000638-36-8 | 91   |
| 4    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 87   |
| 5    |    |   | Pentacosane                         | 352 | C25H52  | 000629-99-2 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092823\  
Data File : BP017427.D  
Acq On : 28 Sep 2023 23:10  
Operator : MA/JU  
Sample : 04417-39  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EZY6

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
Quant Title : SVOA CALIBRATION

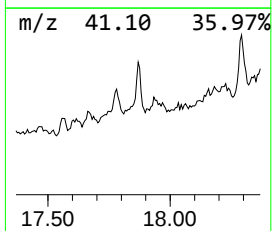
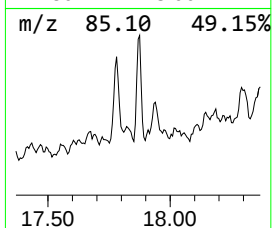
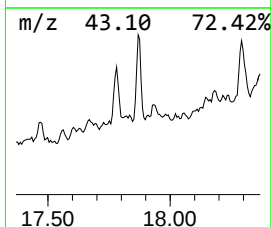
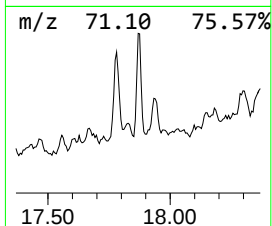
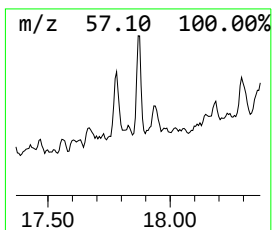
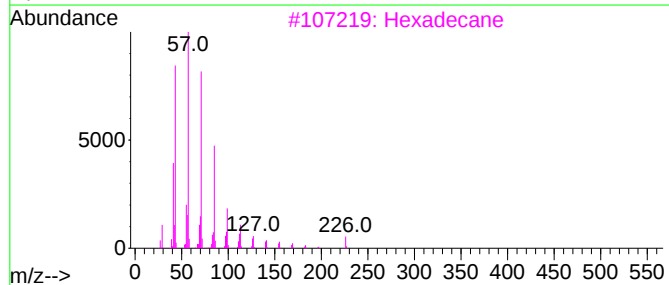
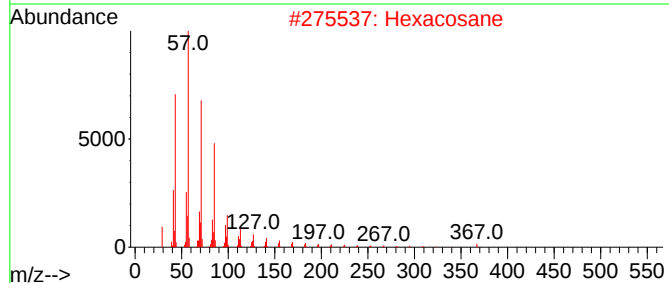
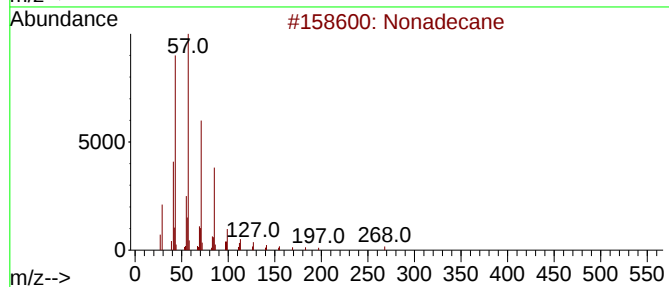
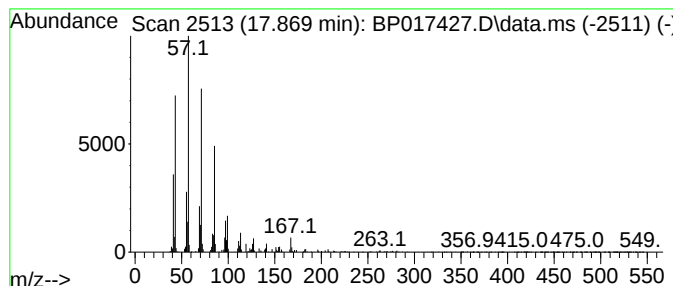
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.869 | 5.63 ng/ul | 465570 | Phenanthrene-d10 | 17.445 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Nonadecane           | 268 | C19H40  | 000629-92-5 | 91   |
| 2    |    |   | Hexacosane           | 366 | C26H54  | 000630-01-3 | 91   |
| 3    |    |   | Hexadecane           | 226 | C16H34  | 000544-76-3 | 87   |
| 4    |    |   | Tridecane, 6-propyl- | 226 | C16H34  | 055045-10-8 | 87   |
| 5    |    |   | Heneicosane          | 296 | C21H44  | 000629-94-7 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092823\  
Data File : BP017427.D  
Acq On : 28 Sep 2023 23:10  
Operator : MA/JU  
Sample : 04417-39  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

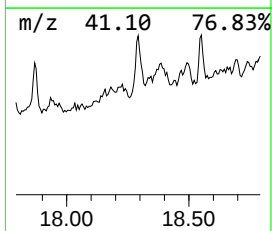
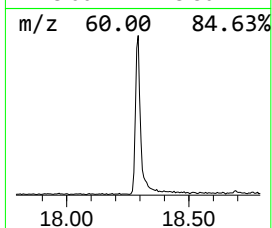
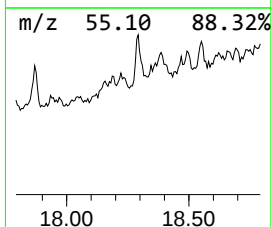
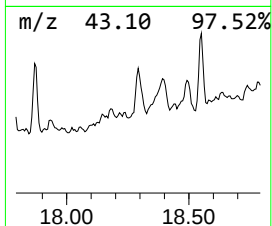
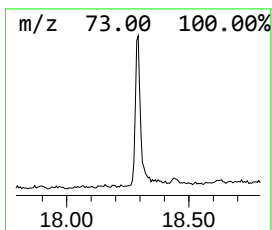
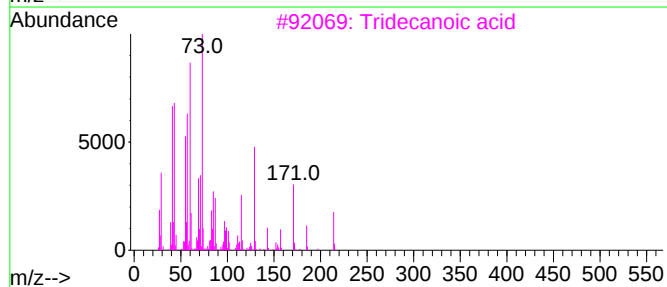
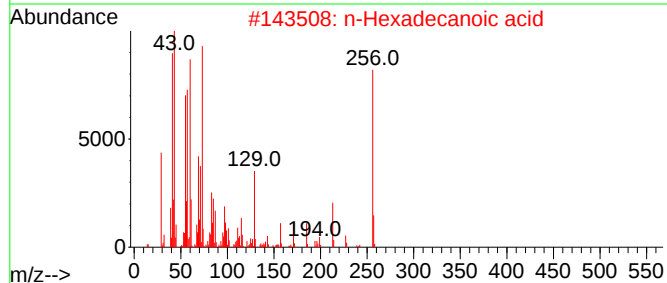
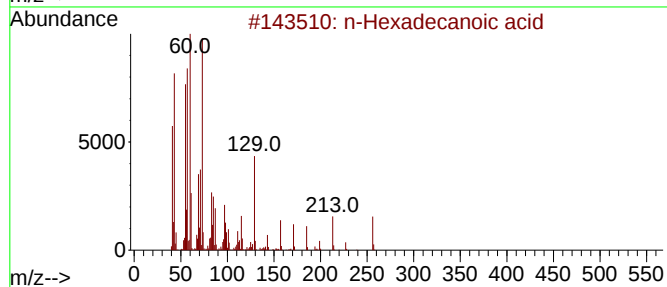
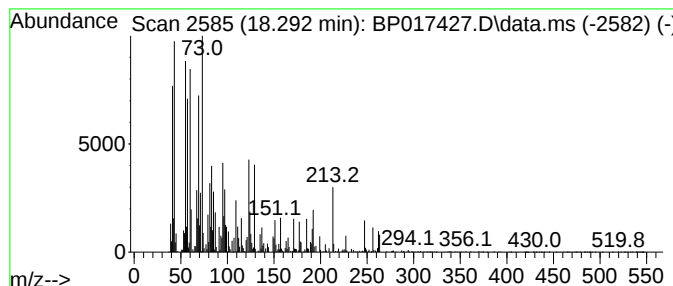
Instrument :  
BNA\_P  
ClientSampleId :  
EZY6

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 16 n-Hexadecanoic acid Concentration Rank 5

| R.T.      | EstConc             | Area   | Relative to ISTD |          | R.T.        |      |
|-----------|---------------------|--------|------------------|----------|-------------|------|
| 18.292    | 10.34 ng/ul         | 855357 | Phenanthrene-d10 |          | 17.445      |      |
| Hit# of 5 | Tentative ID        |        | MW               | MolForm  | CAS#        | Qual |
| 1         | n-Hexadecanoic acid |        | 256              | C16H32O2 | 000057-10-3 | 95   |
| 2         | n-Hexadecanoic acid |        | 256              | C16H32O2 | 000057-10-3 | 60   |
| 3         | Tridecanoic acid    |        | 214              | C13H26O2 | 000638-53-9 | 55   |
| 4         | Tetradecanoic acid  |        | 228              | C14H28O2 | 000544-63-8 | 55   |
| 5         | Tridecanoic acid    |        | 214              | C13H26O2 | 000638-53-9 | 49   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092823\  
Data File : BP017427.D  
Acq On : 28 Sep 2023 23:10  
Operator : MA/JU  
Sample : 04417-39  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EZY6

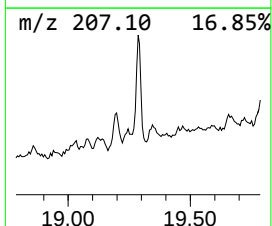
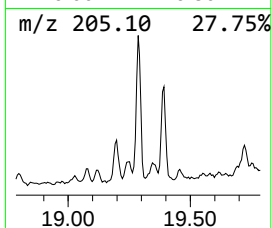
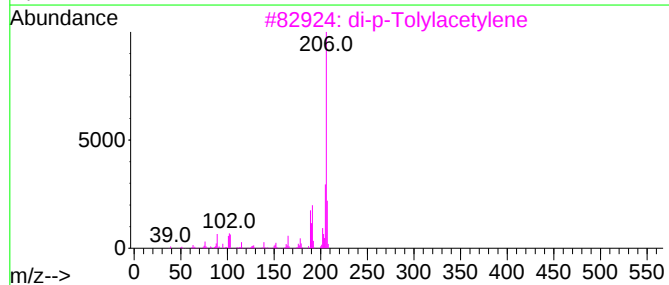
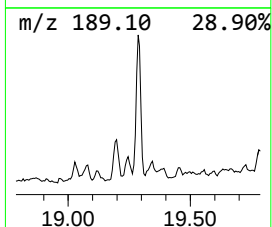
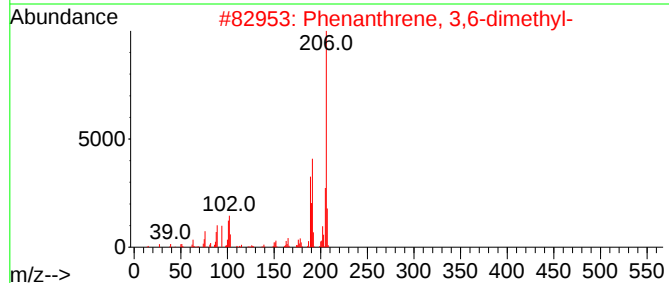
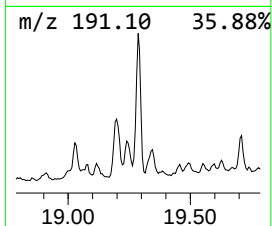
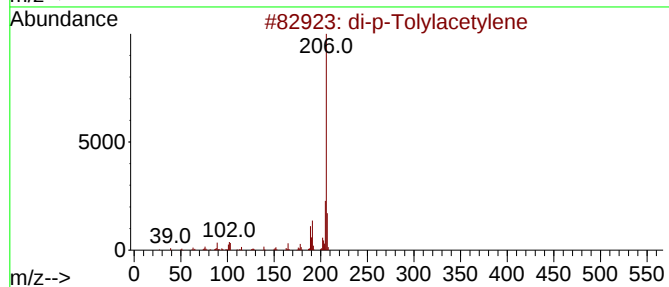
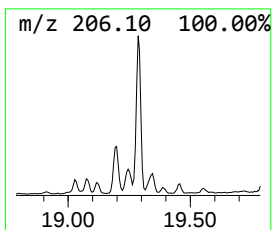
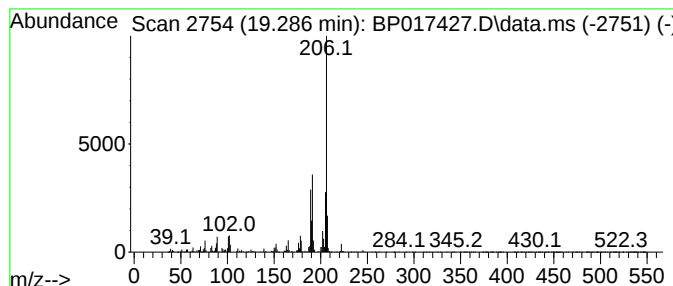
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Quant Title : SVOA CALIBRATION

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

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Peak Number 17 di-p-Tolylacetylene Concentration Rank 3

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 19.286 | 14.63 ng/ul | 1210440 | Phenanthrene-d10 | 17.445 |

| Hit# | of 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|------|-----------------------------|-----|---------|-------------|------|
| 1    |      | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 94   |
| 2    |      | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 93   |
| 3    |      | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 93   |
| 4    |      | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 93   |
| 5    |      | Phenanthrene, 1,7-dimethyl- | 206 | C16H14  | 000483-87-4 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092823\  
Data File : BP017427.D  
Acq On : 28 Sep 2023 23:10  
Operator : MA/JU  
Sample : 04417-39  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EZY6

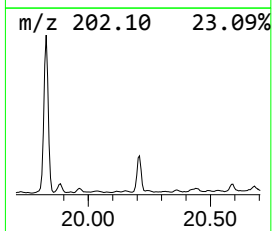
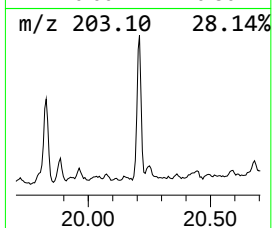
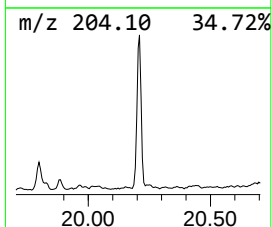
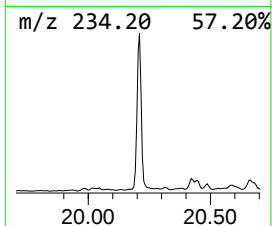
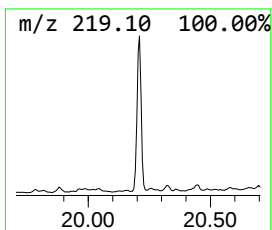
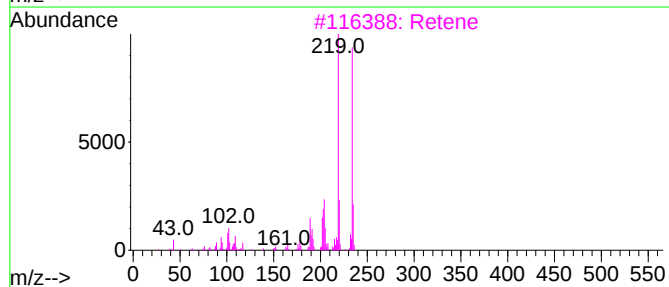
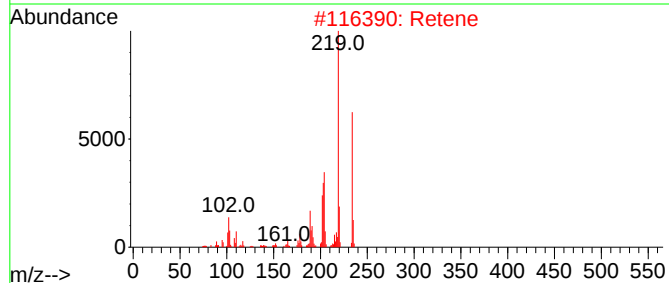
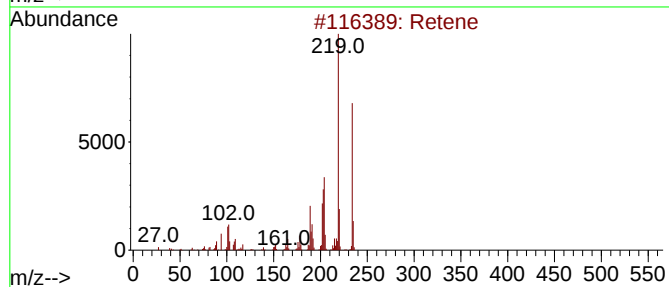
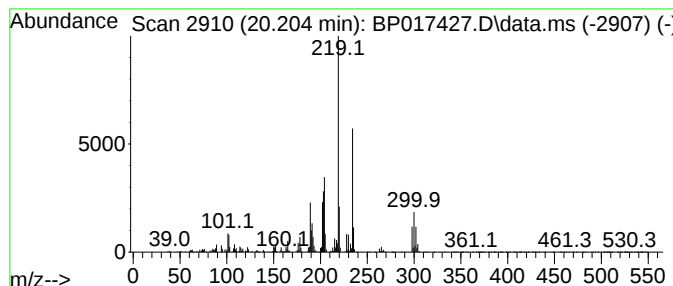
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Quant Title : SVOA CALIBRATION

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

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Peak Number 18 Retene Concentration Rank 1

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 20.204 | 20.00 ng/ul | 2337390 | Chrysene-d12     | 21.580 |

| Hit# | of                                 | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------------------------------------|---|--------------|-----|---------|-------------|------|
| 1    | Retene                             |   |              | 234 | C18H18  | 000483-65-8 | 98   |
| 2    | Retene                             |   |              | 234 | C18H18  | 000483-65-8 | 98   |
| 3    | Retene                             |   |              | 234 | C18H18  | 000483-65-8 | 94   |
| 4    | 2-Isopropyl-10-methylphenanthrene  |   |              | 234 | C18H18  | 066552-97-4 | 93   |
| 5    | Phenanthrene, 2,4,5,7-tetramethyl- |   |              | 234 | C18H18  | 007396-38-5 | 52   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092823\  
 Data File : BP017427.D  
 Acq On : 28 Sep 2023 23:10  
 Operator : MA/JU  
 Sample : 04417-39  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EYZE6

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
 Quant Title : SVOA CALIBRATION

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

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
|                    |        |         |       |          | #                     | RT     | Resp    | Conc |
| Butane, 2,3-dim... | 3.928  | 2.6     | ng/ul | 98086    | 1                     | 7.963  | 751670  | 20.0 |
| (DEL) Alkane: S... | 10.687 | 2.2     | ng/ul | 113573   | 2                     | 10.804 | 1032990 | 20.0 |
| (DEL) Alkane: S... | 12.146 | 2.8     | ng/ul | 143203   | 2                     | 10.804 | 1032990 | 20.0 |
| unknown-01         | 12.322 | 6.9     | ng/ul | 357812   | 2                     | 10.804 | 1032990 | 20.0 |
| (DEL) Alkane: S... | 13.398 | 5.6     | ng/ul | 362435   | 3                     | 14.657 | 1303810 | 20.0 |
| 2,4,7,9-Tetrame... | 13.493 | 15.2    | ng/ul | 989150   | 3                     | 14.657 | 1303810 | 20.0 |
| unknown-02         | 14.016 | 6.1     | ng/ul | 397028   | 3                     | 14.657 | 1303810 | 20.0 |
| unknown-03         | 14.428 | 6.6     | ng/ul | 428840   | 3                     | 14.657 | 1303810 | 20.0 |
| (DEL) Alkane: S... | 14.481 | 3.6     | ng/ul | 233656   | 3                     | 14.657 | 1303810 | 20.0 |
| unknown-04         | 15.257 | 2.3     | ng/ul | 152608   | 3                     | 14.657 | 1303810 | 20.0 |
| unknown-05         | 15.322 | 2.6     | ng/ul | 171291   | 3                     | 14.657 | 1303810 | 20.0 |
| 4-Fluorobenzoic... | 15.398 | 6.2     | ng/ul | 402458   | 3                     | 14.657 | 1303810 | 20.0 |
| (DEL) Alkane: S... | 16.322 | 10.9    | ng/ul | 897917   | 4                     | 17.445 | 1654950 | 20.0 |
| (DEL) Alkane: S... | 17.163 | 8.3     | ng/ul | 688117   | 4                     | 17.445 | 1654950 | 20.0 |
| (DEL) Alkane: S... | 17.869 | 5.6     | ng/ul | 465570   | 4                     | 17.445 | 1654950 | 20.0 |
| n-Hexadecanoic ... | 18.292 | 10.3    | ng/ul | 855357   | 4                     | 17.445 | 1654950 | 20.0 |
| di-p-Tolylacety... | 19.286 | 14.6    | ng/ul | 1210440  | 4                     | 17.445 | 1654950 | 20.0 |
| Retene             | 20.204 | 20.0    | ng/ul | 2337390  | 5                     | 21.580 | 2337390 | 20.0 |