

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\  
 Data File : BP004388.D  
 Acq On : 19 Dec 2020 11:40  
 Operator : CG/JU  
 Sample : L5128-02  
 Misc :  
 ALS Vial : 35 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 CB8C0

## Integration Parameters: LSCINT.P

Integrator: RTE  
 Smoothing : 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\SOM-EPA-BP121820MA.M  
 Title : SVOA CALIBRATION

Signal : TIC

| 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         | 3.352       | 23            | 28          | 42           | rVB      | 765894         | 1124994       | 19.17%          | 1.631%        |
| 2         | 4.116       | 154           | 158         | 163          | rBV      | 125544         | 160054        | 2.73%           | 0.232%        |
| 3         | 4.481       | 215           | 220         | 227          | rVB      | 252665         | 382634        | 6.52%           | 0.555%        |
| 4         | 6.104       | 492           | 496         | 501          | rVB      | 43533          | 71365         | 1.22%           | 0.103%        |
| 5         | 6.775       | 605           | 610         | 619          | rVB3     | 36180          | 70392         | 1.20%           | 0.102%        |
| 6         | 6.992       | 639           | 647         | 662          | rVB2     | 569559         | 1076510       | 18.34%          | 1.561%        |
| 7         | 7.151       | 668           | 674         | 686          | rBV      | 673594         | 1101673       | 18.77%          | 1.597%        |
| 8         | 7.269       | 690           | 694         | 702          | rVB2     | 40948          | 66262         | 1.13%           | 0.096%        |
| 9         | 7.357       | 702           | 709         | 716          | rBV      | 881990         | 1414591       | 24.10%          | 2.051%        |
| 10        | 7.822       | 779           | 788         | 795          | rBV      | 776237         | 1325313       | 22.58%          | 1.922%        |
| 11        | 7.892       | 795           | 800         | 812          | rVB2     | 68966          | 145429        | 2.48%           | 0.211%        |
| 12        | 8.528       | 901           | 908         | 916          | rVV      | 960094         | 1564096       | 26.65%          | 2.268%        |
| 13        | 8.604       | 917           | 921         | 927          | rVB5     | 20463          | 38104         | 0.65%           | 0.055%        |
| 14        | 8.716       | 935           | 940         | 949          | rVB7     | 17123          | 33818         | 0.58%           | 0.049%        |
| 15        | 8.804       | 949           | 955         | 964          | rBV      | 588354         | 992408        | 16.91%          | 1.439%        |
| 16        | 8.922       | 970           | 975         | 978          | rBV      | 52528          | 80707         | 1.38%           | 0.117%        |
| 17        | 8.975       | 978           | 984         | 996          | rVB      | 805485         | 1389949       | 23.68%          | 2.015%        |
| 18        | 9.181       | 1009          | 1019        | 1027         | rVV10    | 30819          | 87861         | 1.50%           | 0.127%        |
| 19        | 9.334       | 1039          | 1045        | 1050         | rBV5     | 61697          | 108784        | 1.85%           | 0.158%        |
| 20        | 9.428       | 1050          | 1061        | 1072         | rVB2     | 119124         | 288396        | 4.91%           | 0.418%        |
| 21        | 9.522       | 1072          | 1077        | 1081         | rBV5     | 26703          | 46178         | 0.79%           | 0.067%        |
| 22        | 9.622       | 1089          | 1094        | 1101         | rVB6     | 24494          | 55549         | 0.95%           | 0.081%        |
| 23        | 9.698       | 1101          | 1107        | 1114         | rBV      | 717779         | 1212705       | 20.66%          | 1.758%        |
| 24        | 9.828       | 1121          | 1129        | 1134         | rBV4     | 56645          | 112041        | 1.91%           | 0.162%        |
| 25        | 9.881       | 1134          | 1138        | 1145         | rVB2     | 98261          | 162234        | 2.76%           | 0.235%        |
| 26        | 10.034      | 1159          | 1164        | 1171         | rVB3     | 39136          | 76394         | 1.30%           | 0.111%        |
| 27        | 10.163      | 1177          | 1186        | 1193         | rVB      | 170478         | 330201        | 5.63%           | 0.479%        |
| 28        | 10.239      | 1193          | 1199        | 1207         | rBV      | 1210177        | 2094984       | 35.69%          | 3.038%        |
| 29        | 10.298      | 1207          | 1209        | 1215         | rVB3     | 40210          | 66258         | 1.13%           | 0.096%        |
| 30        | 10.398      | 1218          | 1226        | 1230         | rBV6     | 31930          | 80756         | 1.38%           | 0.117%        |
| 31        | 10.528      | 1240          | 1248        | 1256         | rBV2     | 54018          | 114389        | 1.95%           | 0.166%        |
| 32        | 10.616      | 1256          | 1263        | 1270         | rVV      | 1134407        | 1921799       | 32.74%          | 2.787%        |
| 33        | 10.669      | 1270          | 1272        | 1276         | rVV2     | 39899          | 53748         | 0.92%           | 0.078%        |
| 34        | 10.751      | 1276          | 1286        | 1300         | rVB      | 998481         | 1887082       | 32.15%          | 2.736%        |

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 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\SOM-EPA-BP121820MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |       |         |         |        |        |
|----|--------|------|------|------|-------|---------|---------|--------|--------|
| 35 | 10.875 | 1300 | 1307 | 1316 | rBV10 | 48533   | 154718  | 2.64%  | 0.224% |
| 36 | 10.963 | 1320 | 1322 | 1328 | rVB7  | 19076   | 34813   | 0.59%  | 0.050% |
| 37 | 11.045 | 1328 | 1336 | 1341 | rBV9  | 35018   | 73562   | 1.25%  | 0.107% |
| 38 | 11.192 | 1354 | 1361 | 1363 | rBV6  | 26918   | 63765   | 1.09%  | 0.092% |
| 39 | 11.281 | 1374 | 1376 | 1387 | rVB5  | 33951   | 72267   | 1.23%  | 0.105% |
| 40 | 11.439 | 1395 | 1403 | 1404 | rBV2  | 37926   | 68239   | 1.16%  | 0.099% |
| 41 | 11.469 | 1405 | 1408 | 1414 | rVB3  | 60286   | 97225   | 1.66%  | 0.141% |
| 42 | 11.533 | 1415 | 1419 | 1423 | rBV4  | 23765   | 43063   | 0.73%  | 0.062% |
| 43 | 11.586 | 1423 | 1428 | 1436 | rVB8  | 58651   | 117991  | 2.01%  | 0.171% |
| 44 | 11.657 | 1436 | 1440 | 1445 | rBV5  | 34863   | 64090   | 1.09%  | 0.093% |
| 45 | 11.780 | 1457 | 1461 | 1465 | rBV4  | 22427   | 36774   | 0.63%  | 0.053% |
| 46 | 11.822 | 1465 | 1468 | 1474 | rVB   | 51940   | 79346   | 1.35%  | 0.115% |
| 47 | 11.969 | 1488 | 1493 | 1505 | rVB   | 60390   | 132041  | 2.25%  | 0.191% |
| 48 | 12.069 | 1505 | 1510 | 1514 | rBV3  | 31683   | 55756   | 0.95%  | 0.081% |
| 49 | 12.145 | 1517 | 1523 | 1527 | rBV   | 214117  | 335421  | 5.71%  | 0.486% |
| 50 | 12.186 | 1528 | 1530 | 1536 | rVV4  | 43889   | 75088   | 1.28%  | 0.109% |
| 51 | 12.269 | 1536 | 1544 | 1550 | rVV4  | 20631   | 66569   | 1.13%  | 0.097% |
| 52 | 12.339 | 1551 | 1556 | 1566 | rVV9  | 45468   | 115696  | 1.97%  | 0.168% |
| 53 | 12.439 | 1567 | 1573 | 1580 | rVB10 | 21018   | 49541   | 0.84%  | 0.072% |
| 54 | 12.504 | 1580 | 1584 | 1589 | rBV6  | 20172   | 36484   | 0.62%  | 0.053% |
| 55 | 12.575 | 1589 | 1596 | 1599 | rBV9  | 15804   | 31981   | 0.54%  | 0.046% |
| 56 | 12.616 | 1599 | 1603 | 1613 | rVV   | 52743   | 98890   | 1.68%  | 0.143% |
| 57 | 12.716 | 1616 | 1620 | 1625 | rVB3  | 30153   | 44471   | 0.76%  | 0.064% |
| 58 | 12.839 | 1634 | 1641 | 1648 | rVV9  | 32620   | 69248   | 1.18%  | 0.100% |
| 59 | 13.010 | 1666 | 1670 | 1675 | rBV6  | 24410   | 43971   | 0.75%  | 0.064% |
| 60 | 13.180 | 1694 | 1699 | 1709 | rBV9  | 35036   | 67293   | 1.15%  | 0.098% |
| 61 | 13.298 | 1717 | 1719 | 1728 | rVB5  | 35297   | 69536   | 1.18%  | 0.101% |
| 62 | 13.504 | 1750 | 1754 | 1761 | rBV3  | 74006   | 144120  | 2.46%  | 0.209% |
| 63 | 13.563 | 1761 | 1764 | 1774 | rVB6  | 39665   | 95505   | 1.63%  | 0.138% |
| 64 | 13.645 | 1774 | 1778 | 1782 | rBV4  | 21218   | 34861   | 0.59%  | 0.051% |
| 65 | 13.704 | 1783 | 1788 | 1793 | rVB5  | 27488   | 48027   | 0.82%  | 0.070% |
| 66 | 13.874 | 1812 | 1817 | 1822 | rBV   | 2072665 | 2790915 | 47.55% | 4.047% |
| 67 | 13.922 | 1822 | 1825 | 1830 | rVB   | 213739  | 280504  | 4.78%  | 0.407% |
| 68 | 14.022 | 1837 | 1842 | 1846 | rBV7  | 28120   | 62074   | 1.06%  | 0.090% |
| 69 | 14.069 | 1848 | 1850 | 1856 | rVB5  | 22905   | 33701   | 0.57%  | 0.049% |
| 70 | 14.163 | 1859 | 1866 | 1874 | rVB   | 2279274 | 3457519 | 58.91% | 5.013% |
| 71 | 14.280 | 1881 | 1886 | 1890 | rBV2  | 78531   | 129670  | 2.21%  | 0.188% |

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 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121820MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |        |
|-----|--------|------|------|------|------|---------|---------|---------|--------|
| 72  | 14.392 | 1899 | 1905 | 1910 | rVB2 | 212440  | 324179  | 5.52%   | 0.470% |
| 73  | 14.469 | 1912 | 1918 | 1927 | rVB  | 1753006 | 2629790 | 44.80%  | 3.813% |
| 74  | 14.551 | 1929 | 1932 | 1936 | rBV4 | 25071   | 36146   | 0.62%   | 0.052% |
| 75  | 14.674 | 1948 | 1953 | 1960 | rBV  | 496643  | 755458  | 12.87%  | 1.095% |
| 76  | 14.827 | 1976 | 1979 | 1983 | rBV5 | 23340   | 31540   | 0.54%   | 0.046% |
| 77  | 15.004 | 2006 | 2009 | 2013 | rVB5 | 24797   | 30782   | 0.52%   | 0.045% |
| 78  | 15.069 | 2017 | 2020 | 2025 | rVB4 | 37457   | 50672   | 0.86%   | 0.073% |
| 79  | 15.216 | 2041 | 2045 | 2050 | rBV2 | 51184   | 115725  | 1.97%   | 0.168% |
| 80  | 15.469 | 2082 | 2088 | 2099 | rVB  | 2839317 | 4137864 | 70.50%  | 6.000% |
| 81  | 15.592 | 2103 | 2109 | 2113 | rBV  | 578071  | 832445  | 14.18%  | 1.207% |
| 82  | 15.721 | 2125 | 2131 | 2138 | rBV3 | 64867   | 160280  | 2.73%   | 0.232% |
| 83  | 15.927 | 2162 | 2166 | 2170 | rBV2 | 48738   | 68095   | 1.16%   | 0.099% |
| 84  | 15.980 | 2171 | 2175 | 2180 | rVV  | 211851  | 289241  | 4.93%   | 0.419% |
| 85  | 16.133 | 2197 | 2201 | 2209 | rVB4 | 58474   | 114202  | 1.95%   | 0.166% |
| 86  | 16.204 | 2209 | 2213 | 2219 | rBV8 | 23997   | 53934   | 0.92%   | 0.078% |
| 87  | 16.368 | 2237 | 2241 | 2245 | rVB  | 39692   | 53458   | 0.91%   | 0.078% |
| 88  | 16.498 | 2259 | 2263 | 2266 | rBV  | 51458   | 70228   | 1.20%   | 0.102% |
| 89  | 16.763 | 2305 | 2308 | 2315 | rVB  | 75248   | 104907  | 1.79%   | 0.152% |
| 90  | 17.227 | 2379 | 2387 | 2396 | rBV  | 2128633 | 3183445 | 54.24%  | 4.616% |
| 91  | 17.327 | 2398 | 2404 | 2414 | rVV  | 3155466 | 4524330 | 77.08%  | 6.560% |
| 92  | 18.127 | 2536 | 2540 | 2549 | rBV  | 218592  | 373973  | 6.37%   | 0.542% |
| 93  | 18.874 | 2663 | 2667 | 2676 | rVB  | 255108  | 365906  | 6.23%   | 0.531% |
| 94  | 19.568 | 2780 | 2785 | 2790 | rBV  | 4001069 | 5465029 | 93.11%  | 7.924% |
| 95  | 19.615 | 2790 | 2793 | 2800 | rVB  | 173376  | 237418  | 4.04%   | 0.344% |
| 96  | 21.039 | 3031 | 3035 | 3042 | rBV2 | 749051  | 1131216 | 19.27%  | 1.640% |
| 97  | 21.321 | 3078 | 3083 | 3091 | rBV  | 2871191 | 3622745 | 61.72%  | 5.253% |
| 98  | 22.551 | 3288 | 3292 | 3300 | rVB  | 1131432 | 1567965 | 26.71%  | 2.274% |
| 99  | 23.527 | 3451 | 3458 | 3475 | rVB2 | 3007210 | 5869564 | 100.00% | 8.511% |
| 100 | 23.680 | 3477 | 3484 | 3495 | rVB  | 1848743 | 3751410 | 63.91%  | 5.440% |

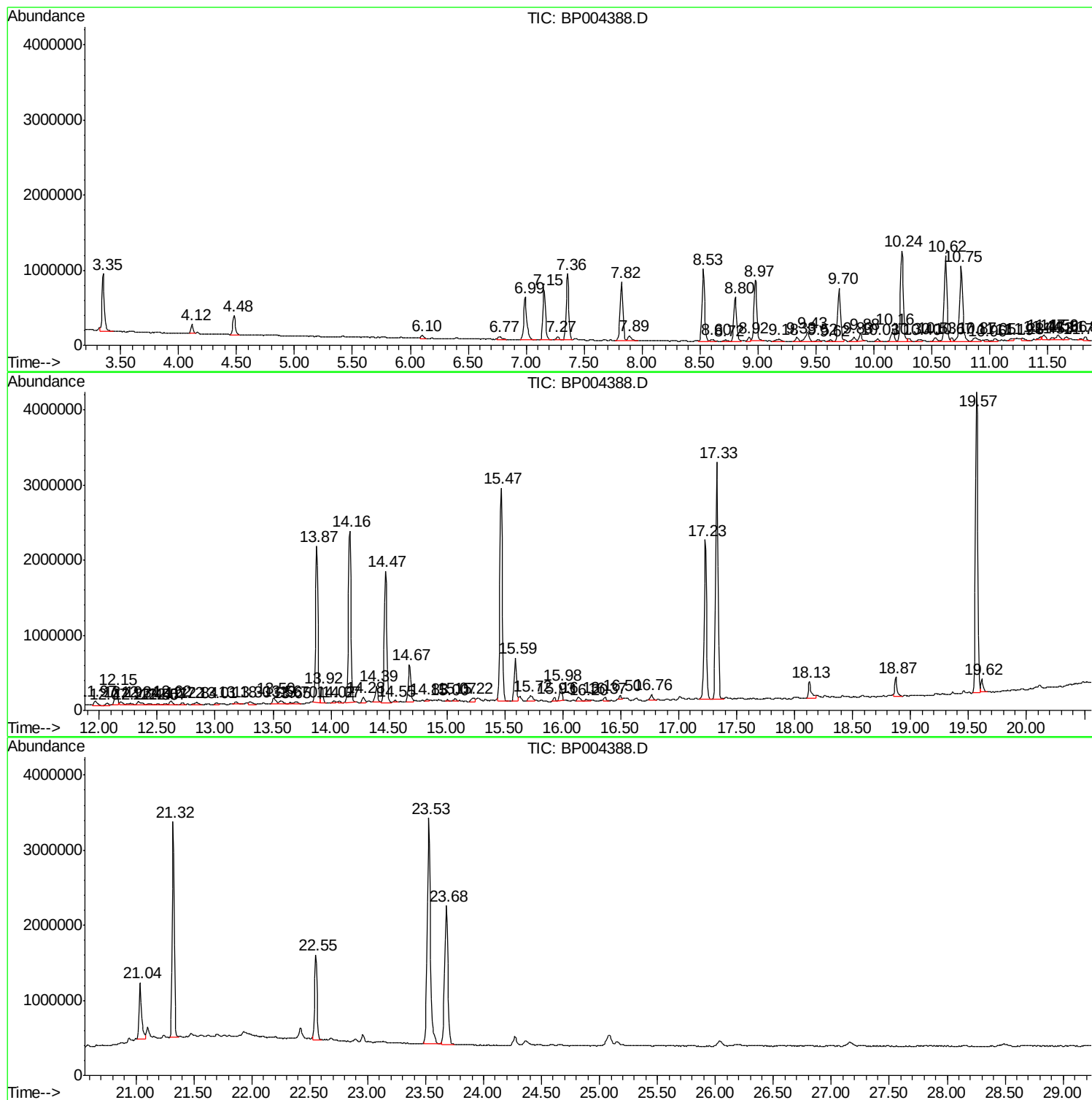
Sum of corrected areas: 68964340

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121820MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P



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Quant Title : SVOA CALIBRATION

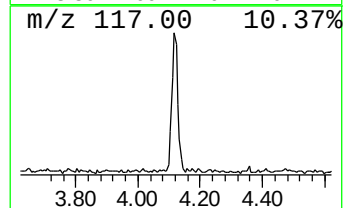
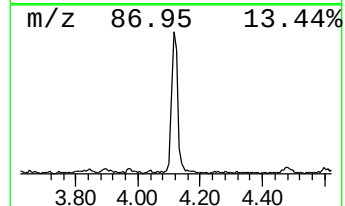
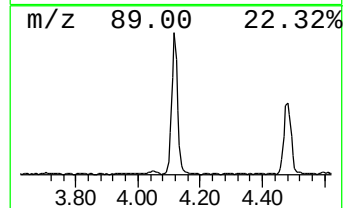
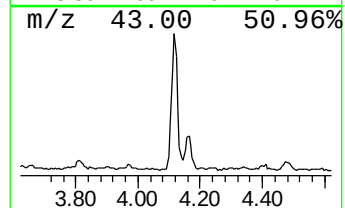
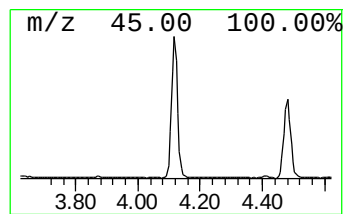
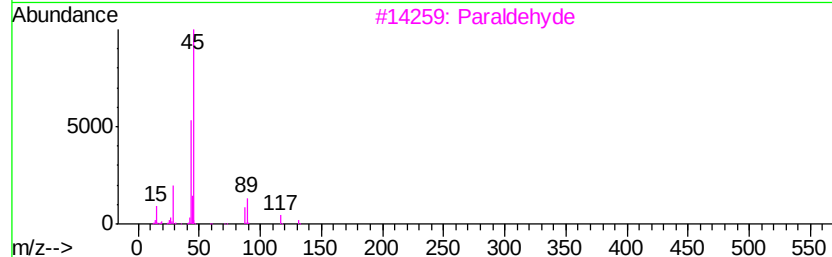
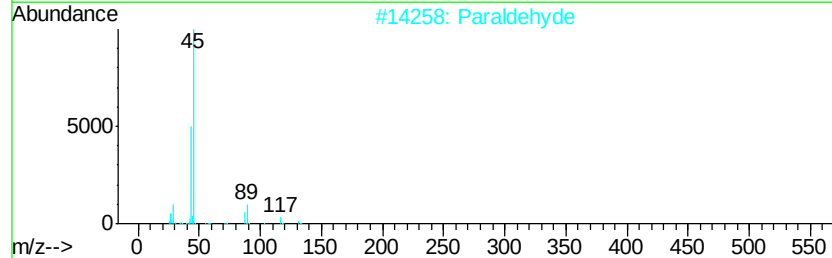
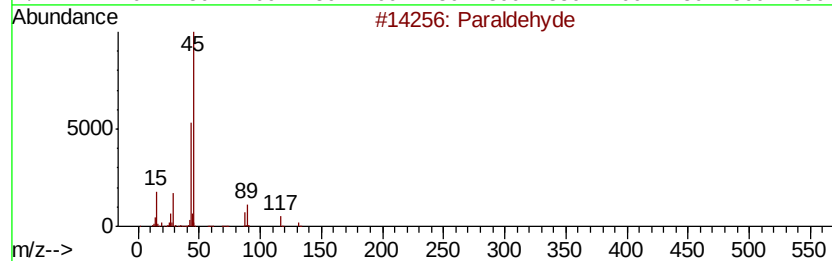
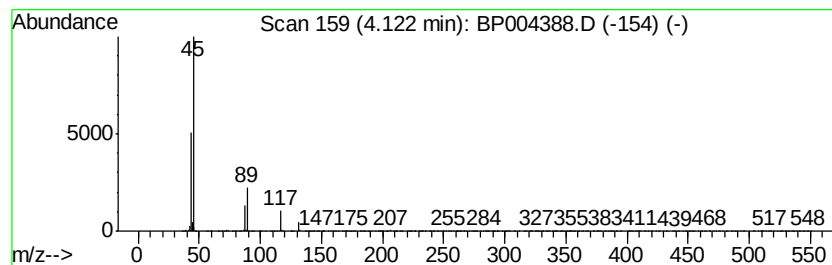
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Paraldehyde Concentration Rank 9

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 4.12 | 2.42 ng/ul | 160054 | 1,4-Dichlorobenzene-d4 | 7.82 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Paraldehyde                         | 132 | C6H12O3  | 000123-63-7 | 90   |
| 2    |    | Paraldehyde                         | 132 | C6H12O3  | 000123-63-7 | 74   |
| 3    |    | Paraldehyde                         | 132 | C6H12O3  | 000123-63-7 | 64   |
| 4    |    | Acetaldehyde, tetramer              | 176 | C8H16O4  | 000108-62-3 | 56   |
| 5    |    | Ethanol, 2-[2-(2-butoxyethoxy)et... | 206 | C10H22O4 | 000143-22-6 | 40   |



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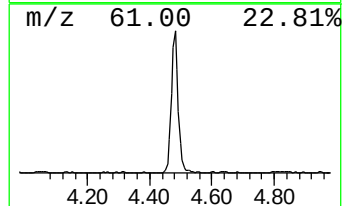
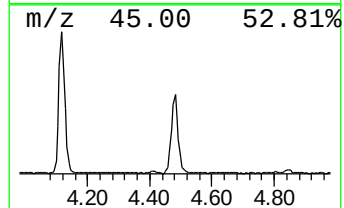
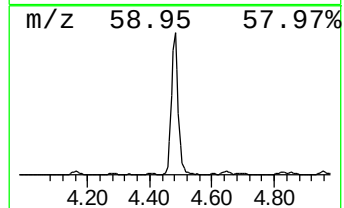
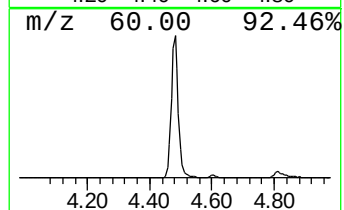
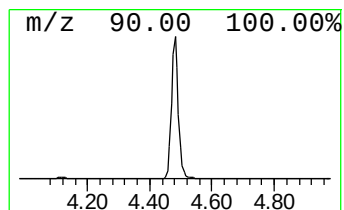
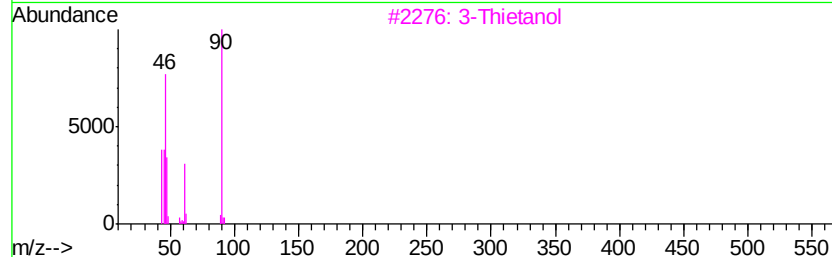
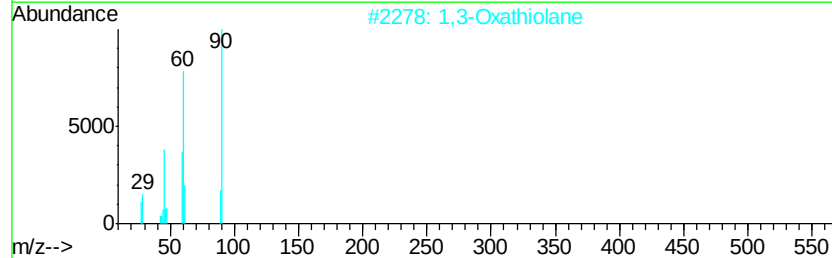
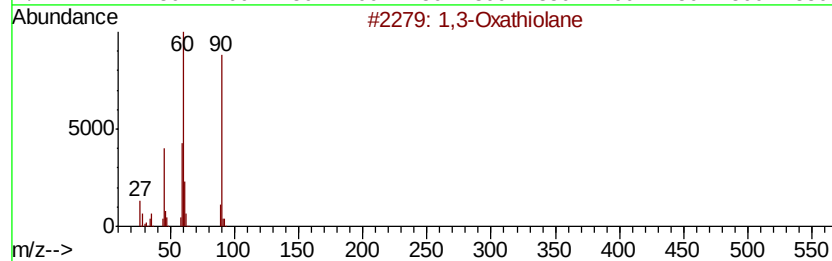
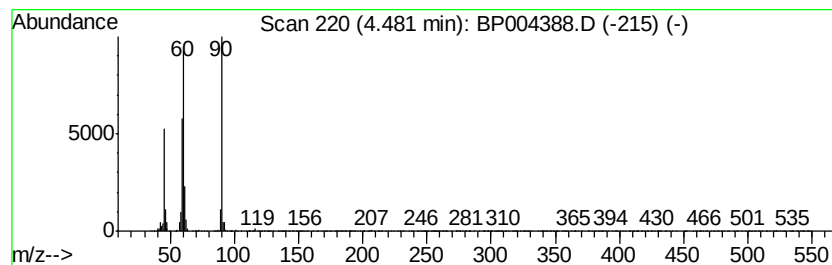
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 1,3-Oxathiolane Concentration Rank 4

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 4.48 | 5.77 ng/ul | 382634 | 1,4-Dichlorobenzene-d4 | 7.82 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1,3-Oxathiolane                     | 90  | C3H6OS  | 002094-97-5 | 90   |
| 2    |    | 1,3-Oxathiolane                     | 90  | C3H6OS  | 002094-97-5 | 78   |
| 3    |    | 3-Thietanol                         | 90  | C3H6OS  | 010304-16-2 | 53   |
| 4    |    | Thiourea, methyl-                   | 90  | C2H6N2S | 000598-52-7 | 50   |
| 5    |    | Heptanoic acid, 2-hydroxy-, meth... | 160 | C8H16O3 | 054340-91-9 | 42   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\  
Data File : BP004388.D  
Acq On : 19 Dec 2020 11:40  
Operator : CG/JU  
Sample : L5128-02  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

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

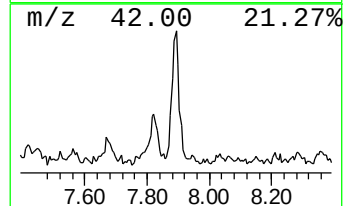
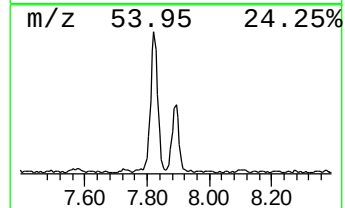
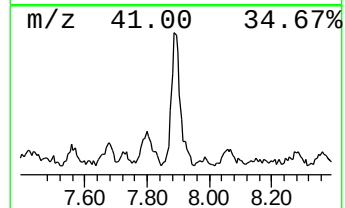
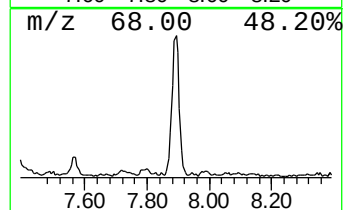
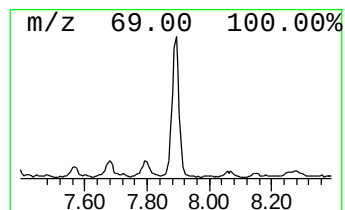
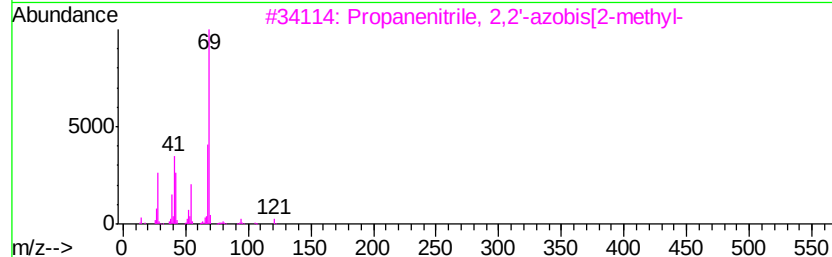
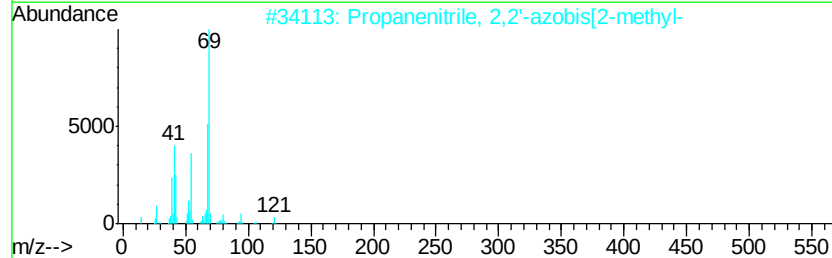
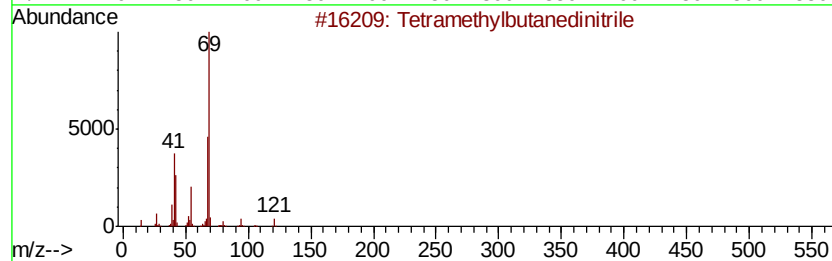
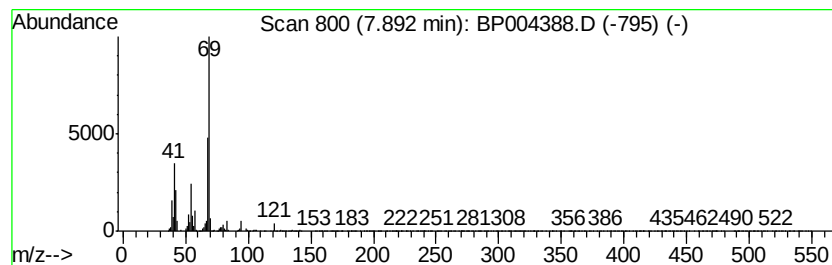
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Tetramethylbutanedinitrile Concentration Rank 12

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.89 | 2.19 ng/ul | 145429 | 1,4-Dichlorobenzene-d4 | 7.82 |

| Hit# | of | Tentative ID                           | MW  | MolForm  | CAS#         | Qual |
|------|----|----------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Tetramethylbutanedinitrile             | 136 | C8H12N2  | 003333-52-6  | 86   |
| 2    |    | Propanenitrile, 2,2'-azobis[2-methyl-] | 164 | C8H12N4  | 000078-67-1  | 78   |
| 3    |    | Propanenitrile, 2,2'-azobis[2-methyl-] | 164 | C8H12N4  | 000078-67-1  | 78   |
| 4    |    | Heptafluorobutyric acid, cyclope...    | 282 | C9H9F7O2 | 1000245-81-9 | 50   |
| 5    |    | 1,5-Pentandiol, 0,0'-di(proparg...     | 268 | C13H16O6 | 1000331-35-4 | 42   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\  
Data File : BP004388.D  
Acq On : 19 Dec 2020 11:40  
Operator : CG/JU  
Sample : L5128-02  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CB8C0

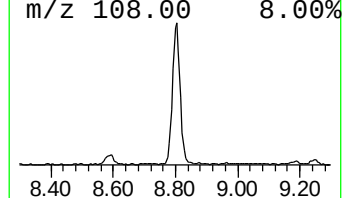
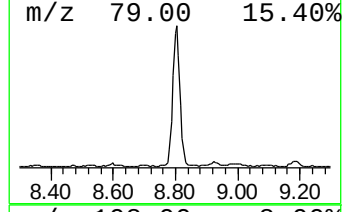
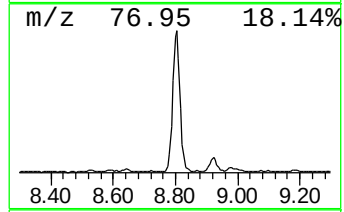
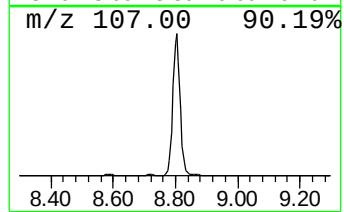
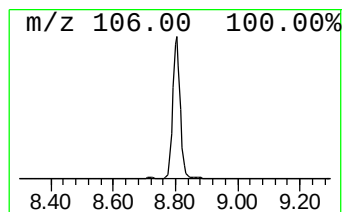
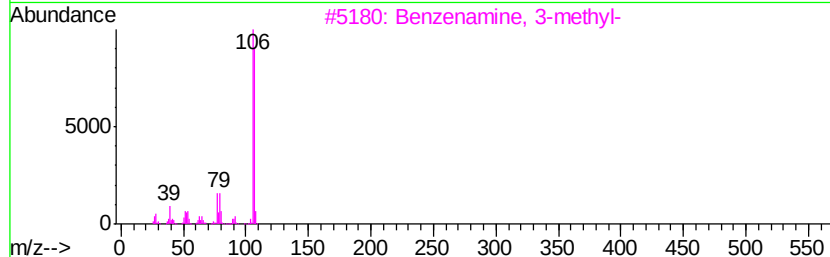
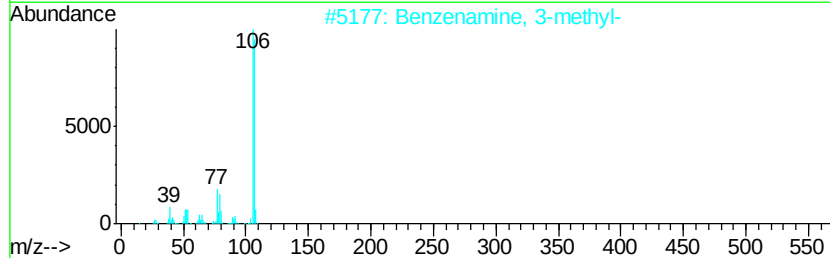
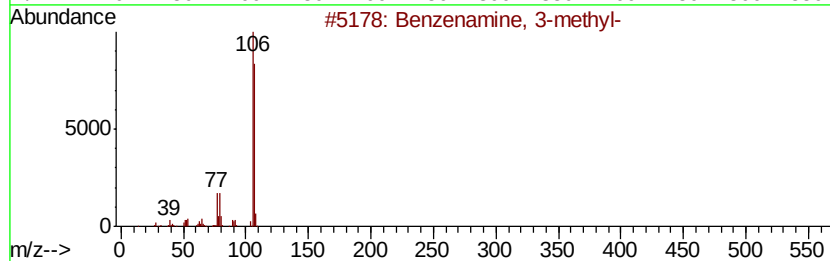
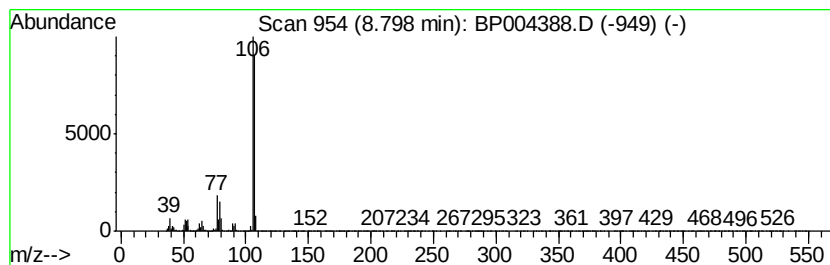
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Benzenamine, 3-methyl- Concentration Rank 1

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 8.80 | 14.98 ng/ul | 992408 | 1,4-Dichlorobenzene-d4 | 7.82 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzenamine, 3-methyl- | 107 | C7H9N   | 000108-44-1 | 97   |
| 2    |    | Benzenamine, 3-methyl- | 107 | C7H9N   | 000108-44-1 | 97   |
| 3    |    | Benzenamine, 3-methyl- | 107 | C7H9N   | 000108-44-1 | 97   |
| 4    |    | Benzenamine, 3-methyl- | 107 | C7H9N   | 000108-44-1 | 95   |
| 5    |    | o-Toluidine            | 107 | C7H9N   | 000095-53-4 | 94   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\  
Data File : BP004388.D  
Acq On : 19 Dec 2020 11:40  
Operator : CG/JU  
Sample : L5128-02  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CB8C0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121820MA.M  
Quant Title : SVOA CALIBRATION

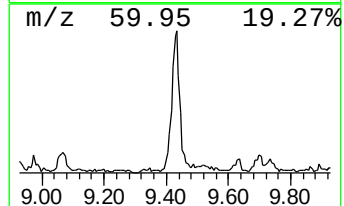
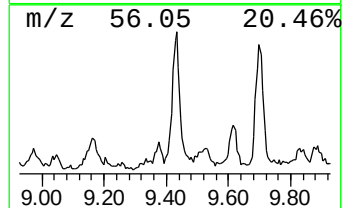
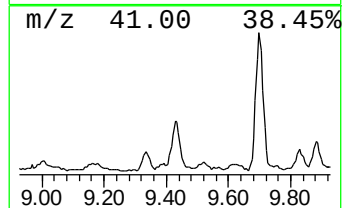
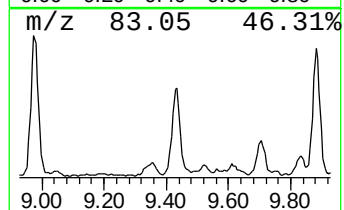
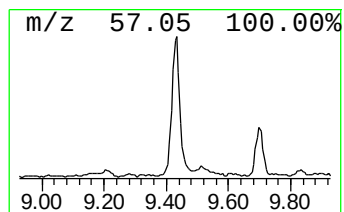
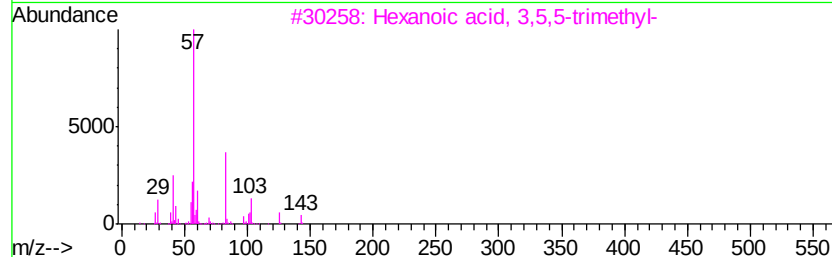
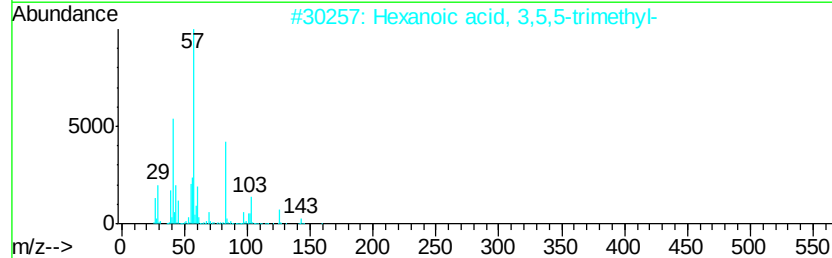
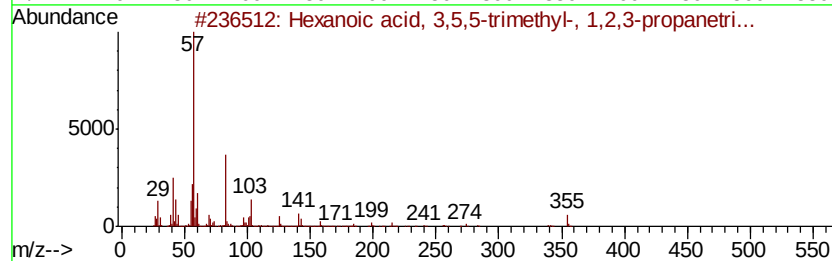
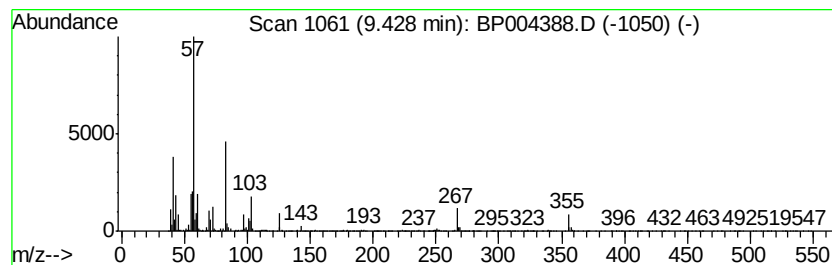
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Hexanoic acid, 3,5,5-trimet... Concentration Rank 7

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.43 | 3.00 ng/ul | 288396 | Naphthalene-d8   | 10.62 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Hexanoic acid, 3,5,5-trimethyl-,... | 512 | C30H56O6 | 056554-53-1 | 83   |
| 2    |    | Hexanoic acid, 3,5,5-trimethyl-     | 158 | C9H18O2  | 003302-10-1 | 72   |
| 3    |    | Hexanoic acid, 3,5,5-trimethyl-     | 158 | C9H18O2  | 003302-10-1 | 64   |
| 4    |    | 3,3-Dimethylheptanoic acid          | 158 | C9H18O2  | 067061-30-7 | 64   |
| 5    |    | Hexanoic acid, 3,5,5-trimethyl-     | 158 | C9H18O2  | 003302-10-1 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\  
Data File : BP004388.D  
Acq On : 19 Dec 2020 11:40  
Operator : CG/JU  
Sample : L5128-02  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

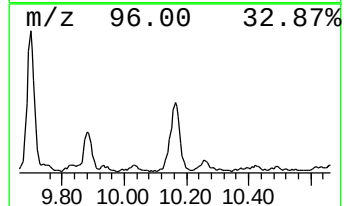
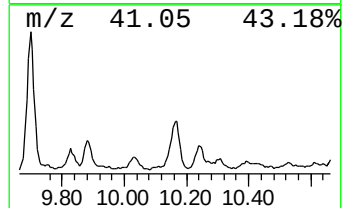
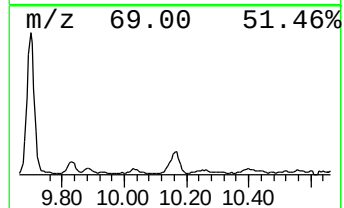
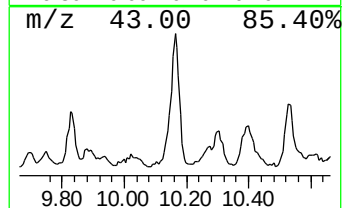
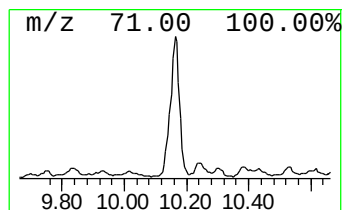
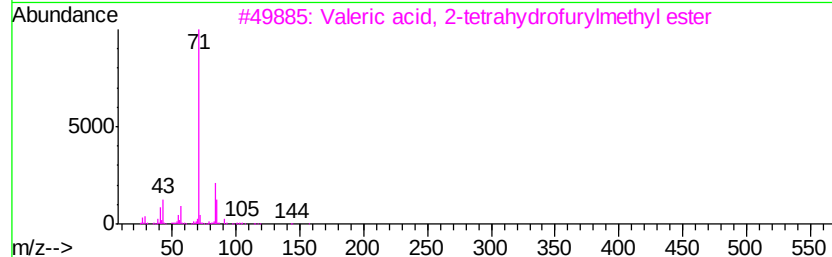
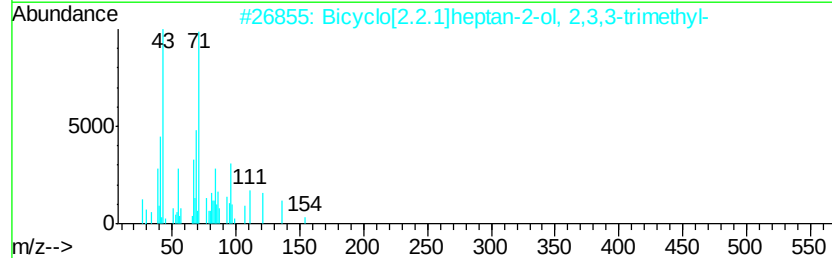
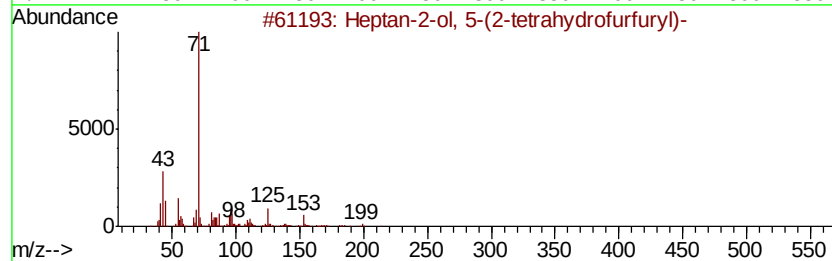
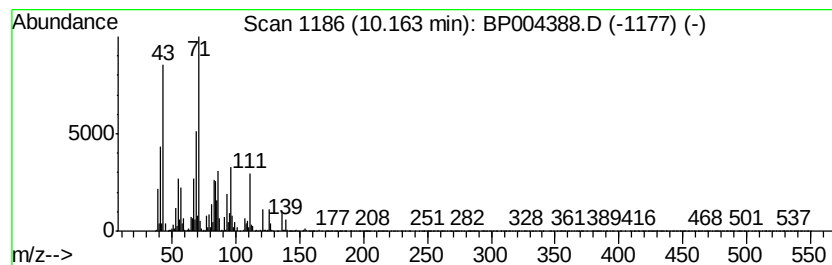
Instrument :  
BNA\_P  
ClientSampleId :  
CB8C0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121820MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 unknown-01 Concentration Rank 6

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.      |
|---------|-------------------------------------|--------------|------------------|-----------|
| 10.16   | 3.44 ng/ul                          | 330201       | Naphthalene-d8   | 10.62     |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Heptan-2-ol, 5-(2-tetrahydrofurf... | 200 C12H24O2 | 281676-71-9      | 46        |
| 2       | Bicyclo[2.2.1]heptan-2-ol, 2,3,3... | 154 C10H18O  | 000465-31-6      | 43        |
| 3       | Valeric acid, 2-tetrahydrofurylm... | 186 C10H18O3 | 005451-86-5      | 38        |
| 4       | 1-Propene, 3-methoxy-2-methyl-      | 86 C5H10O    | 022418-49-1      | 38        |
| 5       | 1,3-Dimethylbutyl butyrate          | 172 C10H20O2 | 005332-88-7      | 38        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\  
Data File : BP004388.D  
Acq On : 19 Dec 2020 11:40  
Operator : CG/JU  
Sample : L5128-02  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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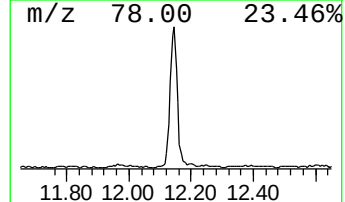
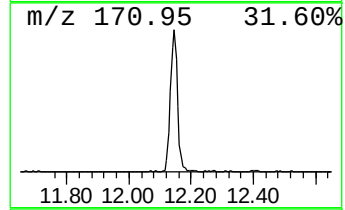
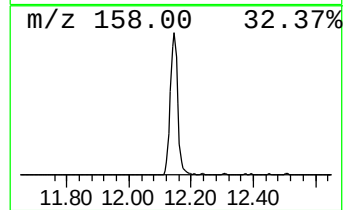
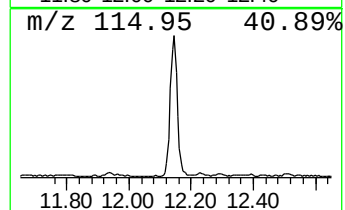
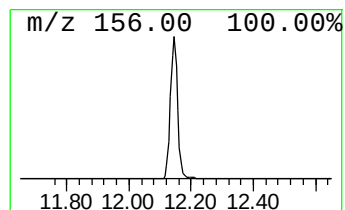
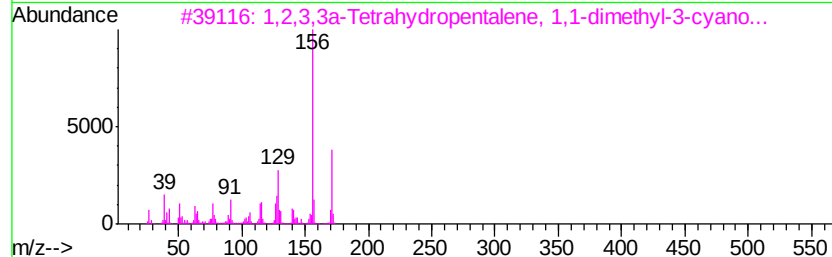
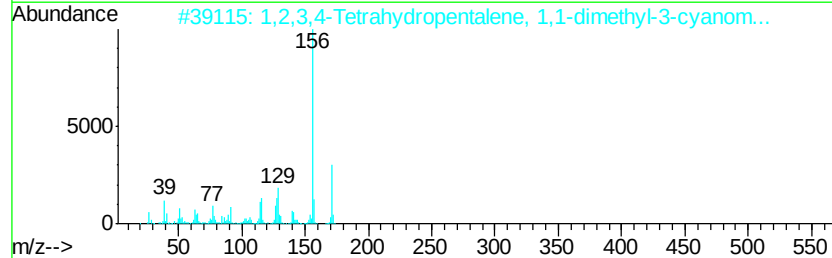
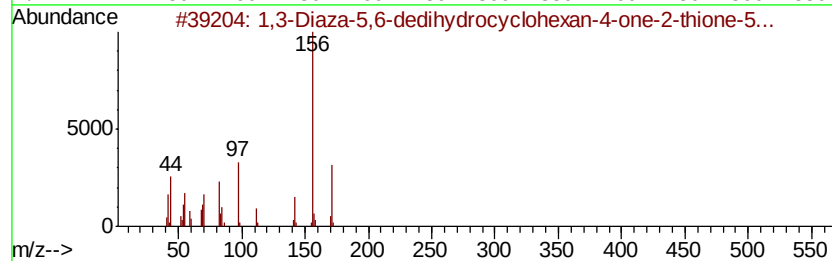
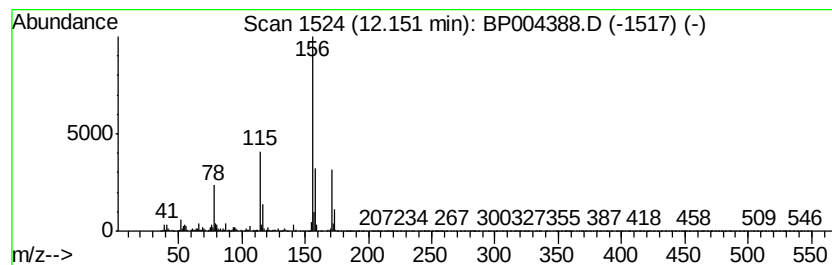
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 unknown-02 Concentration Rank 5

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.15 | 3.49 ng/ul | 335421 | Naphthalene-d8   | 10.62 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 1,3-Diaza-5,6-dedihydrocyclohexa... | 171 | C6H9N3OS | 021263-85-4  | 38   |
| 2    |    | 1,2,3,4-Tetrahydropentalene, 1,1... | 171 | C12H13N  | 1000217-18-6 | 37   |
| 3    |    | 1,2,3,3a-Tetrahydropentalene, 1,... | 171 | C12H13N  | 1000217-17-0 | 37   |
| 4    |    | Bicyclo[3.2.0]hept-2-ene, 4-isop... | 171 | C12H13N  | 1000217-15-6 | 32   |
| 5    |    | 6-Isopropylquinoline                | 171 | C12H13N  | 000135-79-5  | 32   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\  
Data File : BP004388.D  
Acq On : 19 Dec 2020 11:40  
Operator : CG/JU  
Sample : L5128-02  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

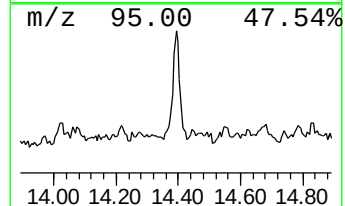
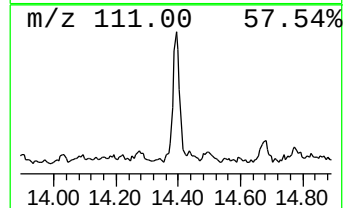
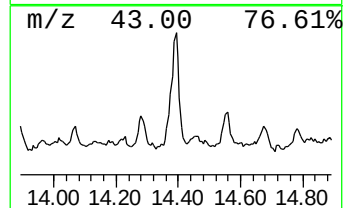
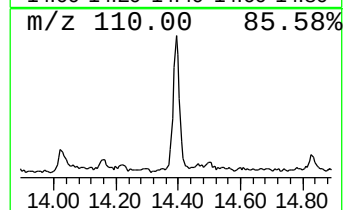
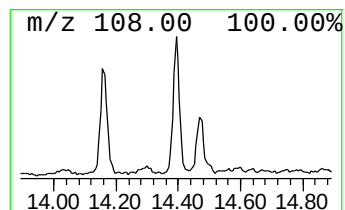
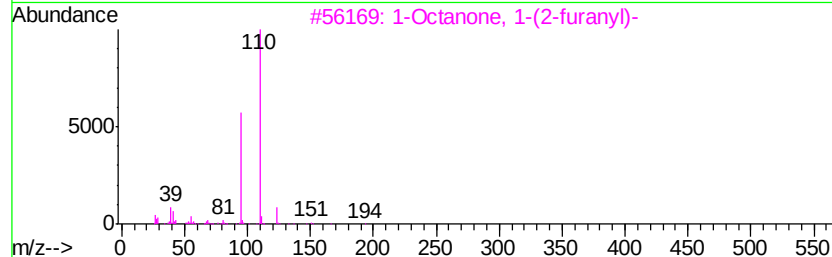
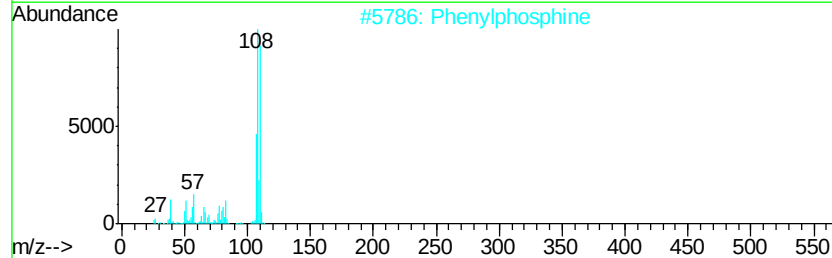
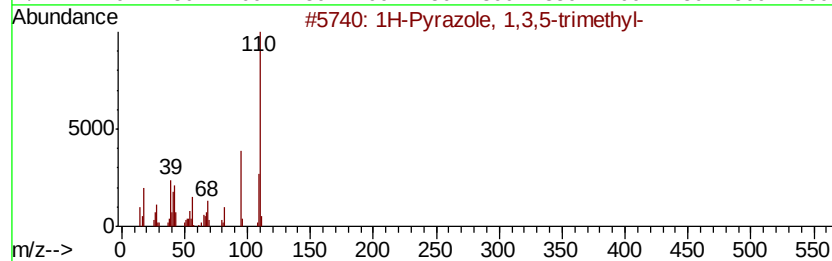
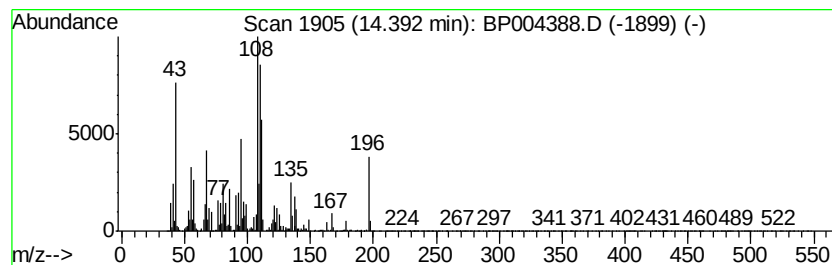
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ClientSampleId :  
CB8C0

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.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.            |
|---------|------------|--------------------------------------|------------------|-----------------|
| 14.39   | 2.47 ng/ul | 324179                               | Acenaphthene-d10 | 14.47           |
| Hit# of | 5          | Tentative ID                         | MW MolForm       | CAS# Qual       |
| 1       |            | 1H-Pyrazole, 1,3,5-trimethyl-        | 110 C6H10N2      | 001072-91-9 30  |
| 2       |            | Phenylphosphine                      | 110 C6H7P        | 000638-21-1 25  |
| 3       |            | 1-Octanone, 1-(2-furanvl)-           | 194 C12H18O2     | 005456-77-9 14  |
| 4       |            | Fumaric acid, butyl 2-methylcyclo... | 280 C16H24O4     | 1000345-15-6 14 |
| 5       |            | 7-Nonen-2-one, 4,8-dimethyl-         | 168 C11H20O      | 003664-64-0 14  |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\  
Data File : BP004388.D  
Acq On : 19 Dec 2020 11:40  
Operator : CG/JU  
Sample : L5128-02  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

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

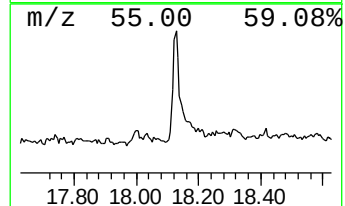
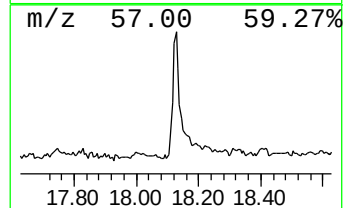
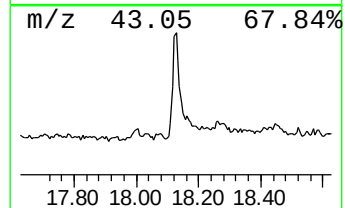
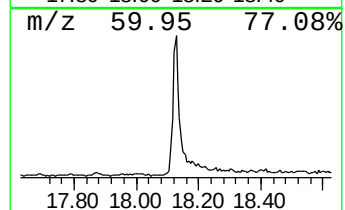
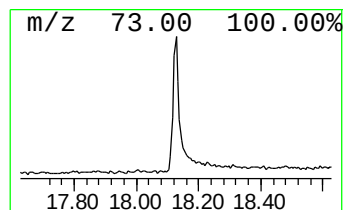
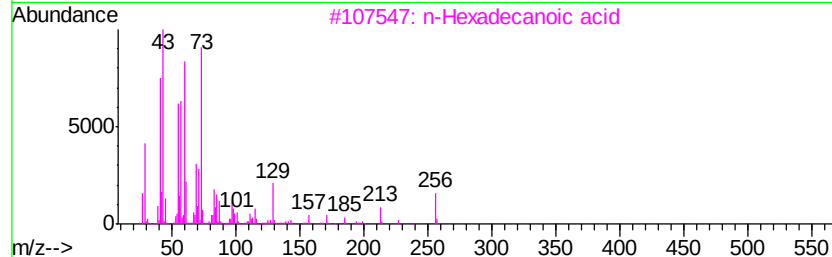
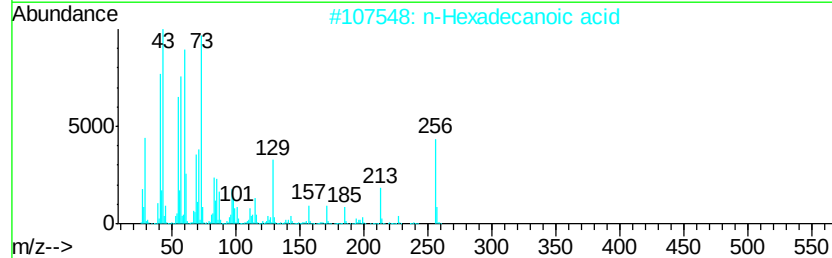
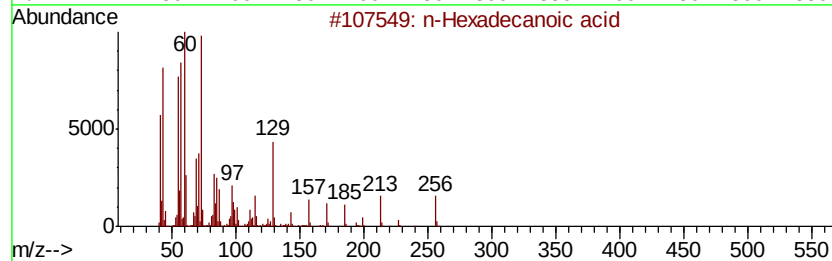
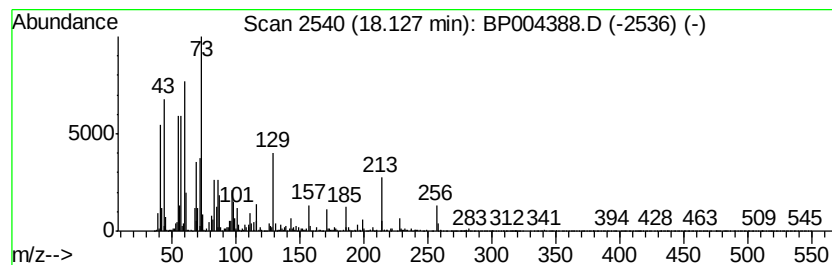
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 n-Hexadecanoic acid Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.13 | 2.35 ng/ul | 373973 | Phenanthrene-d10 | 17.23 |

| Hit# of | 5                   | Tentative ID | MW       | MolForm     | CAS# | Qual |
|---------|---------------------|--------------|----------|-------------|------|------|
| 1       | n-Hexadecanoic acid | 256          | C16H32O2 | 000057-10-3 | 99   |      |
| 2       | n-Hexadecanoic acid | 256          | C16H32O2 | 000057-10-3 | 98   |      |
| 3       | n-Hexadecanoic acid | 256          | C16H32O2 | 000057-10-3 | 95   |      |
| 4       | Pentadecanoic acid  | 242          | C15H30O2 | 001002-84-2 | 91   |      |
| 5       | Pentadecanoic acid  | 242          | C15H30O2 | 001002-84-2 | 87   |      |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\  
Data File : BP004388.D  
Acq On : 19 Dec 2020 11:40  
Operator : CG/JU  
Sample : L5128-02  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CB8C0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121820MA.M  
Quant Title : SVOA CALIBRATION

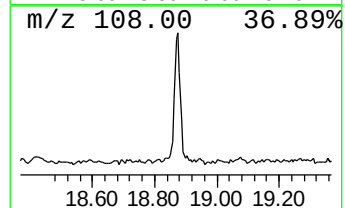
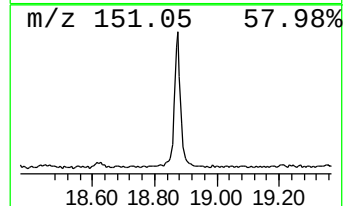
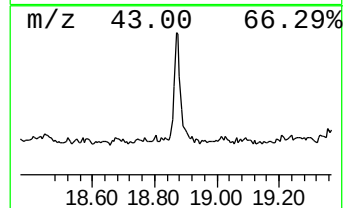
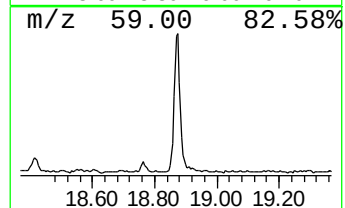
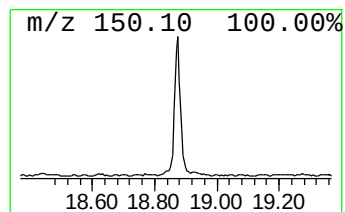
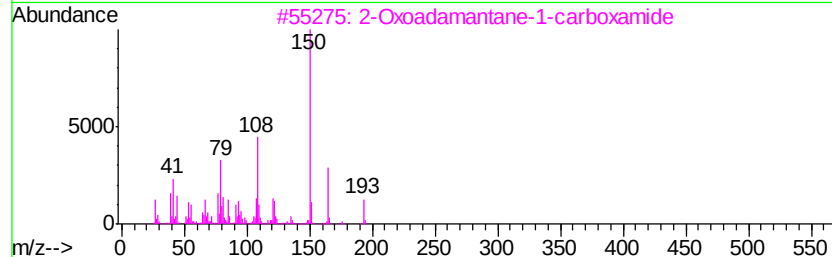
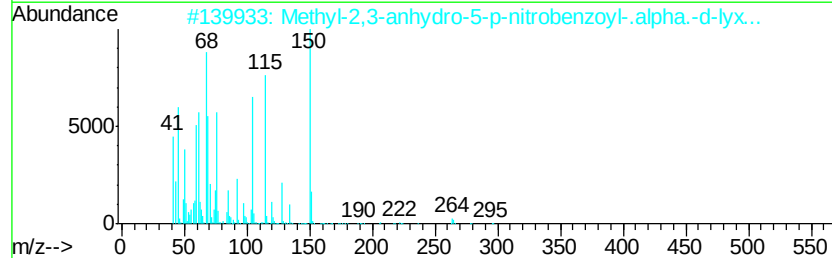
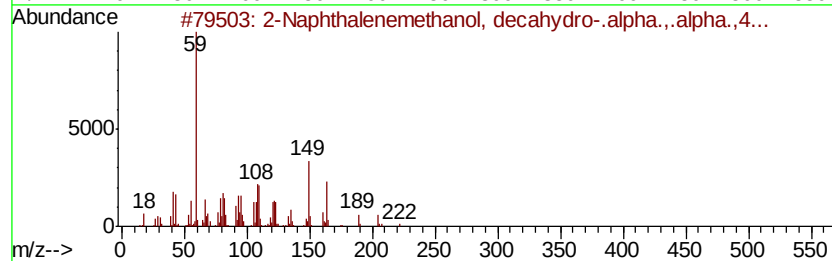
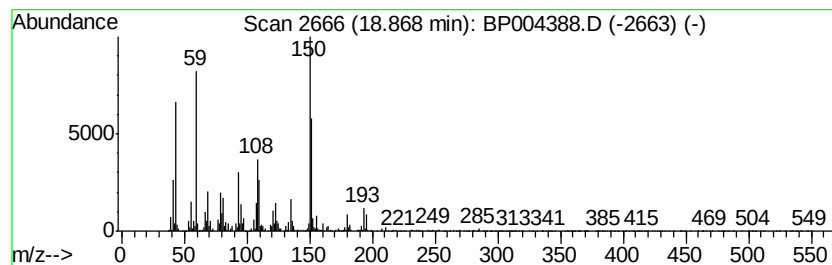
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 unknown-04 Concentration Rank 11

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.87 | 2.30 ng/ul | 365906 | Phenanthrene-d10 | 17.23 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 2-Naphthalenemethanol, decahydro... | 222 | C15H26O      | 000473-15-4  | 38      |      |      |
| 2    | Methyl-2,3-anhydro-5-p-nitrobenz... | 295 | C13H13NO7    | 1000214-61-0 | 35      |      |      |
| 3    | 2-Oxadamantane-1-carboxamide        | 193 | C11H15NO2    | 041171-77-1  | 25      |      |      |
| 4    | 1-Methyl-4(equat)ethynyl-trans-d... | 193 | C12H19NO     | 057581-44-9  | 22      |      |      |
| 5    | 2-Benzothiazolamine                 | 150 | C7H6N2S      | 000136-95-8  | 22      |      |      |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\  
Data File : BP004388.D  
Acq On : 19 Dec 2020 11:40  
Operator : CG/JU  
Sample : L5128-02  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

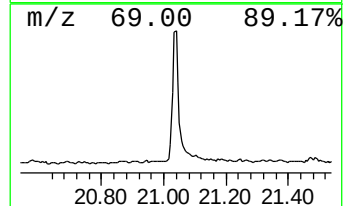
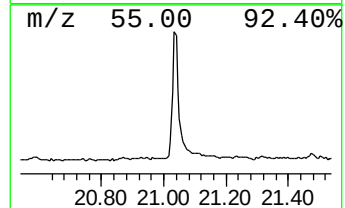
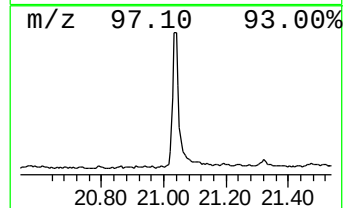
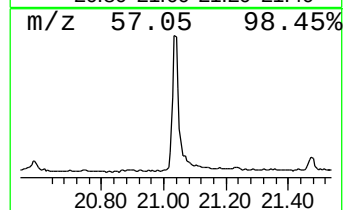
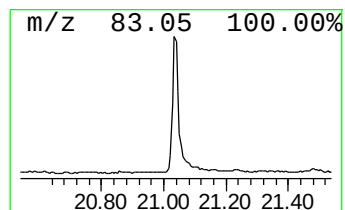
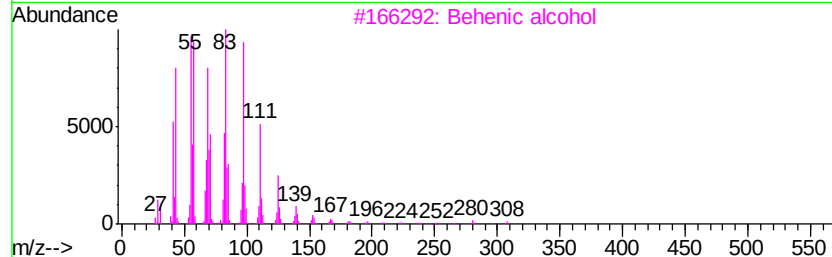
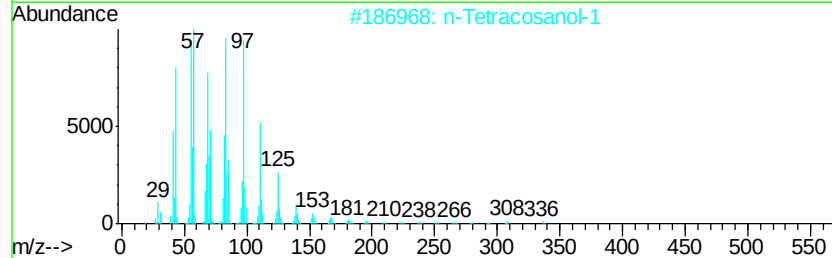
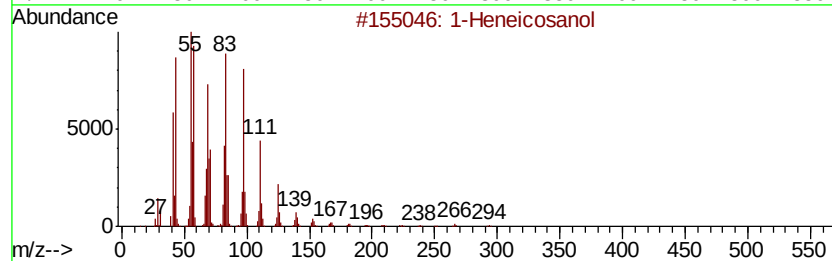
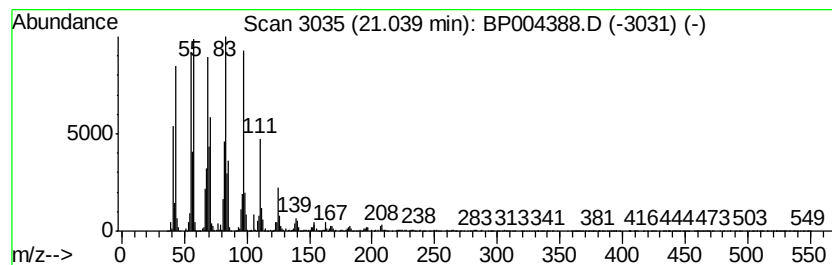
Instrument :  
BNA\_P  
ClientSampleId :  
CB8C0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121820MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 1-Heneicosanol Concentration Rank 3

| R.T.    | EstConc    | Area             | Relative to ISTD | R.T.           |
|---------|------------|------------------|------------------|----------------|
| 21.04   | 6.25 ng/ul | 1131220          | Chrysene-d12     | 21.32          |
| Hit# of | 5          | Tentative ID     | MW MolForm       | CAS# Qual      |
| 1       |            | 1-Heneicosanol   | 312 C21H44O      | 015594-90-8 94 |
| 2       |            | n-Tetracosanol-1 | 354 C24H50O      | 000506-51-4 94 |
| 3       |            | Behenic alcohol  | 326 C22H46O      | 000661-19-8 94 |
| 4       |            | 1-Heptacosanol   | 396 C27H56O      | 002004-39-9 94 |
| 5       |            | n-Nonadecanol-1  | 284 C19H40O      | 001454-84-8 94 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\  
Data File : BP004388.D  
Acq On : 19 Dec 2020 11:40  
Operator : CG/JU  
Sample : L5128-02  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CB8C0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121820MA.M  
Quant Title : SVOA CALIBRATION

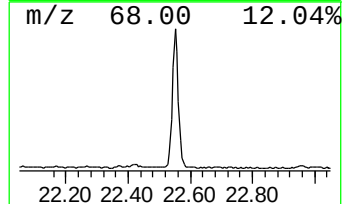
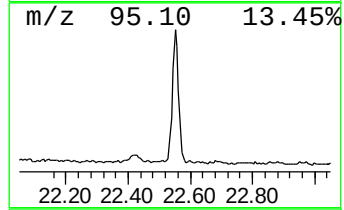
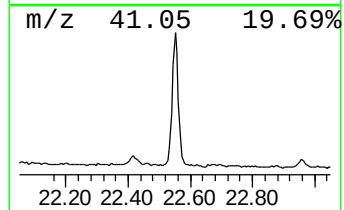
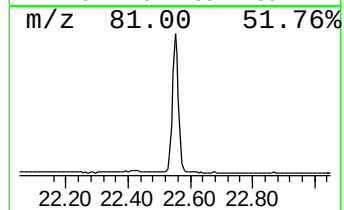
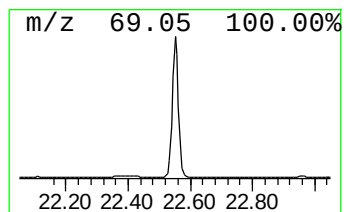
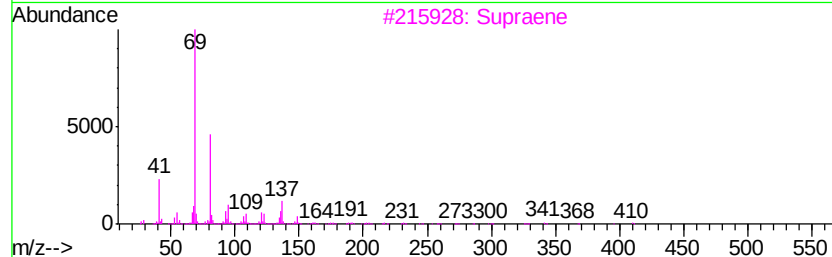
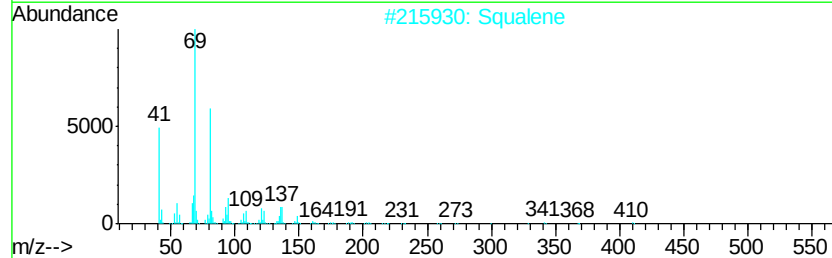
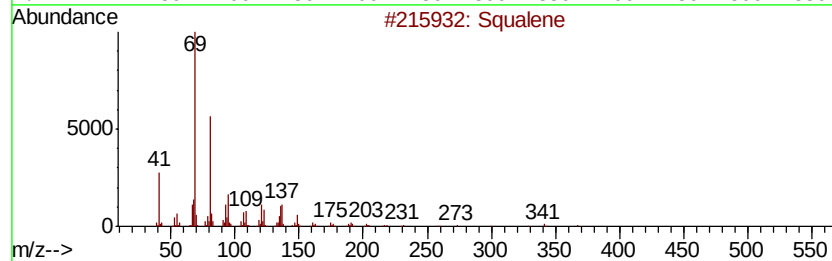
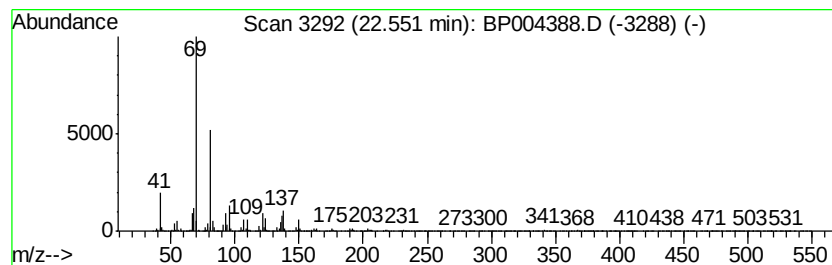
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Squalene Concentration Rank 2

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 22.55 | 8.36 ng/ul | 1567970 | Perylene-d12     | 23.68 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Squalene     | 410 | C30H50  | 000111-02-4 | 94   |
| 2    |    | Squalene     | 410 | C30H50  | 000111-02-4 | 91   |
| 3    |    | Supraene     | 410 | C30H50  | 007683-64-9 | 91   |
| 4    |    | Squalene     | 410 | C30H50  | 000111-02-4 | 90   |
| 5    |    | Squalene     | 410 | C30H50  | 000111-02-4 | 90   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA\_P\DATA\BP121820\  
Data File : BP004388.D  
Acq On : 19 Dec 2020 11:40  
Operator : CG/JU  
Sample : L5128-02  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CB8C0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121820MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| Paraldehyde          | 4.12  | 2.4     | ng/ul | 160054   | 1                     | 7.82  | 1325310 | 20.0 |
| 1,3-Oxathiolane      | 4.48  | 5.8     | ng/ul | 382634   | 1                     | 7.82  | 1325310 | 20.0 |
| Tetramethylbutane... | 7.89  | 2.2     | ng/ul | 145429   | 1                     | 7.82  | 1325310 | 20.0 |
| Benzenamine, 3-me... | 8.80  | 15.0    | ng/ul | 992408   | 1                     | 7.82  | 1325310 | 20.0 |
| Hexanoic acid, 3,... | 9.43  | 3.0     | ng/ul | 288396   | 2                     | 10.62 | 1921800 | 20.0 |
| unknown-01           | 10.16 | 3.4     | ng/ul | 330201   | 2                     | 10.62 | 1921800 | 20.0 |
| unknown-02           | 12.15 | 3.5     | ng/ul | 335421   | 2                     | 10.62 | 1921800 | 20.0 |
| unknown-03           | 14.39 | 2.5     | ng/ul | 324179   | 3                     | 14.47 | 2629790 | 20.0 |
| n-Hexadecanoic acid  | 18.13 | 2.4     | ng/ul | 373973   | 4                     | 17.23 | 3183450 | 20.0 |
| unknown-04           | 18.87 | 2.3     | ng/ul | 365906   | 4                     | 17.23 | 3183450 | 20.0 |
| 1-Heneicosanol       | 21.04 | 6.3     | ng/ul | 1131220  | 5                     | 21.32 | 3622750 | 20.0 |
| Squalene             | 22.55 | 8.4     | ng/ul | 1567970  | 6                     | 23.68 | 3751410 | 20.0 |