

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
 Data File : BP004268.D  
 Acq On : 14 Dec 2020 11:39  
 Operator : CG/JU  
 Sample : L5063-01DL 2X  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0AD5DL

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-BP121020MA.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         | 7.005       | 642           | 649         | 657          | rVV      | 531994         | 1027866       | 5.00%           | 0.258%        |
| 2         | 7.375       | 706           | 712         | 722          | rVV      | 449610         | 787895        | 3.83%           | 0.198%        |
| 3         | 7.840       | 783           | 791         | 797          | rBV      | 1237695        | 2150109       | 10.47%          | 0.540%        |
| 4         | 8.540       | 905           | 910         | 918          | rVB      | 499612         | 831998        | 4.05%           | 0.209%        |
| 5         | 8.951       | 969           | 980         | 984          | rBV4     | 582377         | 1325326       | 6.45%           | 0.333%        |
| 6         | 9.204       | 1018          | 1023        | 1029         | rVB2     | 318193         | 644246        | 3.14%           | 0.162%        |
| 7         | 9.534       | 1074          | 1079        | 1081         | rVV      | 394981         | 648581        | 3.16%           | 0.163%        |
| 8         | 9.563       | 1081          | 1084        | 1090         | rVB      | 431292         | 670262        | 3.26%           | 0.168%        |
| 9         | 9.722       | 1101          | 1111        | 1116         | rBV4     | 479037         | 1092162       | 5.32%           | 0.274%        |
| 10        | 9.828       | 1124          | 1129        | 1137         | rVV3     | 684235         | 1526783       | 7.43%           | 0.383%        |
| 11        | 10.046      | 1160          | 1166        | 1171         | rVV3     | 741747         | 1355528       | 6.60%           | 0.340%        |
| 12        | 10.110      | 1173          | 1177        | 1182         | rVB2     | 361938         | 585645        | 2.85%           | 0.147%        |
| 13        | 10.257      | 1195          | 1202        | 1212         | rVB5     | 689378         | 1722629       | 8.38%           | 0.432%        |
| 14        | 10.346      | 1212          | 1217        | 1221         | rBV      | 426812         | 769108        | 3.74%           | 0.193%        |
| 15        | 10.546      | 1241          | 1251        | 1256         | rBV8     | 295769         | 1136282       | 5.53%           | 0.285%        |
| 16        | 10.634      | 1256          | 1266        | 1272         | rVV2     | 2323673        | 6145860       | 29.91%          | 1.542%        |
| 17        | 10.687      | 1272          | 1275        | 1283         | rVV3     | 1078414        | 2495434       | 12.15%          | 0.626%        |
| 18        | 10.763      | 1283          | 1288        | 1293         | rVV2     | 582985         | 1131442       | 5.51%           | 0.284%        |
| 19        | 10.810      | 1293          | 1296        | 1300         | rVV2     | 495906         | 697697        | 3.40%           | 0.175%        |
| 20        | 10.863      | 1300          | 1305        | 1310         | rVB      | 446401         | 706294        | 3.44%           | 0.177%        |
| 21        | 11.057      | 1334          | 1338        | 1343         | rBV3     | 511011         | 870188        | 4.24%           | 0.218%        |
| 22        | 11.140      | 1348          | 1352        | 1357         | rVB      | 822435         | 1275727       | 6.21%           | 0.320%        |
| 23        | 11.310      | 1377          | 1381        | 1383         | rBV2     | 418188         | 688612        | 3.35%           | 0.173%        |
| 24        | 11.422      | 1395          | 1400        | 1405         | rBV2     | 457999         | 815061        | 3.97%           | 0.205%        |
| 25        | 11.487      | 1406          | 1411        | 1418         | rVV8     | 365994         | 987536        | 4.81%           | 0.248%        |
| 26        | 11.563      | 1418          | 1424        | 1431         | rVB2     | 1045729        | 2199337       | 10.70%          | 0.552%        |
| 27        | 11.681      | 1439          | 1444        | 1446         | rBV2     | 1149457        | 1865953       | 9.08%           | 0.468%        |
| 28        | 11.751      | 1452          | 1456        | 1462         | rVB      | 1105529        | 1765991       | 8.60%           | 0.443%        |
| 29        | 11.840      | 1466          | 1471        | 1478         | rVB3     | 948577         | 1576175       | 7.67%           | 0.396%        |
| 30        | 11.940      | 1484          | 1488        | 1491         | rVB2     | 633300         | 774364        | 3.77%           | 0.194%        |
| 31        | 12.034      | 1500          | 1504        | 1506         | rBV      | 403533         | 627477        | 3.05%           | 0.157%        |
| 32        | 12.075      | 1506          | 1511        | 1514         | rVV3     | 921939         | 1690972       | 8.23%           | 0.424%        |
| 33        | 12.304      | 1541          | 1550        | 1559         | rVB2     | 3379221        | 7612113       | 37.05%          | 1.910%        |
| 34        | 12.404      | 1559          | 1567        | 1570         | rBV4     | 805046         | 2197934       | 10.70%          | 0.552%        |

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 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0AD5DL

## 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-BP121020MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |          |        |        |
|----|--------|------|------|------|------|---------|----------|--------|--------|
| 35 | 12.481 | 1576 | 1580 | 1581 | rBV4 | 498723  | 611831   | 2.98%  | 0.154% |
| 36 | 12.522 | 1582 | 1587 | 1591 | rVV2 | 2435567 | 4294529  | 20.90% | 1.078% |
| 37 | 12.610 | 1599 | 1602 | 1606 | rBV  | 559791  | 697845   | 3.40%  | 0.175% |
| 38 | 12.704 | 1614 | 1618 | 1622 | rBV4 | 628994  | 1021179  | 4.97%  | 0.256% |
| 39 | 13.004 | 1662 | 1669 | 1670 | rBV6 | 783290  | 1539049  | 7.49%  | 0.386% |
| 40 | 13.034 | 1671 | 1674 | 1679 | rVB2 | 1701304 | 2444679  | 11.90% | 0.614% |
| 41 | 13.151 | 1690 | 1694 | 1699 | rBV4 | 807248  | 1515592  | 7.38%  | 0.380% |
| 42 | 13.322 | 1717 | 1723 | 1729 | rVB3 | 1852256 | 3491540  | 16.99% | 0.876% |
| 43 | 13.422 | 1736 | 1740 | 1746 | rVB2 | 2070100 | 3404093  | 16.57% | 0.854% |
| 44 | 13.516 | 1753 | 1756 | 1759 | rBV  | 933183  | 1275773  | 6.21%  | 0.320% |
| 45 | 13.651 | 1774 | 1779 | 1781 | rBV  | 3326340 | 5533458  | 26.93% | 1.389% |
| 46 | 13.816 | 1801 | 1807 | 1812 | rBV  | 6820636 | 11842936 | 57.64% | 2.972% |
| 47 | 13.869 | 1812 | 1816 | 1822 | rBV2 | 4226883 | 6424293  | 31.27% | 1.612% |
| 48 | 13.987 | 1832 | 1836 | 1839 | rVB  | 1485394 | 1730778  | 8.42%  | 0.434% |
| 49 | 14.051 | 1843 | 1847 | 1850 | rVV  | 3025860 | 3982036  | 19.38% | 0.999% |
| 50 | 14.187 | 1866 | 1870 | 1872 | rBV  | 1342879 | 1858190  | 9.04%  | 0.466% |
| 51 | 14.334 | 1891 | 1895 | 1901 | rVB3 | 1368797 | 2588820  | 12.60% | 0.650% |
| 52 | 14.404 | 1902 | 1907 | 1911 | rBV  | 2402072 | 3597920  | 17.51% | 0.903% |
| 53 | 14.492 | 1915 | 1922 | 1927 | rVB3 | 3436153 | 6986488  | 34.01% | 1.753% |
| 54 | 14.581 | 1935 | 1937 | 1941 | rVB2 | 946909  | 1077453  | 5.24%  | 0.270% |
| 55 | 14.628 | 1941 | 1945 | 1951 | rVB  | 1639009 | 2467311  | 12.01% | 0.619% |
| 56 | 14.716 | 1953 | 1960 | 1967 | rVB2 | 3304504 | 9689554  | 47.16% | 2.432% |
| 57 | 14.786 | 1968 | 1972 | 1977 | rBV3 | 1925066 | 3073611  | 14.96% | 0.771% |
| 58 | 14.898 | 1984 | 1991 | 1997 | rVB  | 6628095 | 13112193 | 63.82% | 3.291% |
| 59 | 14.975 | 1998 | 2004 | 2009 | rVB  | 7209391 | 11391197 | 55.44% | 2.859% |
| 60 | 15.028 | 2010 | 2013 | 2020 | rVB5 | 1549544 | 2730807  | 13.29% | 0.685% |
| 61 | 15.116 | 2023 | 2028 | 2033 | rBV  | 6269836 | 11227322 | 54.65% | 2.818% |
| 62 | 15.169 | 2033 | 2037 | 2041 | rVB  | 5025279 | 6834223  | 33.26% | 1.715% |
| 63 | 15.292 | 2050 | 2058 | 2061 | rBV2 | 4097328 | 9558514  | 46.52% | 2.399% |
| 64 | 15.322 | 2061 | 2063 | 2066 | rVV  | 3267241 | 3631477  | 17.68% | 0.911% |
| 65 | 15.357 | 2066 | 2069 | 2073 | rVB2 | 2310271 | 2764283  | 13.45% | 0.694% |
| 66 | 15.451 | 2082 | 2085 | 2090 | rBV4 | 1816687 | 2514916  | 12.24% | 0.631% |
| 67 | 15.539 | 2097 | 2100 | 2104 | rVB2 | 2873464 | 4026671  | 19.60% | 1.011% |
| 68 | 15.598 | 2105 | 2110 | 2114 | rBV3 | 2401600 | 4445165  | 21.64% | 1.116% |
| 69 | 15.669 | 2119 | 2122 | 2126 | rVB3 | 2792122 | 4141201  | 20.16% | 1.039% |
| 70 | 15.763 | 2134 | 2138 | 2145 | rVB2 | 5464573 | 8675703  | 42.23% | 2.177% |
| 71 | 15.828 | 2146 | 2149 | 2153 | rVV3 | 1543966 | 3085554  | 15.02% | 0.774% |

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Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

## 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-BP121020MA.M  
Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |          |         |        |
|-----|--------|------|------|------|------|---------|----------|---------|--------|
| 72  | 15.869 | 2153 | 2156 | 2160 | rVB  | 3589697 | 4745050  | 23.10%  | 1.191% |
| 73  | 15.922 | 2161 | 2165 | 2169 | rBV3 | 2063606 | 3802037  | 18.51%  | 0.954% |
| 74  | 16.004 | 2176 | 2179 | 2185 | rVB2 | 2564675 | 3901946  | 18.99%  | 0.979% |
| 75  | 16.063 | 2186 | 2189 | 2193 | rBV2 | 2402441 | 3230392  | 15.72%  | 0.811% |
| 76  | 16.228 | 2213 | 2217 | 2231 | rVB2 | 9145856 | 20545051 | 100.00% | 5.156% |
| 77  | 16.369 | 2236 | 2241 | 2244 | rBV  | 4785212 | 7145771  | 34.78%  | 1.793% |
| 78  | 16.481 | 2257 | 2260 | 2262 | rBV4 | 1570086 | 2074046  | 10.10%  | 0.521% |
| 79  | 16.610 | 2279 | 2282 | 2287 | rVB3 | 3484384 | 5227704  | 25.45%  | 1.312% |
| 80  | 16.910 | 2330 | 2333 | 2339 | rBV4 | 2147415 | 3725038  | 18.13%  | 0.935% |
| 81  | 17.028 | 2350 | 2353 | 2356 | rVB2 | 2846050 | 3190412  | 15.53%  | 0.801% |
| 82  | 17.075 | 2357 | 2361 | 2370 | rVB5 | 5319874 | 11072189 | 53.89%  | 2.779% |
| 83  | 17.257 | 2388 | 2392 | 2395 | rBV  | 2803951 | 4228342  | 20.58%  | 1.061% |
| 84  | 17.298 | 2395 | 2399 | 2403 | rVB  | 5839969 | 7398022  | 36.01%  | 1.857% |
| 85  | 17.839 | 2488 | 2491 | 2494 | rVB  | 2926758 | 3574293  | 17.40%  | 0.897% |
| 86  | 18.127 | 2536 | 2540 | 2544 | rBV  | 5949285 | 9219863  | 44.88%  | 2.314% |
| 87  | 18.175 | 2544 | 2548 | 2556 | rVB  | 8621085 | 13400272 | 65.22%  | 3.363% |
| 88  | 18.310 | 2567 | 2571 | 2575 | rBV  | 5838171 | 8443657  | 41.10%  | 2.119% |
| 89  | 18.351 | 2575 | 2578 | 2583 | rVB  | 4887924 | 6663541  | 32.43%  | 1.672% |
| 90  | 18.439 | 2589 | 2593 | 2601 | rBV3 | 3356730 | 5299202  | 25.79%  | 1.330% |
| 91  | 18.510 | 2602 | 2605 | 2609 | rVB2 | 3056387 | 4043329  | 19.68%  | 1.015% |
| 92  | 18.610 | 2620 | 2622 | 2626 | rBV4 | 1744177 | 2148304  | 10.46%  | 0.539% |
| 93  | 18.827 | 2655 | 2659 | 2660 | rBV2 | 2611011 | 3280487  | 15.97%  | 0.823% |
| 94  | 18.898 | 2669 | 2671 | 2675 | rVB  | 2129432 | 2347469  | 11.43%  | 0.589% |
| 95  | 19.022 | 2687 | 2692 | 2695 | rBV  | 8363518 | 12468337 | 60.69%  | 3.129% |
| 96  | 19.110 | 2704 | 2707 | 2711 | rVB2 | 3182492 | 3568461  | 17.37%  | 0.896% |
| 97  | 19.157 | 2711 | 2715 | 2720 | rBV2 | 2229598 | 4503853  | 21.92%  | 1.130% |
| 98  | 19.269 | 2731 | 2734 | 2737 | rBV  | 2978945 | 4067492  | 19.80%  | 1.021% |
| 99  | 19.704 | 2804 | 2808 | 2813 | rVB  | 5080938 | 7634455  | 37.16%  | 1.916% |
| 100 | 20.404 | 2924 | 2927 | 2931 | rBV  | 4323392 | 5801467  | 28.24%  | 1.456% |

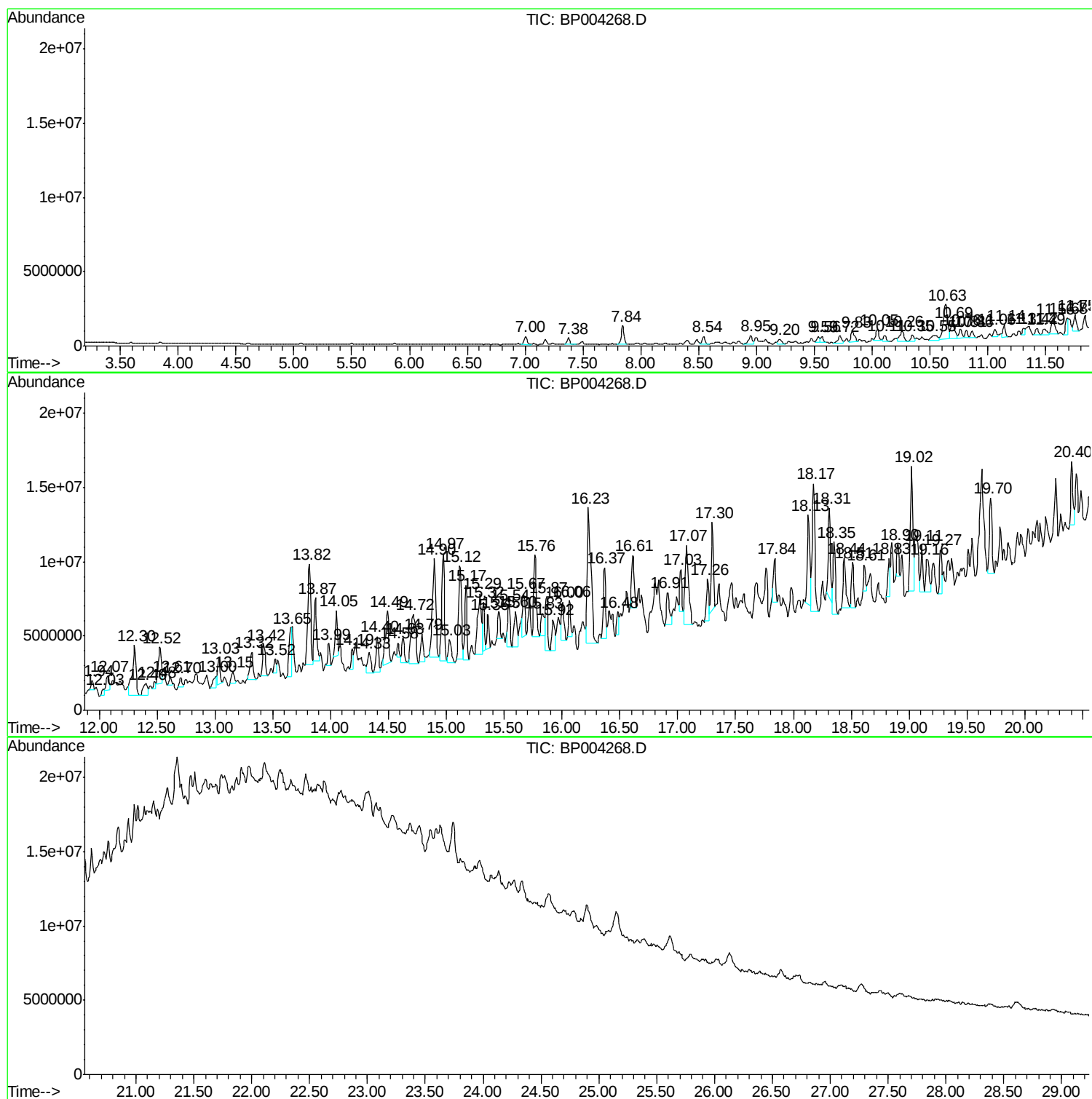
Sum of corrected areas: 398437253

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Instrument :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

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



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
 Quant Title : SVOA CALIBRATION

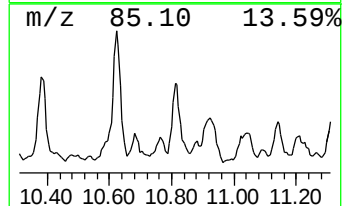
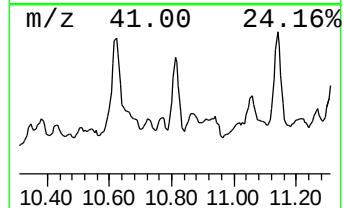
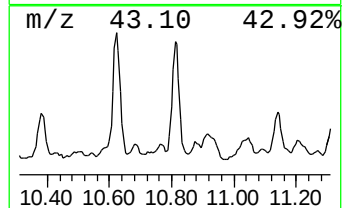
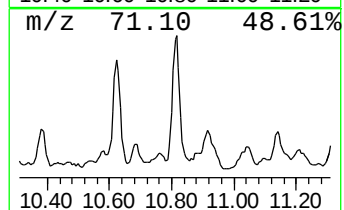
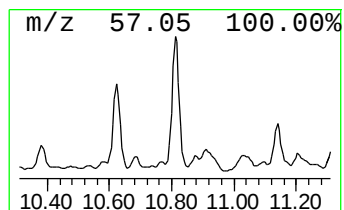
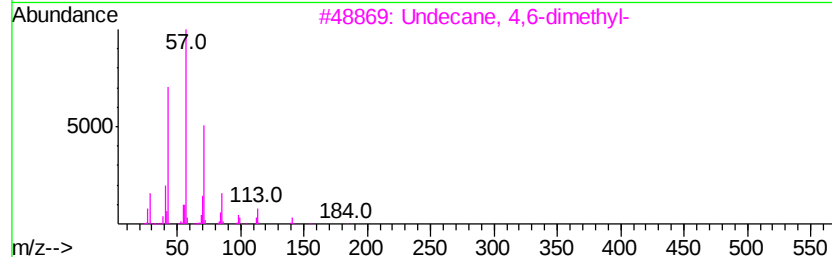
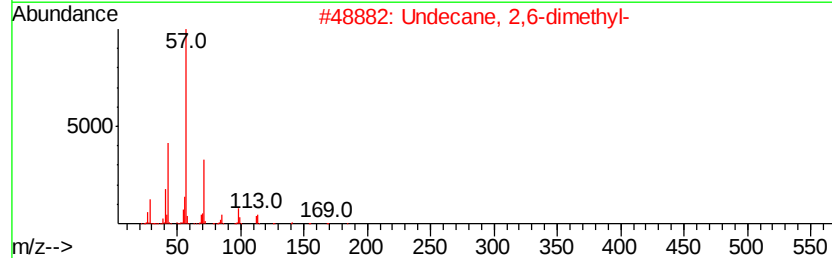
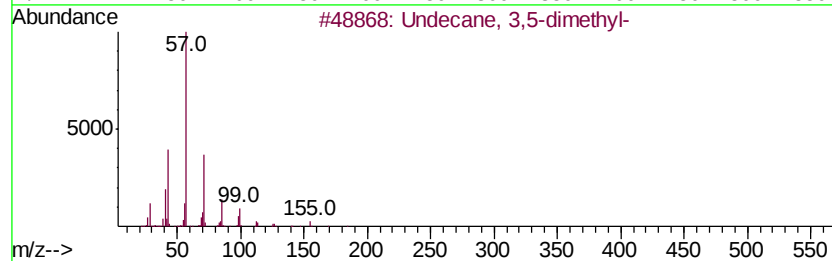
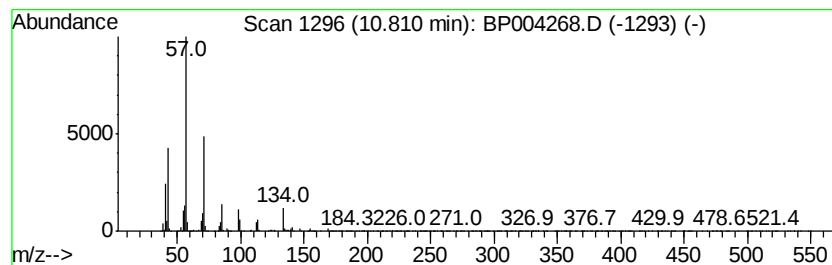
TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
 Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 70

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 10.81 | 2.27 ng/ul | 697697 | Napthalene-d8    | 10.64 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Undecane, 3,5-dimethyl-     | 184 | C13H28  | 017312-81-1 | 72   |
| 2    |    | Undecane, 2,6-dimethyl-     | 184 | C13H28  | 017301-23-4 | 70   |
| 3    |    | Undecane, 4,6-dimethyl-     | 184 | C13H28  | 017312-82-2 | 68   |
| 4    |    | Dodecane, 2,7,10-trimethyl- | 212 | C15H32  | 074645-98-0 | 53   |
| 5    |    | Dodecane, 4,6-dimethyl-     | 198 | C14H30  | 061141-72-8 | 50   |



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Quant Title : SVOA CALIBRATION

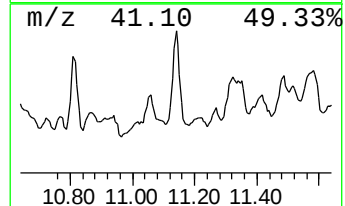
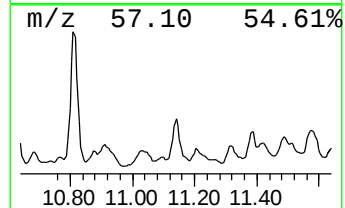
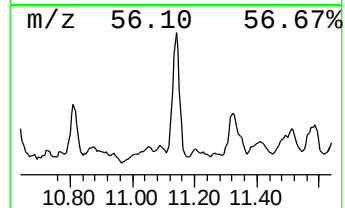
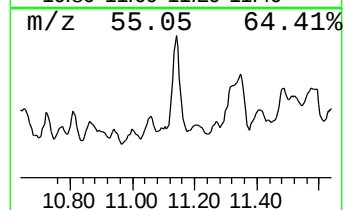
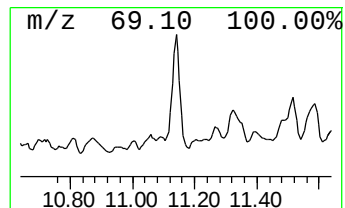
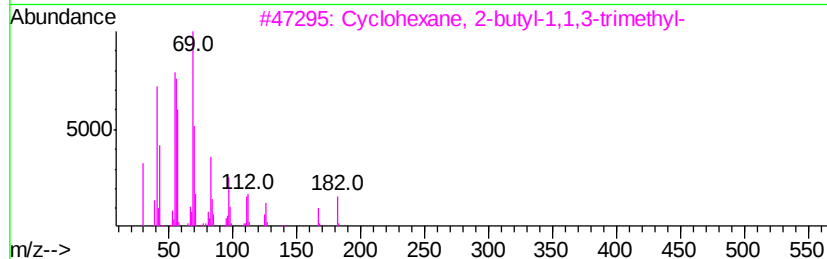
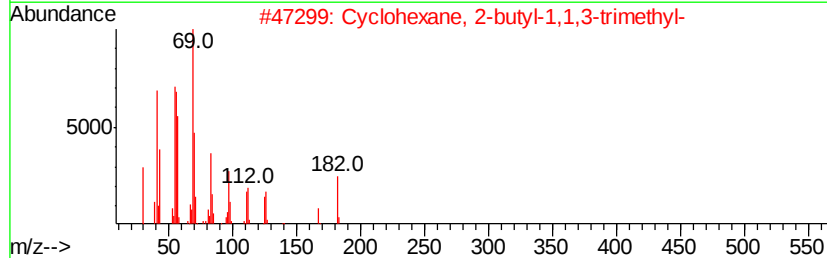
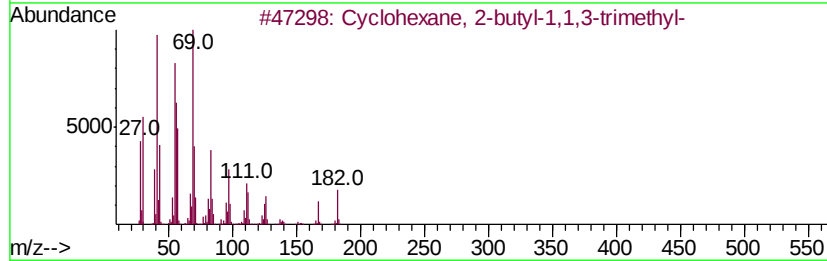
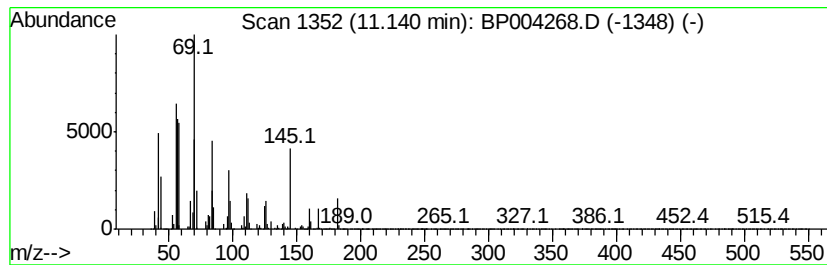
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 (DEL) Alkane: Cyclic11.14 Concentration Rank 60

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 11.14 | 4.15 ng/ul | 1275730 | Naphthalene-d8   | 10.64 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Cyclohexane, 2-butyl-1,1,3-trime... | 182 | C13H26     | 054676-39-0  | 96   |
| 2    |    | Cyclohexane, 2-butyl-1,1,3-trime... | 182 | C13H26     | 054676-39-0  | 89   |
| 3    |    | Cyclohexane, 2-butyl-1,1,3-trime... | 182 | C13H26     | 054676-39-0  | 89   |
| 4    |    | 4-Trifluoroacetoxvtridecane         | 296 | C15H27F3O2 | 1000245-47-3 | 74   |
| 5    |    | Cyclooctane, 1,2-dimethyl-          | 140 | C10H20     | 013151-94-5  | 52   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

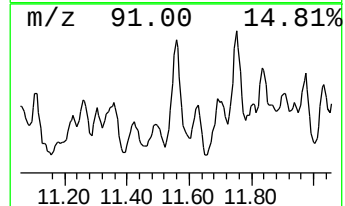
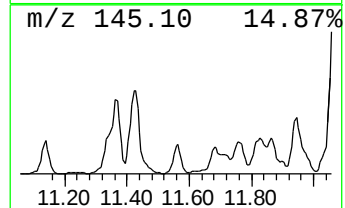
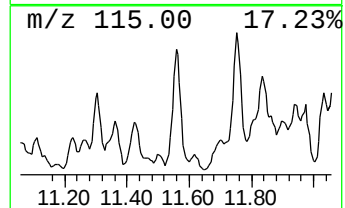
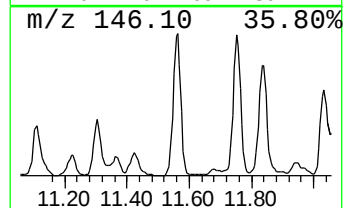
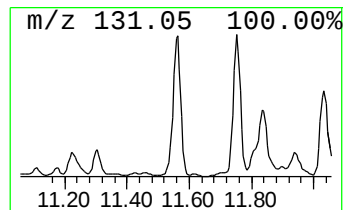
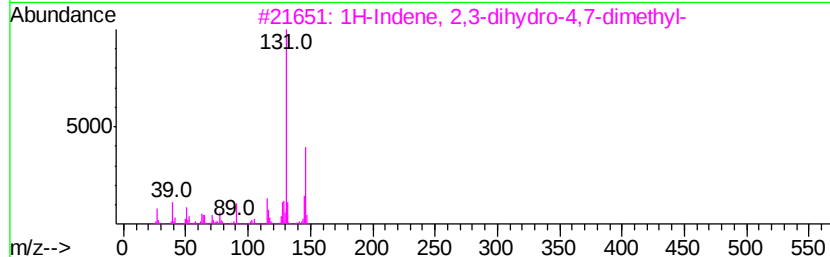
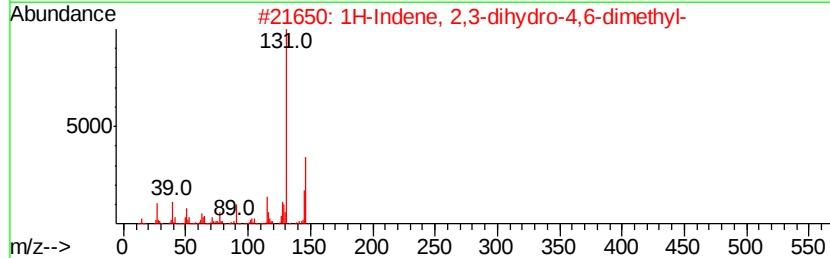
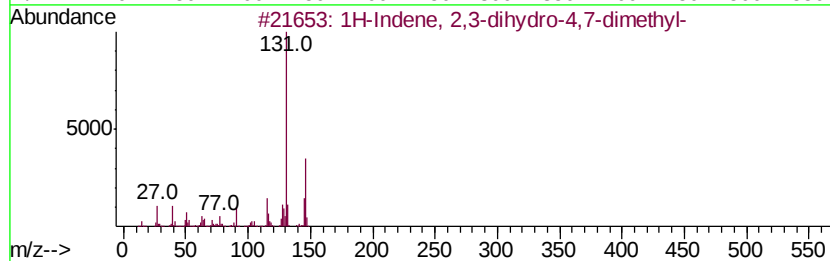
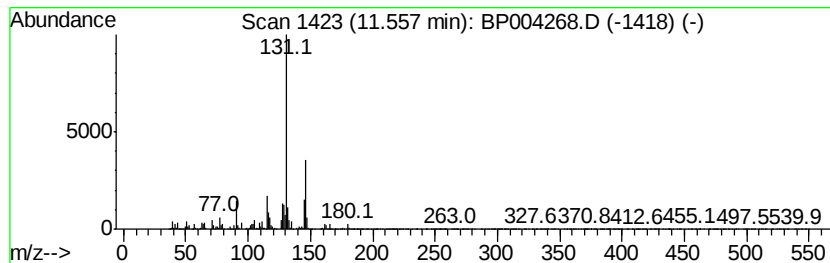
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 47

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 11.56 | 7.16 ng/ul | 2199340 | Napthalene-d8    | 10.64 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 97   |
| 2    |    | 1H-Indene, 2,3-dihydro-4,6-dimet... | 146 | C11H14  | 001685-82-1 | 94   |
| 3    |    | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 94   |
| 4    |    | 1H-Indene, 2,3-dihydro-1,3-dimet... | 146 | C11H14  | 004175-53-5 | 94   |
| 5    |    | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

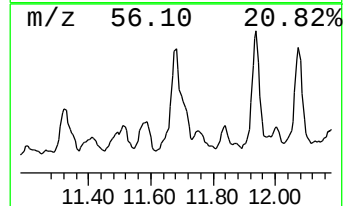
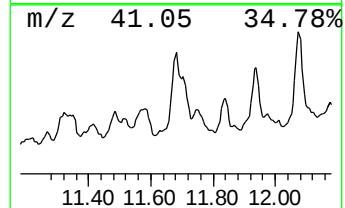
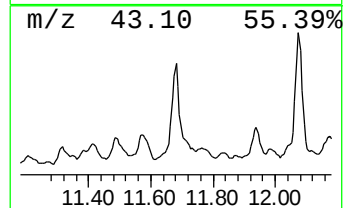
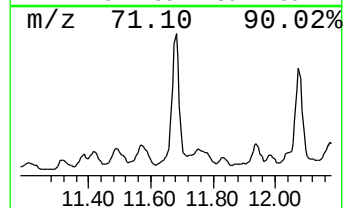
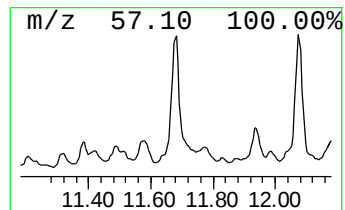
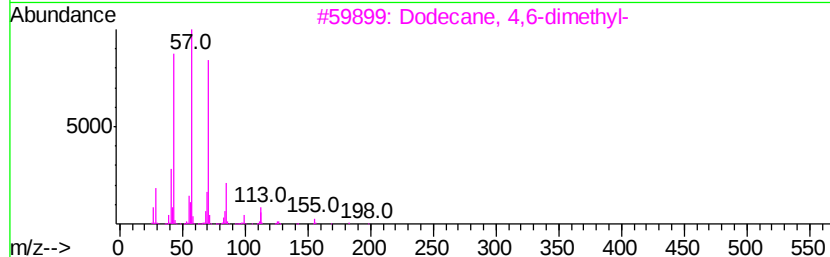
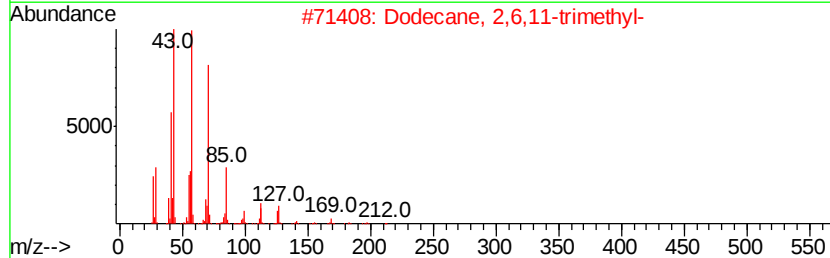
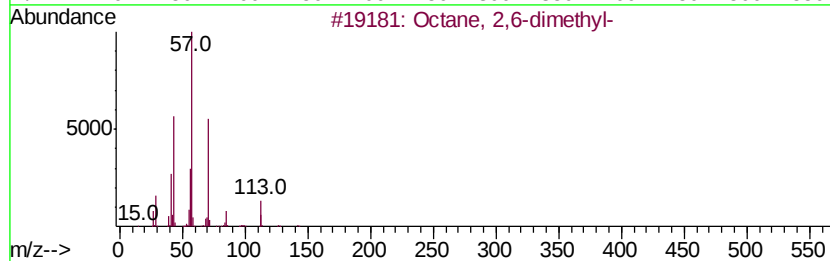
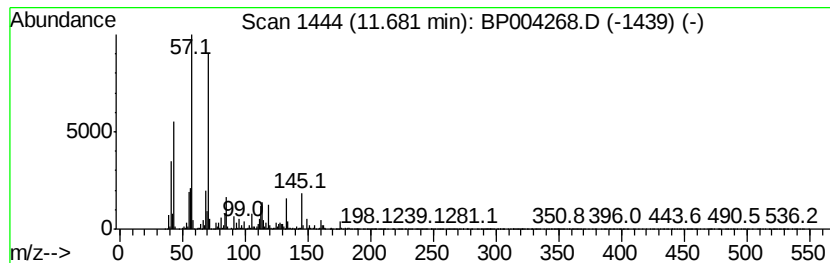
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 51

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 11.68 | 6.07 ng/ul | 1865950 | Naphthalene-d8   | 10.64 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Octane, 2,6-dimethyl-       | 142 | C10H22  | 002051-30-1 | 72   |
| 2    |    |   | Dodecane, 2,6,11-trimethyl- | 212 | C15H32  | 031295-56-4 | 62   |
| 3    |    |   | Dodecane, 4,6-dimethyl-     | 198 | C14H30  | 061141-72-8 | 58   |
| 4    |    |   | Heptane, 3-ethyl-5-methyl-  | 142 | C10H22  | 052896-90-9 | 58   |
| 5    |    |   | Nonane, 2,6-dimethyl-       | 156 | C11H24  | 017302-28-2 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

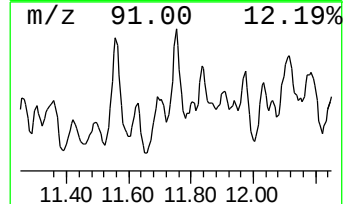
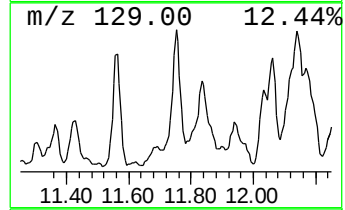
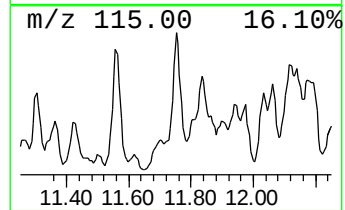
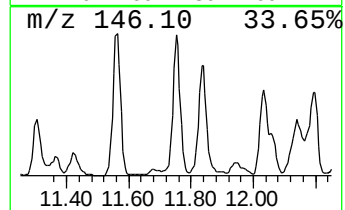
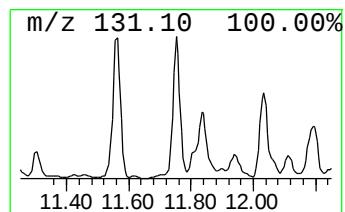
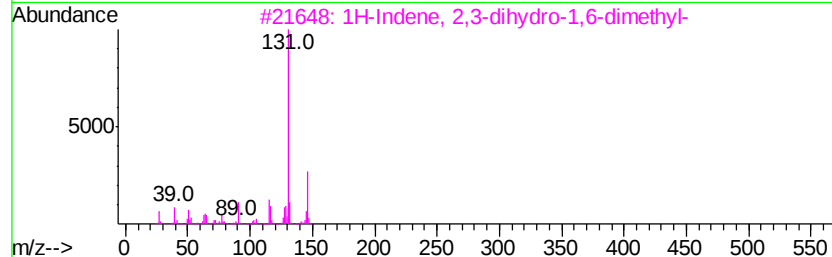
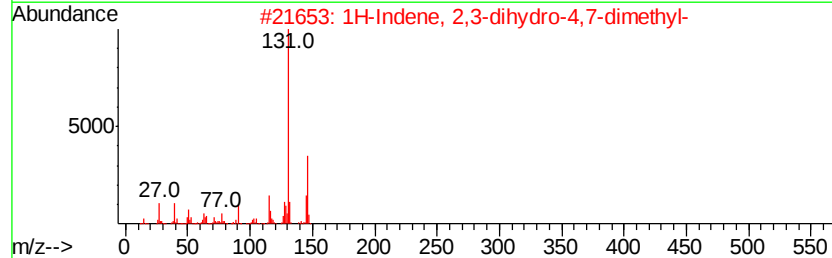
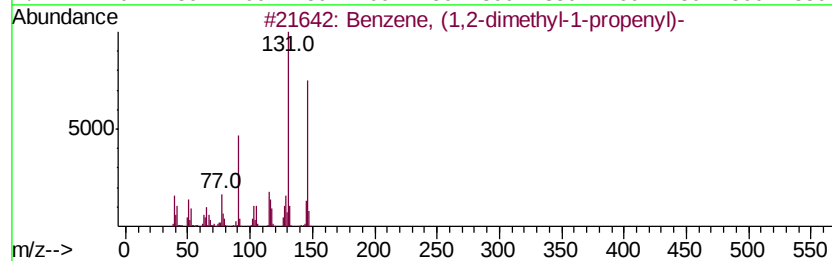
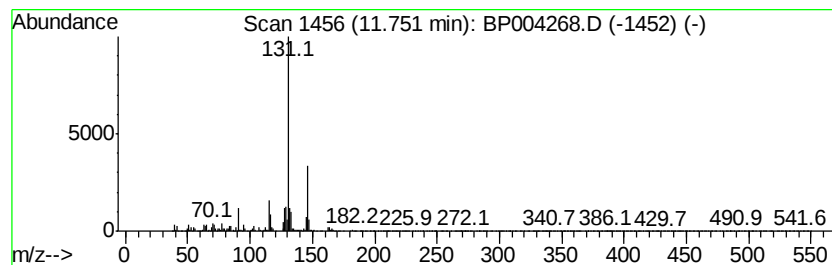
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Benzene, (1,2-dimethyl-1-pr... Concentration Rank 53

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 11.75 | 5.75 ng/ul | 1765990 | Napthalene-d8    | 10.64 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, (1,2-dimethyl-1-propenyl)- | 146 | C11H14  | 000769-57-3 | 93   |
| 2    |    | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 93   |
| 3    |    | 1H-Indene, 2,3-dihydro-1,6-dimet... | 146 | C11H14  | 017059-48-2 | 93   |
| 4    |    | 1H-Indene, 2,3-dihydro-1,3-dimet... | 146 | C11H14  | 004175-53-5 | 90   |
| 5    |    | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

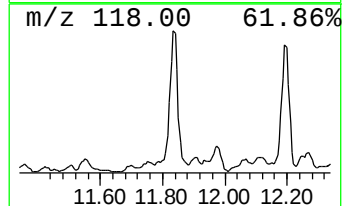
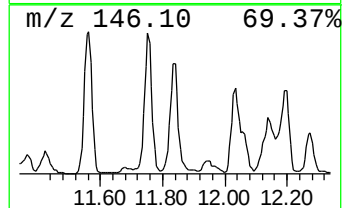
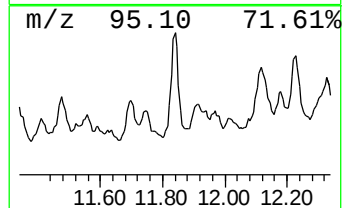
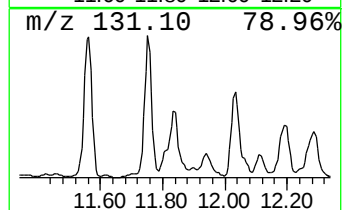
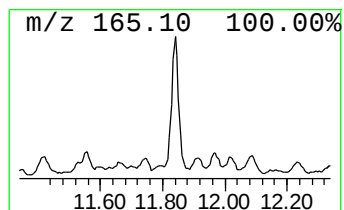
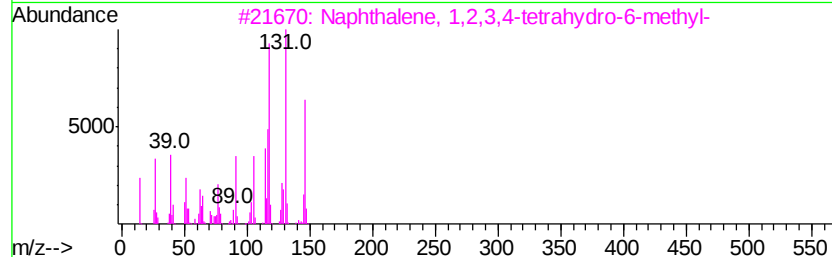
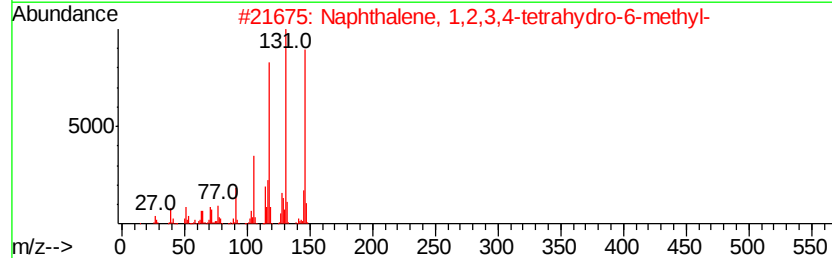
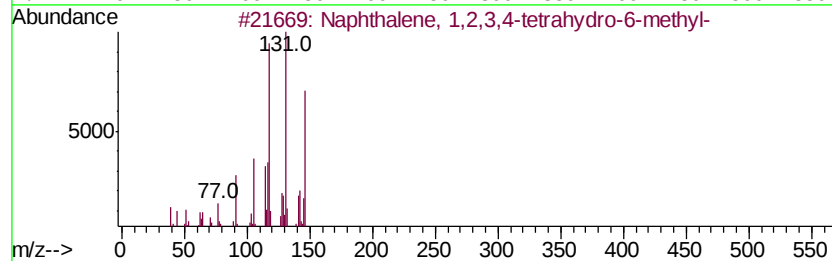
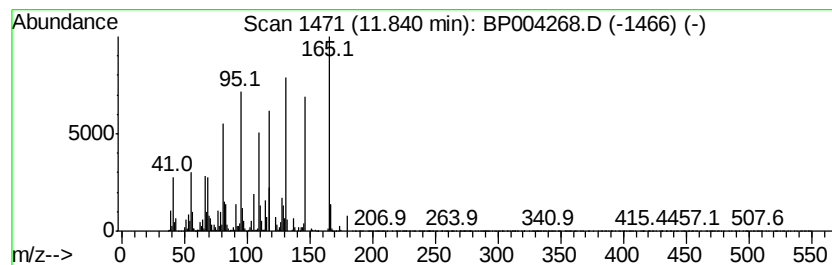
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 55

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 11.84 | 5.13 ng/ul | 1576180 | Naphthalene-d8   | 10.64 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 001680-51-9 | 70   |
| 2    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 001680-51-9 | 64   |
| 3    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 001680-51-9 | 50   |
| 4    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 001680-51-9 | 50   |
| 5    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 002809-64-5 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

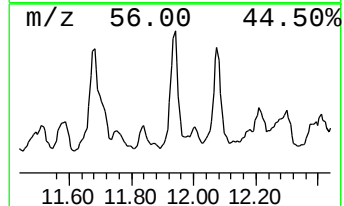
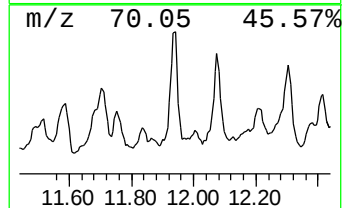
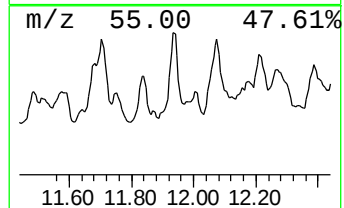
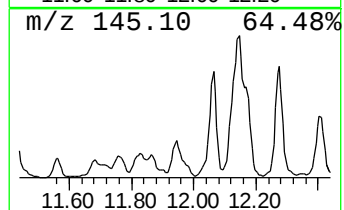
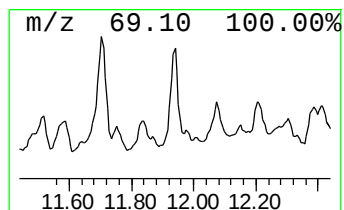
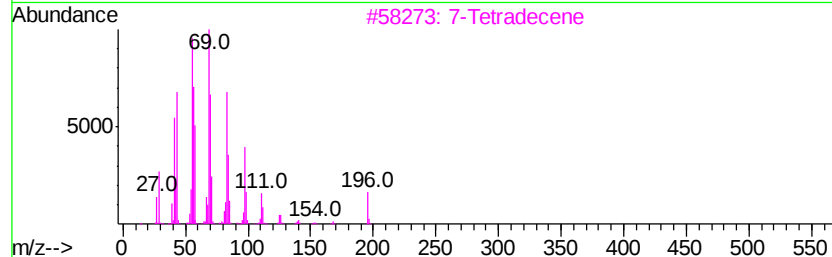
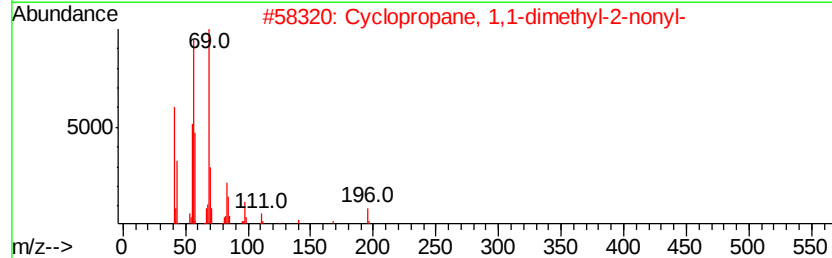
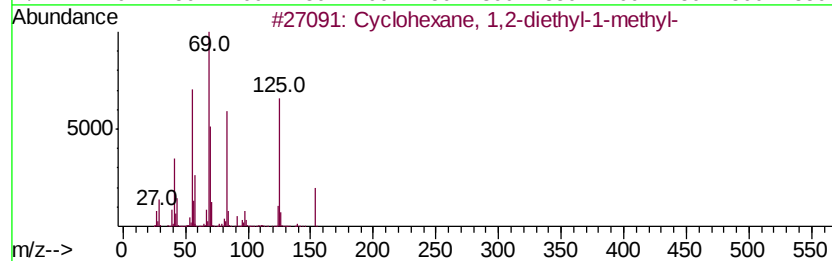
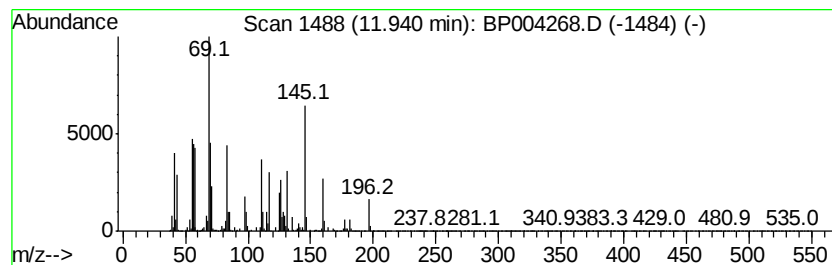
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 (DEL) Alkane: Cyclic11.94 Concentration Rank 67

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.94 | 2.52 ng/ul | 774364 | Naphthalene-d8   | 10.64 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, 1,2-diethyl-1-methyl-  | 154 | C11H22  | 061141-79-5 | 38   |
| 2    |    | Cyclopropane, 1,1-dimethyl-2-nonyl- | 196 | C14H28  | 041977-38-2 | 30   |
| 3    |    | 7-Tetradecene                       | 196 | C14H28  | 010374-74-0 | 22   |
| 4    |    | 9-Oxabicyclo[6.1.0]nonane, 1-met... | 140 | C9H16O  | 052954-47-9 | 22   |
| 5    |    | Cyclohexane, 1,3,5-trimethyl-       | 126 | C9H18   | 001839-63-0 | 22   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

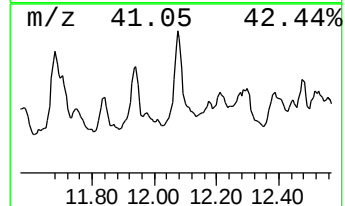
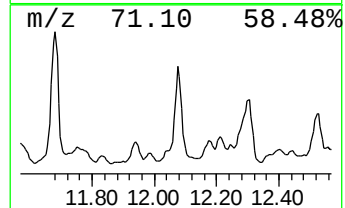
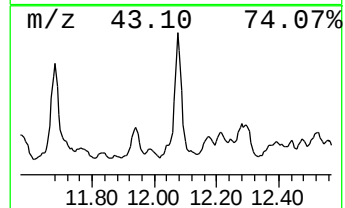
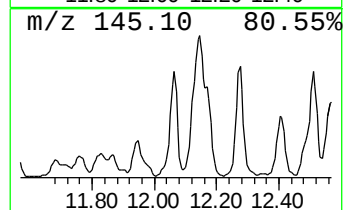
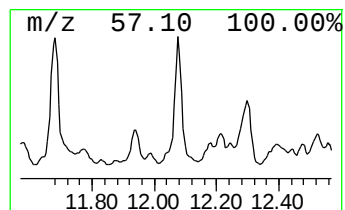
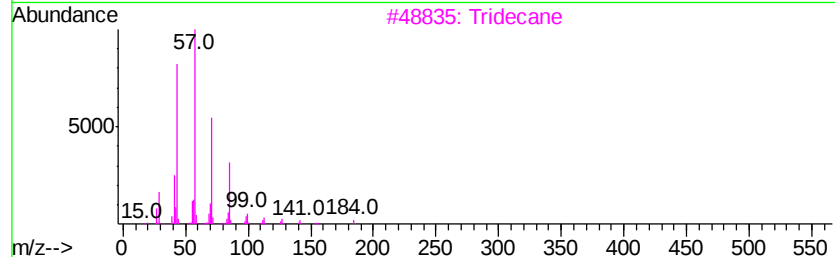
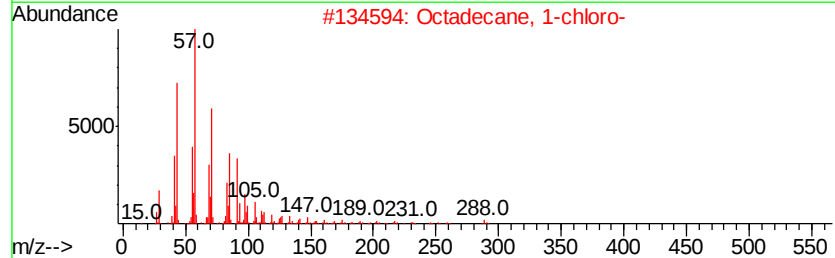
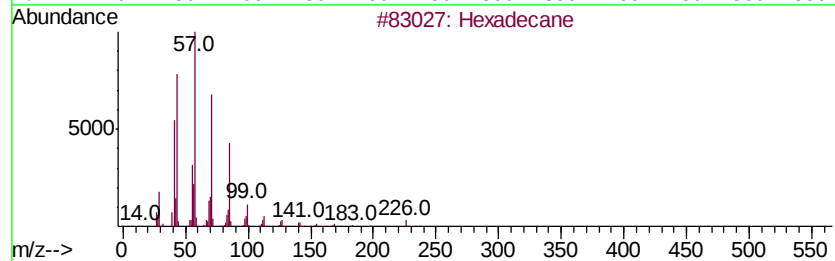
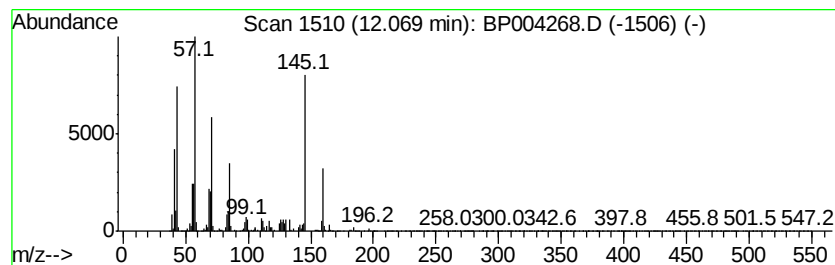
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 54

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 12.07 | 5.50 ng/ul | 1690970 | Napthalene-d8    | 10.64 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm  | CAS#         | Qual |
|------|----|---|---------------------------------|-----|----------|--------------|------|
| 1    |    |   | Hexadecane                      | 226 | C16H34   | 000544-76-3  | 76   |
| 2    |    |   | Octadecane, 1-chloro-           | 288 | C18H37Cl | 003386-33-2  | 76   |
| 3    |    |   | Tridecane                       | 184 | C13H28   | 000629-50-5  | 70   |
| 4    |    |   | Oxalic acid, decyl propyl ester | 272 | C15H28O4 | 1000309-26-3 | 64   |
| 5    |    |   | Hexadecane                      | 226 | C16H34   | 000544-76-3  | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

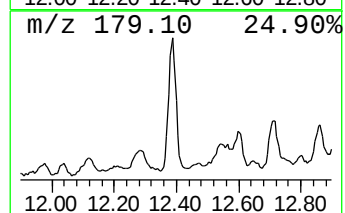
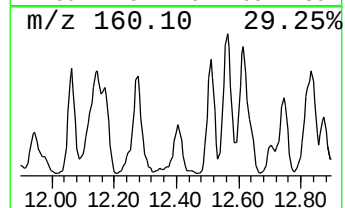
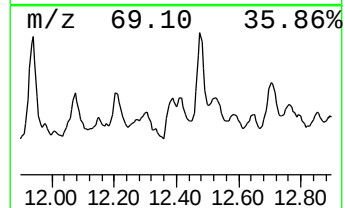
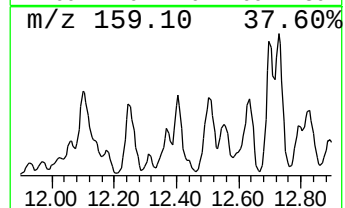
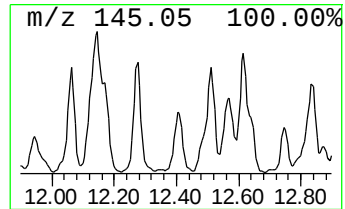
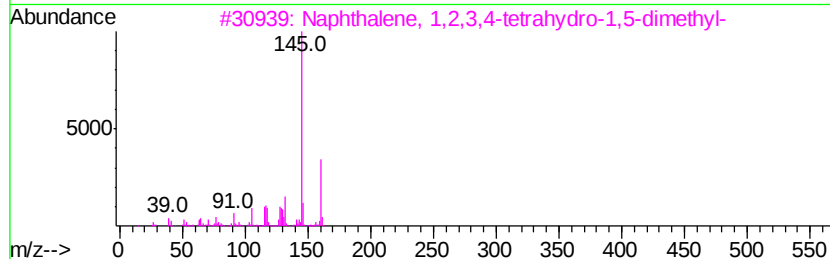
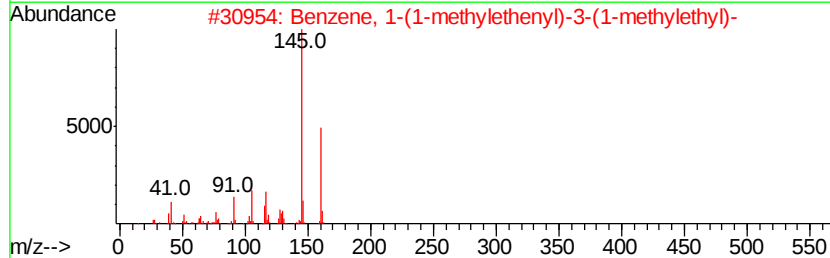
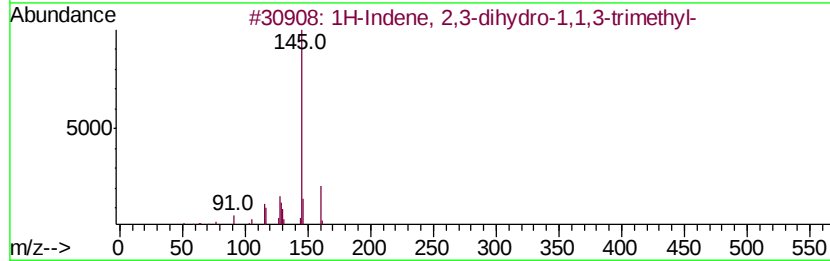
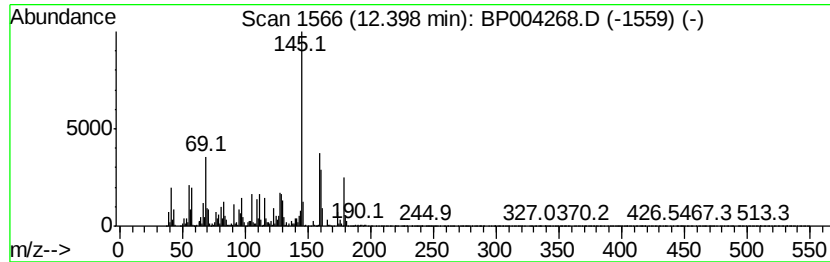
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 48

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 12.40 | 7.15 ng/ul | 2197930 | Naphthalene-d8   | 10.64 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1H-Indene, 2,3-dihydro-1,1,3-tri... | 160 | C12H16  | 002613-76-5 | 70   |
| 2    |    | Benzene, 1-(1-methylethenyl)-3-(... | 160 | C12H16  | 001129-29-9 | 68   |
| 3    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 160 | C12H16  | 021564-91-0 | 64   |
| 4    |    | Benzene, 1-(2-butenyl)-2,3-dimet... | 160 | C12H16  | 054340-85-1 | 64   |
| 5    |    | Benzene, 1,3,5-trimethyl-2-(1-me... | 160 | C12H16  | 014679-13-1 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

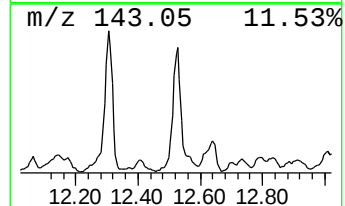
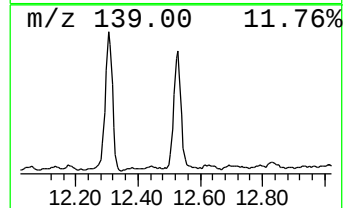
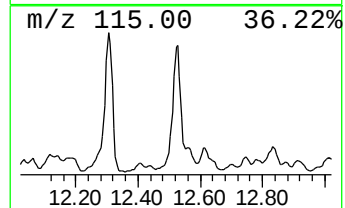
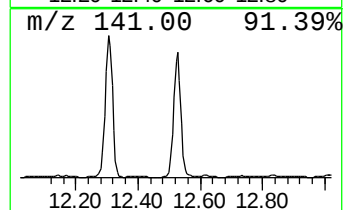
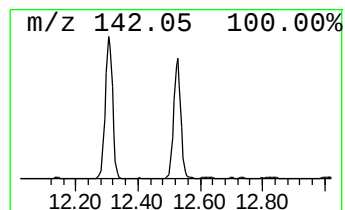
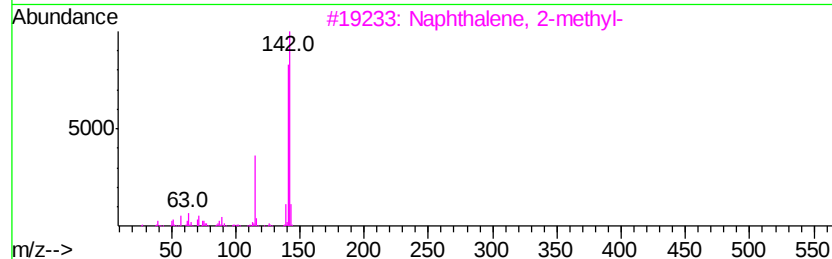
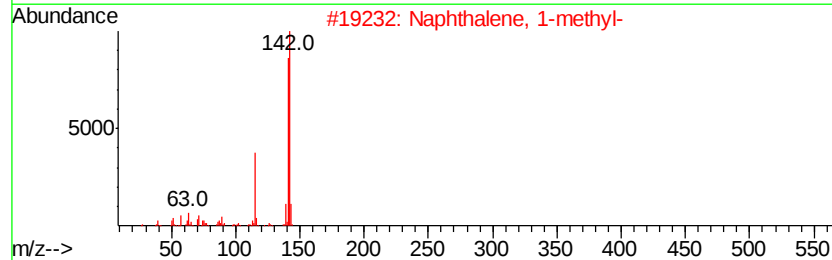
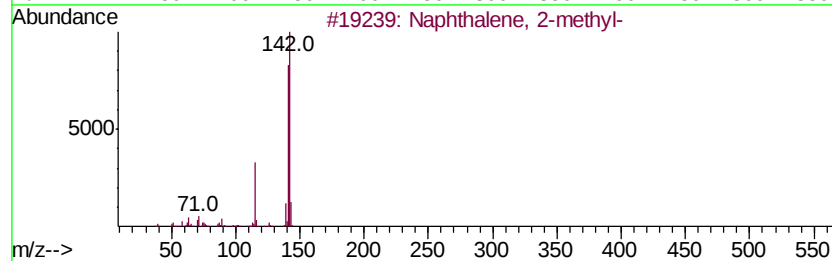
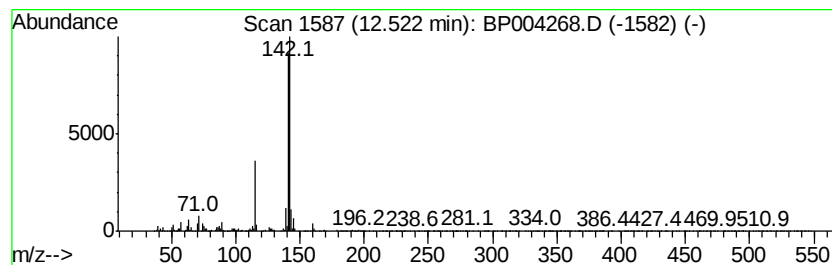
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 23 Naphthalene, 1-methyl- Concentration Rank 31

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 12.52 | 13.98 ng/ul | 4294530 | Naphthalene-d8   | 10.64 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2-methyl- | 142 | C11H10  | 000091-57-6 | 96   |
| 2    |    | Naphthalene, 1-methyl- | 142 | C11H10  | 000090-12-0 | 96   |
| 3    |    | Naphthalene, 2-methyl- | 142 | C11H10  | 000091-57-6 | 96   |
| 4    |    | Naphthalene, 1-methyl- | 142 | C11H10  | 000090-12-0 | 94   |
| 5    |    | Naphthalene, 1-methyl- | 142 | C11H10  | 000090-12-0 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

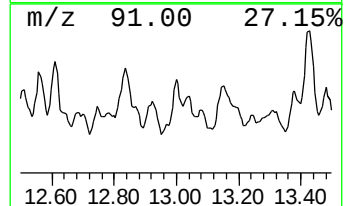
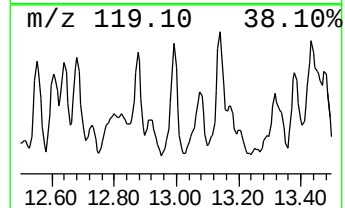
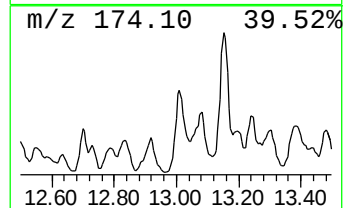
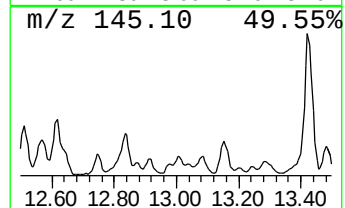
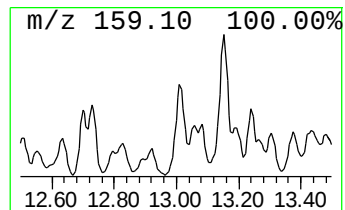
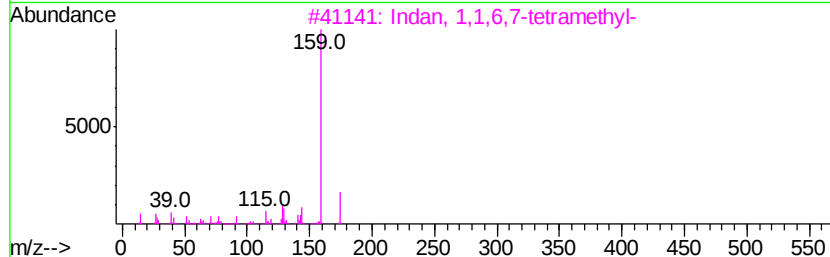
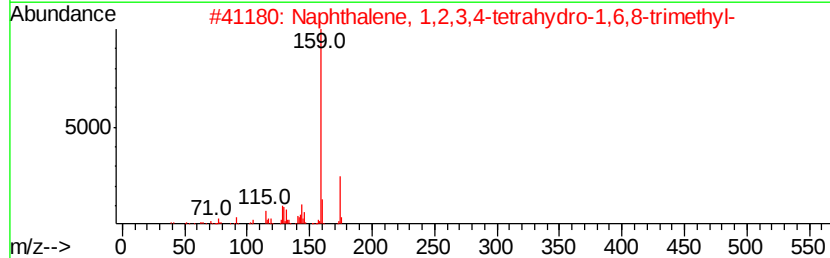
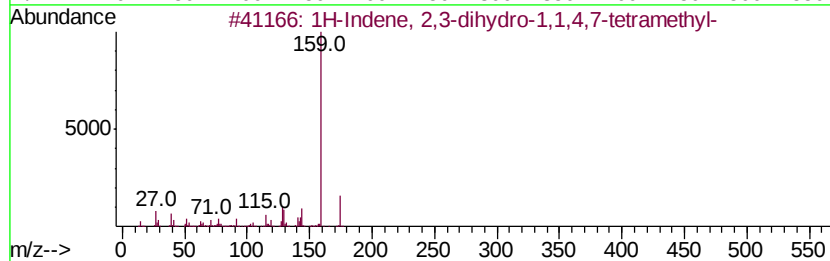
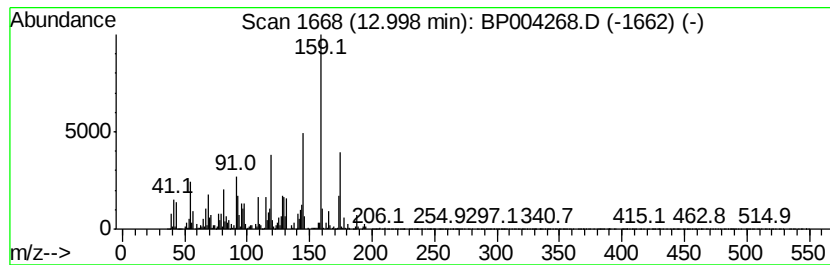
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 25 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 58

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 13.00 | 4.41 ng/ul | 1539050 | Acenaphthene-d10 | 14.49 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1H-Indene, 2,3-dihydro-1,1,4,7-t... | 174 | C13H18  | 001078-04-2 | 92   |
| 2    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 174 | C13H18  | 030316-36-0 | 92   |
| 3    |    | Indan, 1,1,6,7-tetramethyl-         | 174 | C13H18  | 016204-58-3 | 91   |
| 4    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 174 | C13H18  | 021693-51-6 | 86   |
| 5    |    | 1H-Inden-1-one, 2,3-dihydro-3,4,... | 174 | C12H14O | 035322-84-0 | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

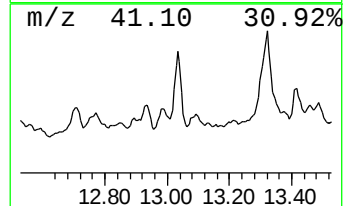
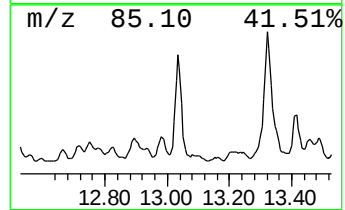
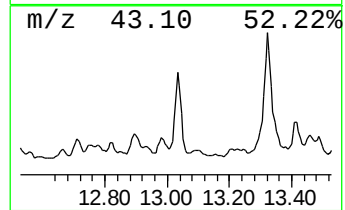
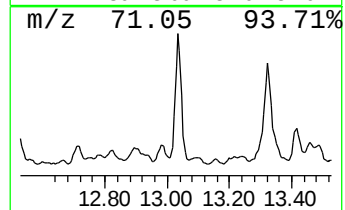
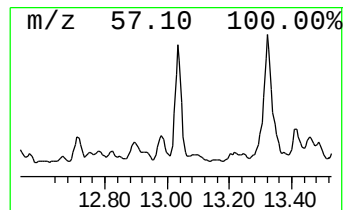
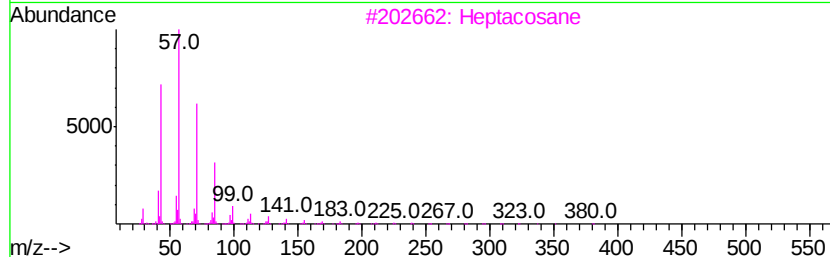
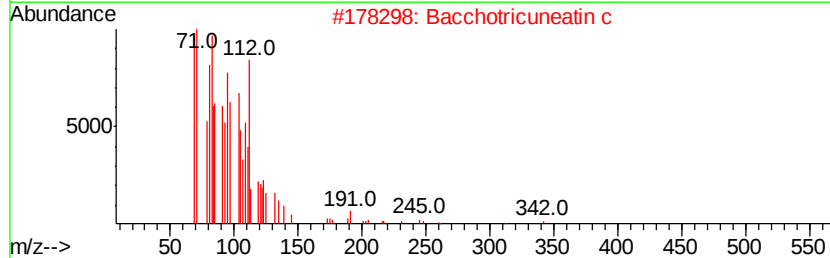
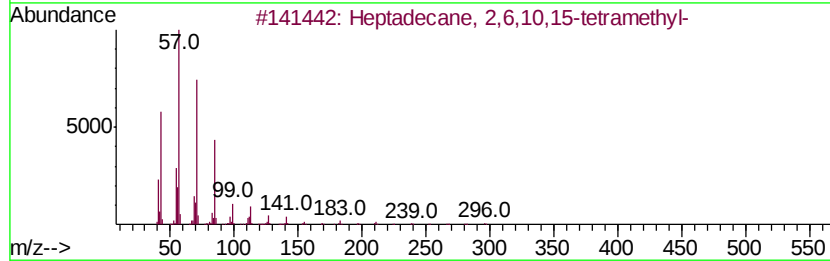
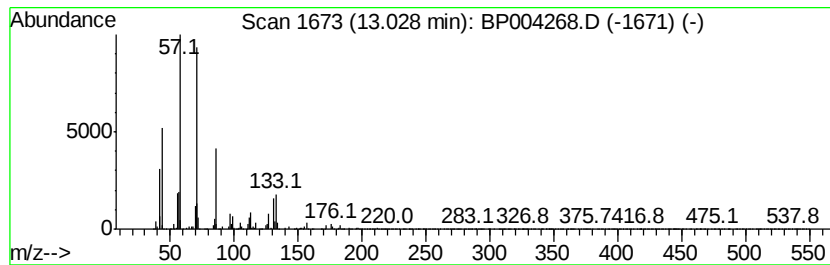
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 50

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 13.03 | 7.00 ng/ul | 2444680 | Acenaphthene-d10 | 14.49 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44   | 054833-48-6 | 81   |
| 2    |    | Bacchotricuneatin c                 | 342 | C20H22O5 | 066563-30-2 | 76   |
| 3    |    | Heptacosane                         | 380 | C27H56   | 000593-49-7 | 74   |
| 4    |    | Heptadecane, 2,6,10,14-tetramethyl- | 296 | C21H44   | 018344-37-1 | 72   |
| 5    |    | Pentadecane, 7-methyl-              | 226 | C16H34   | 006165-40-8 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

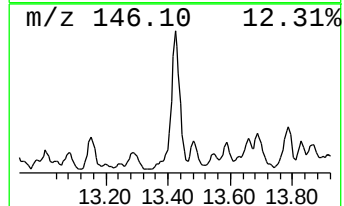
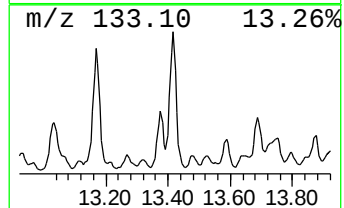
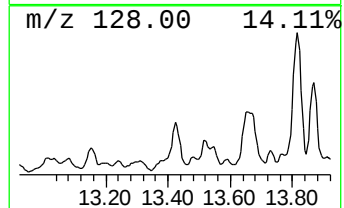
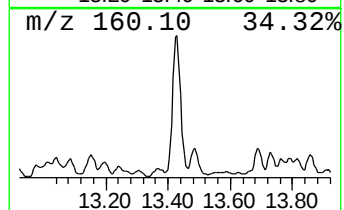
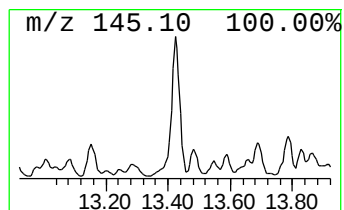
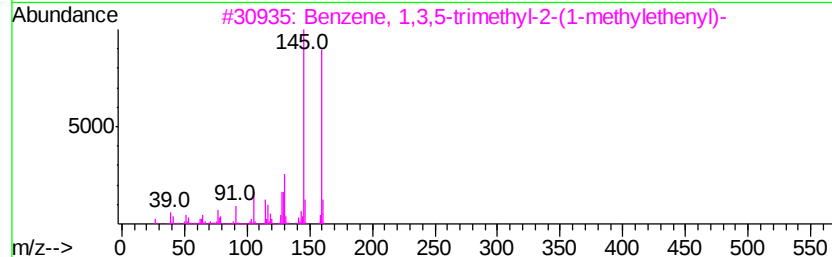
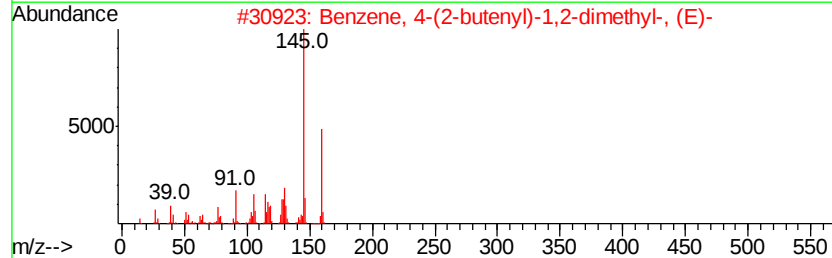
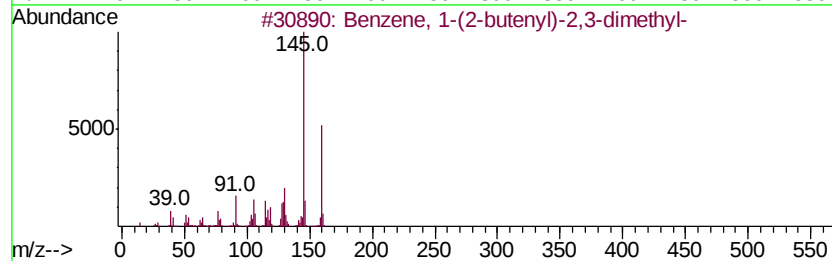
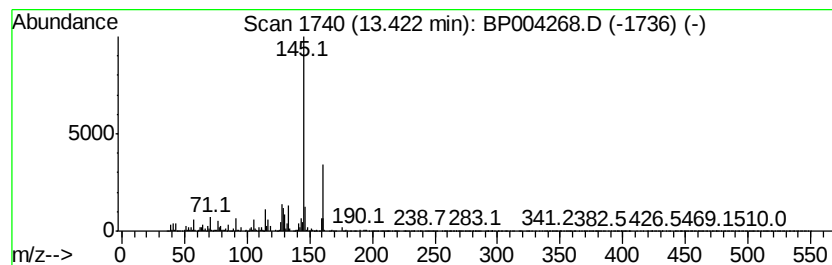
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 28 Benzene, 1-(2-butenyl)-2,3-... Concentration Rank 41

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 13.42 | 9.74 ng/ul | 3404090 | Acenaphthene-d10 | 14.49 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-(2-butenyl)-2,3-dimet... | 160 | C12H16  | 054340-85-1 | 90   |
| 2    |    |   | Benzene, 4-(2-butenyl)-1,2-dimet... | 160 | C12H16  | 054340-86-2 | 90   |
| 3    |    |   | Benzene, 1,3,5-trimethyl-2-(1-me... | 160 | C12H16  | 014679-13-1 | 90   |
| 4    |    |   | 1H-Indene, 2,3-dihydro-4,5,7-tri... | 160 | C12H16  | 006682-06-0 | 90   |
| 5    |    |   | Naphthalene, 1,2,3,4-tetrahydro-... | 160 | C12H16  | 021693-54-9 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

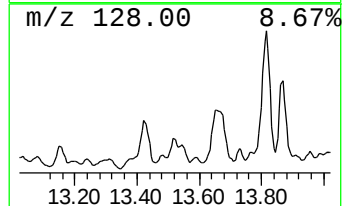
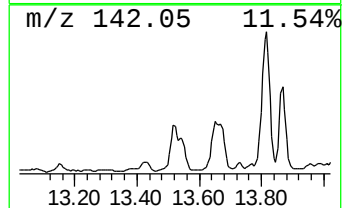
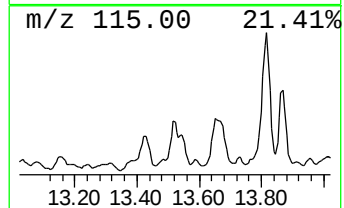
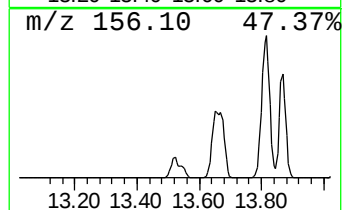
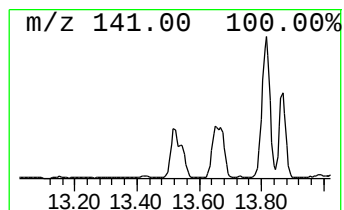
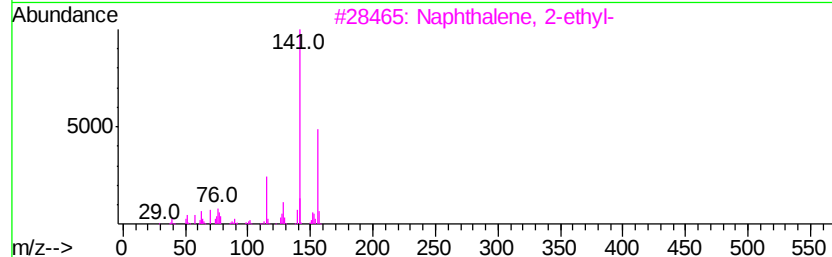
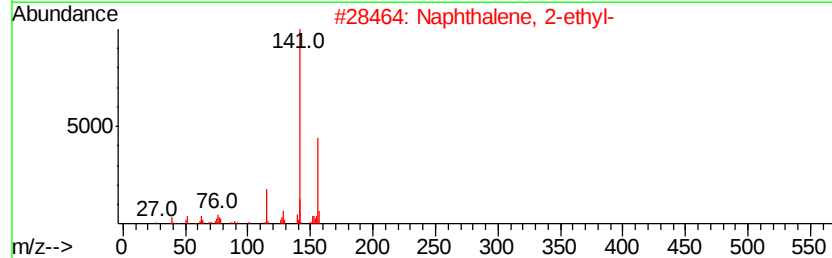
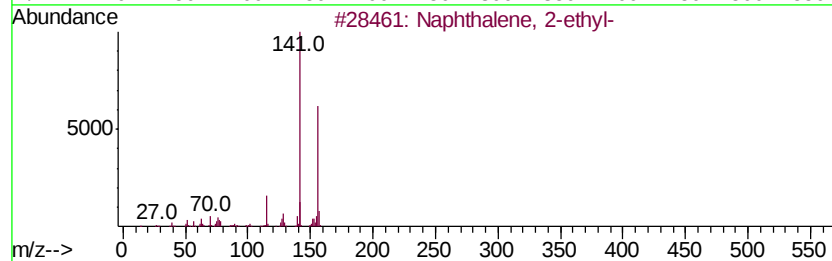
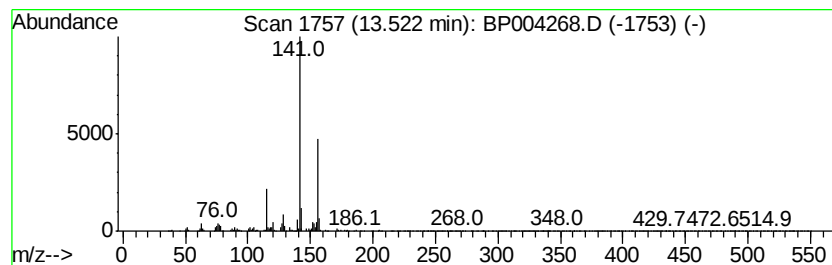
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 29 Naphthalene, 2-ethyl- Concentration Rank 62

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 13.52 | 3.65 ng/ul | 1275770 | Acenaphthene-d10 | 14.49 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2-ethyl- | 156 | C12H12  | 000939-27-5 | 95   |
| 2    |    | Naphthalene, 2-ethyl- | 156 | C12H12  | 000939-27-5 | 95   |
| 3    |    | Naphthalene, 2-ethyl- | 156 | C12H12  | 000939-27-5 | 94   |
| 4    |    | Naphthalene, 1-ethyl- | 156 | C12H12  | 001127-76-0 | 94   |
| 5    |    | Naphthalene, 2-ethyl- | 156 | C12H12  | 000939-27-5 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

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Quant Title : SVOA CALIBRATION

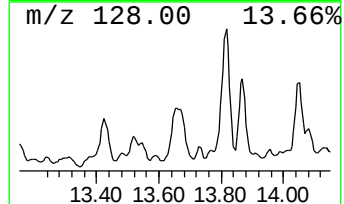
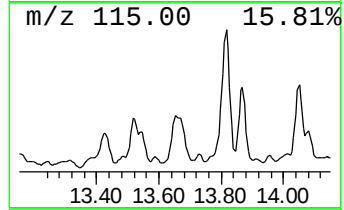
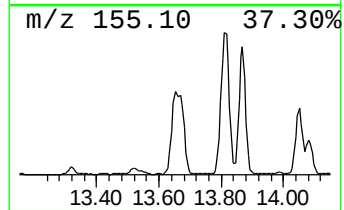
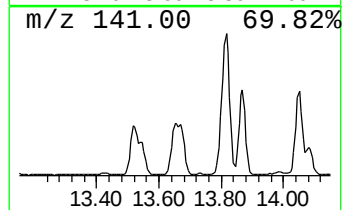
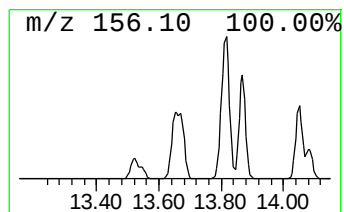
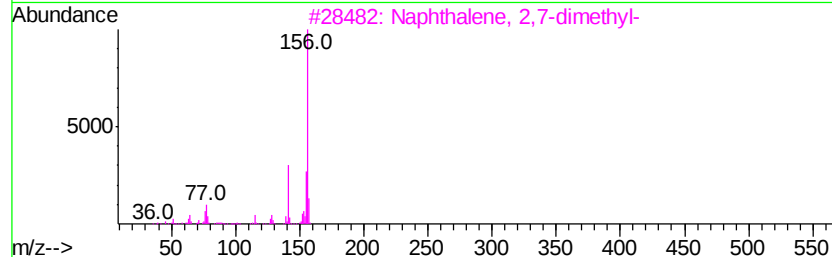
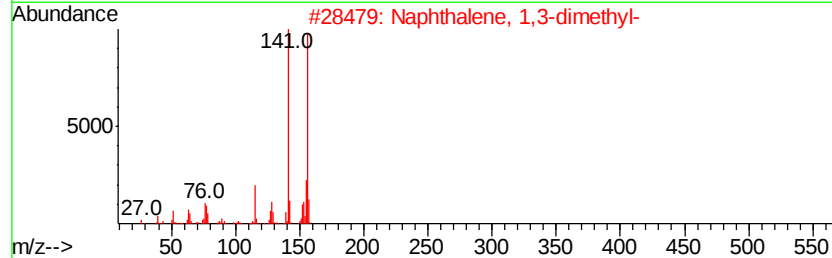
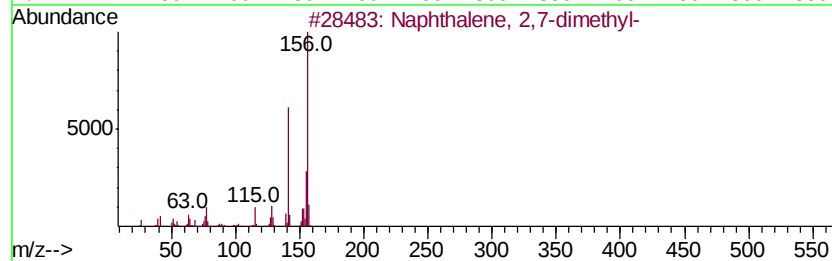
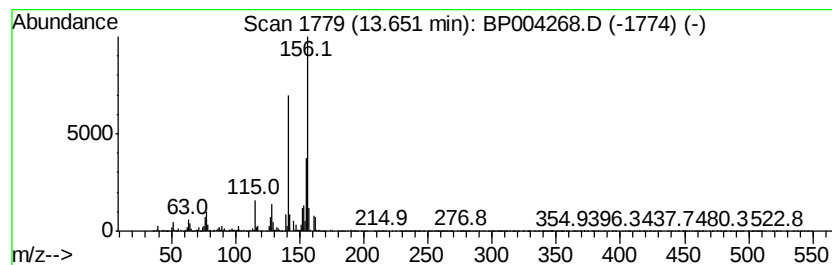
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 30 Naphthalene, 2,7-dimethyl- Concentration Rank 27

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 13.65 | 15.84 ng/ul | 5533460 | Acenaphthene-d10 | 14.49 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 96   |
| 2    |    | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 96   |
| 3    |    | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 96   |
| 4    |    | Naphthalene, 1,2-dimethyl- | 156 | C12H12  | 000573-98-8 | 96   |
| 5    |    | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

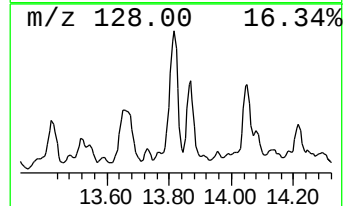
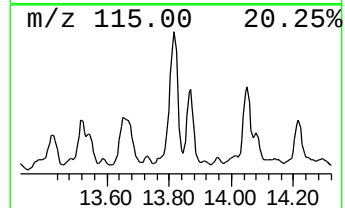
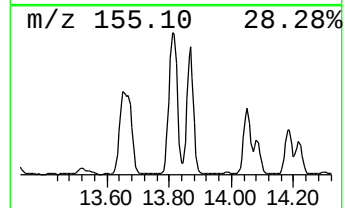
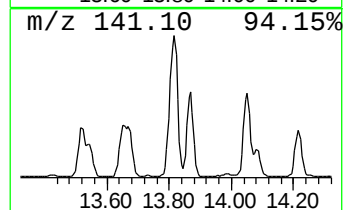
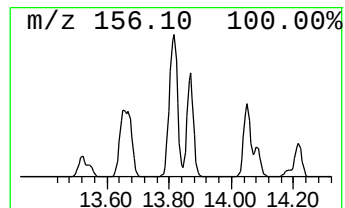
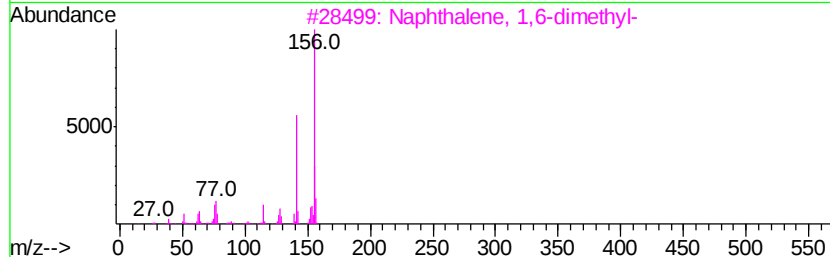
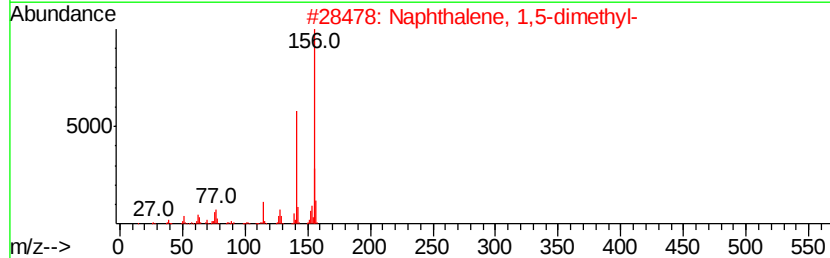
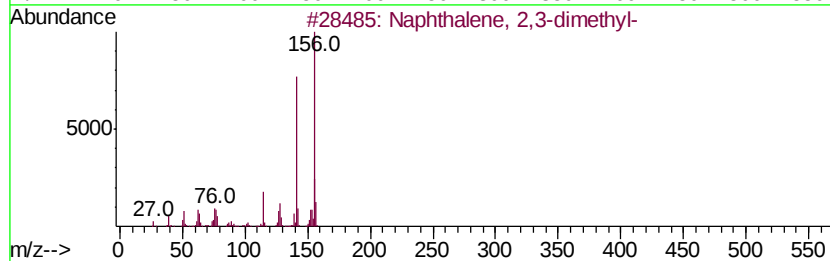
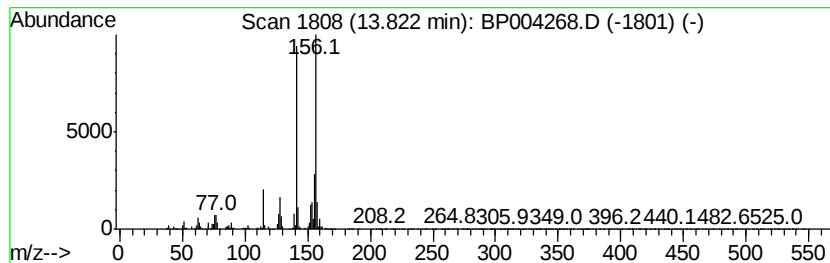
Instrument :  
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ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 31 Naphthalene, 2,3-dimethyl- Concentration Rank 7

| R.T.    | EstConc     | Area                       | Relative to ISTD | R.T.           |
|---------|-------------|----------------------------|------------------|----------------|
| 13.82   | 33.90 ng/ul | 11842900                   | Acenaphthene-d10 | 14.49          |
| Hit# of | 5           | Tentative ID               | MW MolForm       | CAS# Qual      |
| 1       |             | Naphthalene, 2,3-dimethyl- | 156 C12H12       | 000581-40-8 97 |
| 2       |             | Naphthalene, 1,5-dimethyl- | 156 C12H12       | 000571-61-9 97 |
| 3       |             | Naphthalene, 1,6-dimethyl- | 156 C12H12       | 000575-43-9 97 |
| 4       |             | Naphthalene, 1,4-dimethyl- | 156 C12H12       | 000571-58-4 97 |
| 5       |             | Naphthalene, 1,7-dimethyl- | 156 C12H12       | 000575-37-1 97 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

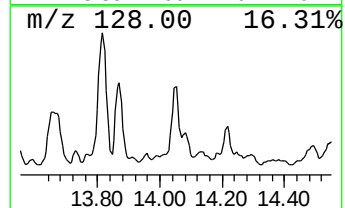
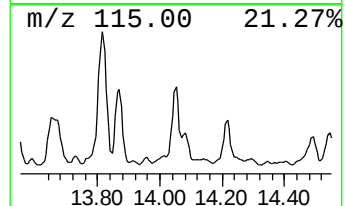
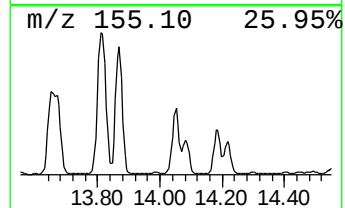
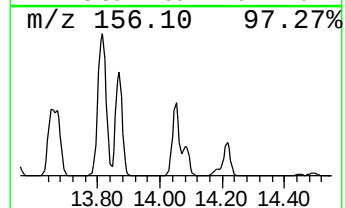
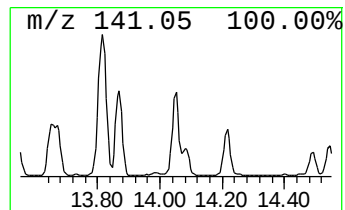
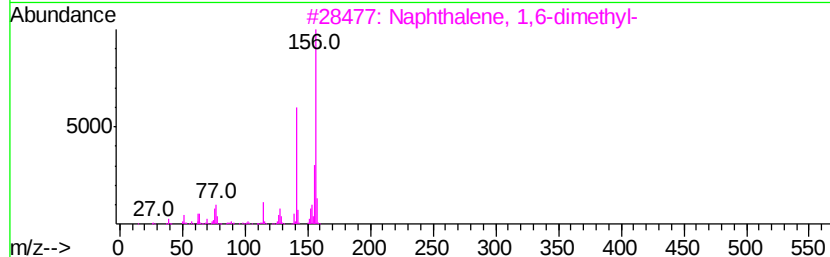
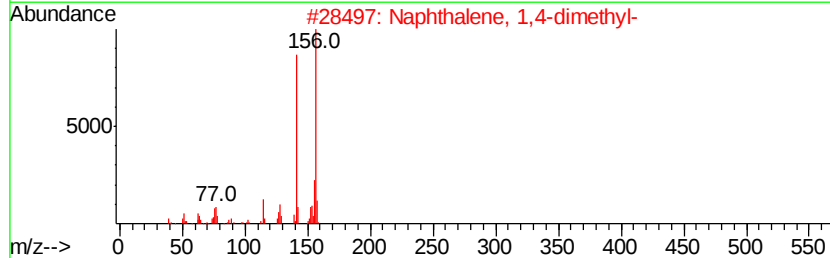
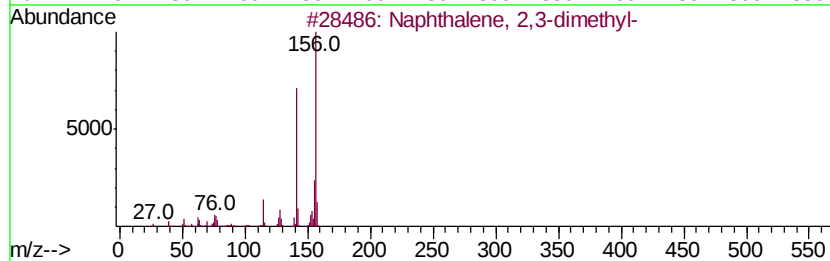
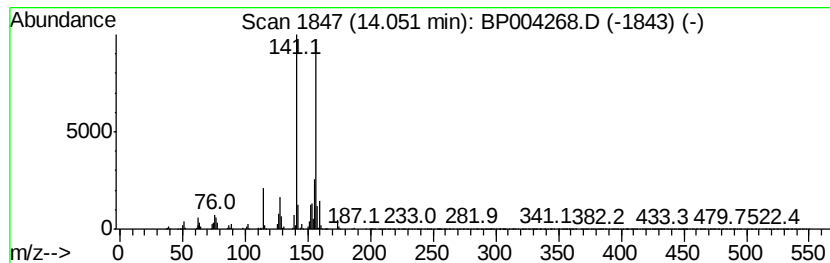
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 32 Naphthalene, 1,4-dimethyl- Concentration Rank 35

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 14.05 | 11.40 ng/ul | 3982040 | Acenaphthene-d10 | 14.49 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |
| 2    |    | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 96   |
| 3    |    | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |
| 4    |    | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 96   |
| 5    |    | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 95   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

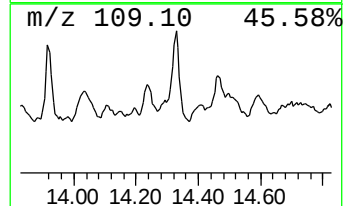
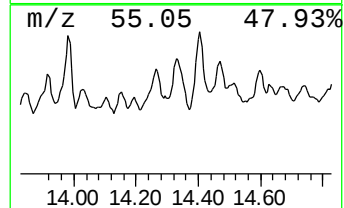
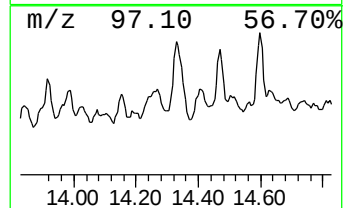
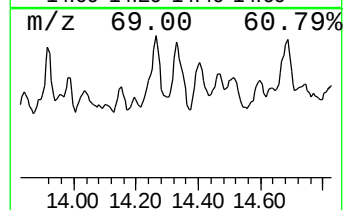
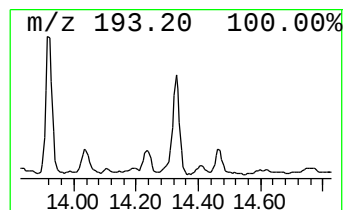
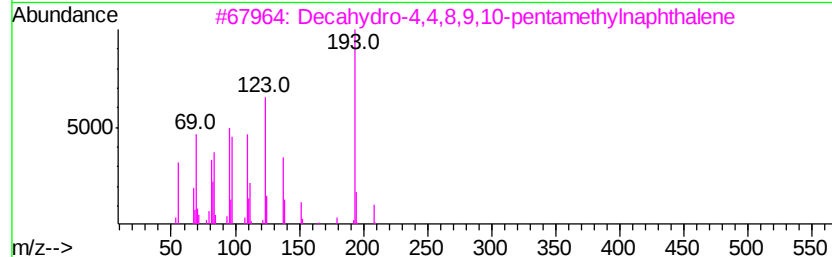
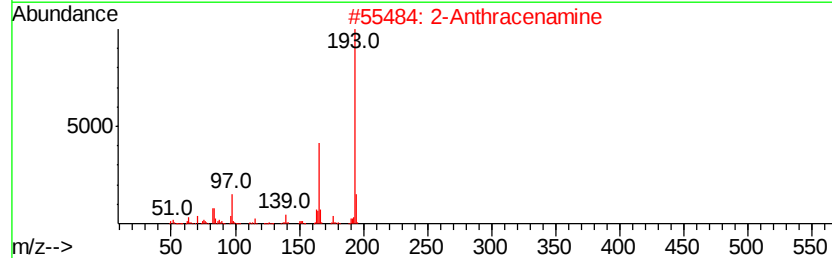
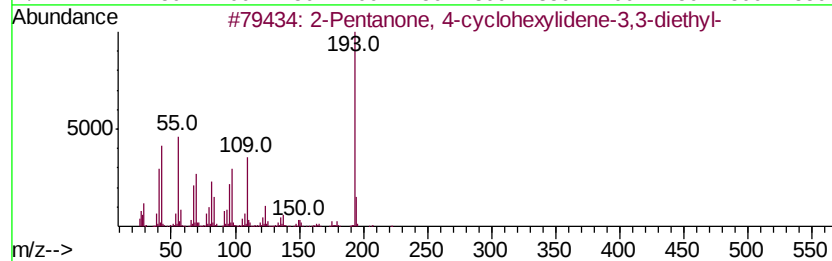
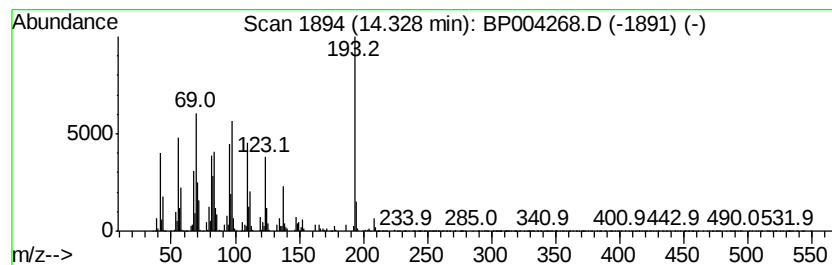
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 33 unknown-01 Concentration Rank 46

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 14.33 | 7.41 ng/ul | 2588820 | Acenaphthene-d10 | 14.49 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-cyclohexylidene-3... | 222 | C15H26O | 313253-65-5 | 38   |
| 2    |    |   | 2-Anthracenamine                    | 193 | C14H11N | 000613-13-8 | 38   |
| 3    |    |   | Decahydro-4,4,8,9,10-pentamethyl... | 208 | C15H28  | 080655-44-3 | 38   |
| 4    |    |   | Cyclohexane, 1,1',1''-(1-ethanyl... | 276 | C20H36  | 055682-86-5 | 35   |
| 5    |    |   | 1H-Indene, octahydro-2,2,4,4,7,7... | 208 | C15H28  | 054832-83-6 | 32   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

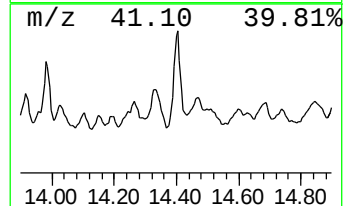
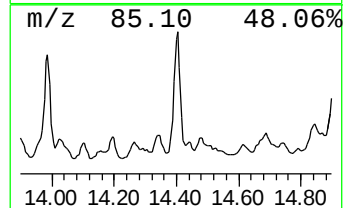
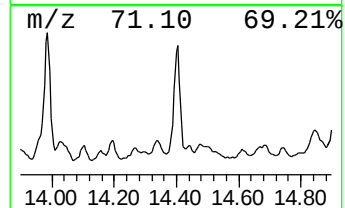
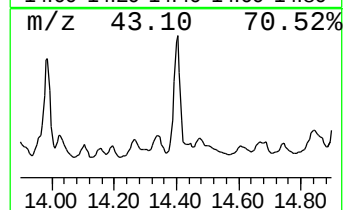
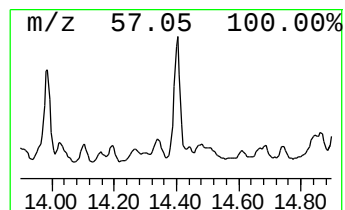
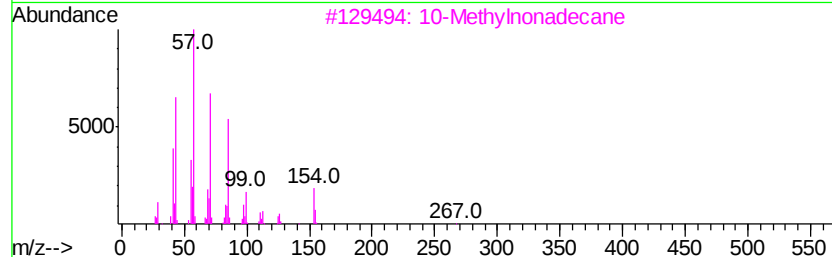
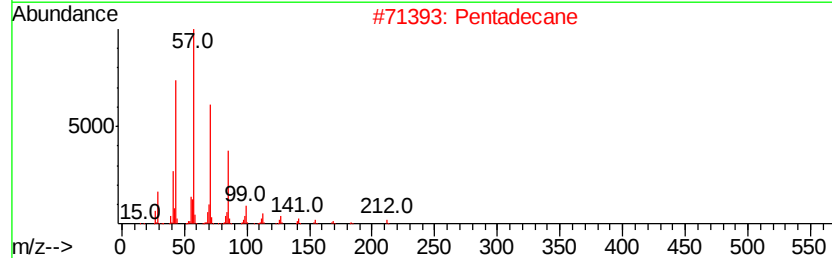
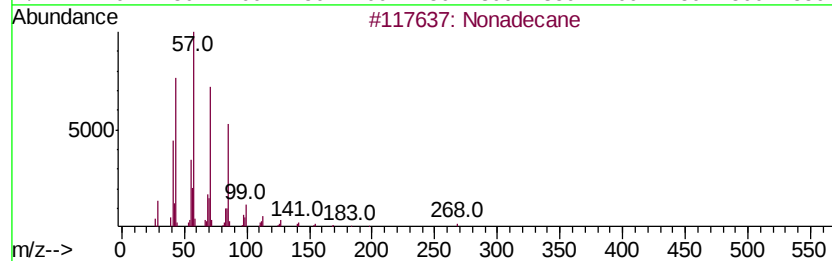
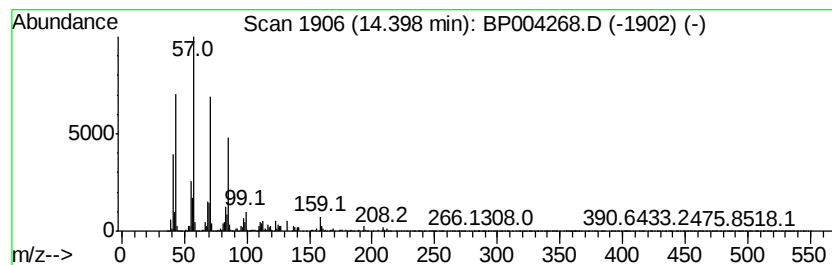
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 34 (DEL) Alkane: Straight-Chai... Concentration Rank 38

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 14.40 | 10.30 ng/ul | 3597920 | Acenaphthene-d10 | 14.49 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------|-----|---------|-------------|------|
| 1    | 5  | Nonadecane          | 268 | C19H40  | 000629-92-5 | 81   |
| 2    |    | Pentadecane         | 212 | C15H32  | 000629-62-9 | 81   |
| 3    |    | 10-Methylnonadecane | 282 | C20H42  | 056862-62-5 | 81   |
| 4    |    | Tetradecane         | 198 | C14H30  | 000629-59-4 | 74   |
| 5    |    | Octacosane          | 394 | C28H58  | 000630-02-4 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

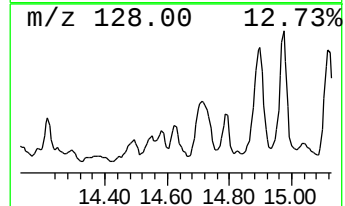
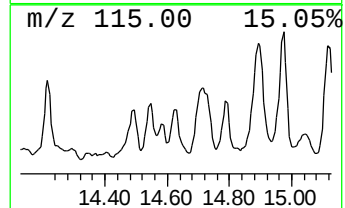
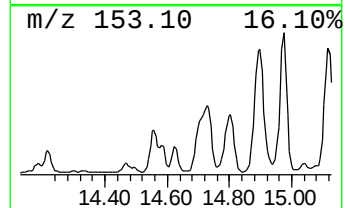
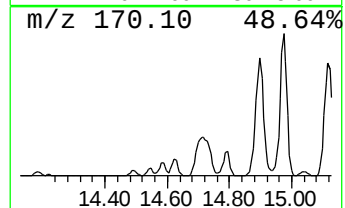
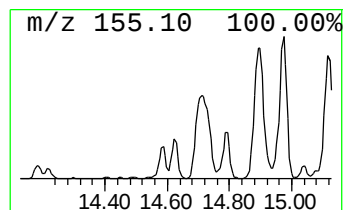
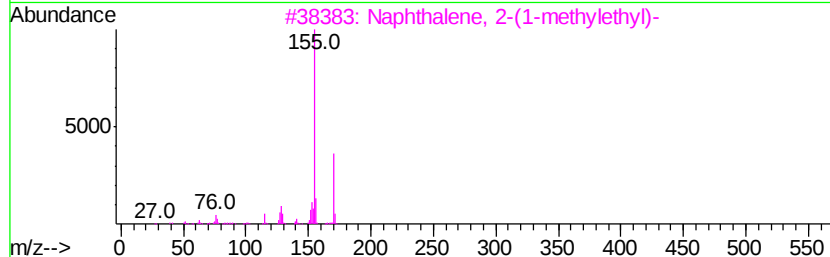
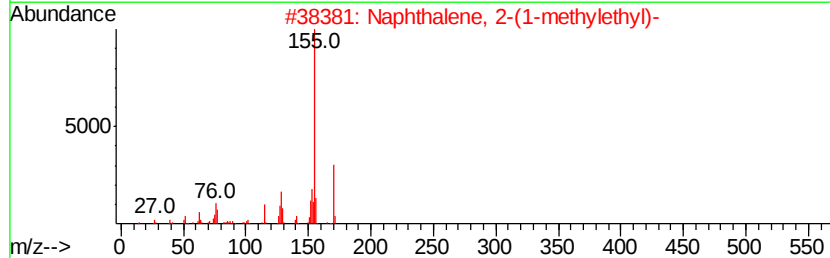
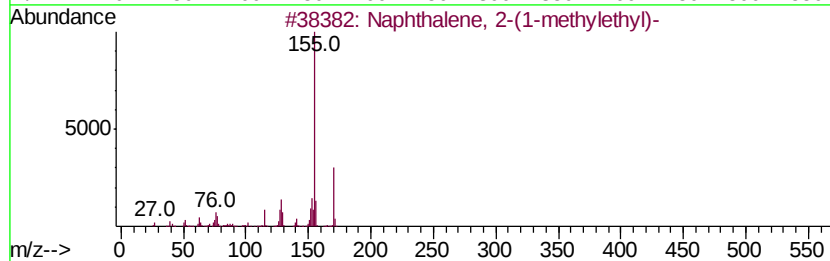
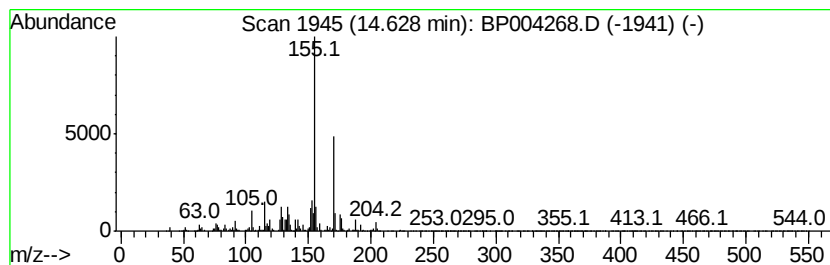
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 35 Naphthalene, 2-(1-methylethyl) Concentration Rank 49

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 14.63 | 7.06 ng/ul | 2467310 | Acenaphthene-d10 | 14.49 |

| Hit# | of | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2-(1-methylethyl)- | 170 | C13H14  | 002027-17-0 | 90   |
| 2    |    | Naphthalene, 2-(1-methylethyl)- | 170 | C13H14  | 002027-17-0 | 90   |
| 3    |    | Naphthalene, 2-(1-methylethyl)- | 170 | C13H14  | 002027-17-0 | 87   |
| 4    |    | Naphthalene, 1,4,6-trimethyl-   | 170 | C13H14  | 002131-42-2 | 86   |
| 5    |    | Naphthalene, 1-(1-methylethyl)- | 170 | C13H14  | 006158-45-8 | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

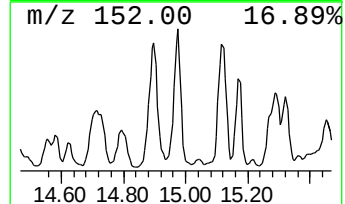
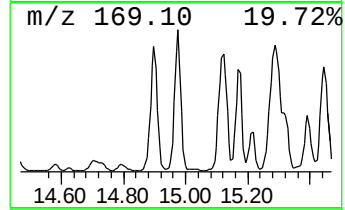
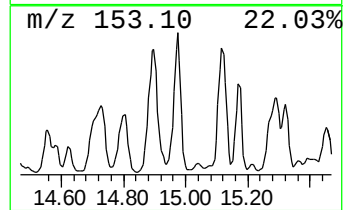
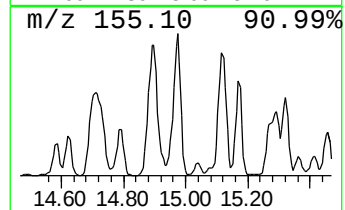
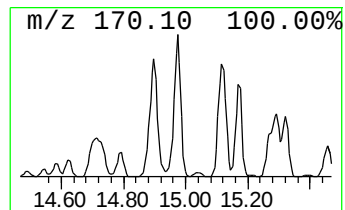
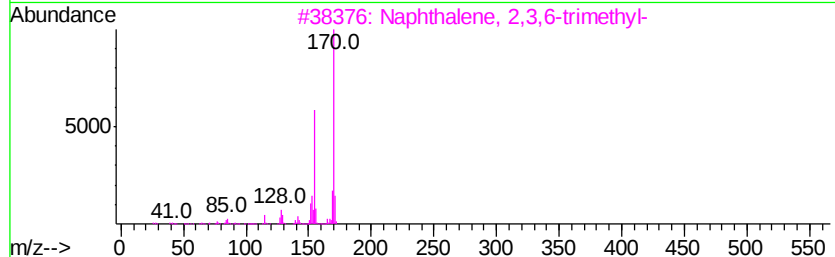
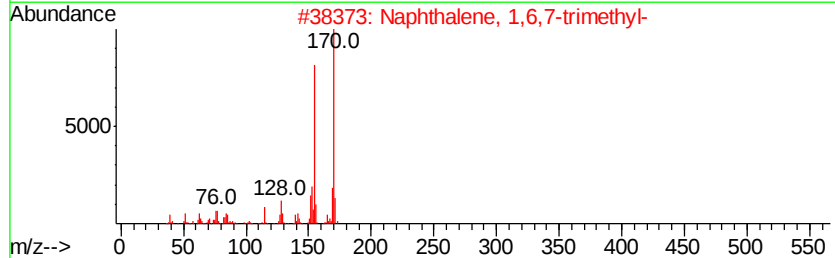
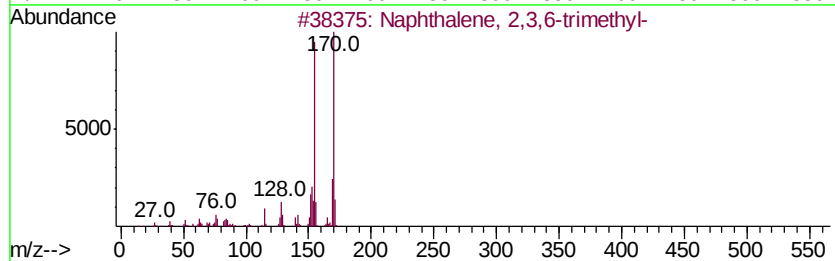
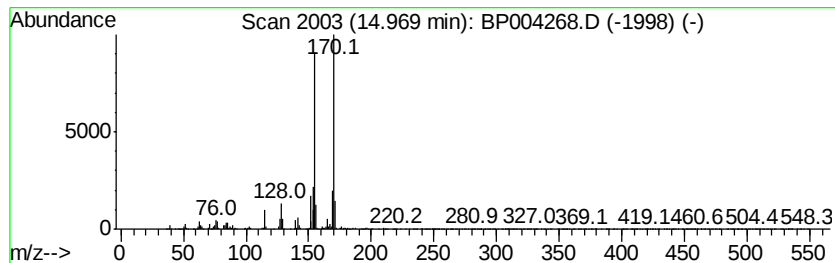
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 38 Naphthalene, 2,3,6-trimethyl- Concentration Rank 9

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 14.97 | 32.61 ng/ul | 11391200 | Acenaphthene-d10 | 14.49 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 98   |
| 2    |    | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 97   |
| 3    |    | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 96   |
| 4    |    | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 96   |
| 5    |    | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 95   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

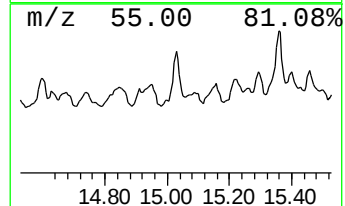
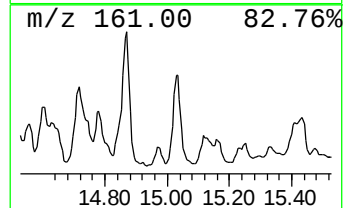
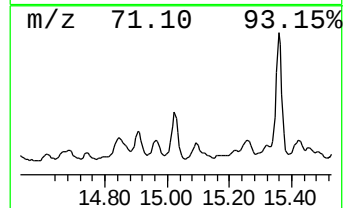
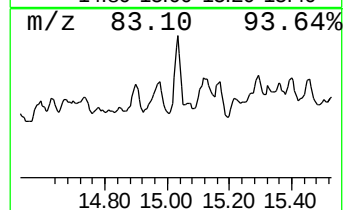
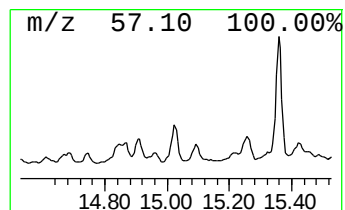
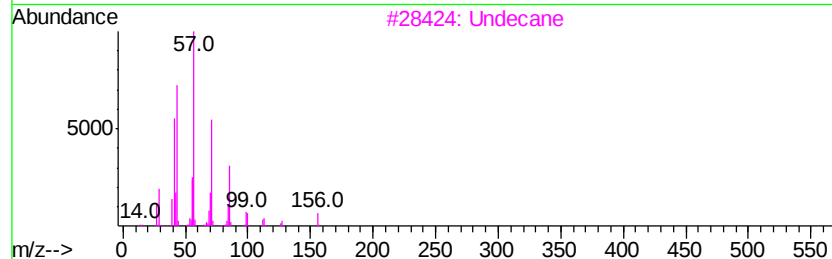
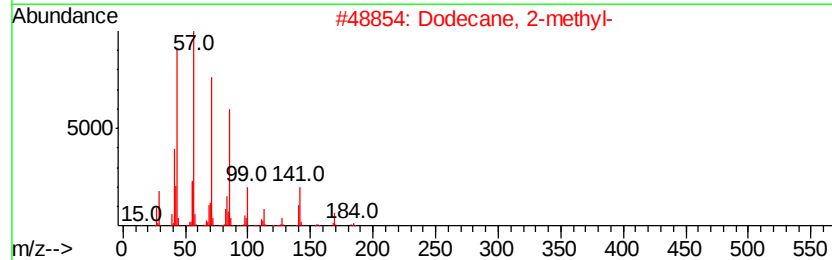
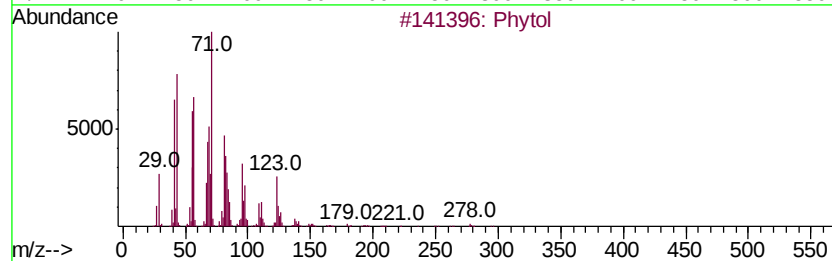
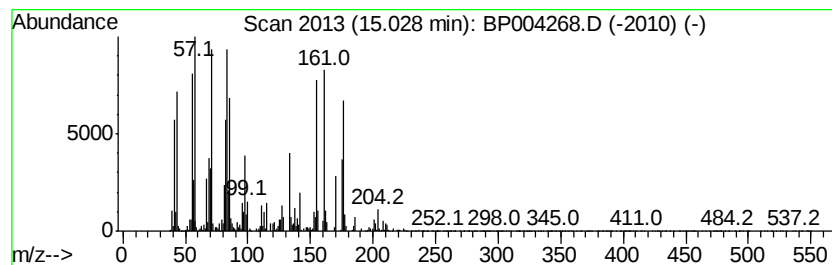
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 39 unknown-02 Concentration Rank 45

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 15.03 | 7.82 ng/ul | 2730810 | Acenaphthene-d10 | 14.49 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Phytol                              | 296 | C20H40O | 000150-86-7 | 25   |
| 2    |    |   | Dodecane, 2-methyl-                 | 184 | C13H28  | 001560-97-0 | 25   |
| 3    |    |   | Undecane                            | 156 | C11H24  | 001120-21-4 | 25   |
| 4    |    |   | Undecane, 4-cyclohexyl-             | 238 | C17H34  | 013151-79-6 | 18   |
| 5    |    |   | Benzofuran, 2,3-dihydro-2,2,4,6-... | 176 | C12H16O | 003698-49-5 | 18   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

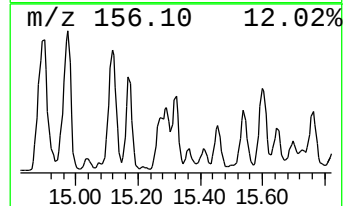
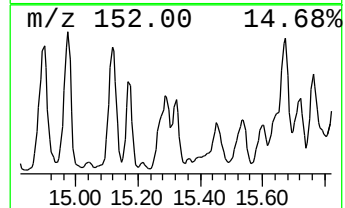
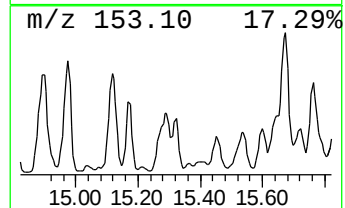
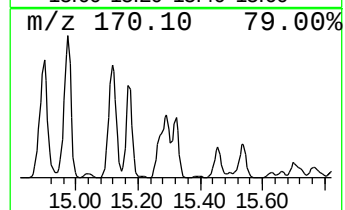
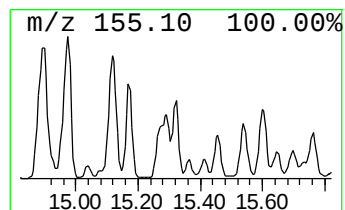
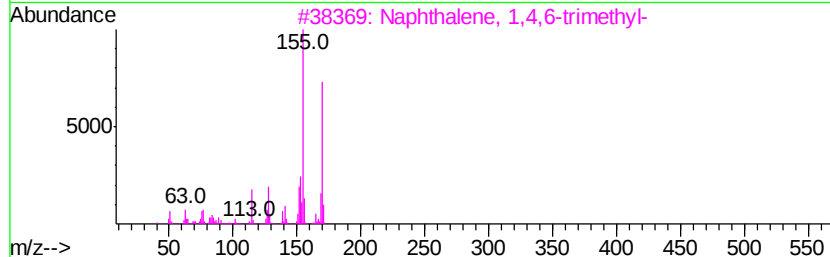
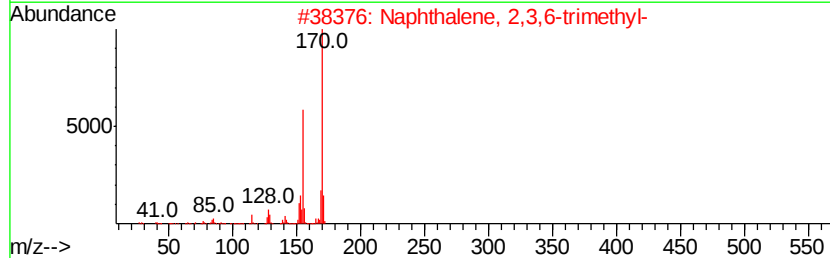
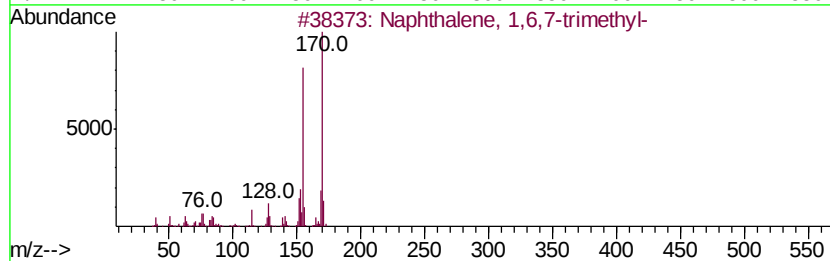
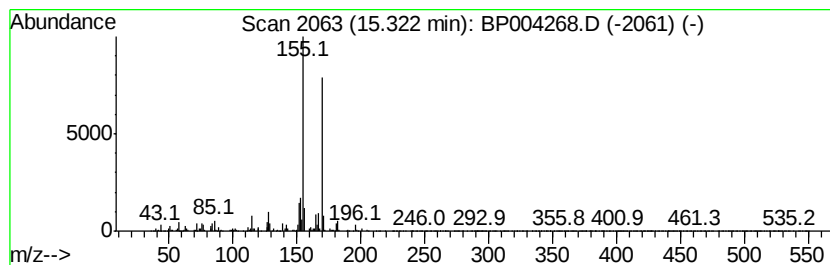
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 43 Naphthalene, 1,6,7-trimethyl- Concentration Rank 37

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.32 | 10.40 ng/ul | 3631480 | Acenaphthene-d10 | 14.49 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 95   |
| 2    |    | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 90   |
| 3    |    | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 87   |
| 4    |    | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 87   |
| 5    |    | 4,6,8-Trimethylazulene        | 170 | C13H14  | 000941-81-1 | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

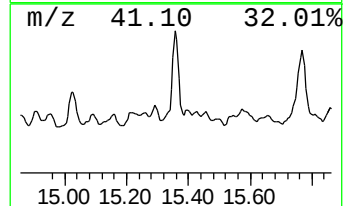
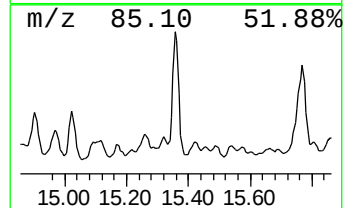
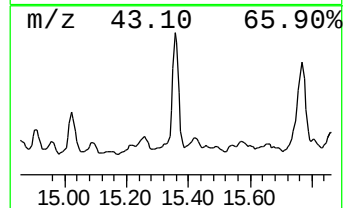
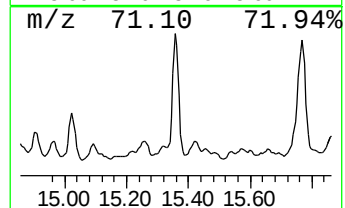
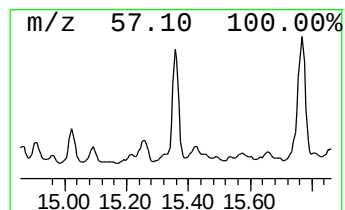
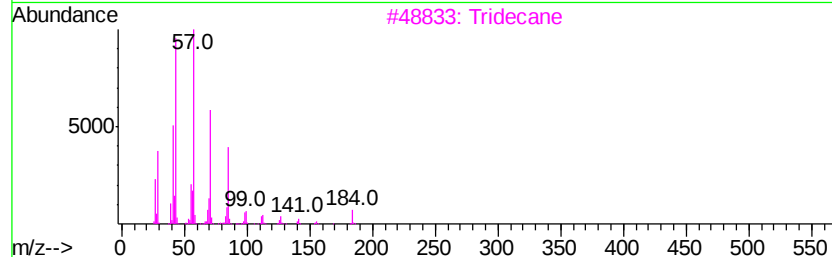
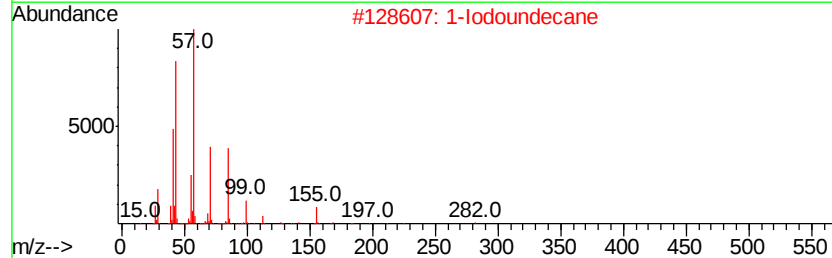
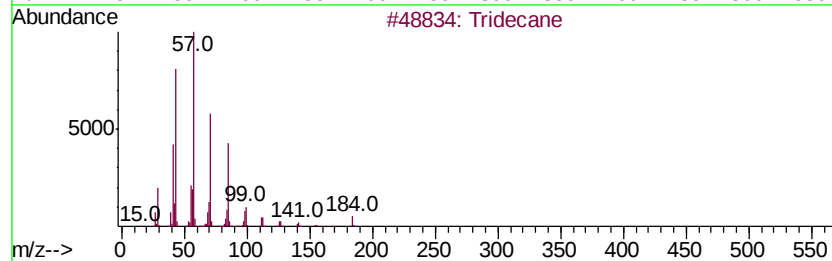
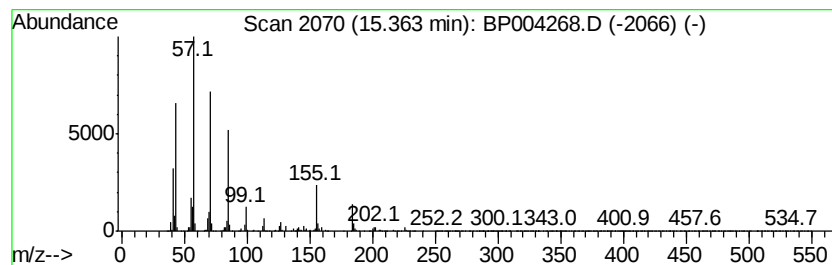
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 15.36 | 7.91 ng/ul | 2764280 | Acenaphthene-d10 | 14.49 |

| Hit# | of | 5 | Tentative ID   | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------|-----|---------|-------------|------|
| 1    |    |   | Tridecane      | 184 | C13H28  | 000629-50-5 | 86   |
| 2    |    |   | 1-Iodoundecane | 282 | C11H23I | 004282-44-4 | 81   |
| 3    |    |   | Tridecane      | 184 | C13H28  | 000629-50-5 | 72   |
| 4    |    |   | Tridecane      | 184 | C13H28  | 000629-50-5 | 72   |
| 5    |    |   | Hexadecane     | 226 | C16H34  | 000544-76-3 | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

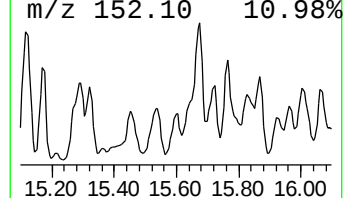
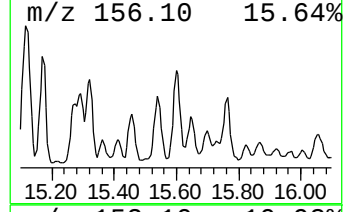
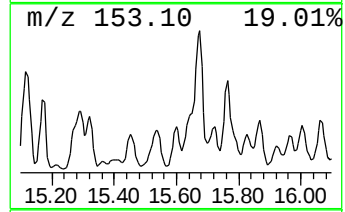
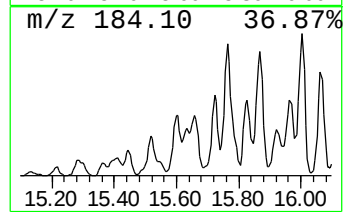
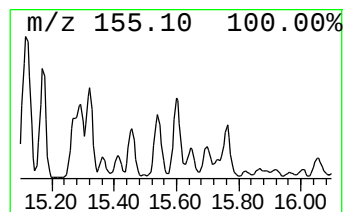
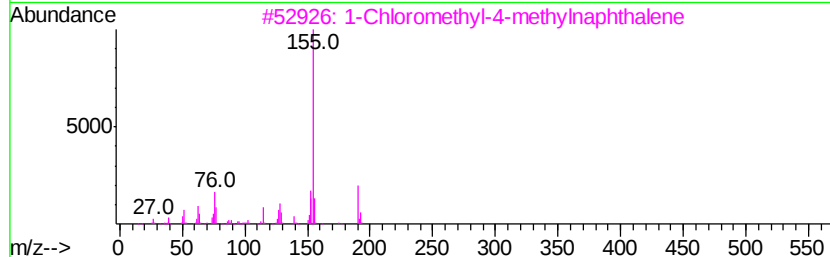
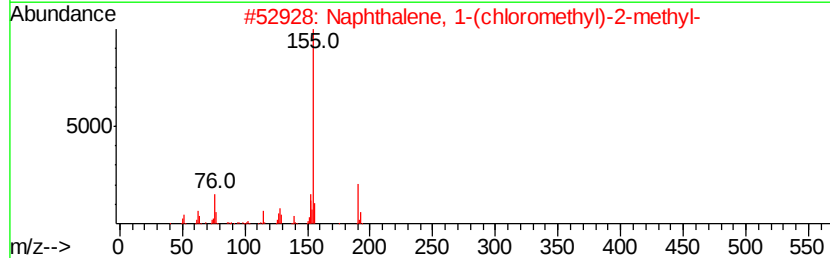
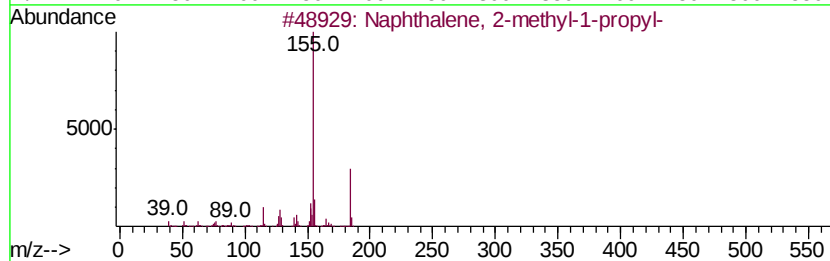
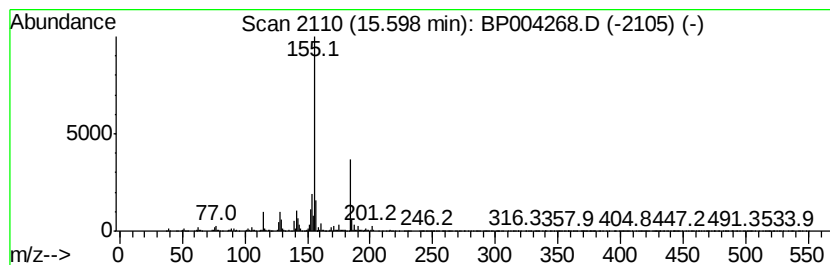
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 45 Naphthalene, 2-methyl-1-pro... Concentration Rank 33

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.60 | 12.73 ng/ul | 4445170 | Acenaphthene-d10 | 14.49 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Naphthalene, 2-methyl-1-propyl-     | 184 | C14H16   | 054774-89-9 | 94   |
| 2    |    | Naphthalene, 1-(chloromethyl)-2-... | 190 | C12H11Cl | 006626-23-9 | 70   |
| 3    |    | 1-Chloromethyl-4-methylnaphthalene  | 190 | C12H11Cl | 005261-50-7 | 45   |
| 4    |    | Naphthalene, 1-(chloromethyl)-2-... | 190 | C12H11Cl | 006626-23-9 | 45   |
| 5    |    | Benzaldehyde, 4-chloro-, oxime, ... | 155 | C7H6ClNO | 003717-24-6 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

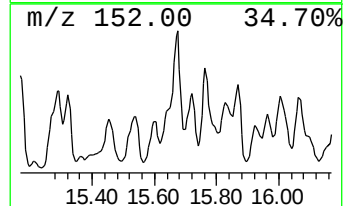
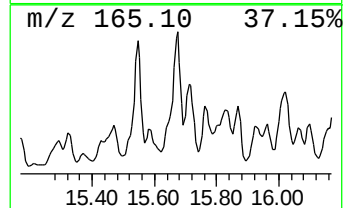
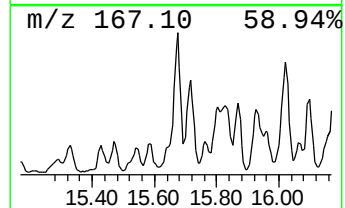
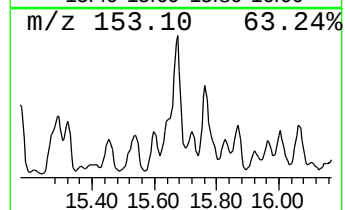
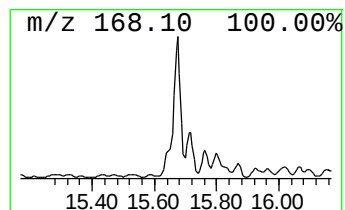
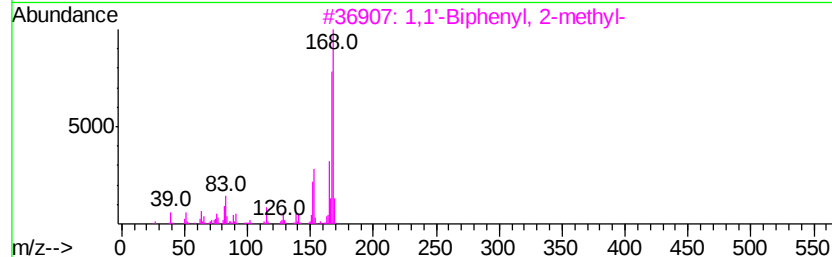
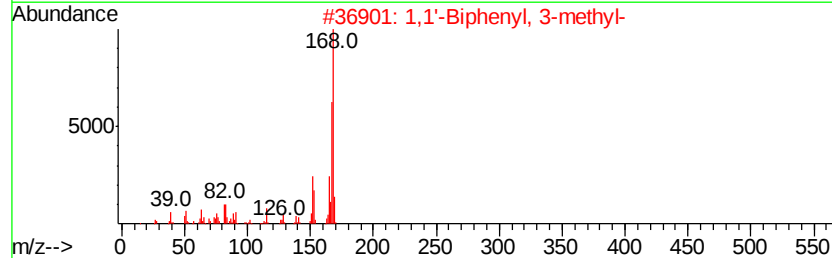
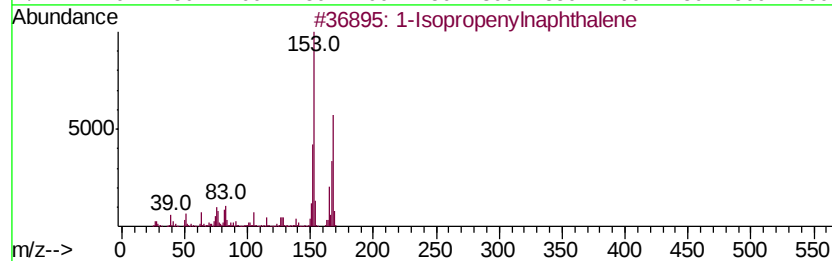
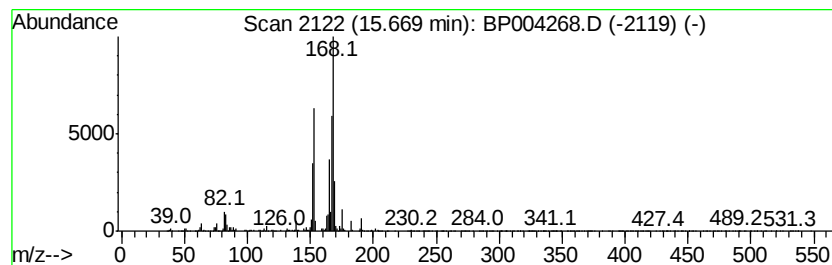
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 46 1-Isopropenylnaphthalene Concentration Rank 34

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.67 | 11.85 ng/ul | 4141200 | Acenaphthene-d10 | 14.49 |

| Hit# | of | Tentative ID             | MW  | MolForm   | CAS#        | Qual |
|------|----|--------------------------|-----|-----------|-------------|------|
| 1    | 5  | 1-Isopropenylnaphthalene | 168 | C13H12    | 001855-47-6 | 74   |
| 2    |    | 1,1'-Biphenyl, 3-methyl- | 168 | C13H12    | 000643-93-6 | 52   |
| 3    |    | 1,1'-Biphenyl, 2-methyl- | 168 | C13H12    | 000643-58-3 | 50   |
| 4    |    | N-Nitrosodiphenylamine   | 198 | C12H10N2O | 000086-30-6 | 49   |
| 5    |    | 1,1'-Biphenyl, 4-methyl- | 168 | C13H12    | 000644-08-6 | 49   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

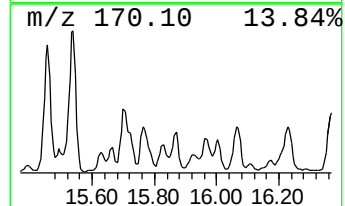
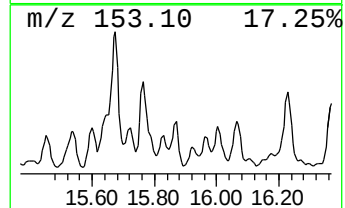
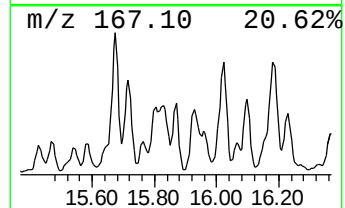
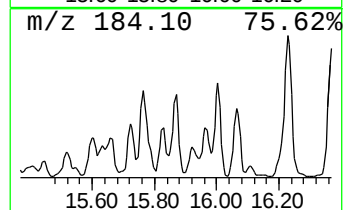
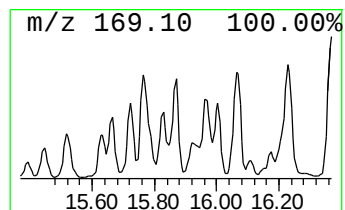
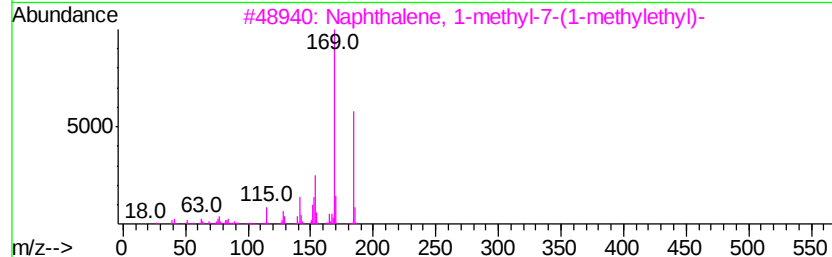
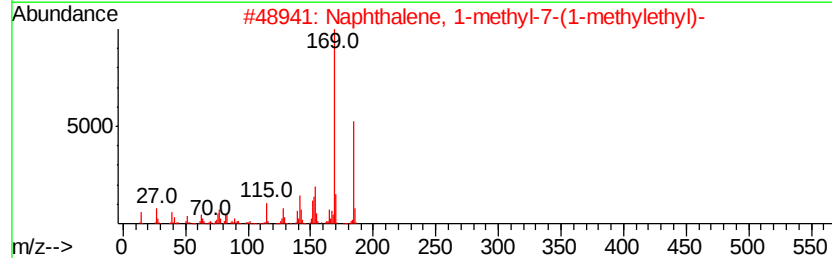
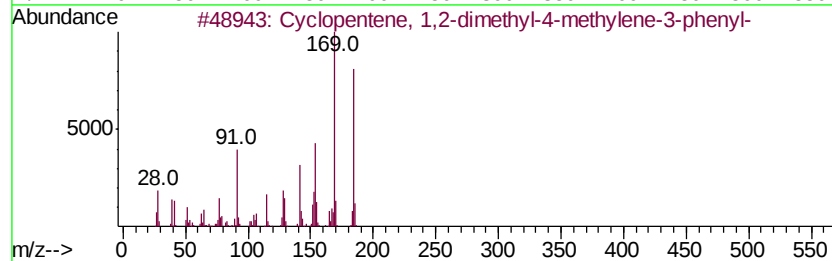
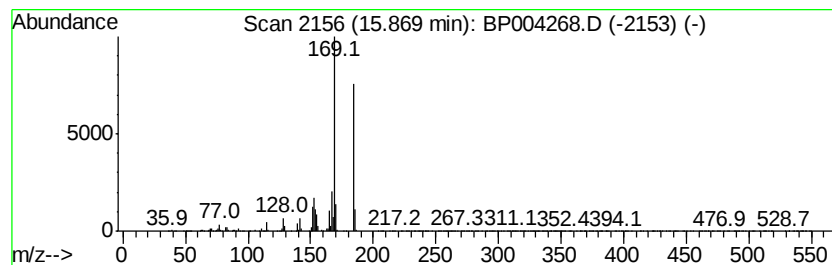
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 49 Cyclopentene, 1,2-dimethyl-... Concentration Rank 32

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.87 | 13.58 ng/ul | 4745050 | Acenaphthene-d10 | 14.49 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Cyclopentene, 1,2-dimethyl-4-met... | 184 | C14H16  | 1000152-22-4 | 90   |
| 2    |    | Naphthalene, 1-methyl-7-(1-methy... | 184 | C14H16  | 000490-65-3  | 90   |
| 3    |    | Naphthalene, 1-methyl-7-(1-methy... | 184 | C14H16  | 000490-65-3  | 87   |
| 4    |    | Naphthalene, 1,2,3,4-tetramethyl-   | 184 | C14H16  | 003031-15-0  | 86   |
| 5    |    | Chamazulene                         | 184 | C14H16  | 000529-05-5  | 76   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

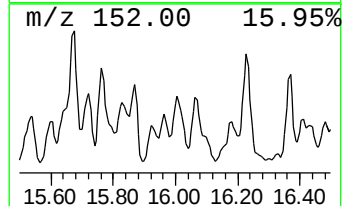
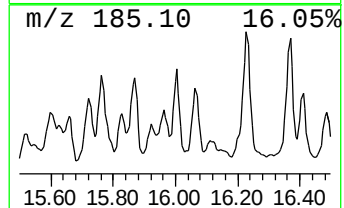
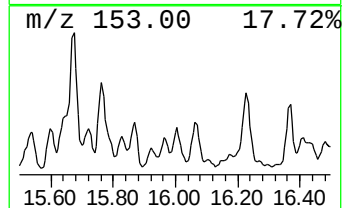
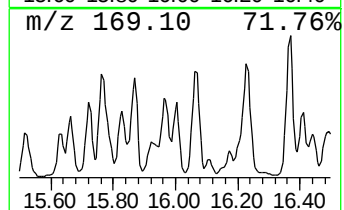
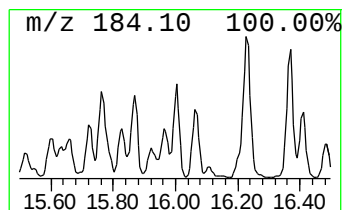
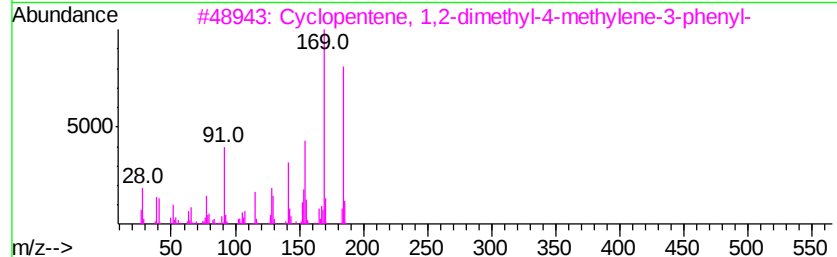
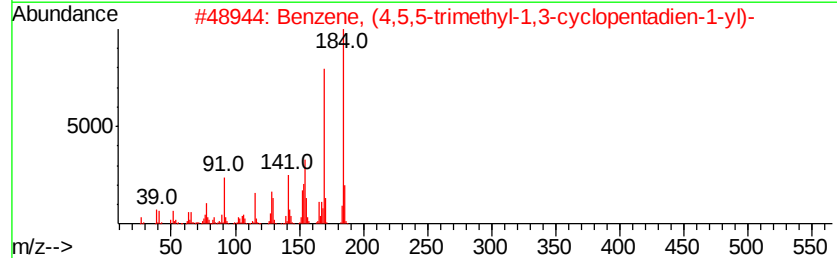
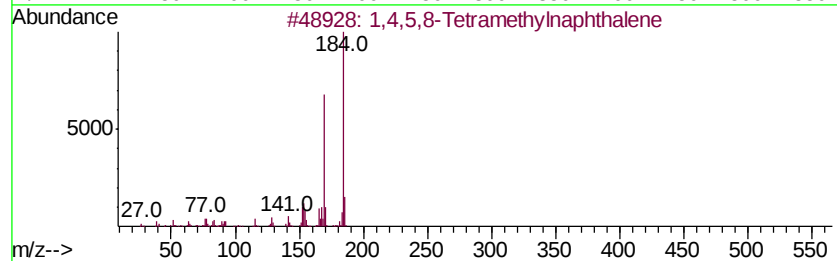
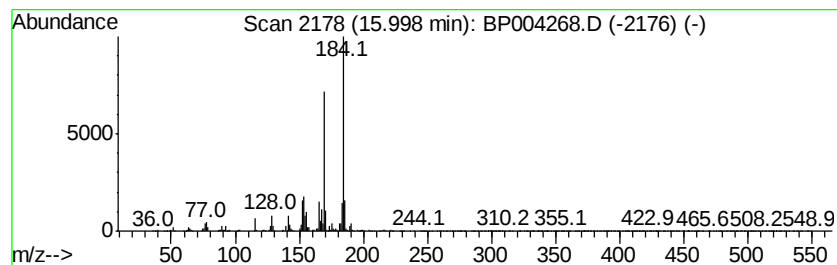
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 51 1,4,5,8-Tetramethylnaphthalene Concentration Rank 22

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 16.00 | 18.46 ng/ul | 3901950 | Phenanthrene-d10 | 17.26 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | 1,4,5,8-Tetramethylnaphthalene      | 184 | C14H16  | 002717-39-7  | 96   |
| 2    |    | Benzene, (4,5,5-trimethyl-1,3-cy... | 184 | C14H16  | 033930-85-7  | 87   |
| 3    |    | Cyclopentene, 1,2-dimethyl-4-met... | 184 | C14H16  | 1000152-22-4 | 80   |
| 4    |    | Naphthalene, 1,2,3,4-tetramethyl-   | 184 | C14H16  | 003031-15-0  | 74   |
| 5    |    | Chamazulene                         | 184 | C14H16  | 000529-05-5  | 68   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

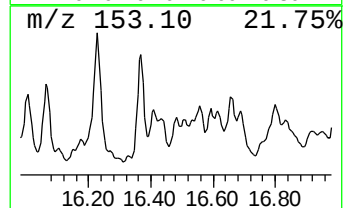
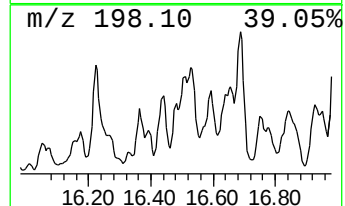
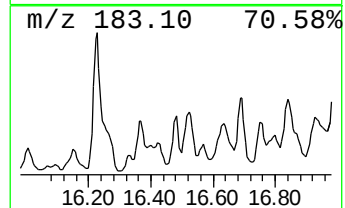
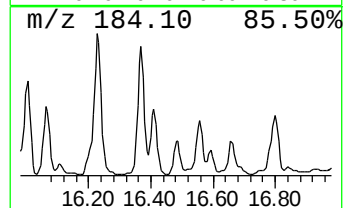
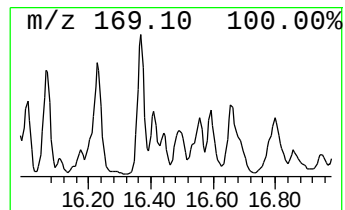
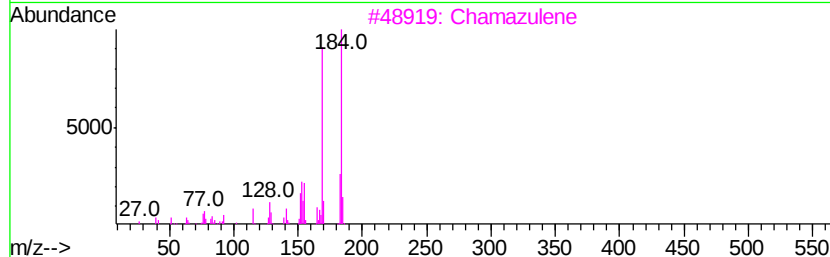
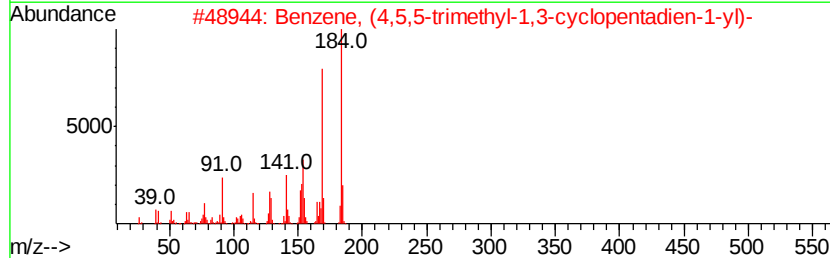
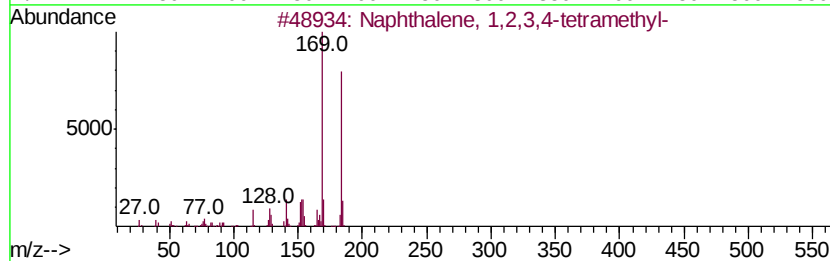
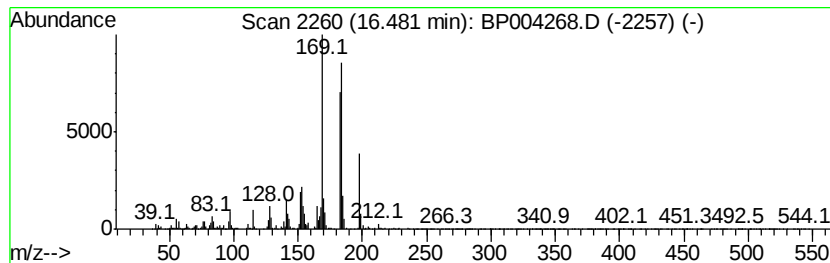
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 55 Naphthalene, 1,2,3,4-tetram... Concentration Rank 40

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 16.48 | 9.81 ng/ul | 2074050 | Phenanthrene-d10 | 17.26 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,2,3,4-tetramethyl-   | 184 | C14H16  | 003031-15-0 | 93   |
| 2    |    |   | Benzene, (4,5,5-trimethyl-1,3-cy... | 184 | C14H16  | 033930-85-7 | 89   |
| 3    |    |   | Chamazulene                         | 184 | C14H16  | 000529-05-5 | 89   |
| 4    |    |   | 1,4,5,8-Tetramethylnaphthalene      | 184 | C14H16  | 002717-39-7 | 86   |
| 5    |    |   | Naphthalene, 1-methyl-7-(1-methy... | 184 | C14H16  | 000490-65-3 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

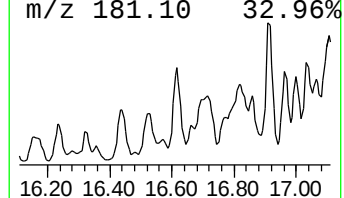
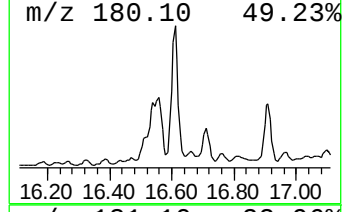
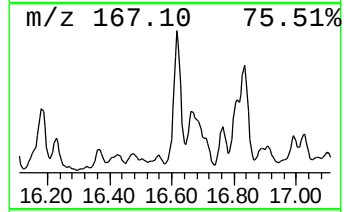
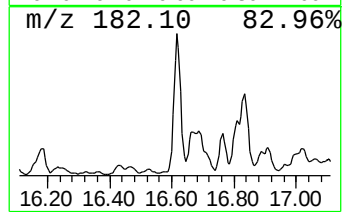
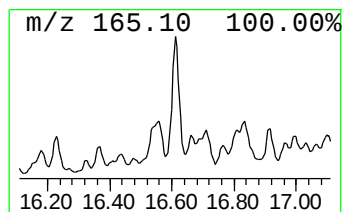
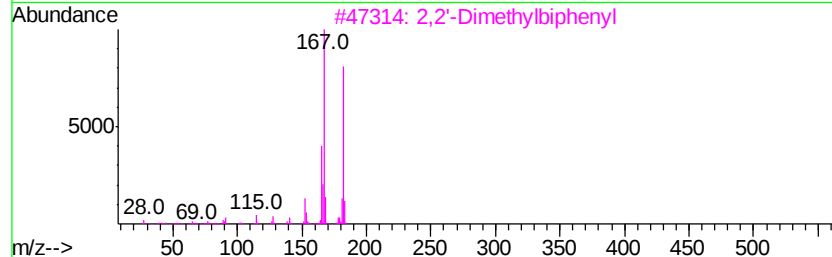
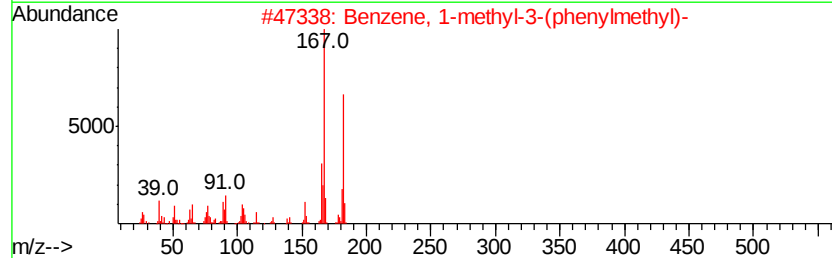
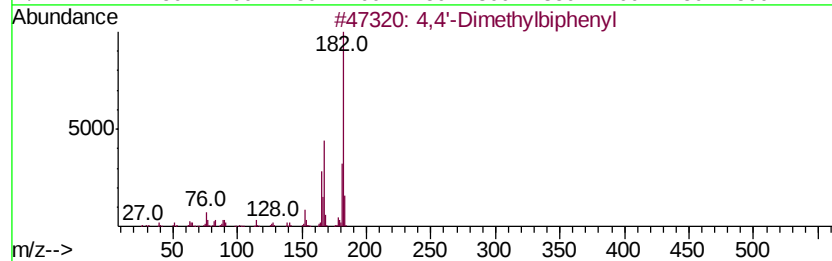
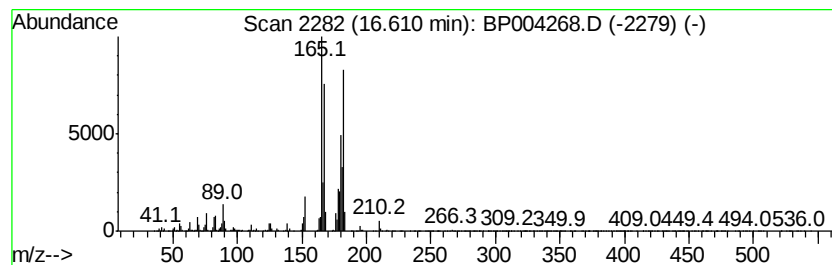
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 56 4,4'-Dimethylbiphenyl Concentration Rank 17

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 16.61 | 24.73 ng/ul | 5227700 | Phenanthrene-d10 | 17.26 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 4,4'-Dimethylbiphenyl               | 182 | C14H14   | 000613-33-2 | 70   |
| 2    |    | Benzene, 1-methyl-3-(phenylmethyl)- | 182 | C14H14   | 000620-47-3 | 46   |
| 3    |    | 2,2'-Dimethylbiphenyl               | 182 | C14H14   | 000605-39-0 | 43   |
| 4    |    | 1-Methyl-1,6-diazaplenalene         | 182 | C12H10N2 | 079691-99-9 | 43   |
| 5    |    | 9H-Fluorene, 2-methyl-              | 180 | C14H12   | 001430-97-3 | 42   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

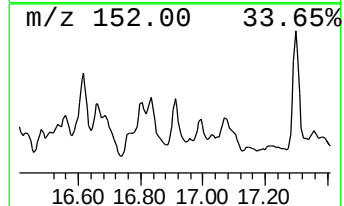
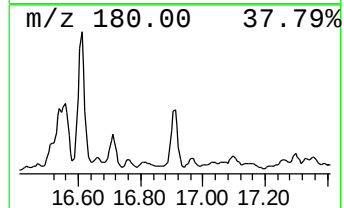
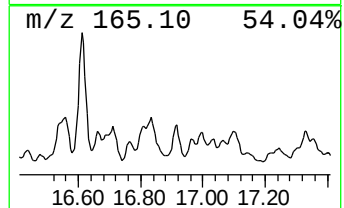
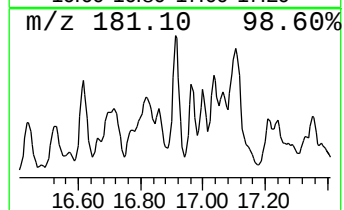
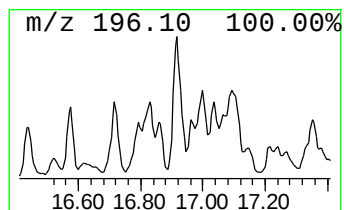
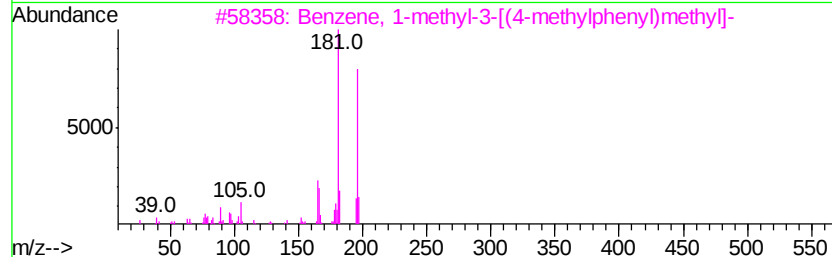
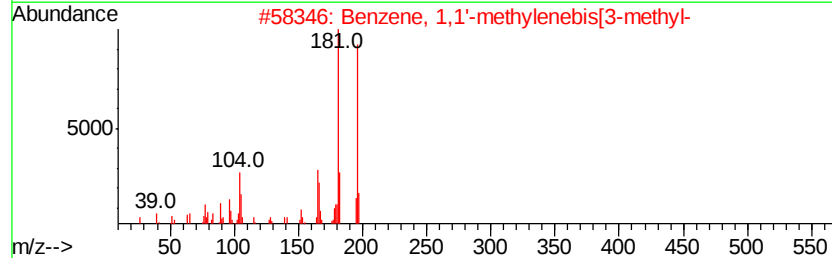
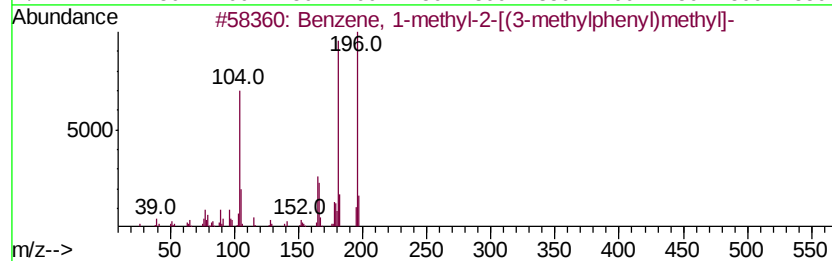
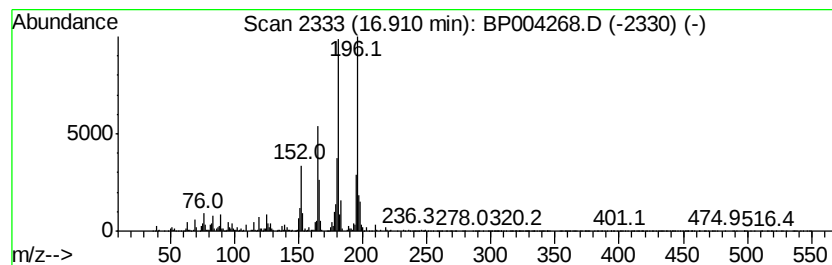
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 57 Benzene, 1-methyl-2-[(3-met... Concentration Rank 24

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 16.91 | 17.62 ng/ul | 3725040 | Phenanthrene-d10 | 17.26 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-2-[(3-methylph... | 196 | C15H16  | 021895-13-6 | 70   |
| 2    |    | Benzene, 1,1'-methylenebis[3-met... | 196 | C15H16  | 021895-14-7 | 68   |
| 3    |    | Benzene, 1-methyl-3-[(4-methylph... | 196 | C15H16  | 021895-16-9 | 58   |
| 4    |    | 2,2-Dimethyl-1-acenaphthenone       | 196 | C14H12O | 018086-43-6 | 52   |
| 5    |    | Methane, di-p-tolyl-                | 196 | C15H16  | 004957-14-6 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

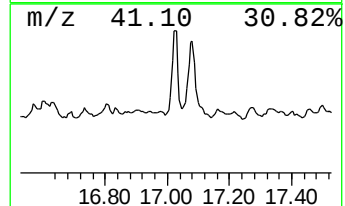
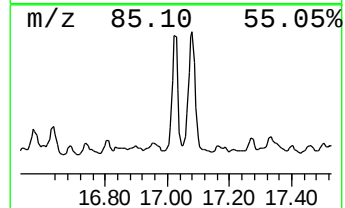
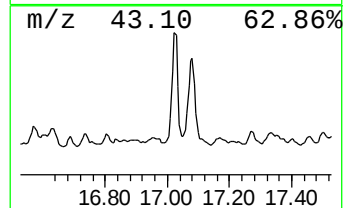
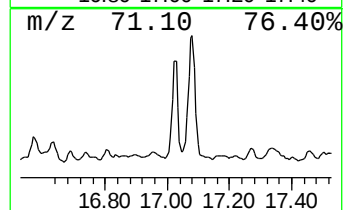
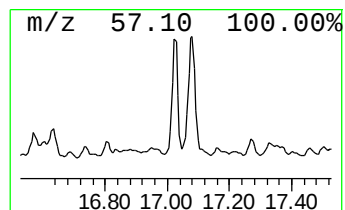
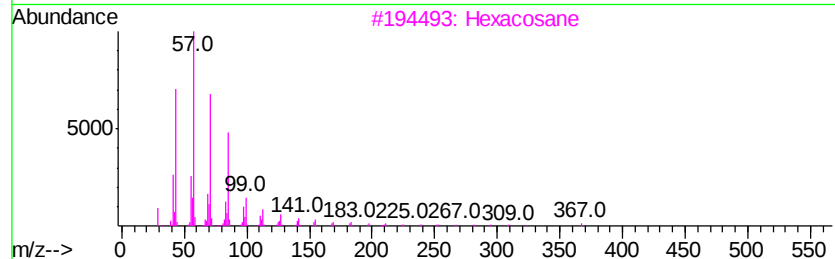
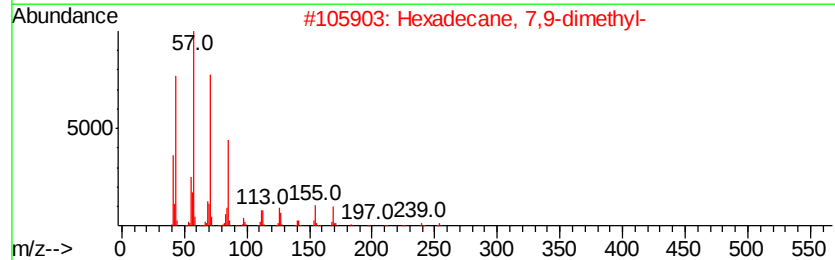
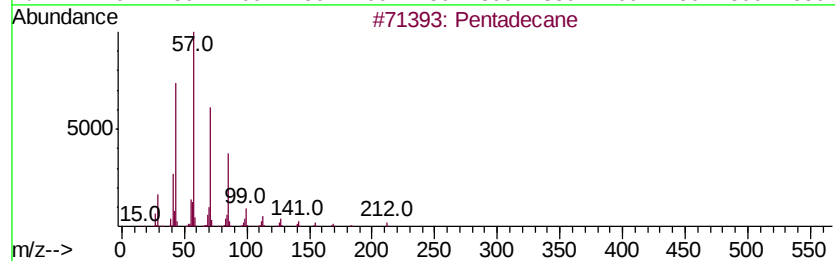
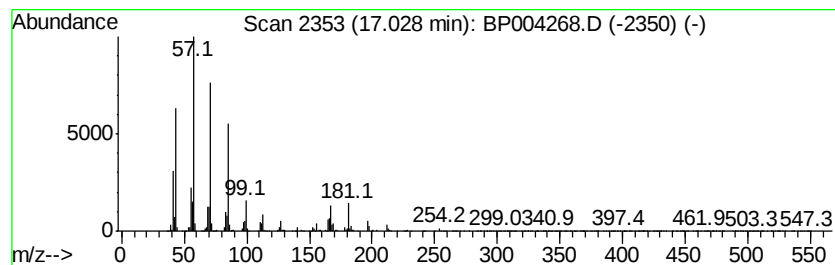
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 58 (DEL) Alkane: Straight-Chai... Concentration Rank 30

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 17.03 | 15.09 ng/ul | 3190410 | Phenanthrene-d10 | 17.26 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Pentadecane               | 212 | C15H32  | 000629-62-9 | 93   |
| 2    |    | Hexadecane, 7,9-dimethyl- | 254 | C18H38  | 021164-95-4 | 83   |
| 3    |    | Hexacosane                | 366 | C26H54  | 000630-01-3 | 81   |
| 4    |    | Pentadecane               | 212 | C15H32  | 000629-62-9 | 78   |
| 5    |    | Heptadecane               | 240 | C17H36  | 000629-78-7 | 74   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

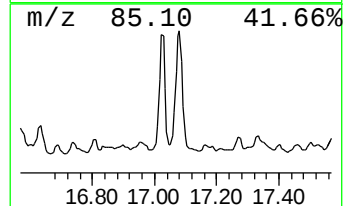
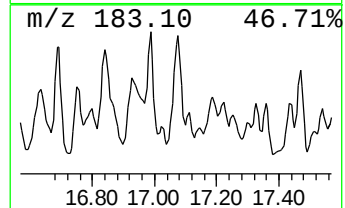
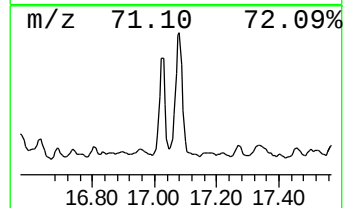
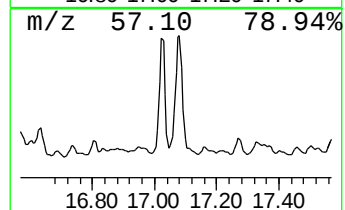
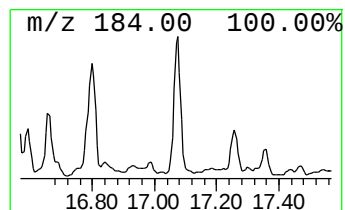
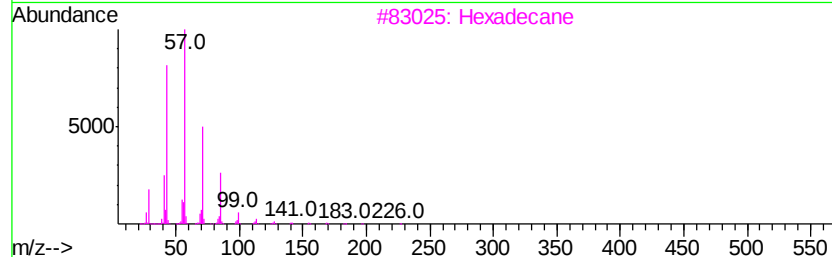
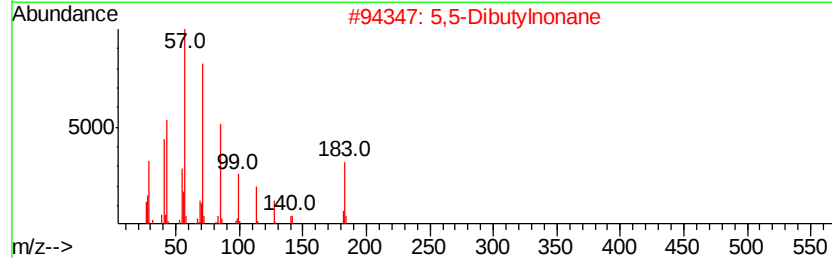
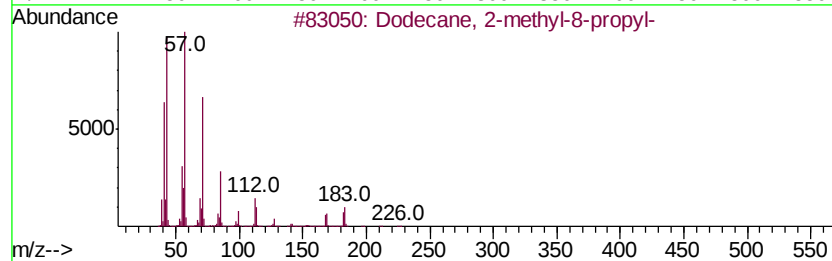
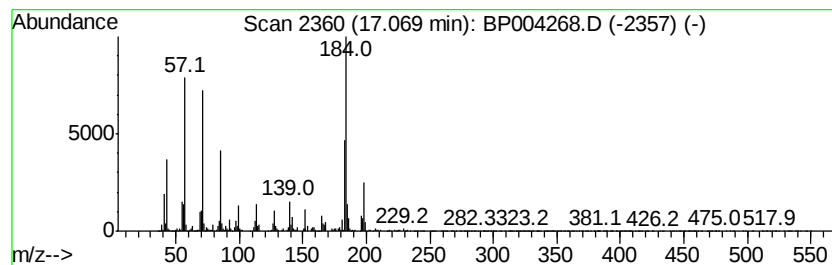
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 17.07 | 52.37 ng/ul | 11072200 | Phenanthrene-d10 | 17.26 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Dodecane, 2-methyl-8-propyl- | 226 | C16H34  | 055045-07-3 | 45   |
| 2    |    | 5,5-Dibutylnonane            | 240 | C17H36  | 006008-17-9 | 43   |
| 3    |    | Hexadecane                   | 226 | C16H34  | 000544-76-3 | 38   |
| 4    |    | Dibenzothiophene             | 184 | C12H8S  | 000132-65-0 | 38   |
| 5    |    | Hexadecane                   | 226 | C16H34  | 000544-76-3 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

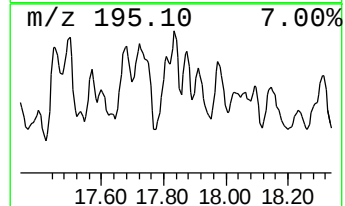
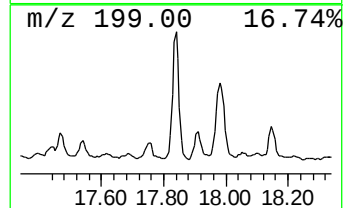
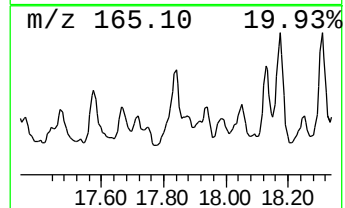
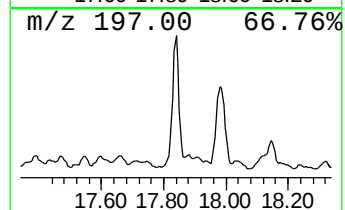
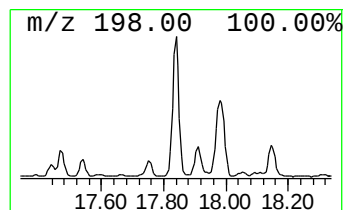
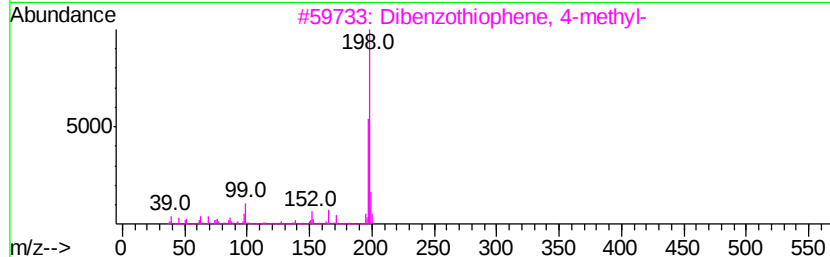
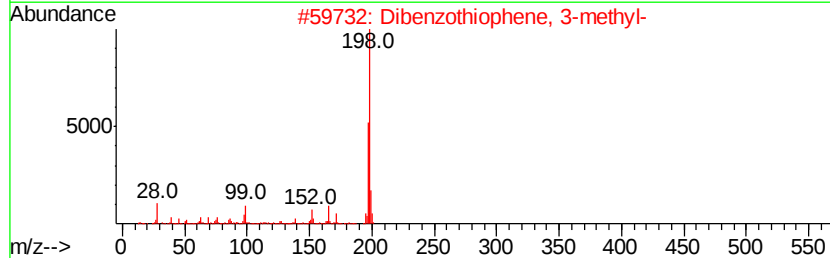
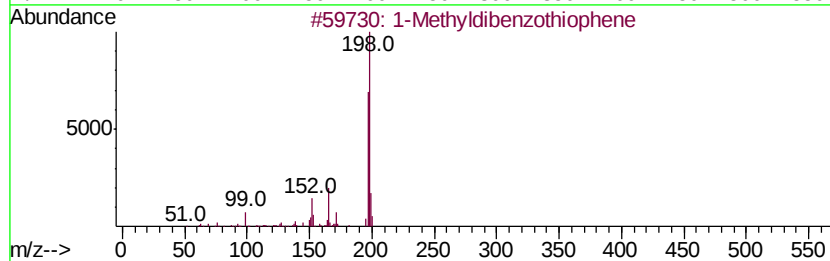
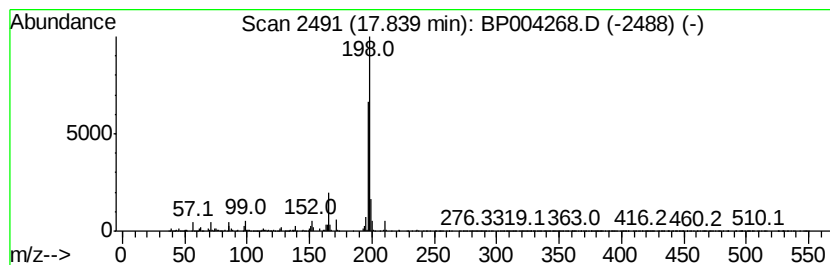
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TIC Integration Parameters: LSCINT.P

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Peak Number 60 1-Methyldibenzothiophene Concentration Rank 25

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 17.84 | 16.91 ng/ul | 3574290 | Phenanthrene-d10 | 17.26 |

| Hit# | of | Tentative ID                | MW  | MolForm  | CAS#        | Qual |
|------|----|-----------------------------|-----|----------|-------------|------|
| 1    | 5  | 1-Methyldibenzothiophene    | 198 | C13H10S  | 031317-07-4 | 87   |
| 2    |    | Dibenzothiophene, 3-methyl- | 198 | C13H10S  | 016587-52-3 | 87   |
| 3    |    | Dibenzothiophene, 4-methyl- | 198 | C13H10S  | 007372-88-5 | 80   |
| 4    |    | Tacrine                     | 198 | C13H14N2 | 000321-64-2 | 64   |
| 5    |    | Tacrine                     | 198 | C13H14N2 | 000321-64-2 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

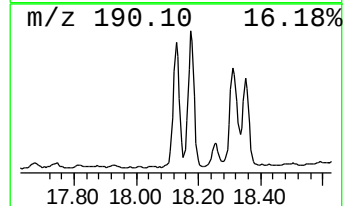
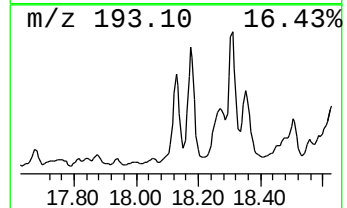
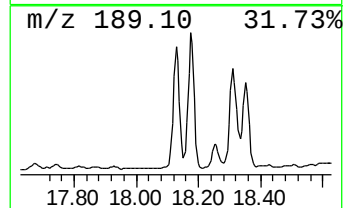
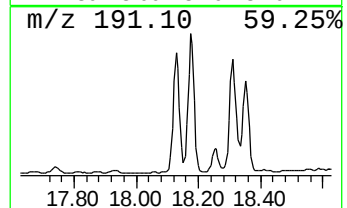
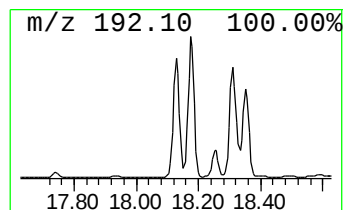
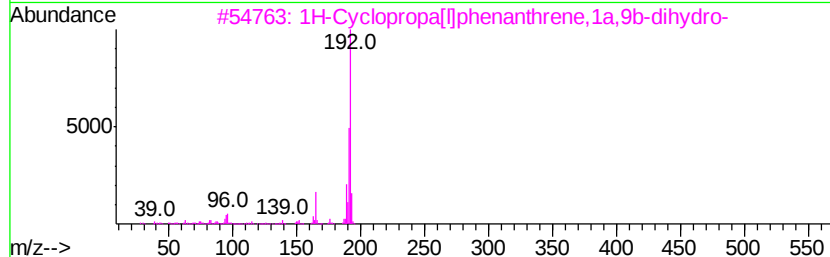
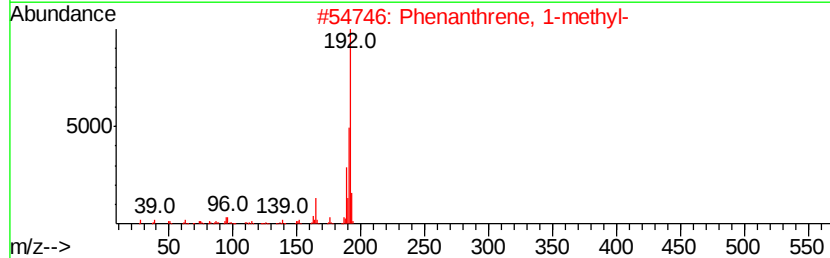
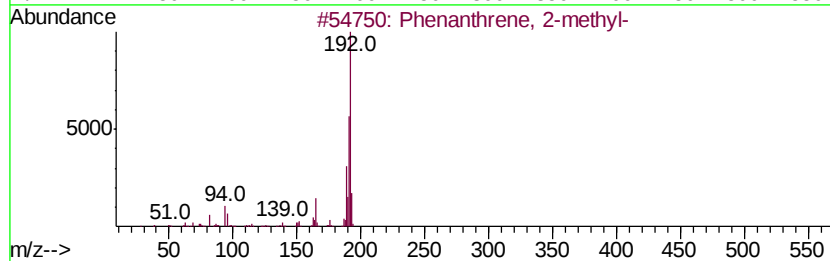
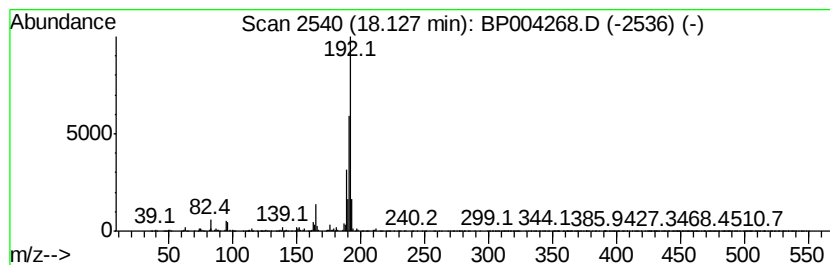
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 61 Phenanthrene, 2-methyl- Concentration Rank 5

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 18.13 | 43.61 ng/ul | 9219860 | Phenanthrene-d10 | 17.26 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 97   |
| 2    |    | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 97   |
| 3    |    | 1H-Cyclopropa[1]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 96   |
| 4    |    | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 96   |
| 5    |    | Anthracene, 2-methyl-               | 192 | C15H12  | 000613-12-7 | 95   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

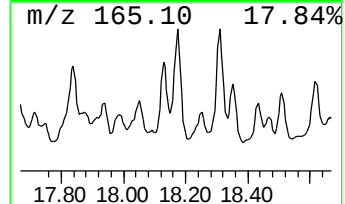
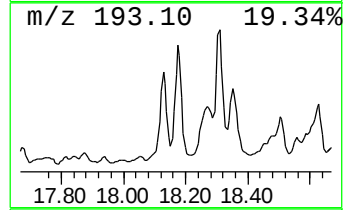
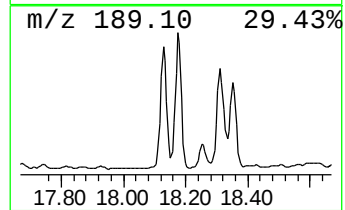
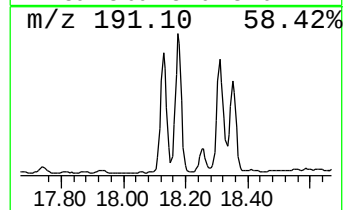
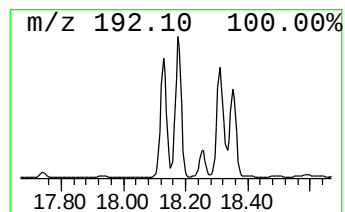
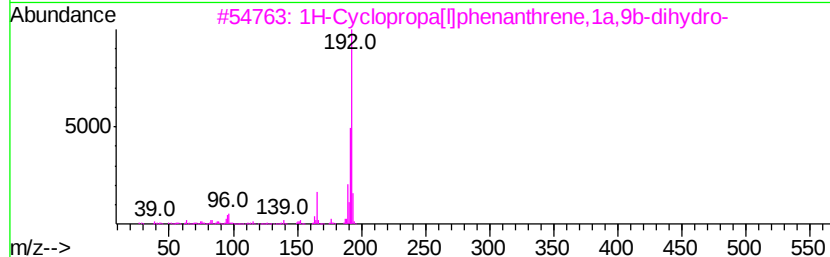
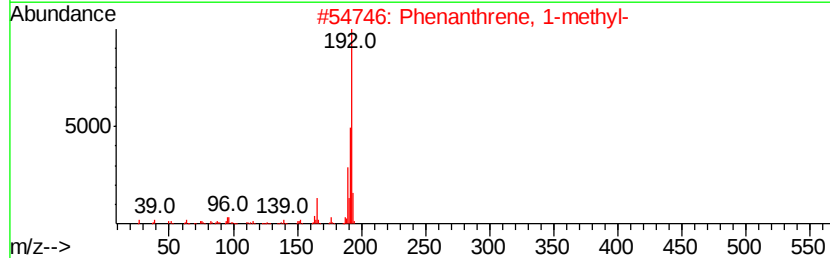
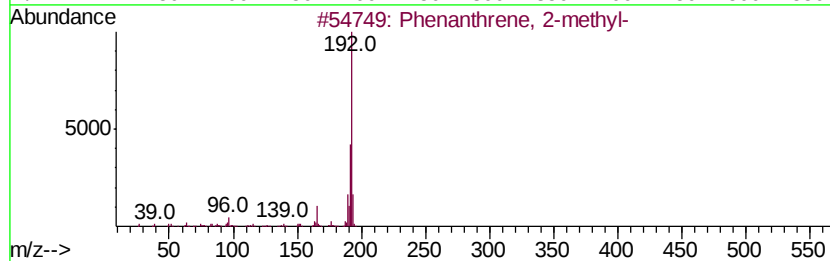
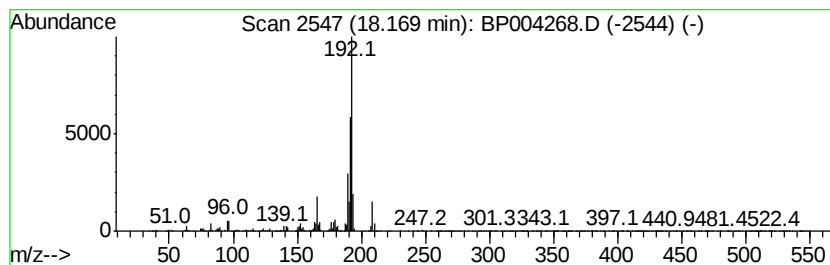
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 62 Phenanthrene, 1-methyl- Concentration Rank 2

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 18.17 | 63.38 ng/ul | 13400300 | Phenanthrene-d10 | 17.26 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 97   |
| 2    |    | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 97   |
| 3    |    | 1H-Cyclopropa[1]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 96   |
| 4    |    | Anthracene, 1-methyl-               | 192 | C15H12  | 000610-48-0 | 96   |
| 5    |    | 1H-Indene, 2-phenyl-                | 192 | C15H12  | 004505-48-0 | 95   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

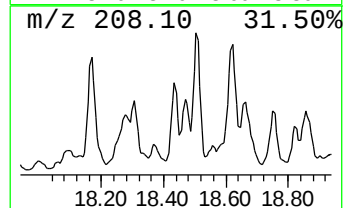
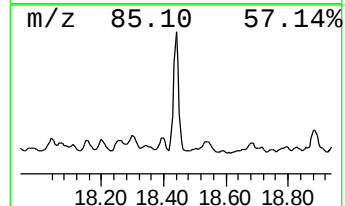
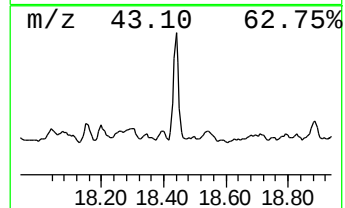
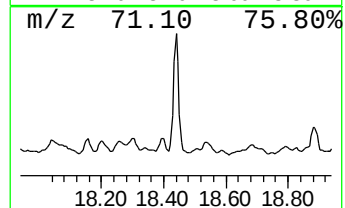
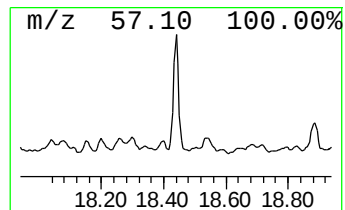
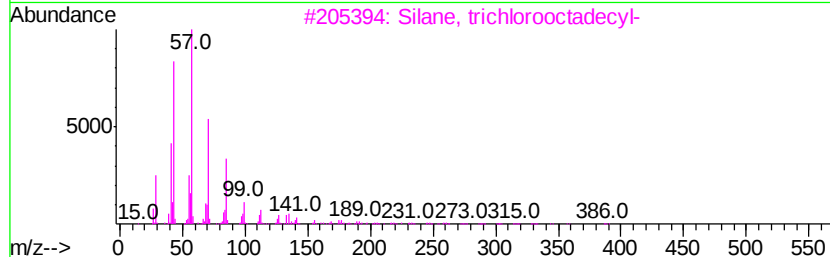
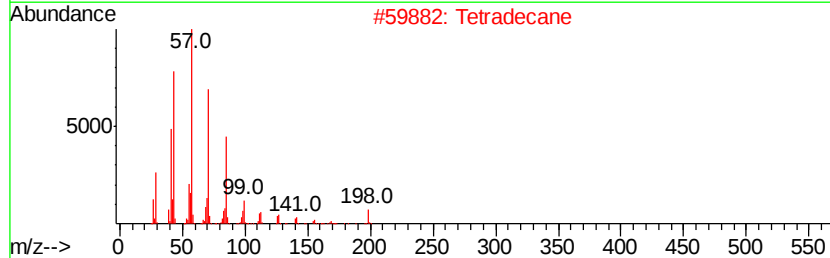
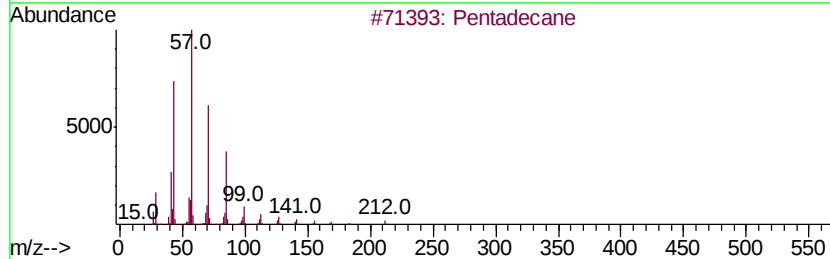
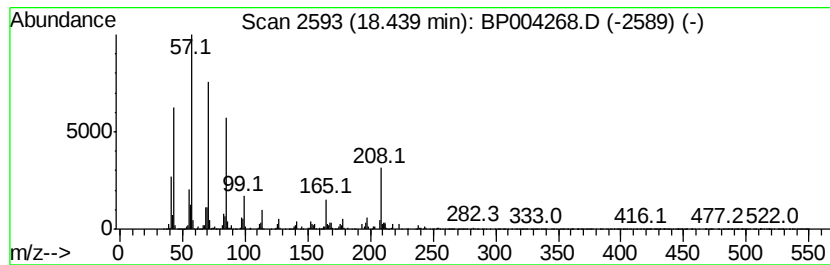
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 65 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 18.44 | 25.07 ng/ul | 5299200 | Phenanthrene-d10 | 17.26 |

| Hit# | of | Tentative ID                | MW  | MolForm     | CAS#        | Qual |
|------|----|-----------------------------|-----|-------------|-------------|------|
| 1    | 5  | Pentadecane                 | 212 | C15H32      | 000629-62-9 | 95   |
| 2    |    | Tetradecane                 | 198 | C14H30      | 000629-59-4 | 81   |
| 3    |    | Silane, trichlorooctadecyl- | 386 | C18H37Cl3Si | 000112-04-9 | 74   |
| 4    |    | Nonadecane                  | 268 | C19H40      | 000629-92-5 | 74   |
| 5    |    | Hexadecane                  | 226 | C16H34      | 000544-76-3 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

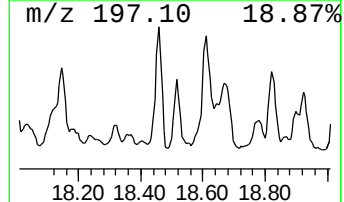
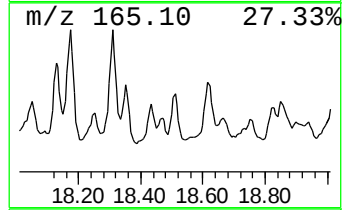
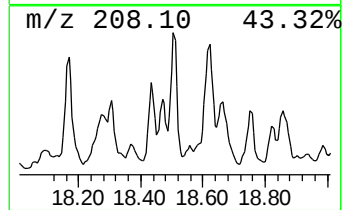
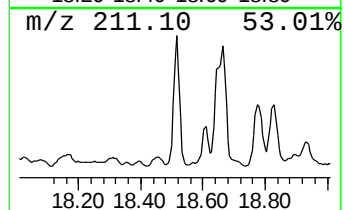
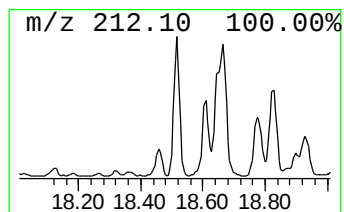
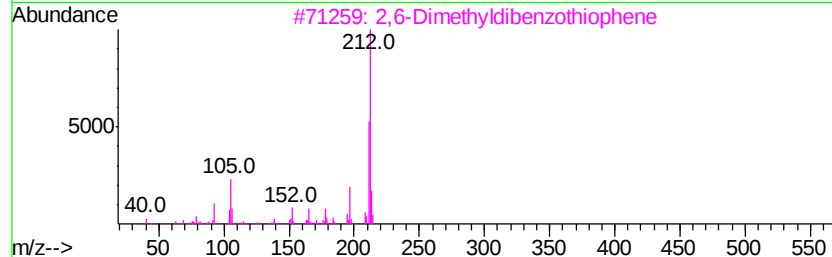
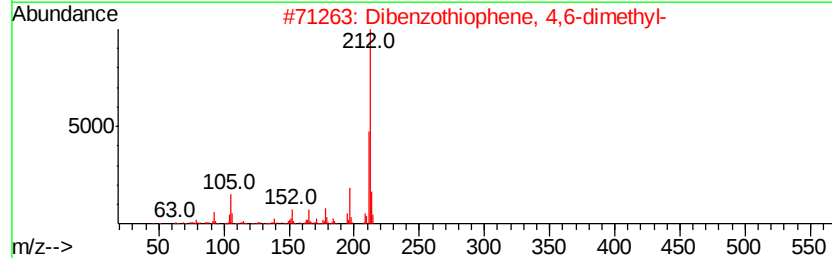
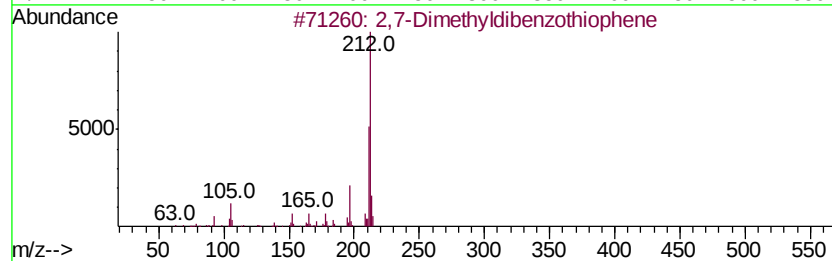
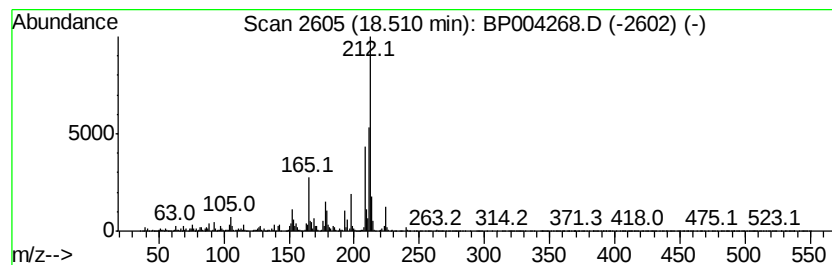
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 66 2,7-Dimethyldibenzothiophene Concentration Rank 21

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 18.51 | 19.12 ng/ul | 4043330 | Phenanthrene-d10 | 17.26 |

| Hit# | of | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2,7-Dimethyldibenzothiophene      | 212 | C14H12S | 031317-19-8 | 86   |
| 2    |    | Dibenzothiophene, 4,6-dimethyl-   | 212 | C14H12S | 001207-12-1 | 70   |
| 3    |    | 2,6-Dimethyldibenzothiophene      | 212 | C14H12S | 089816-75-1 | 70   |
| 4    |    | 2,8-Dimethyldibenzo(b,d)thiophene | 212 | C14H12S | 001207-15-4 | 62   |
| 5    |    | 1,7-Dimethyldibenzothiophene      | 212 | C14H12S | 089816-53-5 | 58   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
Data File : BP004268.D  
Acq On : 14 Dec 2020 11:39  
Operator : CG/JU  
Sample : L5063-01DL 2X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION

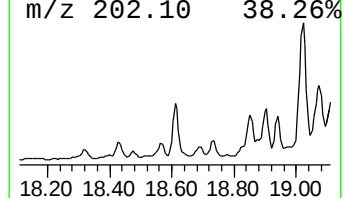
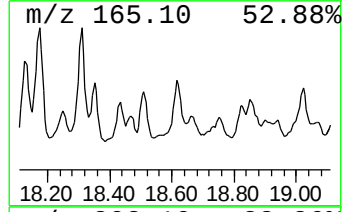
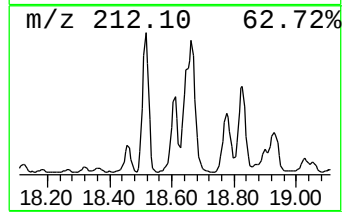
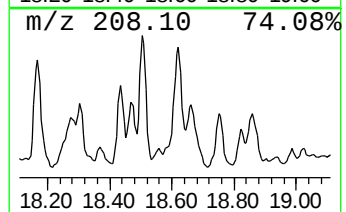
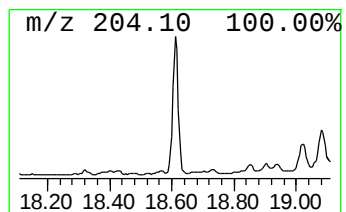
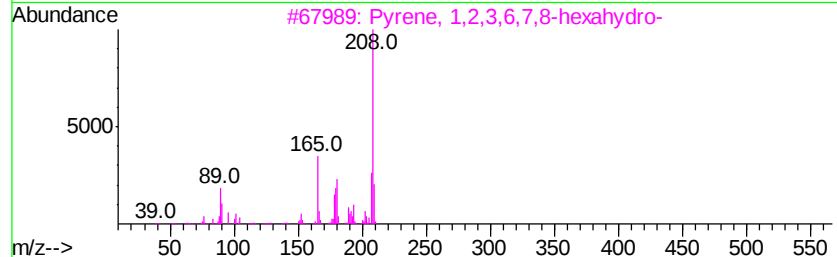
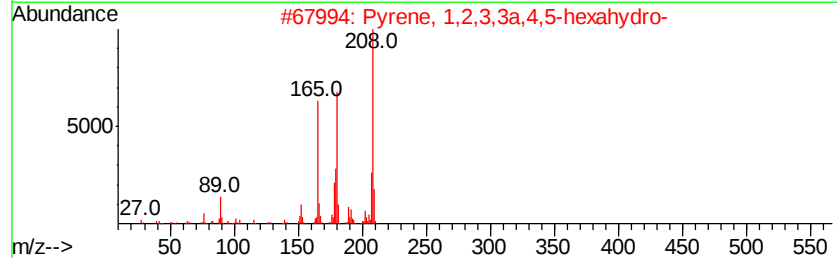
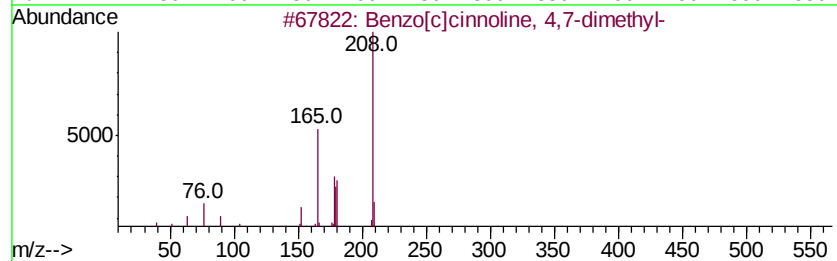
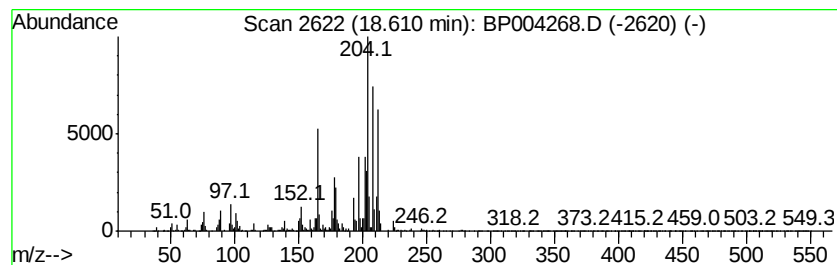
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 67 Benzo[c]cinnoline, 4,7-dime... Concentration Rank 39

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 18.61 | 10.16 ng/ul | 2148300 | Phenanthrene-d10 | 17.26 |

| Hit# | of | Tentative ID                     | MW  | MolForm  | CAS#        | Qual |
|------|----|----------------------------------|-----|----------|-------------|------|
| 1    | 5  | Benzo[c]cinnoline, 4,7-dimethyl- | 208 | C14H12N2 | 020684-54-2 | 80   |
| 2    |    | Pyrene, 1,2,3,3a,4,5-hexahydro-  | 208 | C16H16   | 005385-37-5 | 42   |
| 3    |    | Pyrene, 1,2,3,6,7,8-hexahydro-   | 208 | C16H16   | 001732-13-4 | 38   |
| 4    |    | Anthracene, 1-methoxy-           | 208 | C15H12O  | 054458-84-3 | 38   |
| 5    |    | Pyrene, 1,2,3,6,7,8-hexahydro-   | 208 | C16H16   | 001732-13-4 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_P\DATA\BP121420\  
 Data File : BP004268.D  
 Acq On : 14 Dec 2020 11:39  
 Operator : CG/JU  
 Sample : L5063-01DL 2X  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0AD5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.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 |
| (DEL) Alkane: Str...  | 10.81 | 2.3     | ng/ul | 697697   | 2                     | 10.64 | 6145860 | 20.0 |
| (DEL) Alkane: Cyc...  | 11.14 | 4.2     | ng/ul | 1275730  | 2                     | 10.64 | 6145860 | 20.0 |
| 1H-Indene, 2,3-di...  | 11.56 | 7.2     | ng/ul | 2199340  | 2                     | 10.64 | 6145860 | 20.0 |
| (DEL) Alkane: Str...  | 11.68 | 6.1     | ng/ul | 1865950  | 2                     | 10.64 | 6145860 | 20.0 |
| Benzene, (1,2-dim...  | 11.75 | 5.8     | ng/ul | 1765990  | 2                     | 10.64 | 6145860 | 20.0 |
| Naphthalene, 1,2,...  | 11.84 | 5.1     | ng/ul | 1576180  | 2                     | 10.64 | 6145860 | 20.0 |
| (DEL) Alkane: Cyc...  | 11.94 | 2.5     | ng/ul | 774364   | 2                     | 10.64 | 6145860 | 20.0 |
| (DEL) Alkane: Str...  | 12.07 | 5.5     | ng/ul | 1690970  | 2                     | 10.64 | 6145860 | 20.0 |
| 1H-Indene, 2,3-di...  | 12.40 | 7.2     | ng/ul | 2197930  | 2                     | 10.64 | 6145860 | 20.0 |
| Naphthalene, 1-me...  | 12.52 | 14.0    | ng/ul | 4294530  | 2                     | 10.64 | 6145860 | 20.0 |
| 1H-Indene, 2,3-di...  | 13.00 | 4.4     | ng/ul | 1539050  | 3                     | 14.49 | 6986490 | 20.0 |
| (DEL) Alkane: Str...  | 13.03 | 7.0     | ng/ul | 2444680  | 3                     | 14.49 | 6986490 | 20.0 |
| Benzene, 1-(2-but...  | 13.42 | 9.7     | ng/ul | 3404090  | 3                     | 14.49 | 6986490 | 20.0 |
| Naphthalene, 2-et...  | 13.52 | 3.6     | ng/ul | 1275770  | 3                     | 14.49 | 6986490 | 20.0 |
| Naphthalene, 2,7-...  | 13.65 | 15.8    | ng/ul | 5533460  | 3                     | 14.49 | 6986490 | 20.0 |
| Naphthalene, 2,3-...  | 13.82 | 33.9    | ng/ul | 11842900 | 3                     | 14.49 | 6986490 | 20.0 |
| Naphthalene, 1,4-...  | 14.05 | 11.4    | ng/ul | 3982040  | 3                     | 14.49 | 6986490 | 20.0 |
| unknown-01            | 14.33 | 7.4     | ng/ul | 2588820  | 3                     | 14.49 | 6986490 | 20.0 |
| (DEL) Alkane: Str...  | 14.40 | 10.3    | ng/ul | 3597920  | 3                     | 14.49 | 6986490 | 20.0 |
| Naphthalene, 2-(1...  | 14.63 | 7.1     | ng/ul | 2467310  | 3                     | 14.49 | 6986490 | 20.0 |
| Naphthalene, 2,3,...  | 14.97 | 32.6    | ng/ul | 11391200 | 3                     | 14.49 | 6986490 | 20.0 |
| unknown-02            | 15.03 | 7.8     | ng/ul | 2730810  | 3                     | 14.49 | 6986490 | 20.0 |
| Naphthalene, 1,6,...  | 15.32 | 10.4    | ng/ul | 3631480  | 3                     | 14.49 | 6986490 | 20.0 |
| (DEL) Alkane: Str...  | 15.36 | 7.9     | ng/ul | 2764280  | 3                     | 14.49 | 6986490 | 20.0 |
| Naphthalene, 2-me...  | 15.60 | 12.7    | ng/ul | 4445170  | 3                     | 14.49 | 6986490 | 20.0 |
| 1-Isopropenyl naph... | 15.67 | 11.9    | ng/ul | 4141200  | 3                     | 14.49 | 6986490 | 20.0 |
| Cyclopentene, 1,2...  | 15.87 | 13.6    | ng/ul | 4745050  | 3                     | 14.49 | 6986490 | 20.0 |
| 1,4,5,8-Tetrameth...  | 16.00 | 18.5    | ng/ul | 3901950  | 4                     | 17.26 | 4228340 | 20.0 |
| Chamazulene           | 16.06 | 15.3    | ng/ul | 3230390  | 4                     | 17.26 | 4228340 | 20.0 |
| Naphthalene, 1,2,...  | 16.48 | 9.8     | ng/ul | 2074050  | 4                     | 17.26 | 4228340 | 20.0 |
| 4,4'-Dimethylbiph...  | 16.61 | 24.7    | ng/ul | 5227700  | 4                     | 17.26 | 4228340 | 20.0 |
| Benzene, 1-methyl...  | 16.91 | 17.6    | ng/ul | 3725040  | 4                     | 17.26 | 4228340 | 20.0 |
| (DEL) Alkane: Str...  | 17.03 | 15.1    | ng/ul | 3190410  | 4                     | 17.26 | 4228340 | 20.0 |
| (DEL) Alkane: Str...  | 17.07 | 52.4    | ng/ul | 11072200 | 4                     | 17.26 | 4228340 | 20.0 |
| 1-Methyldibenzoth...  | 17.84 | 16.9    | ng/ul | 3574290  | 4                     | 17.26 | 4228340 | 20.0 |
| Phenanthrene, 2-m...  | 18.13 | 43.6    | ng/ul | 9219860  | 4                     | 17.26 | 4228340 | 20.0 |
| Phenanthrene, 1-m...  | 18.17 | 63.4    | ng/ul | 13400300 | 4                     | 17.26 | 4228340 | 20.0 |
| (DEL) Alkane: Str...  | 18.44 | 25.1    | ng/ul | 5299200  | 4                     | 17.26 | 4228340 | 20.0 |
| 2,7-Dimethyldiben...  | 18.51 | 19.1    | ng/ul | 4043330  | 4                     | 17.26 | 4228340 | 20.0 |
| Benzo[c]cinnoline...  | 18.61 | 10.2    | ng/ul | 2148300  | 4                     | 17.26 | 4228340 | 20.0 |