

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
 Data File : BM013819.D  
 Acq On : 25 Jan 2018 23:11  
 Operator : SJ/JU  
 Sample : J1222-08DL 20X  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 F6A81DL

## Integration Parameters: LSCINT.P

Integrator: RTE  
 Smoothing : OFF  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0

Filtering: 5  
 Min Area: 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:\HPCHEM1\BNA\_M\Methods\SOM-EPA-BM011018.M  
 Title : SVOA CALIBRATION

Signal : TIC

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|--------|----------|------------|----------|-----------|-------|-------------|------------|--------------|------------|
| 1      | 4.757    | 297        | 301      | 310       | rVB   | 96196       | 143644     | 4.17%        | 0.378%     |
| 2      | 5.628    | 443        | 449      | 452       | rBV5  | 44036       | 64258      | 1.86%        | 0.169%     |
| 3      | 5.951    | 498        | 504      | 516       | rBV   | 153440      | 269222     | 7.81%        | 0.708%     |
| 4      | 6.822    | 644        | 652      | 656       | rVB4  | 26095       | 61716      | 1.79%        | 0.162%     |
| 5      | 7.210    | 712        | 718      | 721       | rBV2  | 88326       | 137319     | 3.98%        | 0.361%     |
| 6      | 7.592    | 772        | 783      | 792       | rBV   | 512344      | 879186     | 25.49%       | 2.312%     |
| 7      | 7.834    | 817        | 824      | 832       | rBV4  | 38652       | 79856      | 2.32%        | 0.210%     |
| 8      | 8.233    | 885        | 892      | 897       | rBV5  | 46031       | 95193      | 2.76%        | 0.250%     |
| 9      | 8.339    | 904        | 910      | 914       | rBV7  | 31777       | 58227      | 1.69%        | 0.153%     |
| 10     | 8.804    | 985        | 989      | 998       | rVB   | 181773      | 271074     | 7.86%        | 0.713%     |
| 11     | 9.486    | 1099       | 1105     | 1110      | rVB9  | 22386       | 52820      | 1.53%        | 0.139%     |
| 12     | 9.792    | 1149       | 1157     | 1161      | rBV6  | 44319       | 93232      | 2.70%        | 0.245%     |
| 13     | 10.057   | 1195       | 1202     | 1212      | rVB9  | 18532       | 63602      | 1.84%        | 0.167%     |
| 14     | 10.363   | 1243       | 1254     | 1259      | rVV2  | 640031      | 1546940    | 44.86%       | 4.068%     |
| 15     | 10.416   | 1259       | 1263     | 1270      | rVV   | 242225      | 413249     | 11.98%       | 1.087%     |
| 16     | 10.486   | 1270       | 1275     | 1278      | rVV7  | 33099       | 55860      | 1.62%        | 0.147%     |
| 17     | 10.522   | 1278       | 1281     | 1289      | rVB3  | 74343       | 127882     | 3.71%        | 0.336%     |
| 18     | 10.657   | 1299       | 1304     | 1313      | rVB9  | 29277       | 61840      | 1.79%        | 0.163%     |
| 19     | 11.057   | 1362       | 1372     | 1376      | rBV10 | 32484       | 79276      | 2.30%        | 0.208%     |
| 20     | 11.198   | 1390       | 1396     | 1399      | rBV5  | 36533       | 59387      | 1.72%        | 0.156%     |
| 21     | 11.274   | 1404       | 1409     | 1416      | rVB6  | 71965       | 135488     | 3.93%        | 0.356%     |
| 22     | 11.386   | 1423       | 1428     | 1433      | rBV3  | 136085      | 233686     | 6.78%        | 0.615%     |
| 23     | 11.469   | 1438       | 1442     | 1447      | rVB4  | 58135       | 95779      | 2.78%        | 0.252%     |
| 24     | 11.792   | 1488       | 1497     | 1502      | rBV   | 448874      | 690949     | 20.04%       | 1.817%     |
| 25     | 12.039   | 1533       | 1539     | 1547      | rVB   | 282985      | 486414     | 14.10%       | 1.279%     |
| 26     | 12.257   | 1567       | 1576     | 1585      | rVV   | 180306      | 365074     | 10.59%       | 0.960%     |
| 27     | 12.439   | 1601       | 1607     | 1611      | rBV7  | 40320       | 84291      | 2.44%        | 0.222%     |
| 28     | 12.498   | 1611       | 1617     | 1623      | rVV5  | 53108       | 136979     | 3.97%        | 0.360%     |
| 29     | 12.557   | 1623       | 1627     | 1634      | rVV8  | 39531       | 91820      | 2.66%        | 0.241%     |
| 30     | 12.627   | 1634       | 1639     | 1649      | rVV2  | 84083       | 157135     | 4.56%        | 0.413%     |
| 31     | 12.710   | 1649       | 1653     | 1659      | rVV6  | 64320       | 114698     | 3.33%        | 0.302%     |
| 32     | 12.763   | 1659       | 1662     | 1668      | rVV2  | 153652      | 234738     | 6.81%        | 0.617%     |
| 33     | 13.063   | 1705       | 1713     | 1720      | rVV2  | 663527      | 994061     | 28.82%       | 2.614%     |
| 34     | 13.268   | 1744       | 1748     | 1750      | rBV   | 65139       | 88832      | 2.58%        | 0.234%     |

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 Max Peaks: 100  
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
 Peak separation: 5

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 Title : SVOA CALIBRATION

|    |        |      |      |      |       |         |         |         |        |
|----|--------|------|------|------|-------|---------|---------|---------|--------|
| 35 | 13.410 | 1765 | 1772 | 1782 | rBV4  | 131865  | 368424  | 10.68%  | 0.969% |
| 36 | 13.557 | 1791 | 1797 | 1802 | rBV   | 208995  | 403850  | 11.71%  | 1.062% |
| 37 | 13.616 | 1802 | 1807 | 1810 | rVV2  | 94848   | 144202  | 4.18%   | 0.379% |
| 38 | 13.727 | 1818 | 1826 | 1829 | rVV4  | 227557  | 357539  | 10.37%  | 0.940% |
| 39 | 13.768 | 1829 | 1833 | 1836 | rVV2  | 90077   | 127582  | 3.70%   | 0.336% |
| 40 | 13.798 | 1836 | 1838 | 1841 | rVV3  | 60559   | 64013   | 1.86%   | 0.168% |
| 41 | 13.845 | 1841 | 1846 | 1849 | rVB4  | 47178   | 69259   | 2.01%   | 0.182% |
| 42 | 13.933 | 1858 | 1861 | 1864 | rBV2  | 63004   | 85259   | 2.47%   | 0.224% |
| 43 | 13.968 | 1864 | 1867 | 1871 | rVB3  | 48022   | 64227   | 1.86%   | 0.169% |
| 44 | 14.063 | 1879 | 1883 | 1891 | rVB10 | 47113   | 84171   | 2.44%   | 0.221% |
| 45 | 14.145 | 1891 | 1897 | 1902 | rBV   | 793299  | 1052052 | 30.51%  | 2.767% |
| 46 | 14.239 | 1902 | 1913 | 1920 | rBV3  | 990723  | 2273763 | 65.93%  | 5.980% |
| 47 | 14.304 | 1920 | 1924 | 1933 | rVV   | 265065  | 432510  | 12.54%  | 1.137% |
| 48 | 14.386 | 1934 | 1938 | 1945 | rVB5  | 96221   | 156865  | 4.55%   | 0.413% |
| 49 | 14.457 | 1945 | 1950 | 1959 | rVB3  | 86266   | 239006  | 6.93%   | 0.629% |
| 50 | 14.539 | 1959 | 1964 | 1967 | rBV6  | 37821   | 74784   | 2.17%   | 0.197% |
| 51 | 14.604 | 1970 | 1975 | 1978 | rBV5  | 74593   | 153303  | 4.45%   | 0.403% |
| 52 | 14.651 | 1979 | 1983 | 1989 | rVB   | 289310  | 410356  | 11.90%  | 1.079% |
| 53 | 14.727 | 1990 | 1996 | 2000 | rVB3  | 76092   | 139577  | 4.05%   | 0.367% |
| 54 | 14.774 | 2000 | 2004 | 2008 | rBV4  | 145744  | 230303  | 6.68%   | 0.606% |
| 55 | 14.886 | 2017 | 2023 | 2027 | rBV4  | 126813  | 251241  | 7.29%   | 0.661% |
| 56 | 15.015 | 2040 | 2045 | 2052 | rBV8  | 71398   | 192493  | 5.58%   | 0.506% |
| 57 | 15.104 | 2052 | 2060 | 2065 | rVB   | 766568  | 972869  | 28.21%  | 2.558% |
| 58 | 15.239 | 2080 | 2083 | 2088 | rVB3  | 79972   | 109672  | 3.18%   | 0.288% |
| 59 | 15.298 | 2088 | 2093 | 2097 | rBV   | 241401  | 348520  | 10.11%  | 0.917% |
| 60 | 15.351 | 2098 | 2102 | 2108 | rVB5  | 58386   | 101516  | 2.94%   | 0.267% |
| 61 | 15.415 | 2109 | 2113 | 2118 | rVB7  | 40028   | 62278   | 1.81%   | 0.164% |
| 62 | 15.515 | 2124 | 2130 | 2134 | rBV2  | 212592  | 332630  | 9.65%   | 0.875% |
| 63 | 15.604 | 2141 | 2145 | 2152 | rVB7  | 99523   | 186598  | 5.41%   | 0.491% |
| 64 | 15.668 | 2152 | 2156 | 2159 | rBV3  | 77861   | 127521  | 3.70%   | 0.335% |
| 65 | 15.715 | 2159 | 2164 | 2175 | rVB   | 652751  | 993482  | 28.81%  | 2.613% |
| 66 | 15.809 | 2176 | 2180 | 2184 | rBV7  | 45177   | 69226   | 2.01%   | 0.182% |
| 67 | 15.968 | 2202 | 2207 | 2215 | rBV   | 761493  | 1545559 | 44.82%  | 4.065% |
| 68 | 16.039 | 2215 | 2219 | 2223 | rVB2  | 326821  | 418979  | 12.15%  | 1.102% |
| 69 | 16.309 | 2262 | 2265 | 2269 | rBV5  | 75662   | 92609   | 2.69%   | 0.244% |
| 70 | 16.651 | 2317 | 2323 | 2333 | rBV   | 2434831 | 3448647 | 100.00% | 9.069% |
| 71 | 16.762 | 2338 | 2342 | 2347 | rVV   | 685943  | 893980  | 25.92%  | 2.351% |

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Method : Z:\HPCHEM1\BNA\_M\Methods\SOM-EPA-BM011018.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |        |        |
|-----|--------|------|------|------|------|---------|---------|--------|--------|
| 72  | 16.815 | 2347 | 2351 | 2362 | rVB2 | 239222  | 430088  | 12.47% | 1.131% |
| 73  | 16.904 | 2362 | 2366 | 2374 | rBV2 | 94219   | 164145  | 4.76%  | 0.432% |
| 74  | 16.998 | 2376 | 2382 | 2386 | rBV  | 1109319 | 1586155 | 45.99% | 4.171% |
| 75  | 17.039 | 2386 | 2389 | 2395 | rVV  | 702805  | 975731  | 28.29% | 2.566% |
| 76  | 17.098 | 2395 | 2399 | 2402 | rVV  | 103237  | 174422  | 5.06%  | 0.459% |
| 77  | 17.133 | 2402 | 2405 | 2410 | rVB  | 115243  | 163648  | 4.75%  | 0.430% |
| 78  | 17.227 | 2417 | 2421 | 2426 | rVB4 | 120041  | 208982  | 6.06%  | 0.550% |
| 79  | 17.298 | 2429 | 2433 | 2440 | rVB9 | 50233   | 91095   | 2.64%  | 0.240% |
| 80  | 17.356 | 2440 | 2443 | 2447 | rBV6 | 41475   | 53328   | 1.55%  | 0.140% |
| 81  | 17.421 | 2449 | 2454 | 2462 | rVV7 | 78300   | 203583  | 5.90%  | 0.535% |
| 82  | 17.503 | 2463 | 2468 | 2474 | rVV  | 529936  | 693466  | 20.11% | 1.824% |
| 83  | 17.580 | 2478 | 2481 | 2490 | rVB8 | 60967   | 132256  | 3.84%  | 0.348% |
| 84  | 17.786 | 2512 | 2516 | 2525 | rVB8 | 42797   | 75739   | 2.20%  | 0.199% |
| 85  | 17.874 | 2528 | 2531 | 2534 | rVV  | 47602   | 62330   | 1.81%  | 0.164% |
| 86  | 17.927 | 2534 | 2540 | 2546 | rVB3 | 99737   | 228929  | 6.64%  | 0.602% |
| 87  | 18.009 | 2550 | 2554 | 2558 | rBV7 | 42935   | 63745   | 1.85%  | 0.168% |
| 88  | 18.068 | 2560 | 2564 | 2567 | rBV3 | 80970   | 112115  | 3.25%  | 0.295% |
| 89  | 18.192 | 2581 | 2585 | 2594 | rVB  | 422774  | 529704  | 15.36% | 1.393% |
| 90  | 18.386 | 2614 | 2618 | 2623 | rBV6 | 35637   | 59737   | 1.73%  | 0.157% |
| 91  | 18.539 | 2642 | 2644 | 2651 | rVB6 | 42996   | 63399   | 1.84%  | 0.167% |
| 92  | 18.833 | 2690 | 2694 | 2699 | rVB  | 373858  | 437387  | 12.68% | 1.150% |
| 93  | 19.062 | 2728 | 2733 | 2742 | rVB  | 338341  | 514230  | 14.91% | 1.352% |
| 94  | 19.421 | 2786 | 2794 | 2804 | rBV3 | 466476  | 952775  | 27.63% | 2.506% |
| 95  | 19.956 | 2881 | 2885 | 2889 | rVB2 | 286278  | 341692  | 9.91%  | 0.899% |
| 96  | 20.450 | 2966 | 2969 | 2973 | rBV  | 214754  | 239912  | 6.96%  | 0.631% |
| 97  | 20.915 | 3045 | 3048 | 3055 | rBV3 | 180084  | 228158  | 6.62%  | 0.600% |
| 98  | 21.197 | 3091 | 3096 | 3101 | rBV  | 1394964 | 1840921 | 53.38% | 4.841% |
| 99  | 21.356 | 3120 | 3123 | 3127 | rVB2 | 167314  | 187077  | 5.42%  | 0.492% |
| 100 | 23.391 | 3463 | 3469 | 3476 | rVB  | 856090  | 1607927 | 46.62% | 4.229% |

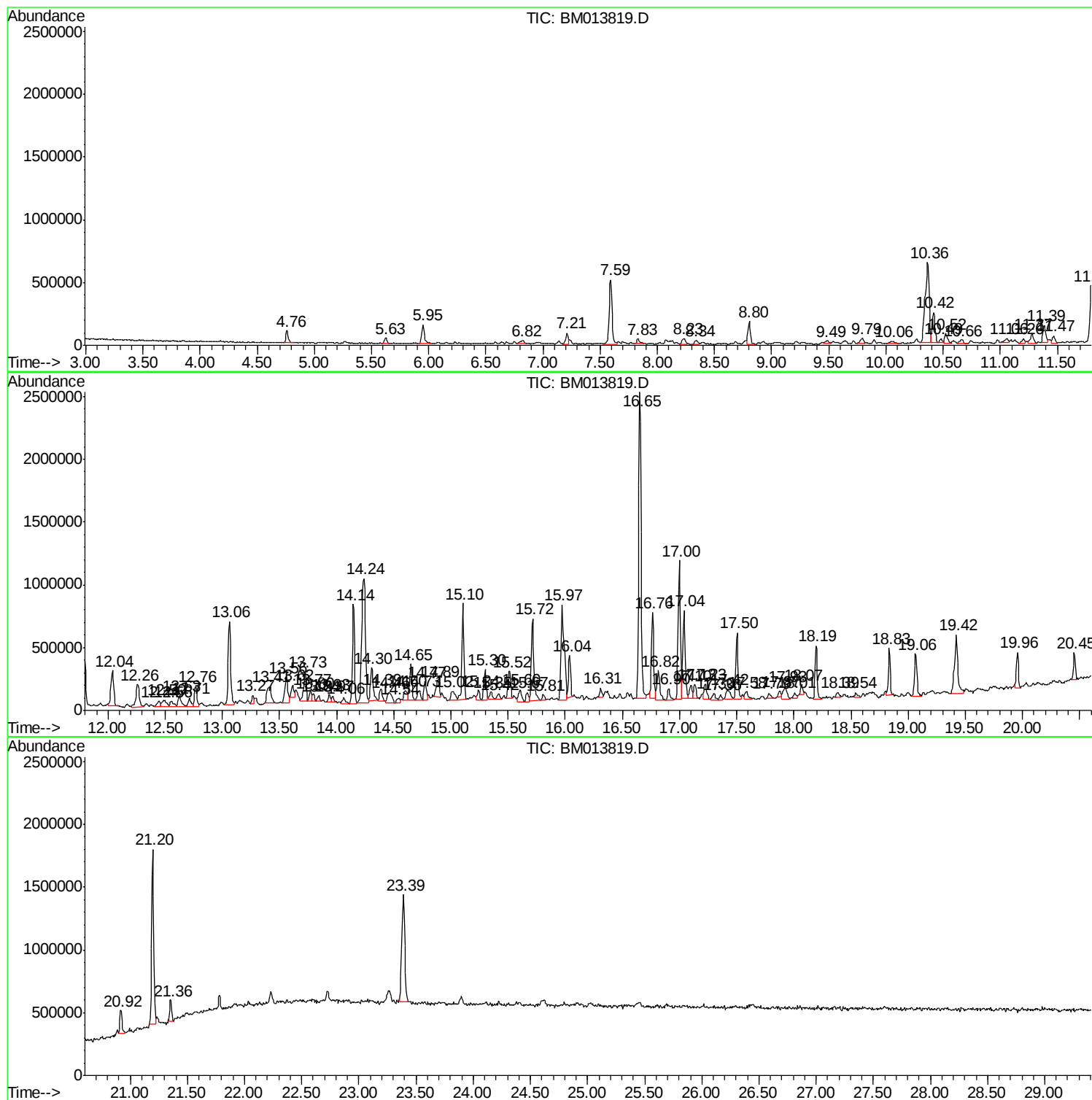
Sum of corrected areas: 38025268

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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

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



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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

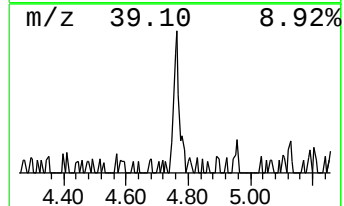
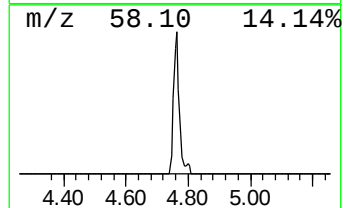
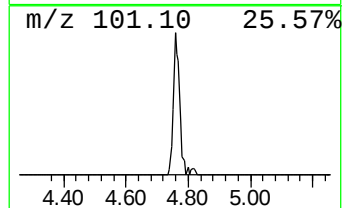
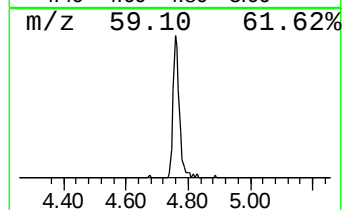
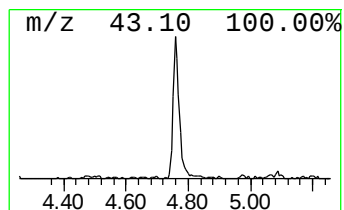
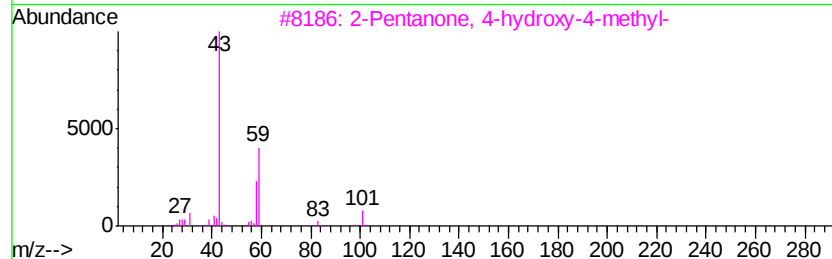
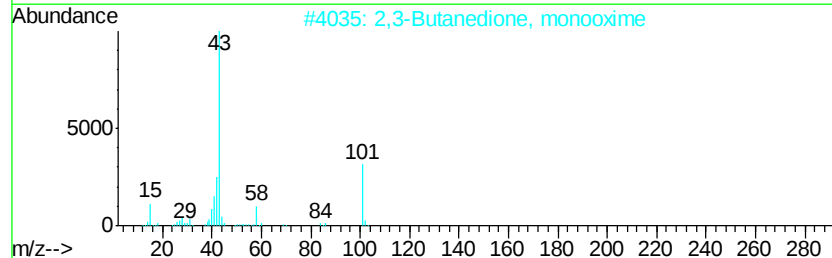
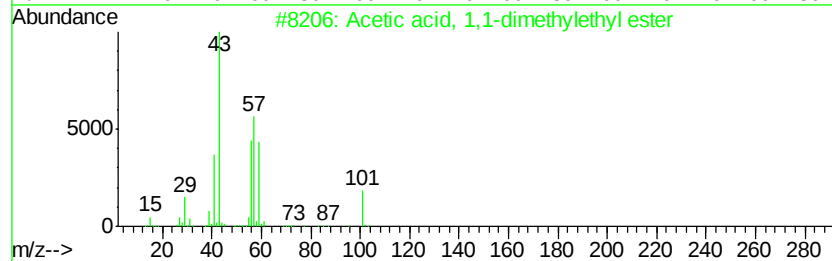
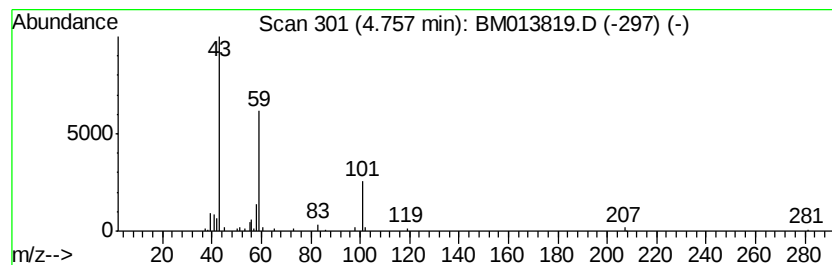
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 unknown-01 Concentration Rank 16

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 4.76 | 3.27 ng/ul | 143644 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2 | 000540-88-5  | 38   |
| 2    |    | 2,3-Butanedione, monooxime          | 101 | C4H7NO2 | 000057-71-6  | 38   |
| 3    |    | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2 | 000123-42-2  | 38   |
| 4    |    | (+)-4-Amino-4,5-dihydro-2(3H)-f...  | 101 | C4H7NO2 | 016504-58-8  | 37   |
| 5    |    | n-Propyl t-butyl ether              | 116 | C7H16O  | 1000334-81-9 | 33   |



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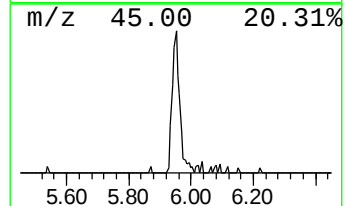
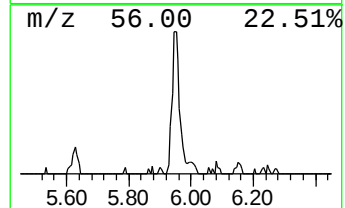
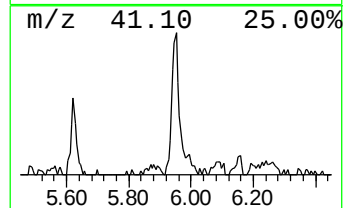
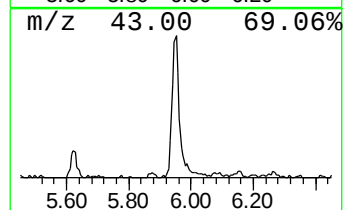
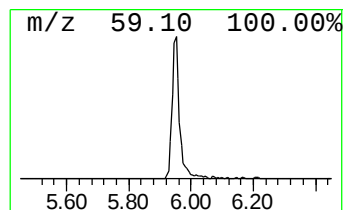
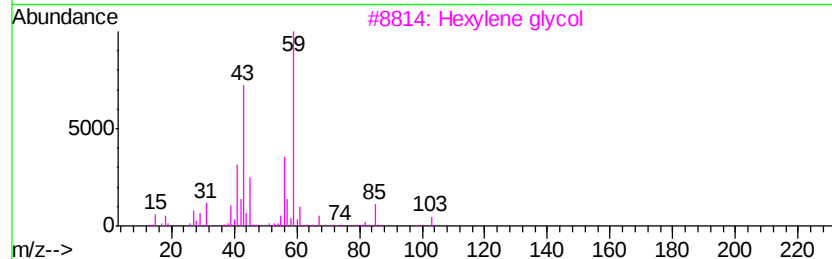
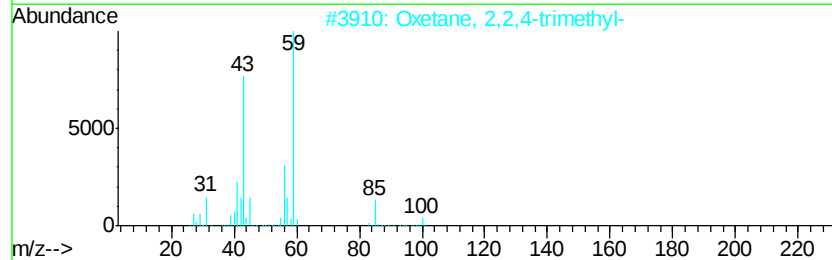
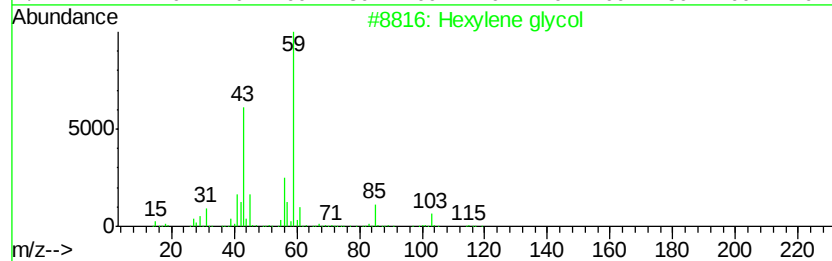
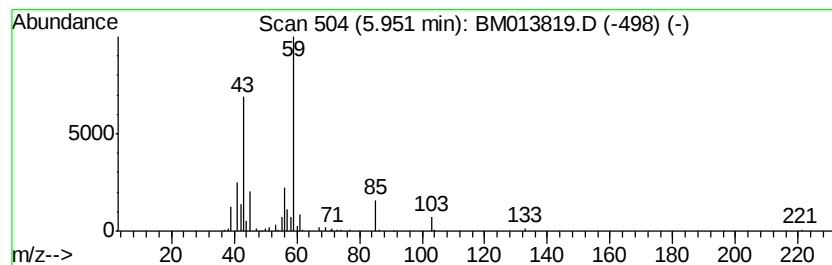
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Hexylene glycol Concentration Rank 9

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 5.95 | 6.12 ng/ul | 269222 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexylene glycol           | 118 | C6H14O2 | 000107-41-5 | 78   |
| 2    |    | Oxetane, 2,2,4-trimethyl- | 100 | C6H12O  | 023120-44-7 | 72   |
| 3    |    | Hexylene glycol           | 118 | C6H14O2 | 000107-41-5 | 56   |
| 4    |    | Hexylene glycol           | 118 | C6H14O2 | 000107-41-5 | 53   |
| 5    |    | Hexylene glycol           | 118 | C6H14O2 | 000107-41-5 | 50   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

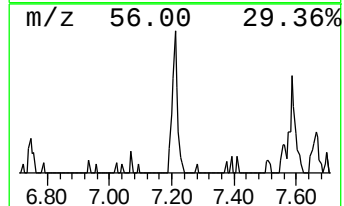
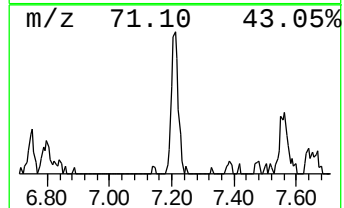
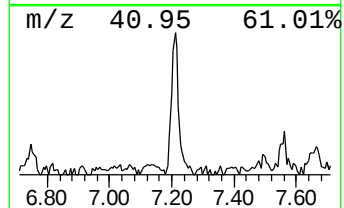
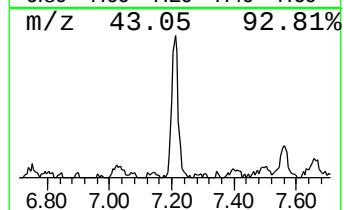
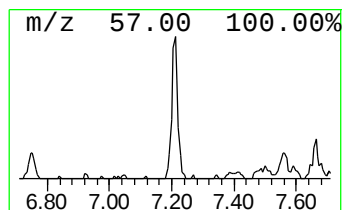
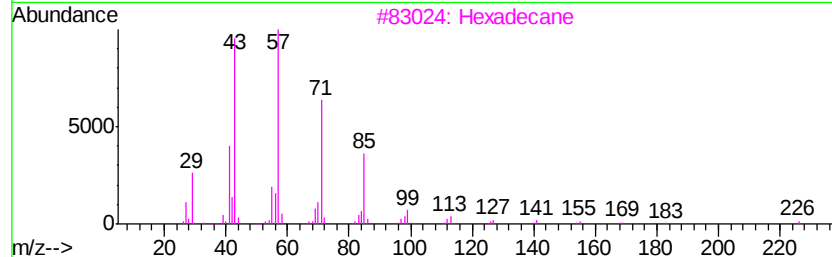
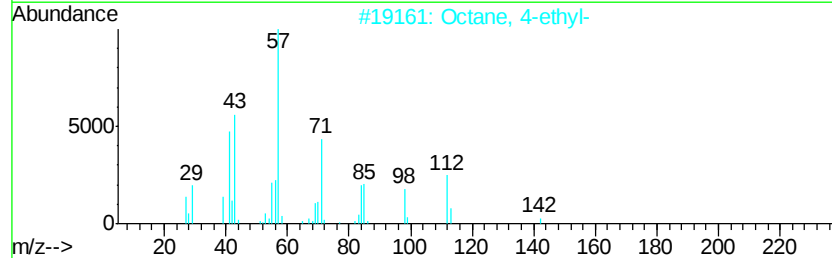
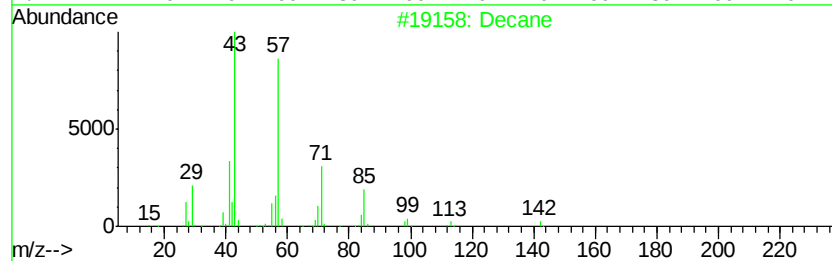
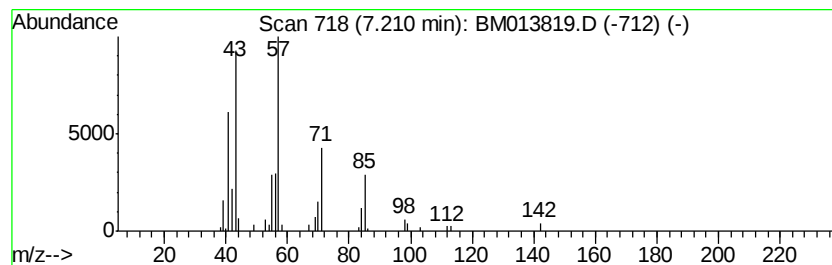
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 19

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.21 | 3.12 ng/ul | 137319 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# | of                | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|-------------------|---|--------------|-----|---------|-------------|------|
| 1    | Decane            |   |              | 142 | C10H22  | 000124-18-5 | 72   |
| 2    | Octane, 4-ethyl-  |   |              | 142 | C10H22  | 015869-86-0 | 72   |
| 3    | Hexadecane        |   |              | 226 | C16H34  | 000544-76-3 | 64   |
| 4    | Undecane          |   |              | 156 | C11H24  | 001120-21-4 | 53   |
| 5    | Decane, 2-methyl- |   |              | 156 | C11H24  | 006975-98-0 | 53   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleID :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

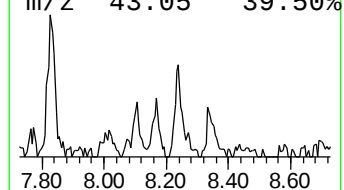
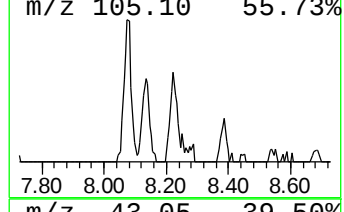
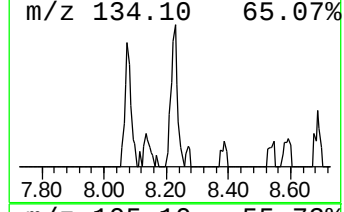
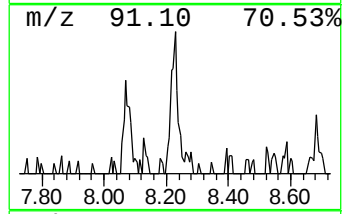
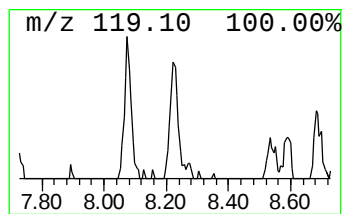
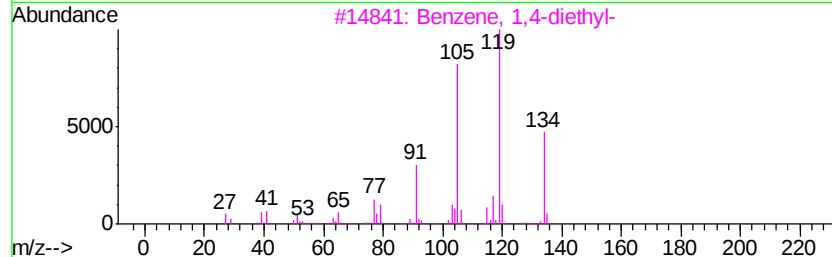
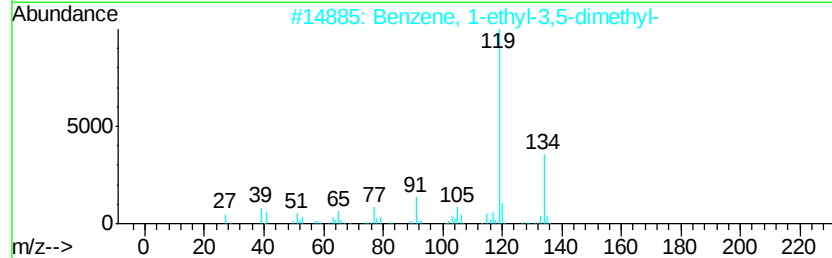
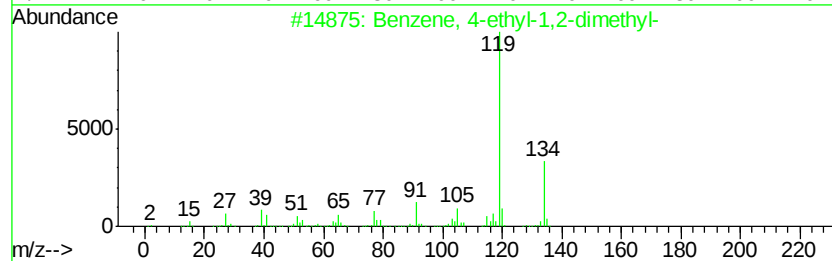
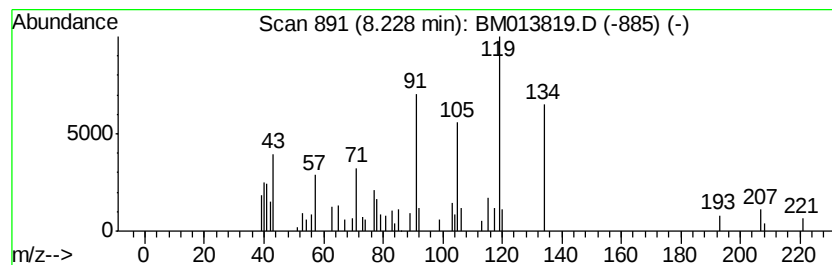
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 unknown-02 Concentration Rank 27

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 8.23 | 2.17 ng/ul | 95193 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 38   |
| 2    |    | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 35   |
| 3    |    | Benzene, 1,4-diethyl-          | 134 | C10H14  | 000105-05-5 | 30   |
| 4    |    | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 30   |
| 5    |    | Benzene, 1,2-diethyl-          | 134 | C10H14  | 000135-01-3 | 30   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

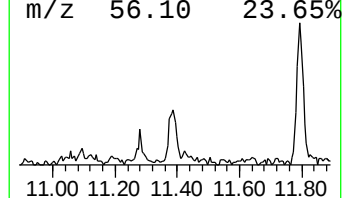
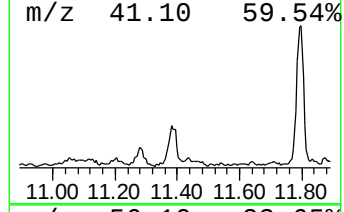
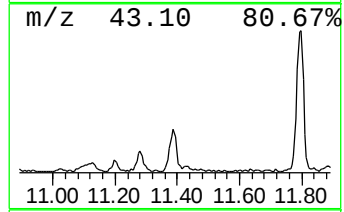
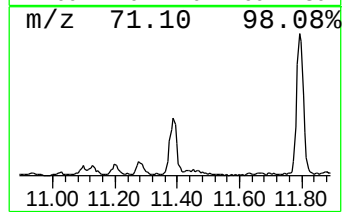
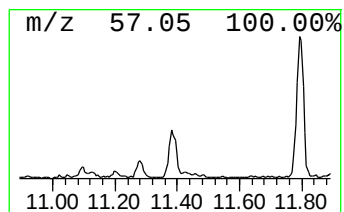
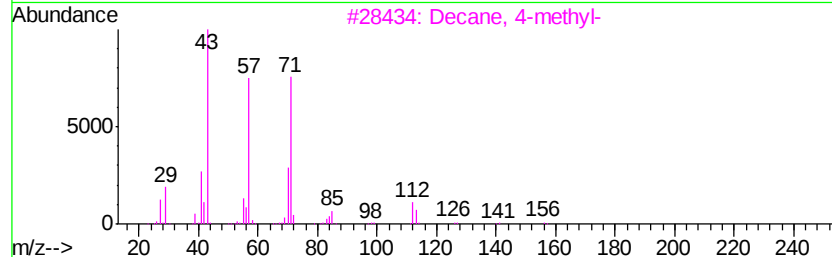
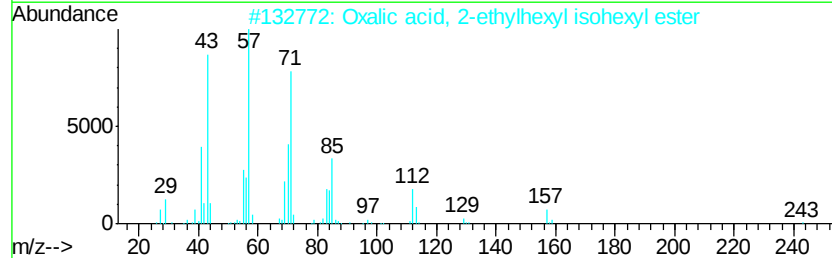
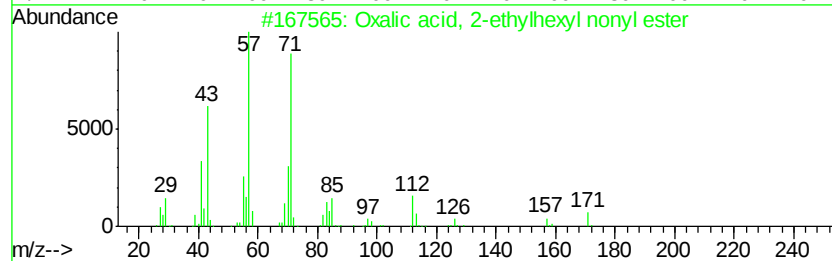
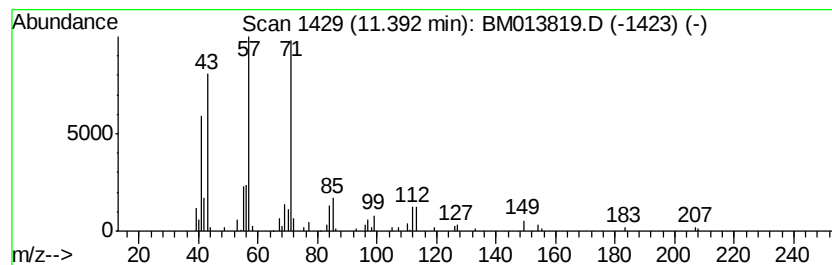
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Oxalic acid, 2-ethylhexyl n... Concentration Rank 20

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.39 | 3.02 ng/ul | 233686 | Naphthalene-d8   | 10.36 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Oxalic acid, 2-ethylhexyl nonyl ... | 328 | C19H36O4  | 1000309-39-2 | 72   |
| 2    |    |   | Oxalic acid, 2-ethylhexyl isohex... | 286 | C16H30O4  | 1000309-38-8 | 72   |
| 3    |    |   | Decane, 4-methyl-                   | 156 | C11H24    | 002847-72-5  | 64   |
| 4    |    |   | Sulfurous acid, 2-ethylhexyl tet... | 390 | C22H46O3S | 1000309-19-7 | 64   |
| 5    |    |   | Sulfurous acid, 2-ethylhexyl hex... | 418 | C24H50O3S | 1000309-19-9 | 64   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

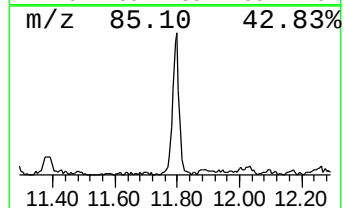
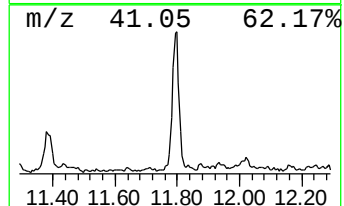
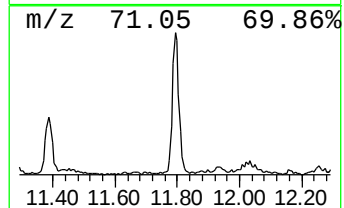
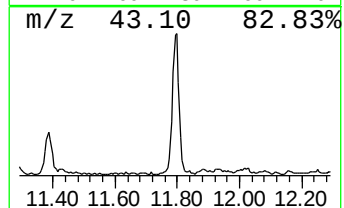
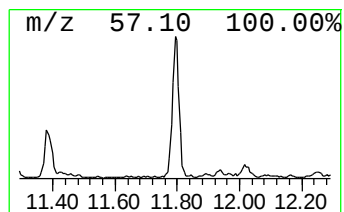
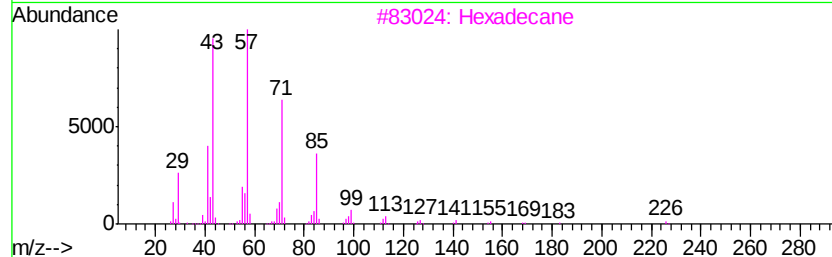
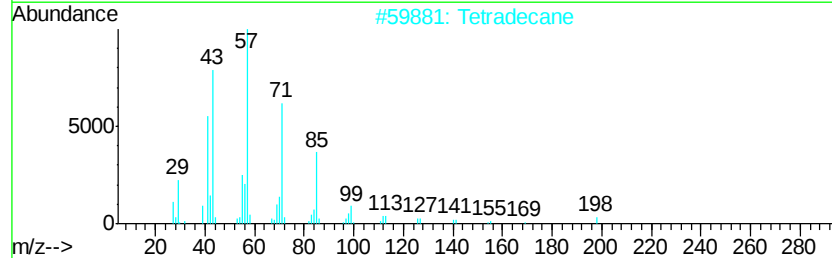
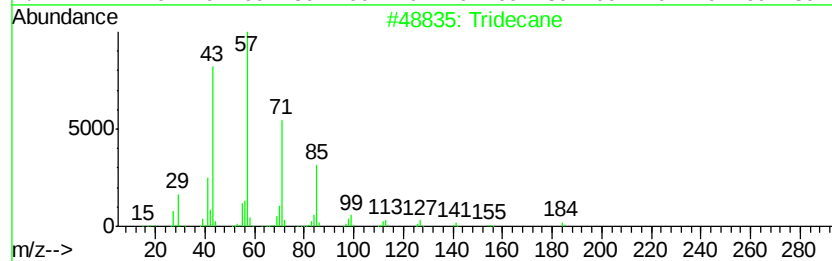
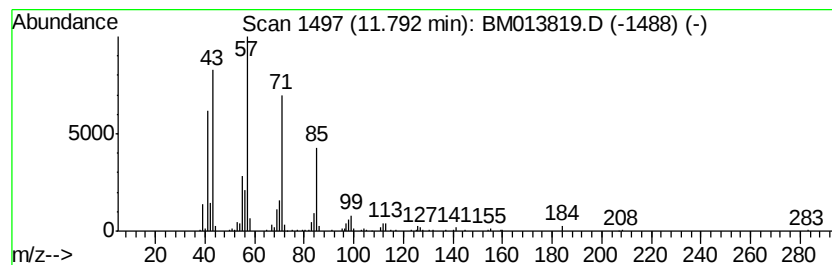
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 5

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.79 | 8.93 ng/ul | 690949 | Napthalene-d8    | 10.36 |

| Hit# of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|-----------|--------------|-----|---------|-------------|------|
| 1         | Tridecane    | 184 | C13H28  | 000629-50-5 | 93   |
| 2         | Tetradecane  | 198 | C14H30  | 000629-59-4 | 91   |
| 3         | Hexadecane   | 226 | C16H34  | 000544-76-3 | 90   |
| 4         | Dodecane     | 170 | C12H26  | 000112-40-3 | 90   |
| 5         | Hexadecane   | 226 | C16H34  | 000544-76-3 | 90   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

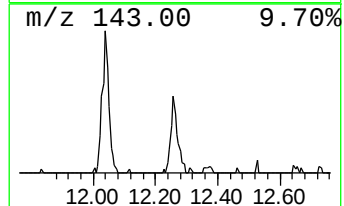
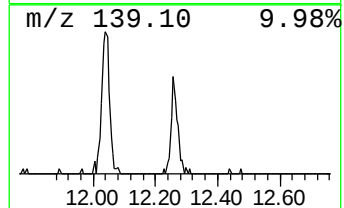
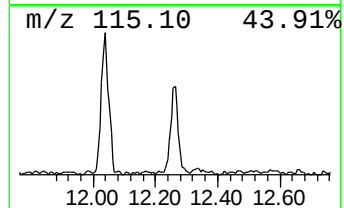
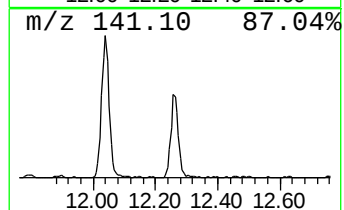
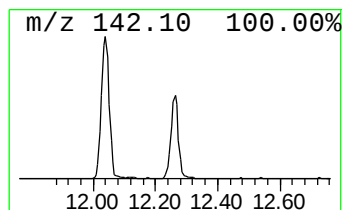
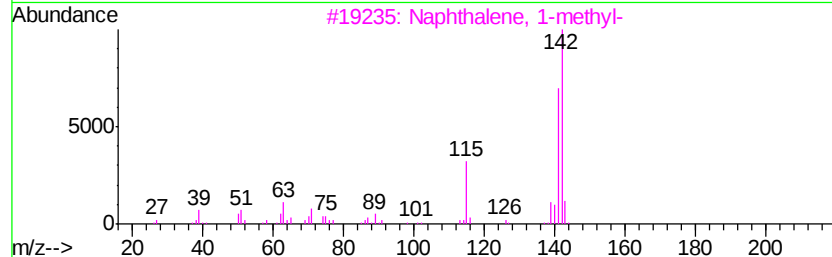
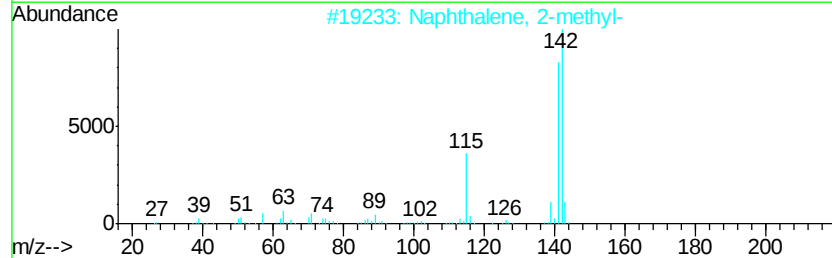
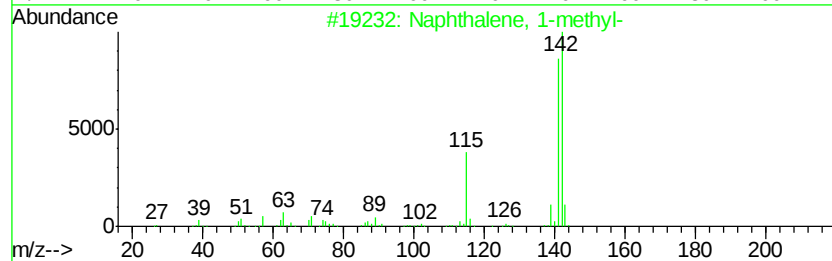
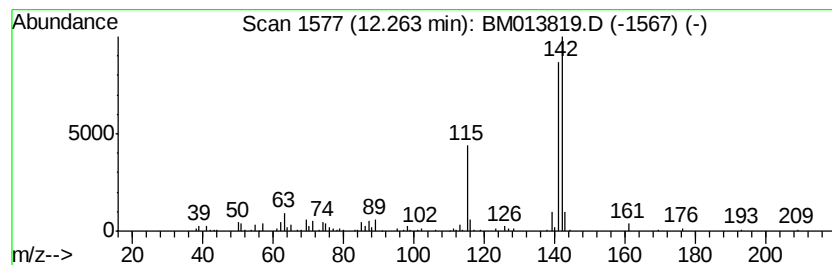
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Naphthalene, 1-methyl- Concentration Rank 13

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.26 | 4.72 ng/ul | 365074 | Naphthalene-d8   | 10.36 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 97   |
| 2    |    | Naphthalene, 2-methyl-              | 142 | C11H10  | 000091-57-6 | 95   |
| 3    |    | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 95   |
| 4    |    | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 94   |
| 5    |    | 1,4-Methanonaphthalene, 1,4-dihy... | 142 | C11H10  | 004453-90-1 | 94   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

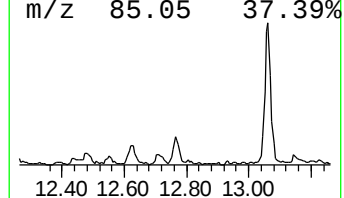
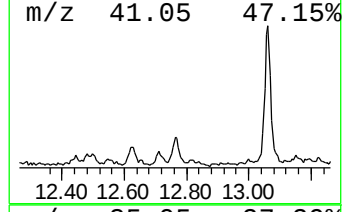
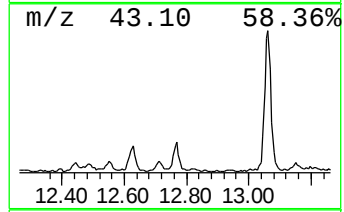
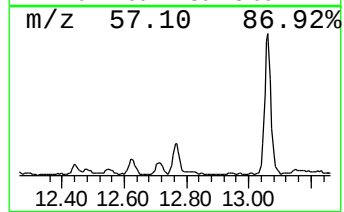
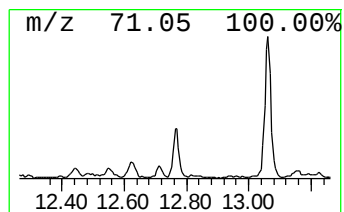
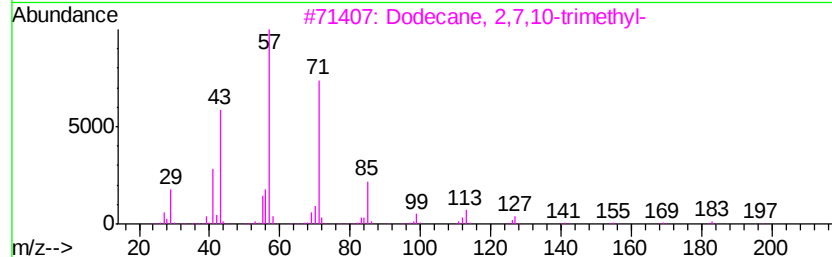
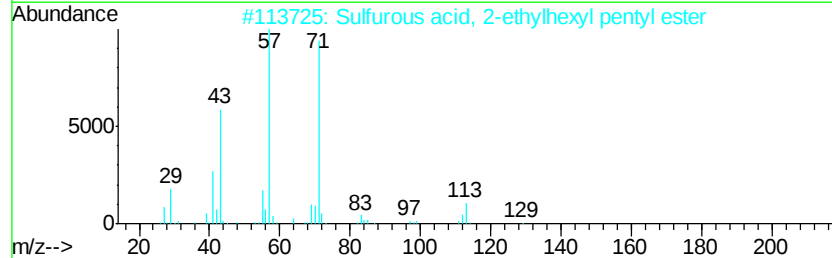
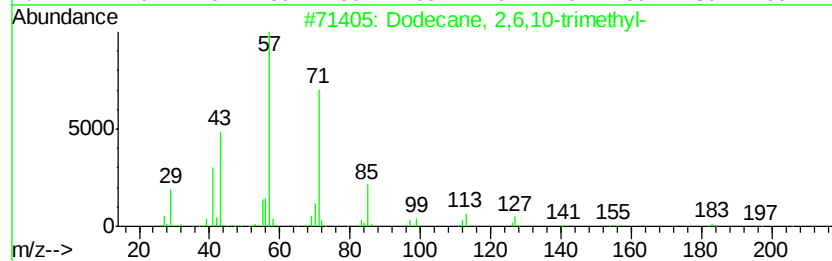
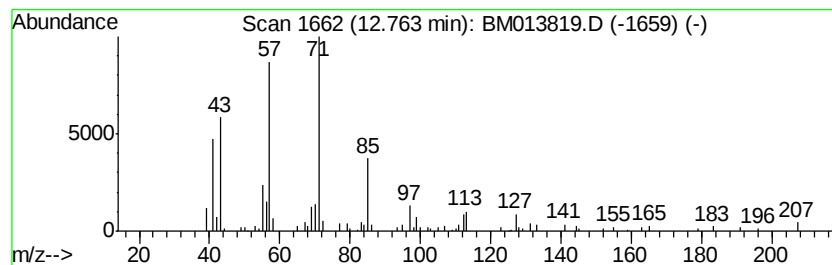
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 30

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.76 | 2.06 ng/ul | 234738 | Acenaphthene-d10 | 14.24 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Dodecane, 2,6,10-trimethyl-         | 212 | C15H32    | 003891-98-3  | 64   |
| 2    |    | Sulfurous acid, 2-ethylhexyl pen... | 264 | C13H28O3S | 1000309-18-9 | 53   |
| 3    |    | Dodecane, 2,7,10-trimethyl-         | 212 | C15H32    | 074645-98-0  | 52   |
| 4    |    | l-Threitol, 2-O-nonyl-              | 248 | C13H28O4  | 163776-15-6  | 50   |
| 5    |    | Sulfurous acid, nonyl pentyl ester  | 278 | C14H30O3S | 1000309-14-2 | 50   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

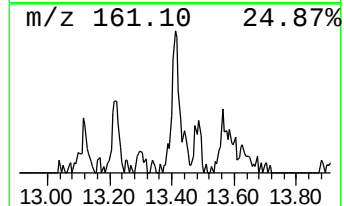
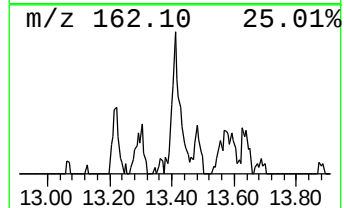
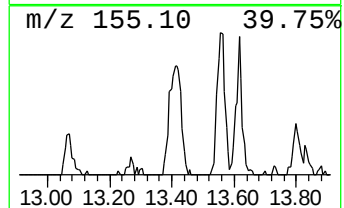
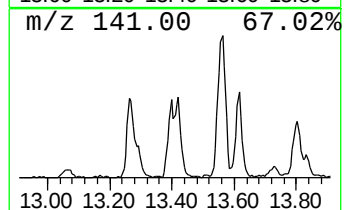
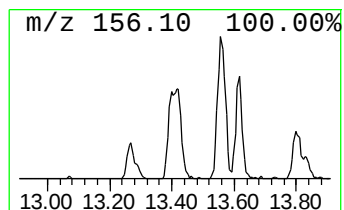
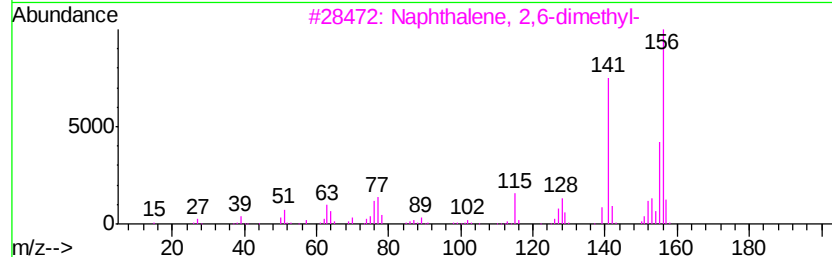
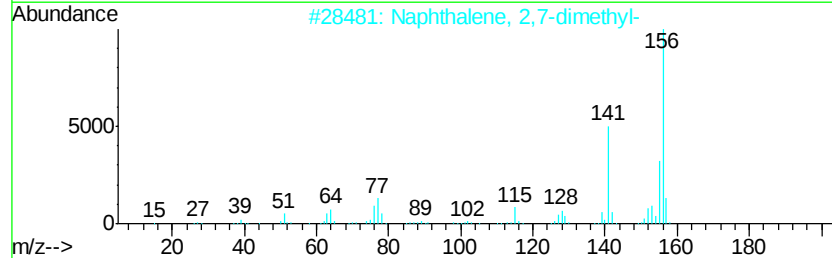
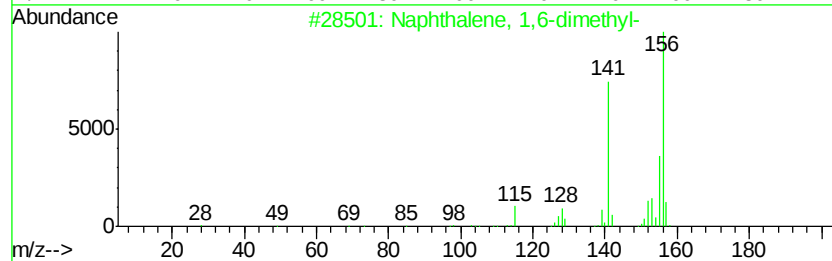
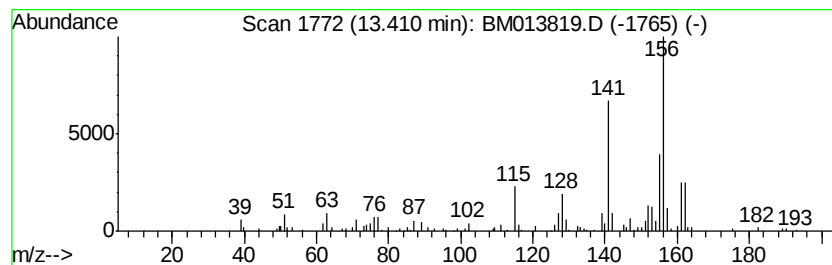
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Naphthalene, 1,6-dimethyl- Concentration Rank 17

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.41 | 3.24 ng/ul | 368424 | Acenaphthene-d10 | 14.24 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 95   |
| 2    |    | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 94   |
| 3    |    | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 93   |
| 4    |    | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 92   |
| 5    |    | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 90   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

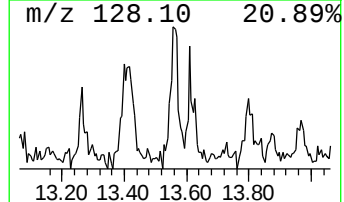
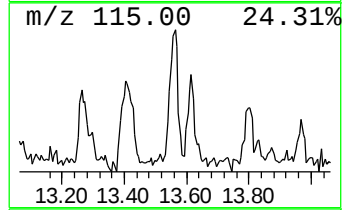
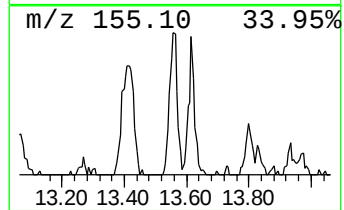
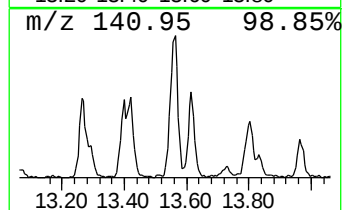
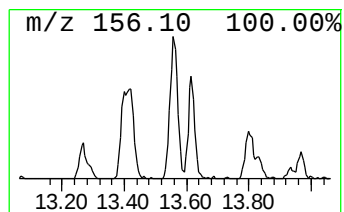
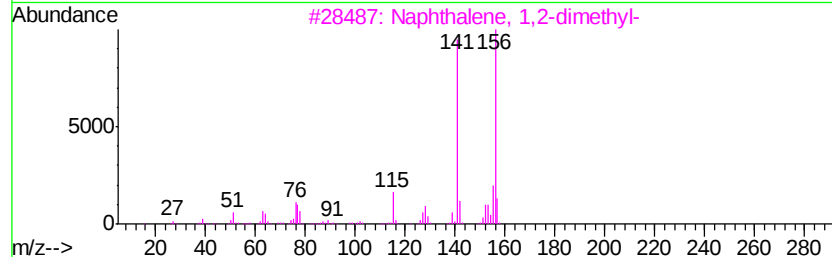
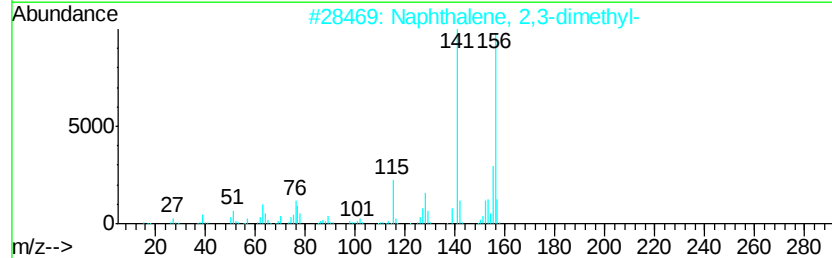
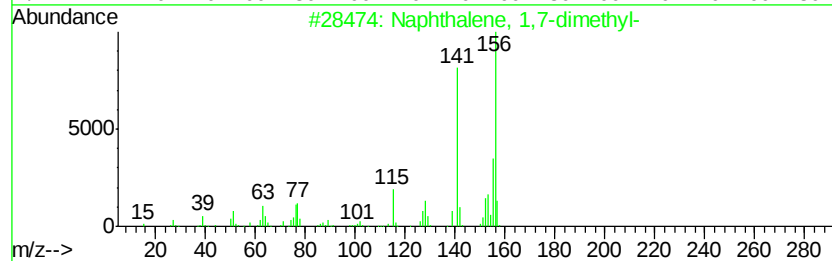
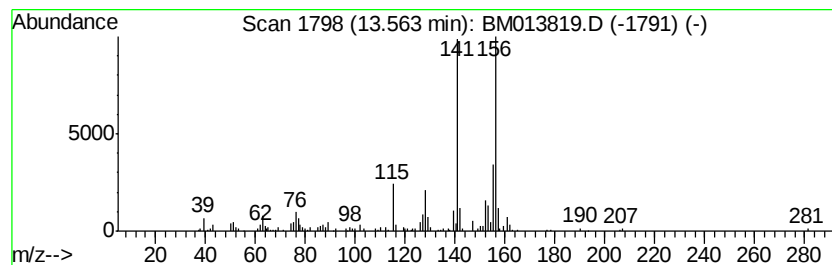
Instrument :  
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ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 10 Naphthalene, 1,7-dimethyl- Concentration Rank 15

| R.T.    | EstConc    | Area                       | Relative to ISTD | R.T.           |
|---------|------------|----------------------------|------------------|----------------|
| 13.56   | 3.55 ng/ul | 403850                     | Acenaphthene-d10 | 14.24          |
| Hit# of | 5          | Tentative ID               | MW MolForm       | CAS# Qual      |
| 1       |            | Naphthalene, 1,7-dimethyl- | 156 C12H12       | 000575-37-1 96 |
| 2       |            | Naphthalene, 2,3-dimethyl- | 156 C12H12       | 000581-40-8 95 |
| 3       |            | Naphthalene, 1,2-dimethyl- | 156 C12H12       | 000573-98-8 94 |
| 4       |            | Naphthalene, 1,8-dimethyl- | 156 C12H12       | 000569-41-5 94 |
| 5       |            | Naphthalene, 1,3-dimethyl- | 156 C12H12       | 000575-41-7 94 |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

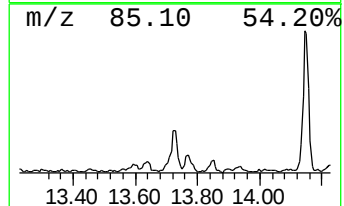
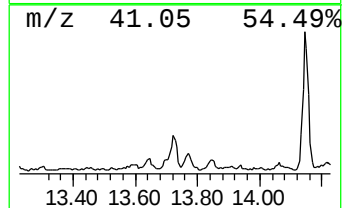
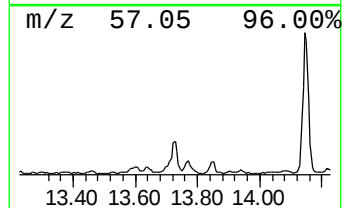
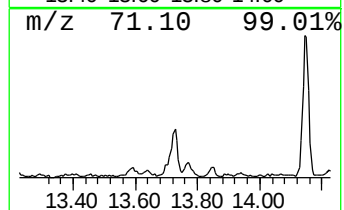
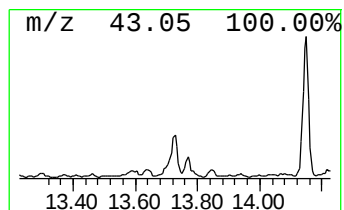
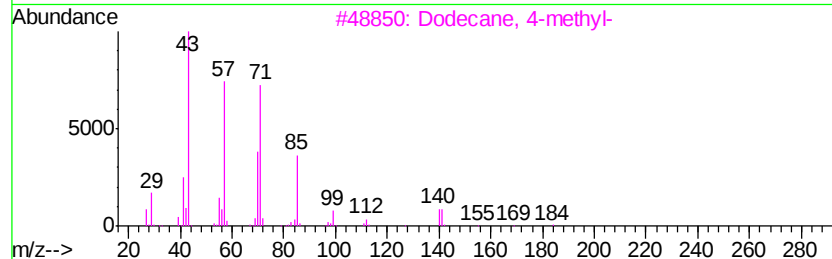
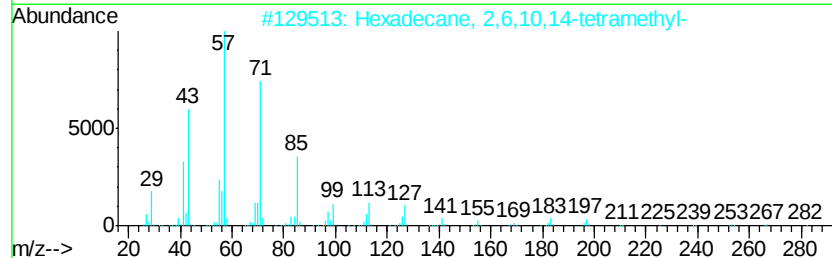
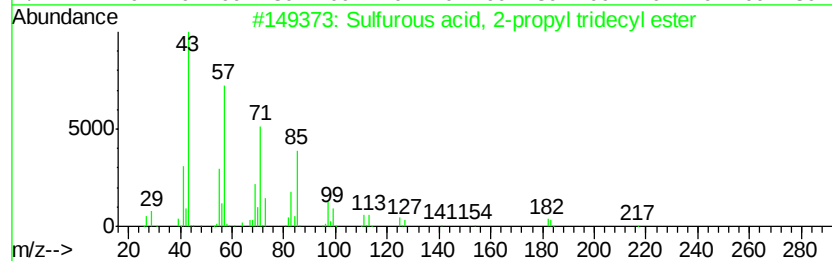
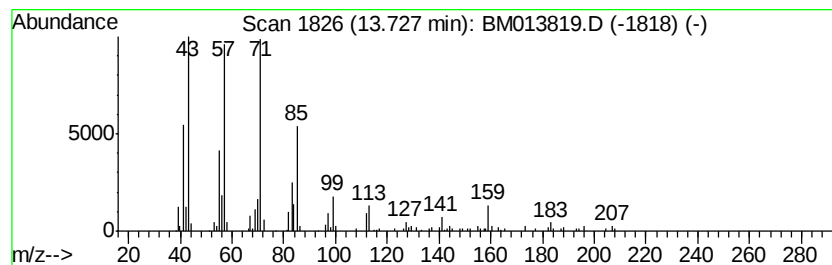
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Sulfurous acid, 2-propyl tr... Concentration Rank 18

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.73 | 3.14 ng/ul | 357539 | Acenaphthene-d10 | 14.24 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Sulfurous acid, 2-propyl tridecy... | 306 | C16H34O3S | 1000309-12-4 | 80   |
| 2    |    |   | Hexadecane, 2,6,10,14-tetramethyl-  | 282 | C20H42    | 000638-36-8  | 64   |
| 3    |    |   | Dodecane, 4-methyl-                 | 184 | C13H28    | 006117-97-1  | 64   |
| 4    |    |   | Hexadecane, 7,9-dimethyl-           | 254 | C18H38    | 021164-95-4  | 64   |
| 5    |    |   | Dodecane, 4-methyl-                 | 184 | C13H28    | 006117-97-1  | 58   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

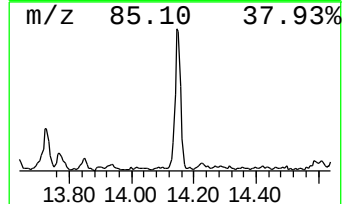
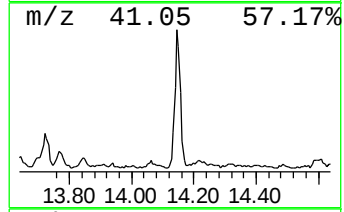
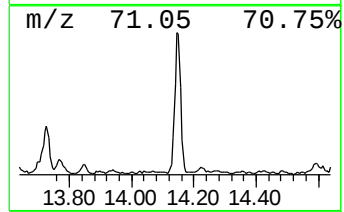
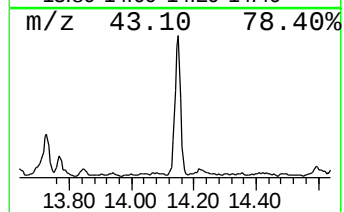
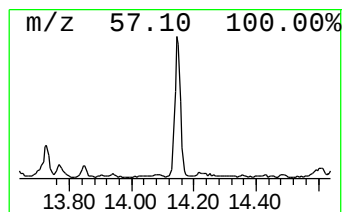
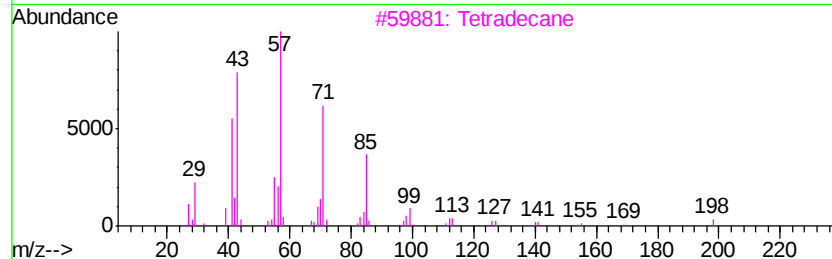
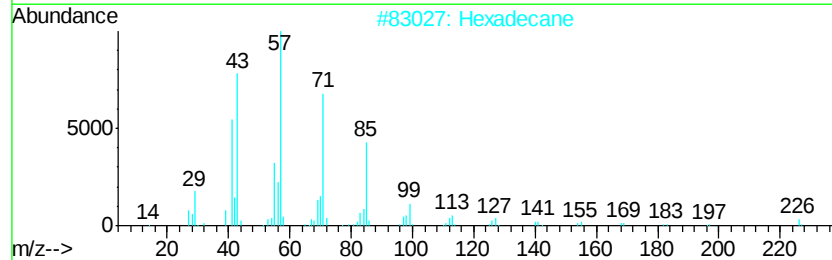
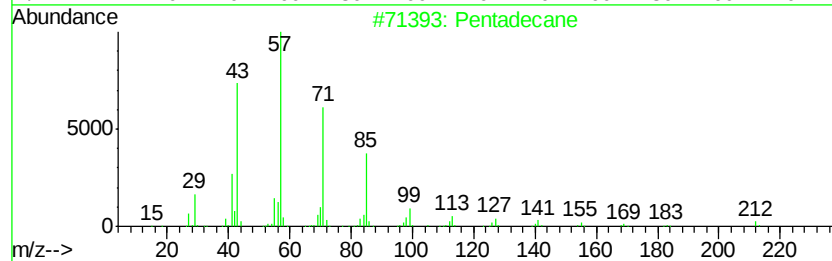
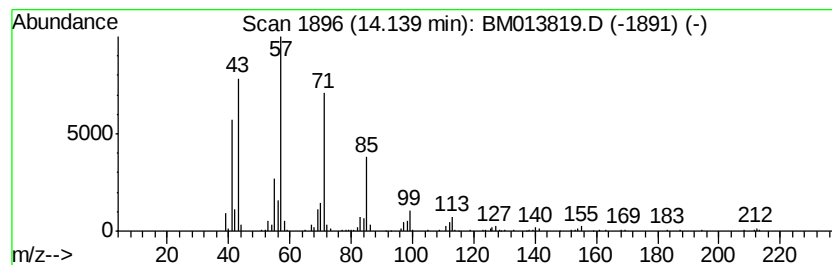
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 14.14 | 9.25 ng/ul | 1052050 | Acenaphthene-d10 | 14.24 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Pentadecane  | 212 | C15H32  | 000629-62-9 | 93   |
| 2    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 91   |
| 3    |    |   | Tetradecane  | 198 | C14H30  | 000629-59-4 | 91   |
| 4    |    |   | Nonadecane   | 268 | C19H40  | 000629-92-5 | 91   |
| 5    |    |   | Nonadecane   | 268 | C19H40  | 000629-92-5 | 90   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

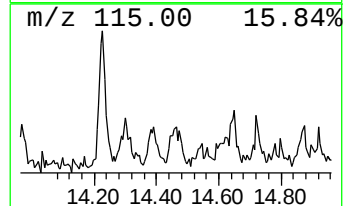
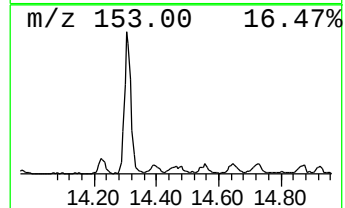
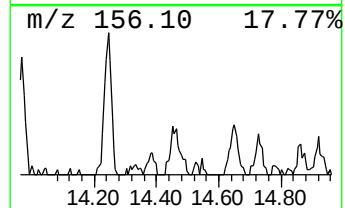
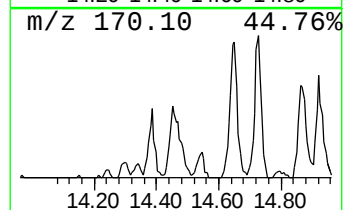
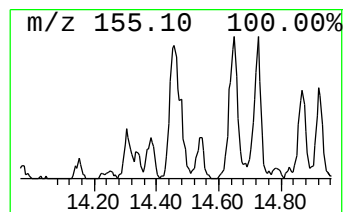
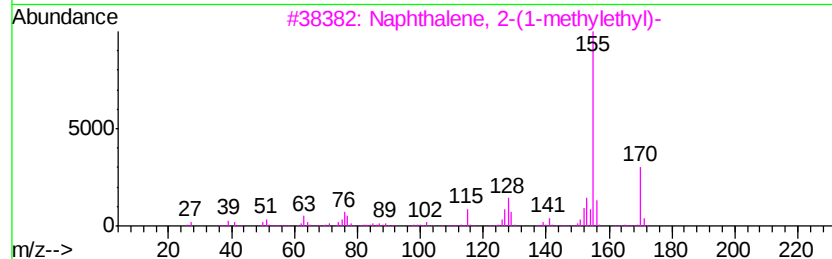
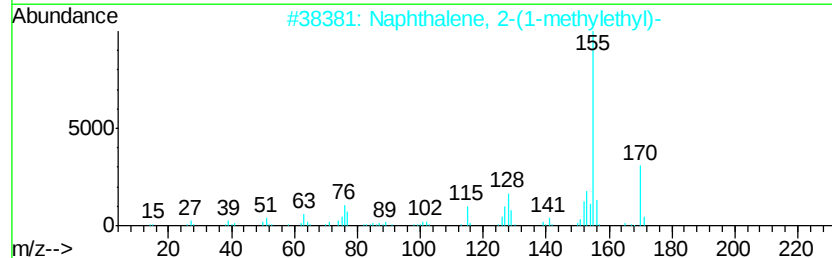
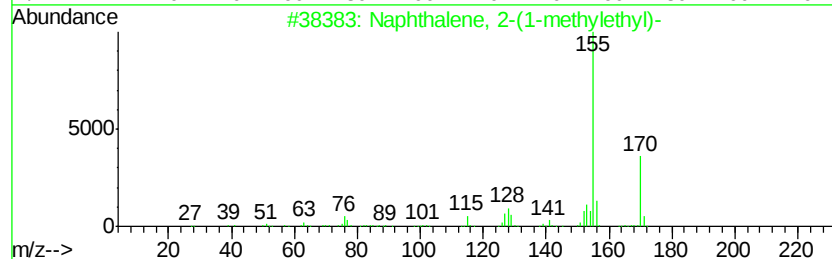
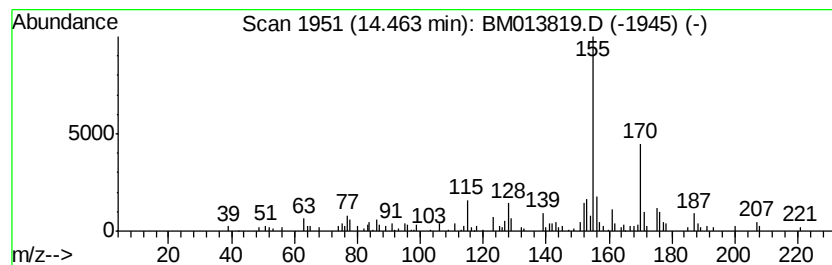
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 Naphthalene, 2-(1-methyleth... Concentration Rank 28

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.46 | 2.10 ng/ul | 239006 | Acenaphthene-d10 | 14.24 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2-(1-methylethyl)- | 170 | C13H14  | 002027-17-0 | 87   |
| 2    |    |   | Naphthalene, 2-(1-methylethyl)- | 170 | C13H14  | 002027-17-0 | 80   |
| 3    |    |   | Naphthalene, 2-(1-methylethyl)- | 170 | C13H14  | 002027-17-0 | 74   |
| 4    |    |   | Naphthalene, 1-(1-methylethyl)- | 170 | C13H14  | 006158-45-8 | 72   |
| 5    |    |   | Naphthalene, 1,4,5-trimethyl-   | 170 | C13H14  | 002131-41-1 | 72   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

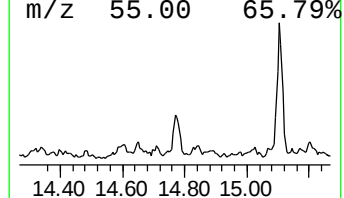
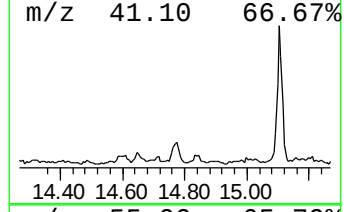
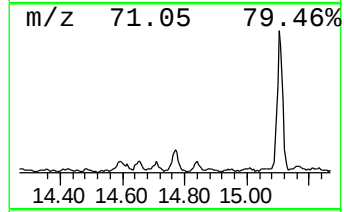
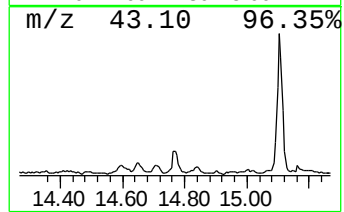
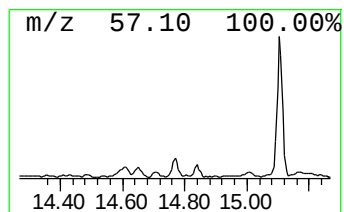
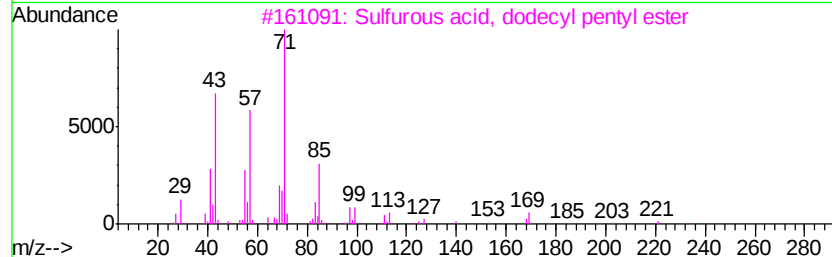
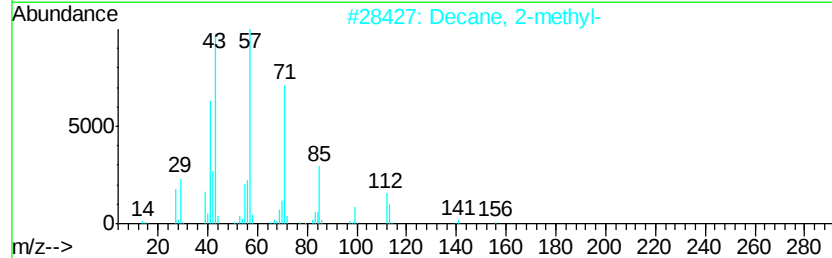
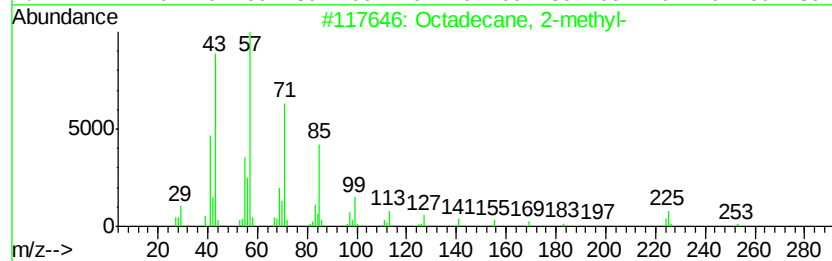
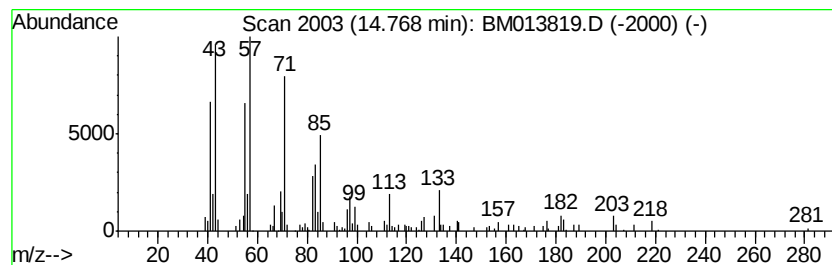
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 32

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.77 | 2.03 ng/ul | 230303 | Acenaphthene-d10 | 14.24 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Octadecane, 2-methyl-               | 268 | C19H40    | 001560-88-9  | 41   |
| 2    |    |   | Decane, 2-methyl-                   | 156 | C11H24    | 006975-98-0  | 38   |
| 3    |    |   | Sulfurous acid, dodecyl pentyl e... | 320 | C17H36O3S | 1000309-14-5 | 35   |
| 4    |    |   | Sulfurous acid, decyl pentyl ester  | 292 | C15H32O3S | 1000309-14-3 | 35   |
| 5    |    |   | Tridecane, 1-iodo-                  | 310 | C13H27I   | 035599-77-0  | 35   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

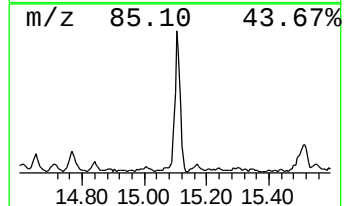
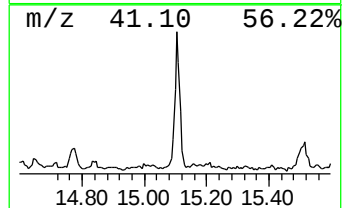
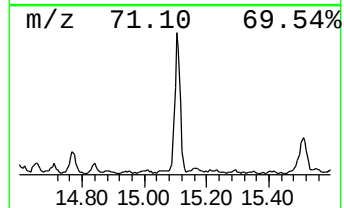
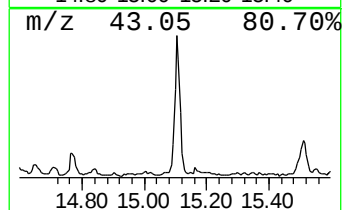
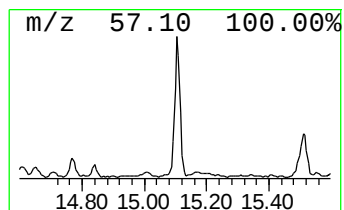
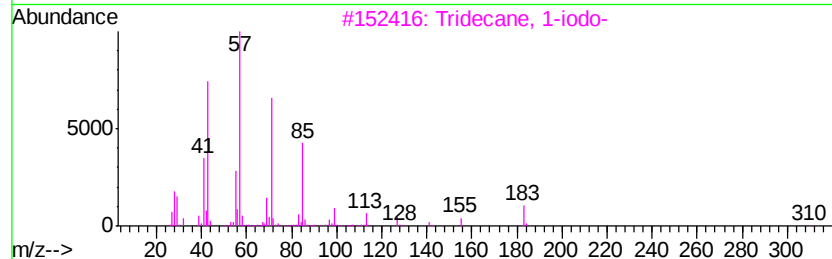
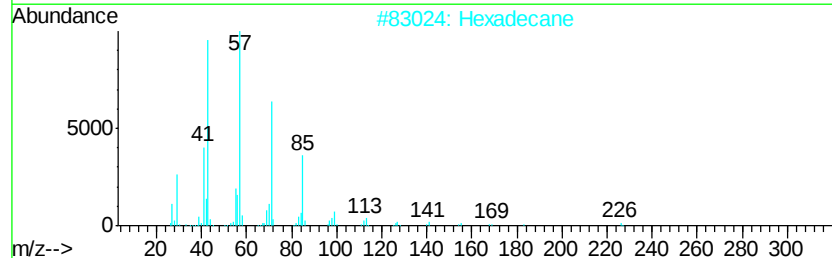
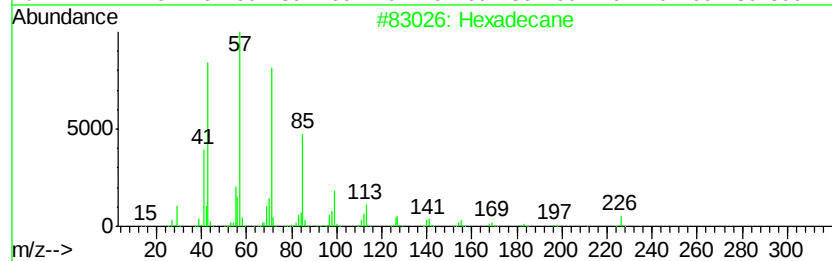
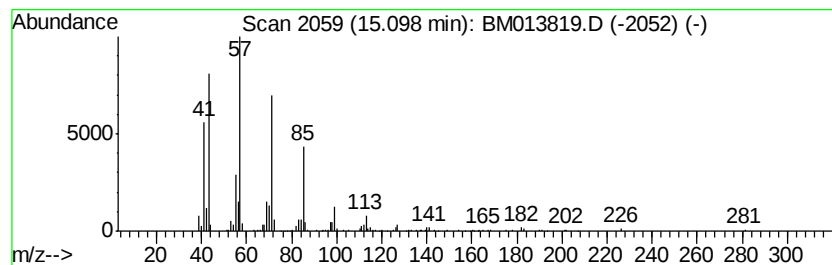
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.10 | 8.56 ng/ul | 972869 | Acenaphthene-d10 | 14.24 |

| Hit# | of | 5 | Tentative ID       | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane         | 226 | C16H34  | 000544-76-3 | 94   |
| 2    |    |   | Hexadecane         | 226 | C16H34  | 000544-76-3 | 94   |
| 3    |    |   | Tridecane, 1-iodo- | 310 | C13H27I | 035599-77-0 | 91   |
| 4    |    |   | Hexadecane         | 226 | C16H34  | 000544-76-3 | 91   |
| 5    |    |   | Pentadecane        | 212 | C15H32  | 000629-62-9 | 90   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

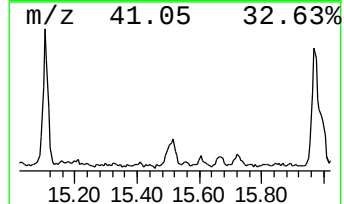
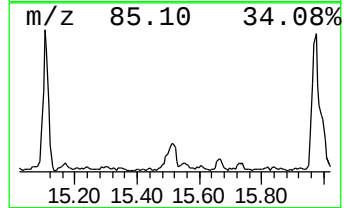
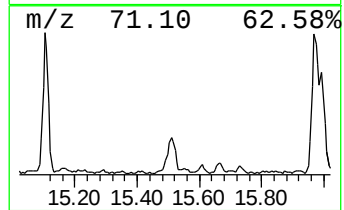
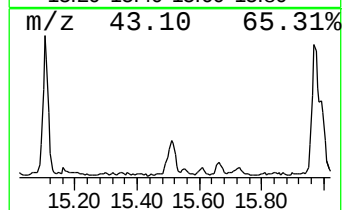
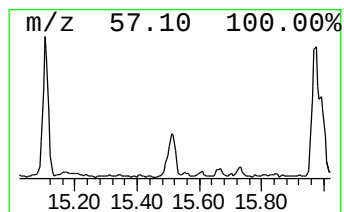
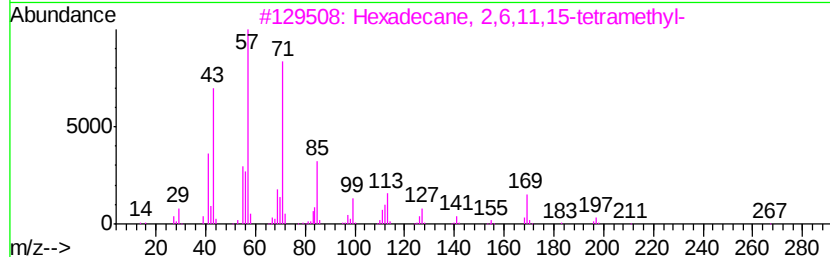
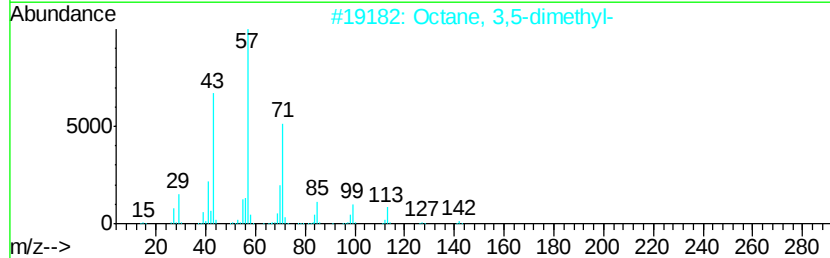
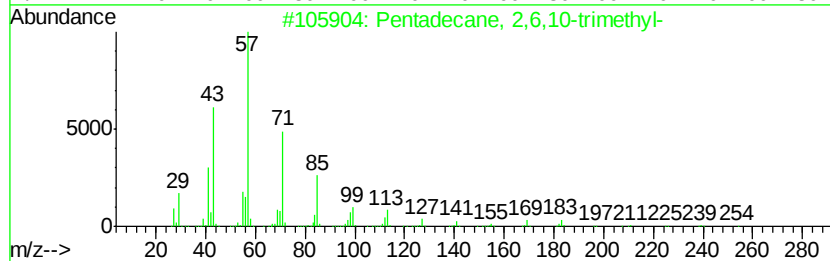
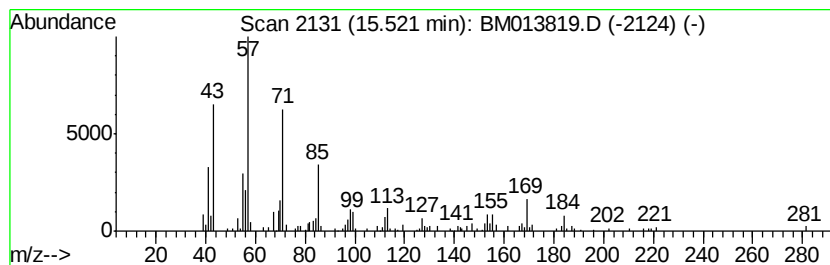
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 21

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.52 | 2.93 ng/ul | 332630 | Acenaphthene-d10 | 14.24 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm     | CAS#        | Qual |
|------|----|---|------------------------------------|-----|-------------|-------------|------|
| 1    |    |   | Pentadecane, 2,6,10-trimethyl-     | 254 | C18H38      | 003892-00-0 | 72   |
| 2    |    |   | Octane, 3,5-dimethyl-              | 142 | C10H22      | 015869-93-9 | 70   |
| 3    |    |   | Hexadecane, 2,6,11,15-tetramethyl- | 282 | C20H42      | 000504-44-9 | 68   |
| 4    |    |   | Silane, trichlorooctadecyl-        | 386 | C18H37Cl3Si | 000112-04-9 | 64   |
| 5    |    |   | Decane, 2-methyl-                  | 156 | C11H24      | 006975-98-0 | 62   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

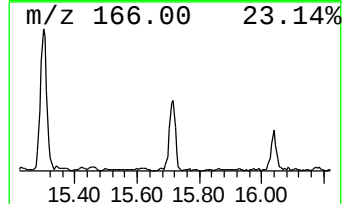
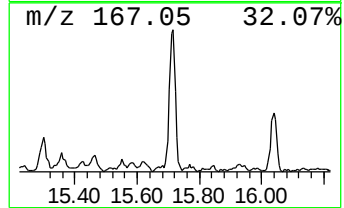
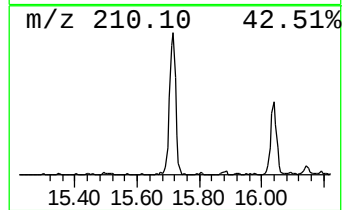
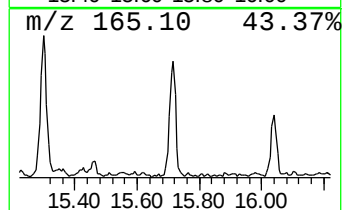
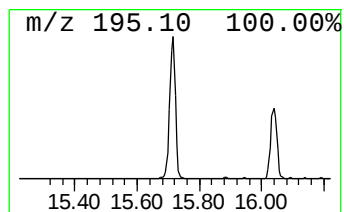
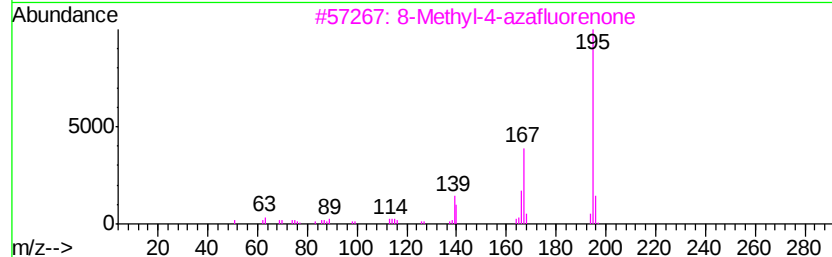
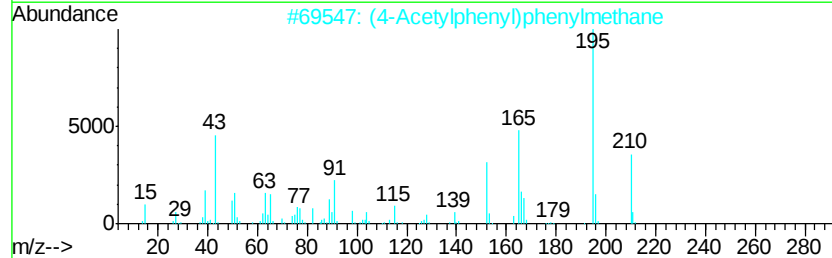
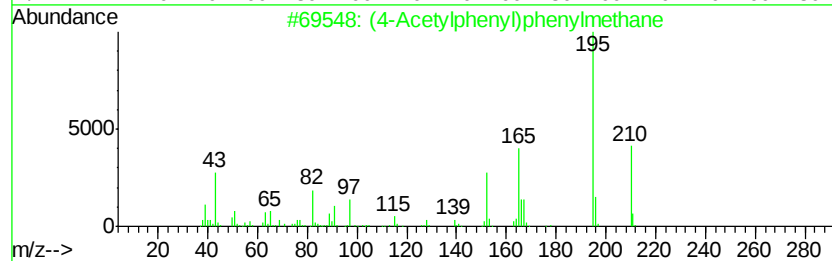
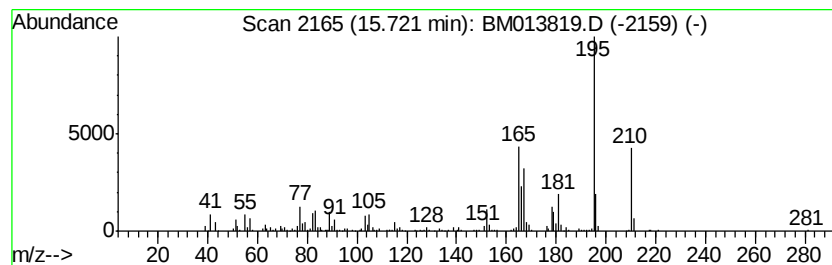
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 (4-Acetylphenyl)phenylmethane Concentration Rank 2

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 15.72 | 12.53 ng/ul | 993482 | Phenanthrene-d10 | 17.00 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | (4-Acetylphenyl)phenylmethane       | 210 | C15H14O  | 000782-92-3 | 68   |
| 2    |    |   | (4-Acetylphenyl)phenylmethane       | 210 | C15H14O  | 000782-92-3 | 62   |
| 3    |    |   | 8-Methyl-4-azafluorenone            | 195 | C13H9NO  | 064292-03-1 | 52   |
| 4    |    |   | 2',3',4',5',6'-Pentafluoroacetop... | 210 | C8H3F5O  | 000652-29-9 | 52   |
| 5    |    |   | 3H-Pyrrolo[3,2-H]quinoline, 2,3,... | 210 | C14H14N2 | 083958-38-7 | 50   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

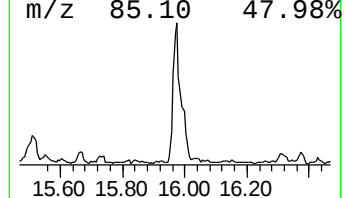
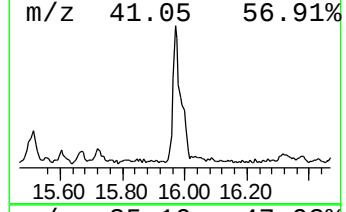
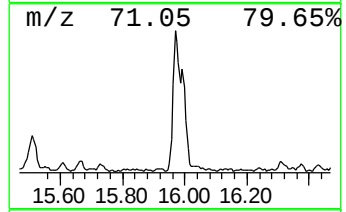
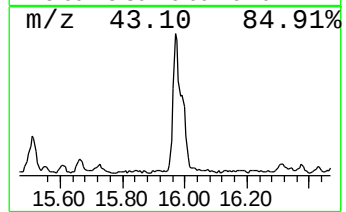
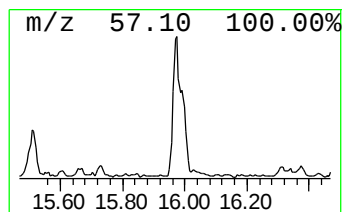
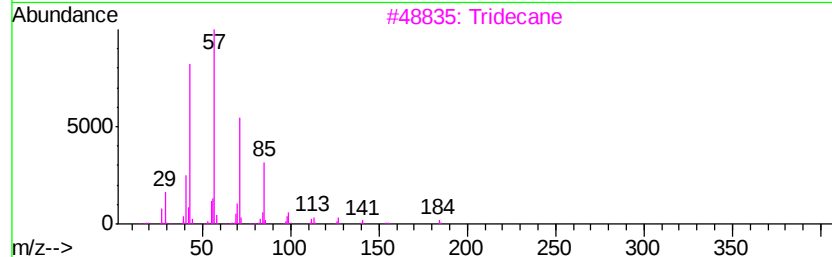
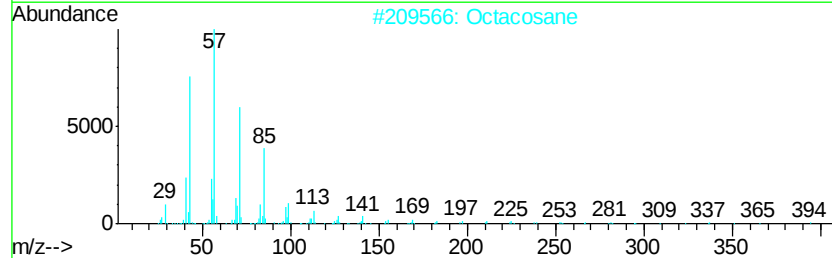
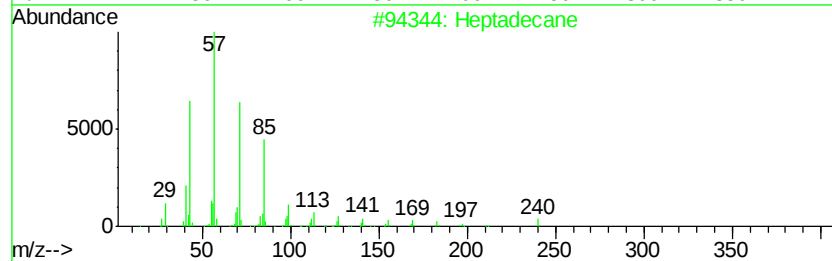
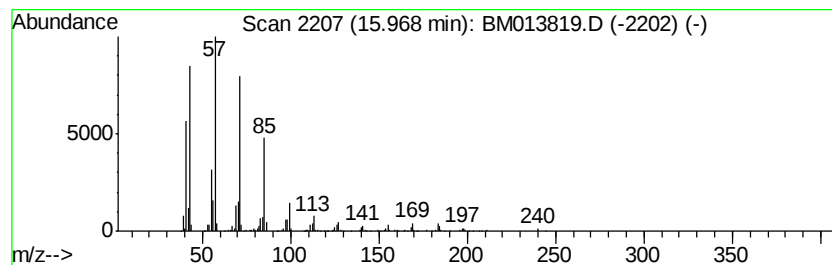
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 1

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.97 | 19.49 ng/ul | 1545560 | Phenanthrene-d10 | 17.00 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane  | 240 | C17H36  | 000629-78-7 | 91   |
| 2    |    |   | Octacosane   | 394 | C28H58  | 000630-02-4 | 90   |
| 3    |    |   | Tridecane    | 184 | C13H28  | 000629-50-5 | 89   |
| 4    |    |   | Tetradecane  | 198 | C14H30  | 000629-59-4 | 87   |
| 5    |    |   | Heptadecane  | 240 | C17H36  | 000629-78-7 | 87   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

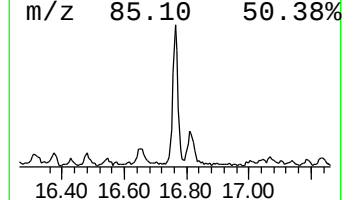
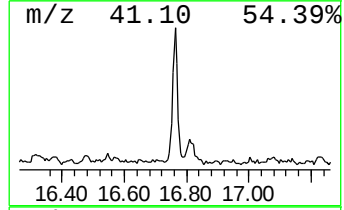
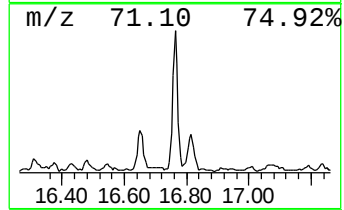
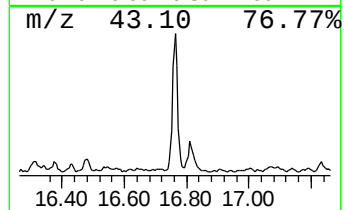
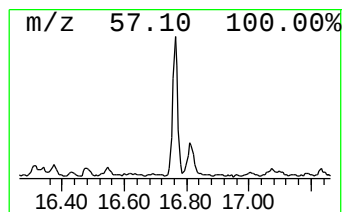
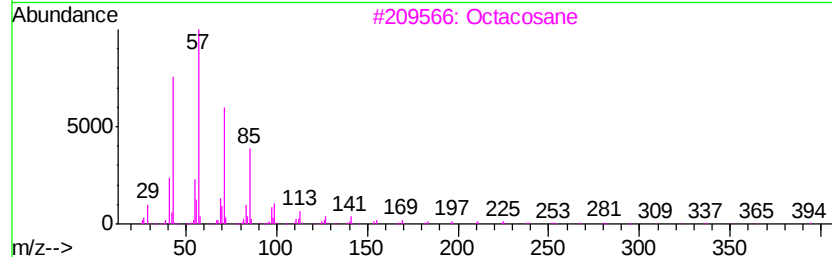
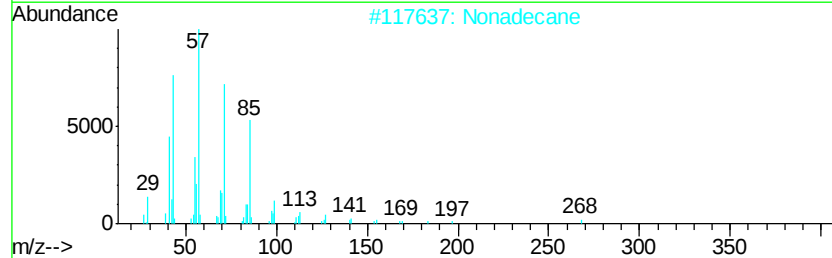
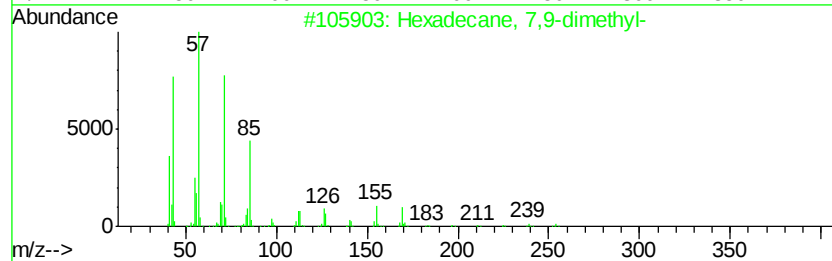
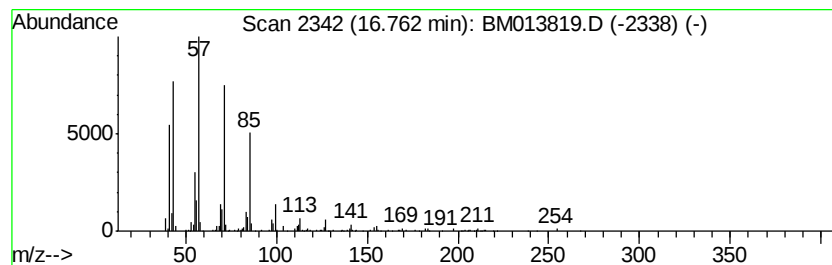
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 3

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 16.76 | 11.27 ng/ul | 893980 | Phenanthrene-d10 | 17.00 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane, 7,9-dimethyl- | 254 | C18H38  | 021164-95-4 | 91   |
| 2    |    | Nonadecane                | 268 | C19H40  | 000629-92-5 | 91   |
| 3    |    | Octacosane                | 394 | C28H58  | 000630-02-4 | 91   |
| 4    |    | Pentadecane               | 212 | C15H32  | 000629-62-9 | 91   |
| 5    |    | Eicosane                  | 282 | C20H42  | 000112-95-8 | 91   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

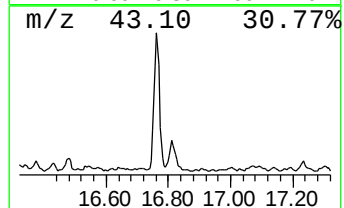
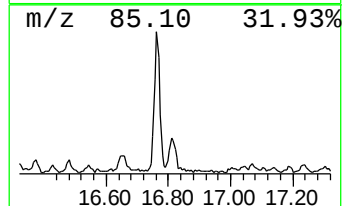
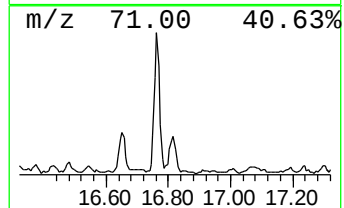
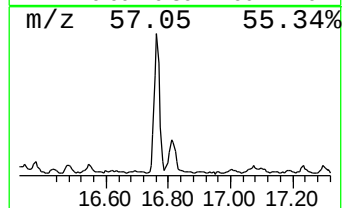
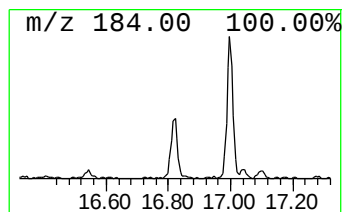
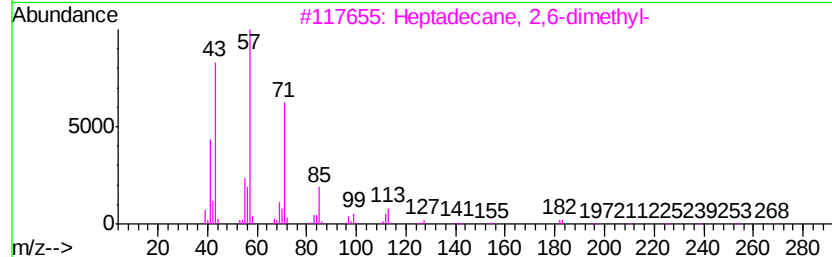
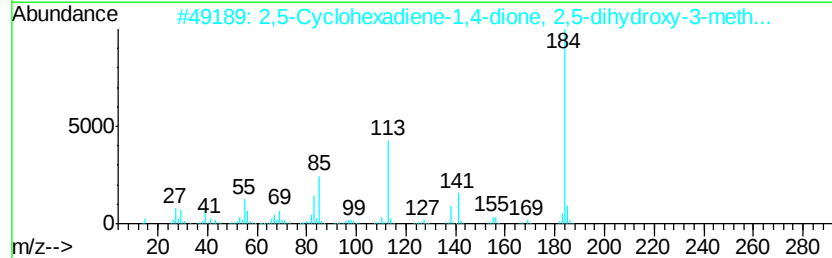
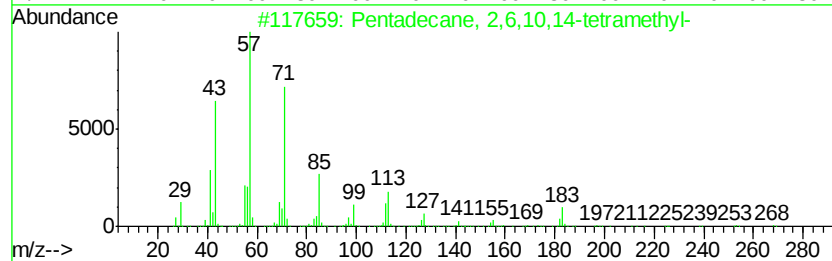
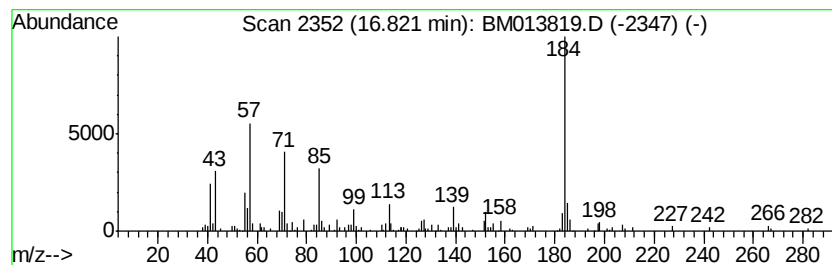
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.82 | 5.42 ng/ul | 430088 | Phenanthrene-d10 | 17.00 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40  | 001921-70-6 | 64   |
| 2    |    |   | 2,5-Cyclohexadiene-1,4-dione, 2,... | 184 | C8H8O5  | 000085-23-4 | 60   |
| 3    |    |   | Heptadecane, 2,6-dimethyl-          | 268 | C19H40  | 054105-67-8 | 50   |
| 4    |    |   | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40  | 001921-70-6 | 50   |
| 5    |    |   | Hexadecane, 2,6,10-trimethyl-       | 268 | C19H40  | 055000-52-7 | 50   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

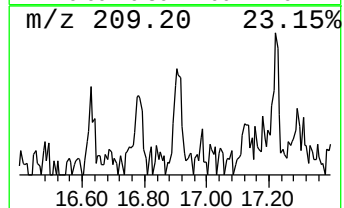
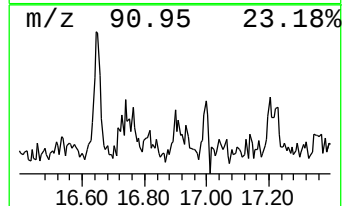
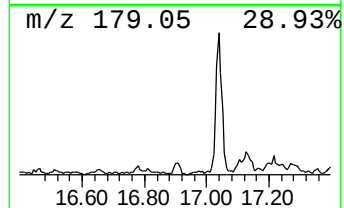
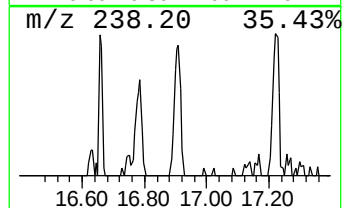
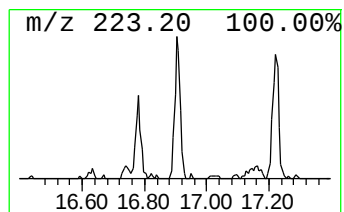
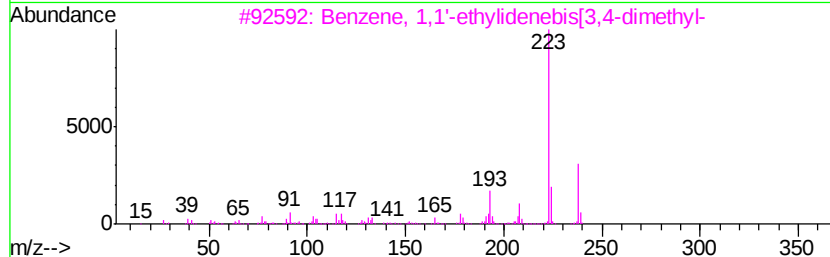
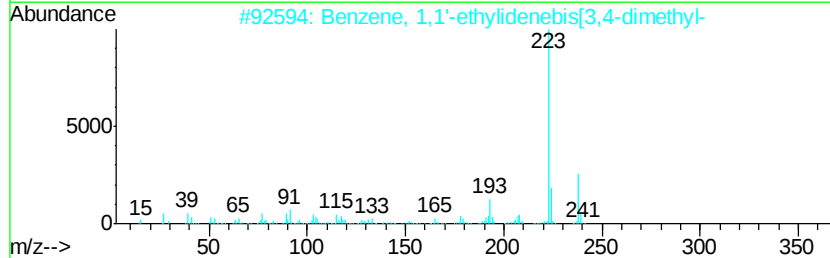
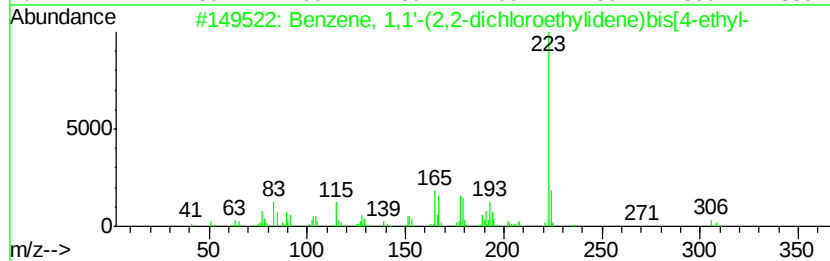
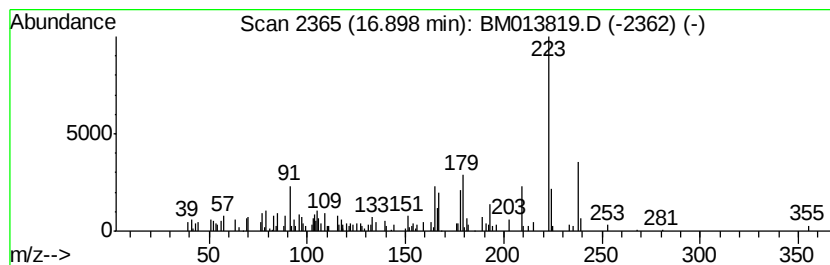
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 22 Benzene, 1,1'-(2,2-dichloro... Concentration Rank 29

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.90 | 2.07 ng/ul | 164145 | Phenanthrene-d10 | 17.00 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | Benzene, 1,1'-(2,2-dichloroethyl... | 306 | C18H20Cl2 | 000072-56-0 | 64   |
| 2    |    |   | Benzene, 1,1'-ethylidenebis[3,4-... | 238 | C18H22    | 001742-14-9 | 62   |
| 3    |    |   | Benzene, 1,1'-ethylidenebis[3,4-... | 238 | C18H22    | 001742-14-9 | 62   |
| 4    |    |   | 3,4'-Diisopropylbiphenyl            | 238 | C18H22    | 061434-46-6 | 60   |
| 5    |    |   | 4,4'-Diisopropylbiphenyl            | 238 | C18H22    | 018970-30-4 | 53   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

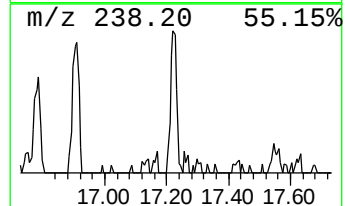
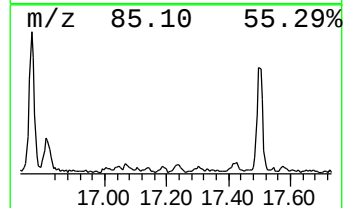
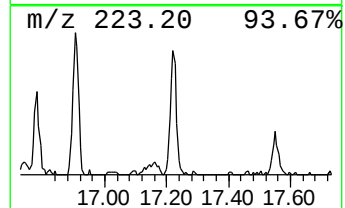
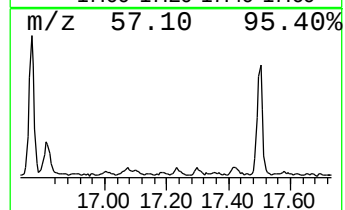
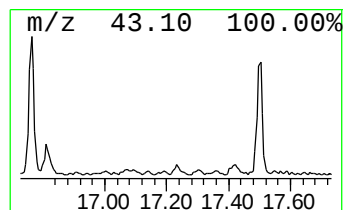
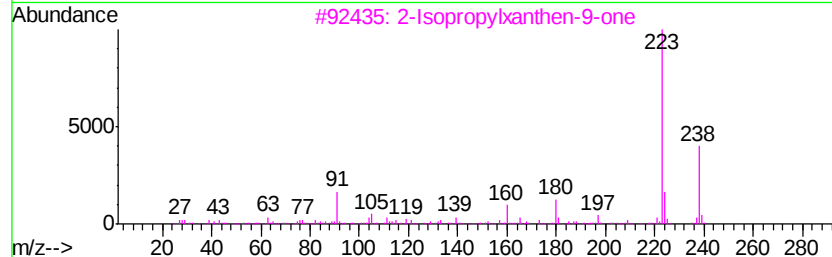
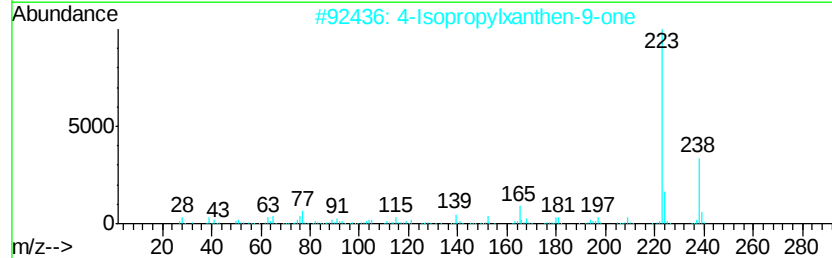
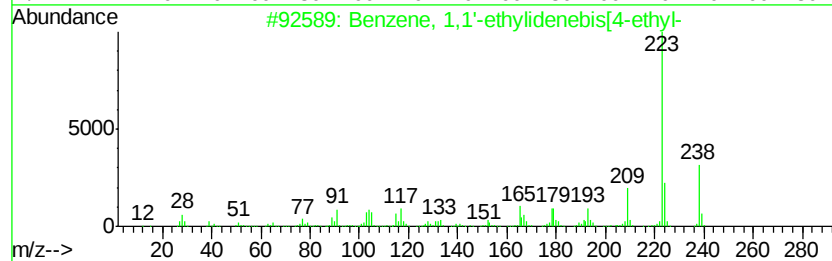
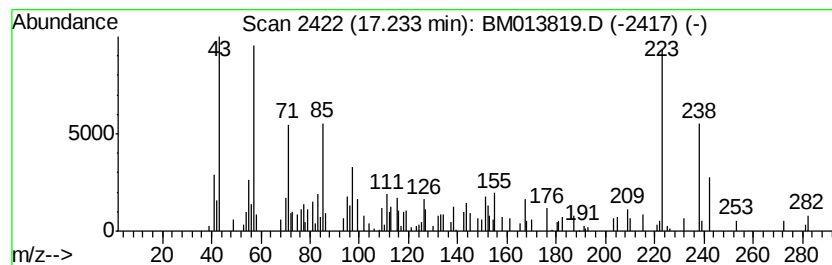
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 23 Benzene, 1,1'-ethylidenebis... Concentration Rank 23

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.23 | 2.64 ng/ul | 208982 | Phenanthrene-d10 | 17.00 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Benzene, 1,1'-ethylidenebis[4-et... | 238 | C18H22   | 010224-91-6  | 90   |
| 2    |    | 4-Isopropylxanthen-9-one            | 238 | C16H14O2 | 1000210-33-1 | 52   |
| 3    |    | 2-Isopropylxanthen-9-one            | 238 | C16H14O2 | 069737-66-2  | 49   |
| 4    |    | Anthracene, 9,10-dimethoxy-         | 238 | C16H14O2 | 002395-97-3  | 49   |
| 5    |    | Cyclopent[a]indene, 3,8-dihydro-... | 238 | C18H22   | 017384-72-4  | 46   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

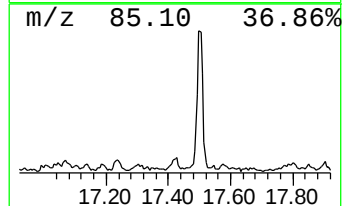
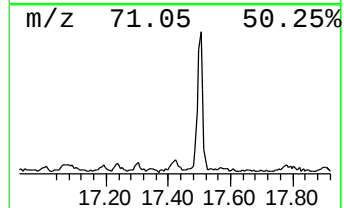
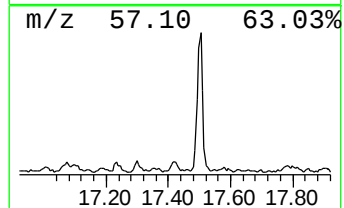
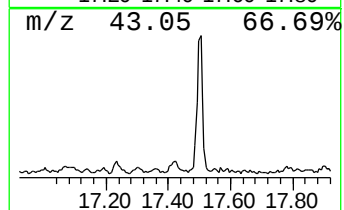
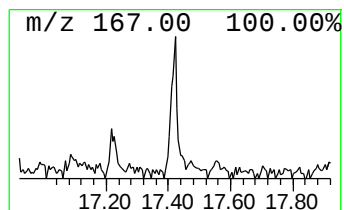
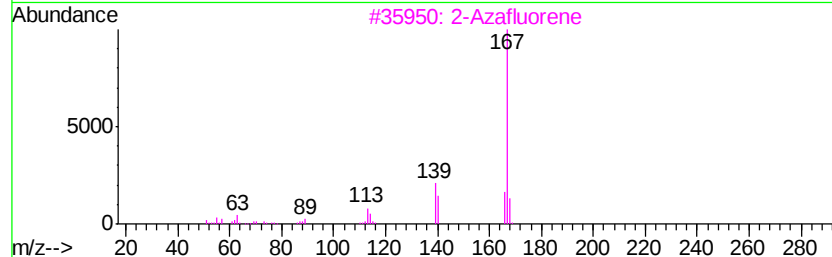
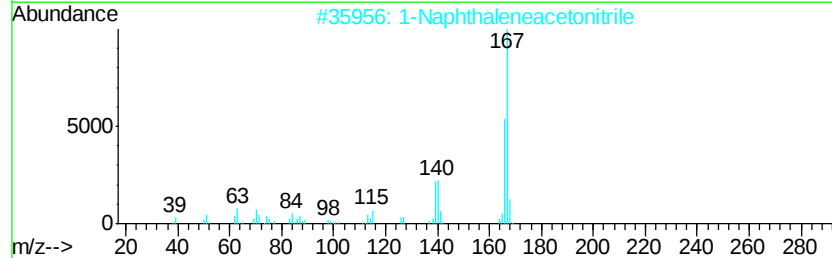
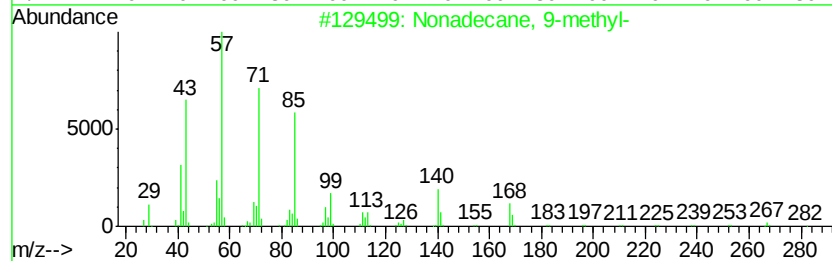
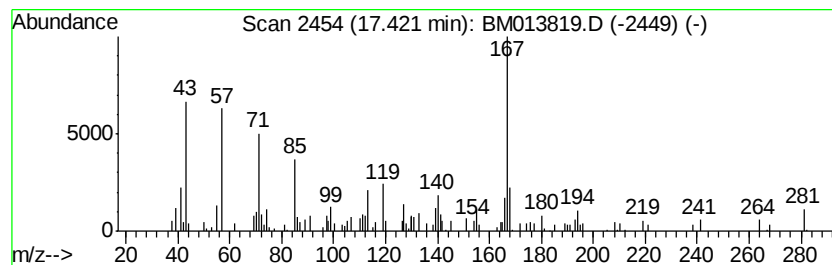
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 25

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.42 | 2.57 ng/ul | 203583 | Phenanthrene-d10 | 17.00 |

| Hit# | of | Tentative ID                       | MW  | MolForm  | CAS#        | Qual |
|------|----|------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Nonadecane, 9-methyl-              | 282 | C20H42   | 013287-24-6 | 35   |
| 2    |    | 1-Naphthaleneacetonitrile          | 167 | C12H9N   | 000132-75-2 | 30   |
| 3    |    | 2-Azafluorene                      | 167 | C12H9N   | 000244-40-6 | 27   |
| 4    |    | 2-Naphthylacetonitrile             | 167 | C12H9N   | 007498-57-9 | 27   |
| 5    |    | 2-Chloro-6-methylphenyl isocyanate | 167 | C8H6ClNO | 019241-34-0 | 25   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

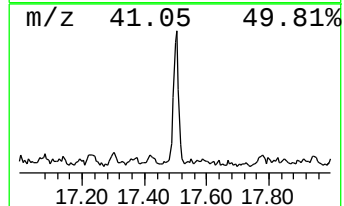
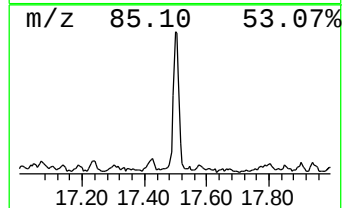
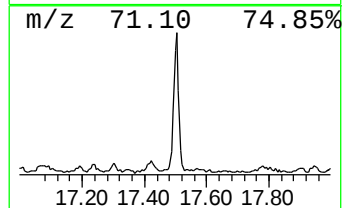
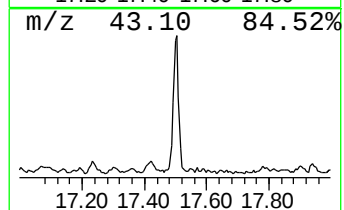
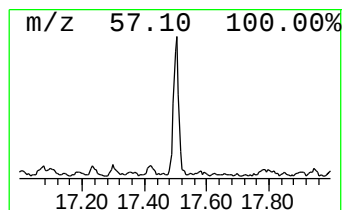
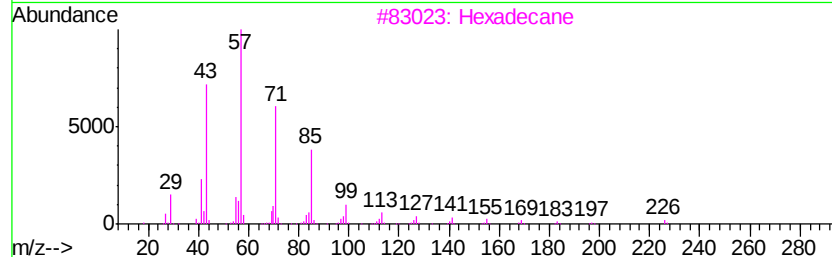
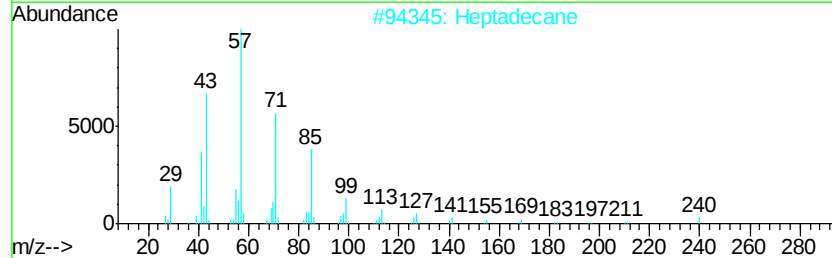
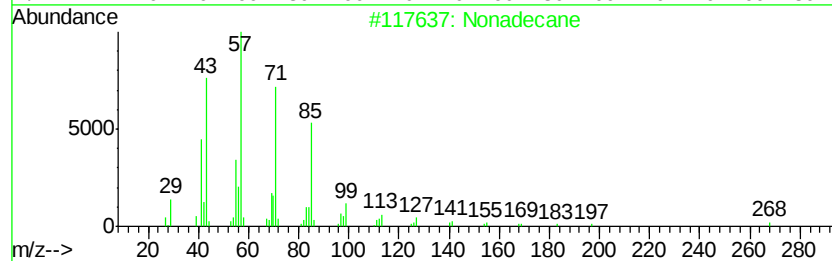
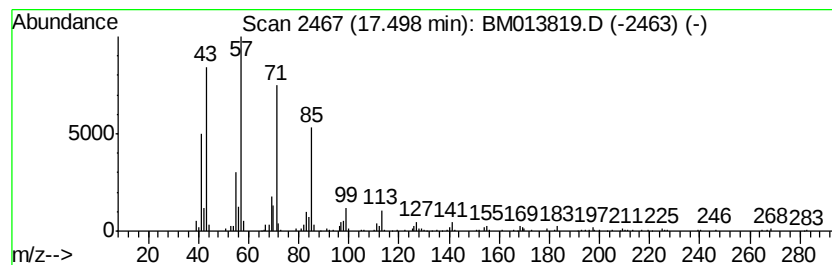
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.50 | 8.74 ng/ul | 693466 | Phenanthrene-d10 | 17.00 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Nonadecane   | 268 | C19H40  | 000629-92-5 | 97   |
| 2    |    |   | Heptadecane  | 240 | C17H36  | 000629-78-7 | 91   |
| 3    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 91   |
| 4    |    |   | Heneicosane  | 296 | C21H44  | 000629-94-7 | 90   |
| 5    |    |   | Docosane     | 310 | C22H46  | 000629-97-0 | 90   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

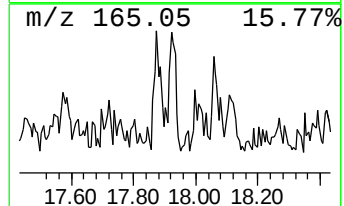
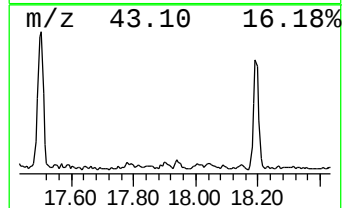
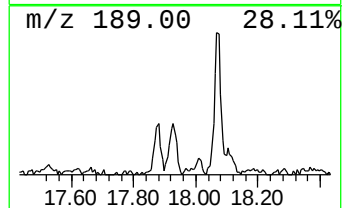
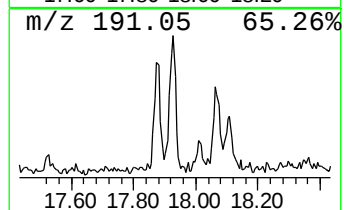
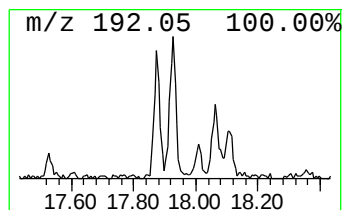
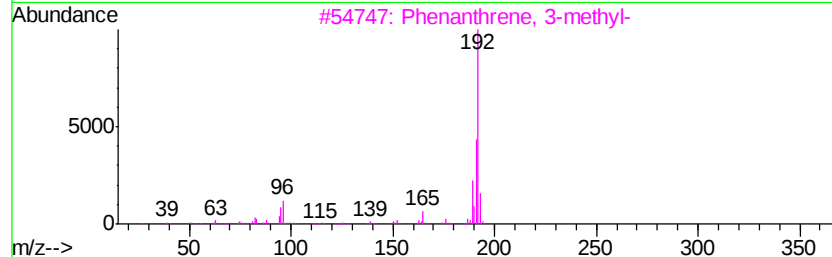
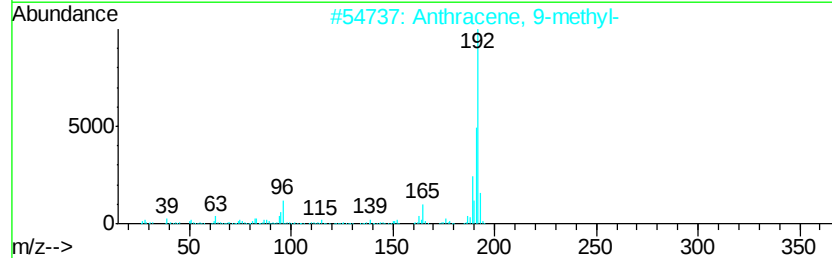
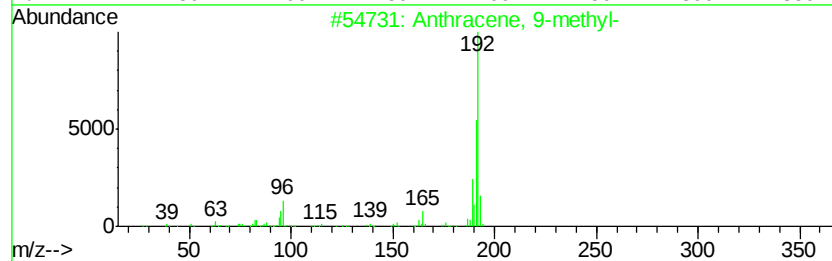
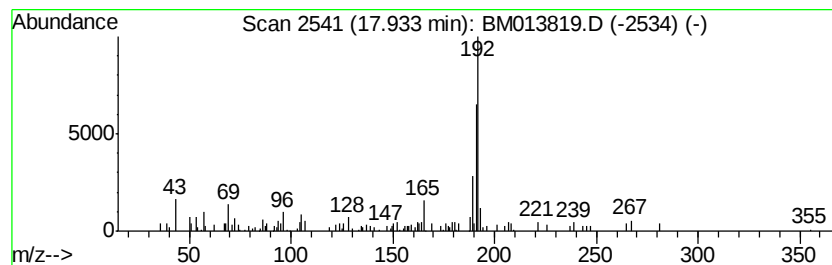
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 26 Anthracene, 9-methyl- Concentration Rank 22

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.93 | 2.89 ng/ul | 228929 | Phenanthrene-d10 | 17.00 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 76   |
| 2    |    |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 64   |
| 3    |    |   | Phenanthrene, 3-methyl- | 192 | C15H12  | 000832-71-3 | 62   |
| 4    |    |   | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 59   |
| 5    |    |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 59   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

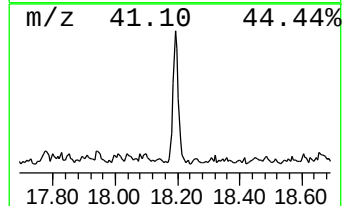
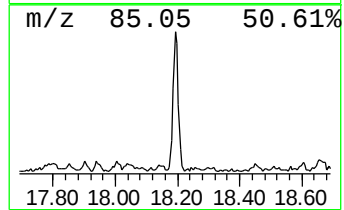
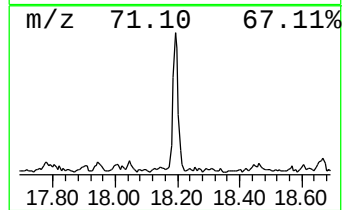
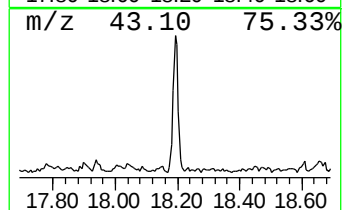
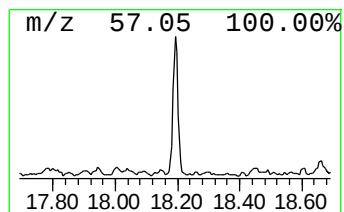
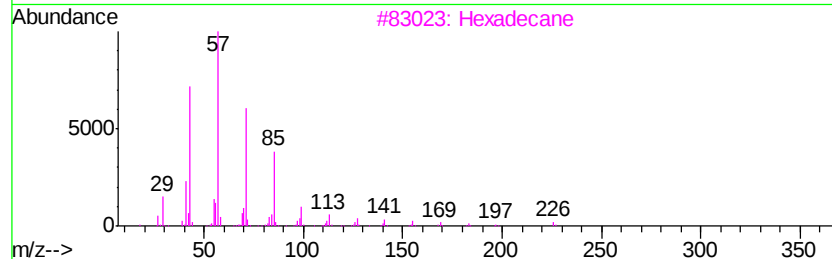
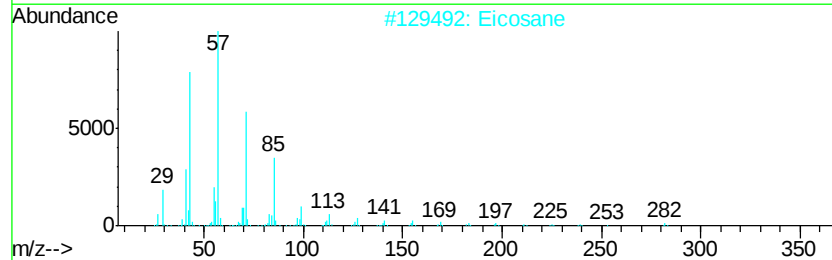
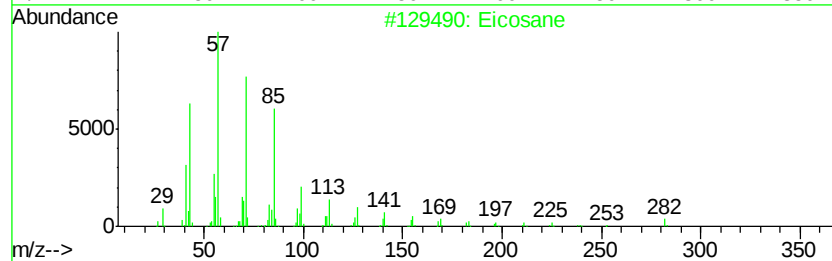
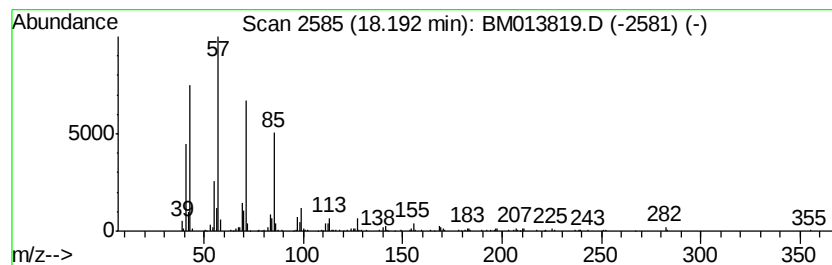
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.19 | 6.68 ng/ul | 529704 | Phenanthrene-d10 | 17.00 |

| Hit# | of                   | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----------------------|---|--------------|-----|---------|-------------|------|
| 1    | Eicosane             |   |              | 282 | C20H42  | 000112-95-8 | 97   |
| 2    | Eicosane             |   |              | 282 | C20H42  | 000112-95-8 | 96   |
| 3    | Hexadecane           |   |              | 226 | C16H34  | 000544-76-3 | 94   |
| 4    | Hexadecane           |   |              | 226 | C16H34  | 000544-76-3 | 94   |
| 5    | Tridecane, 6-propyl- |   |              | 226 | C16H34  | 055045-10-8 | 93   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

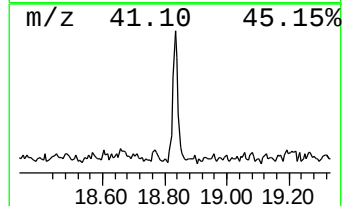
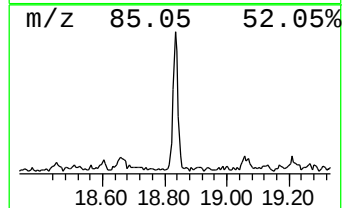
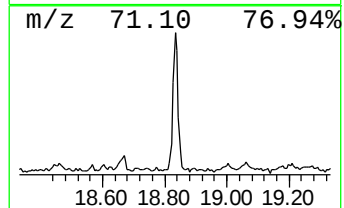
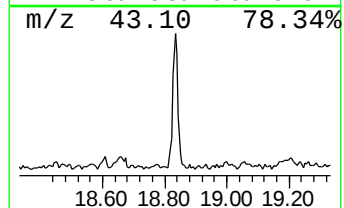
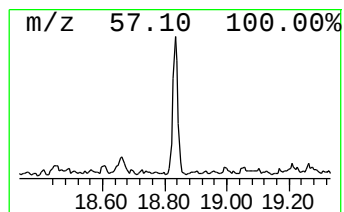
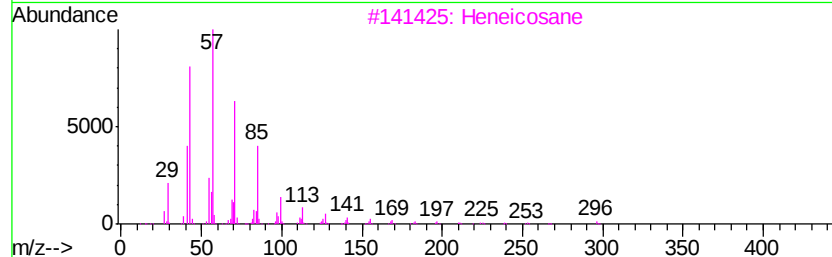
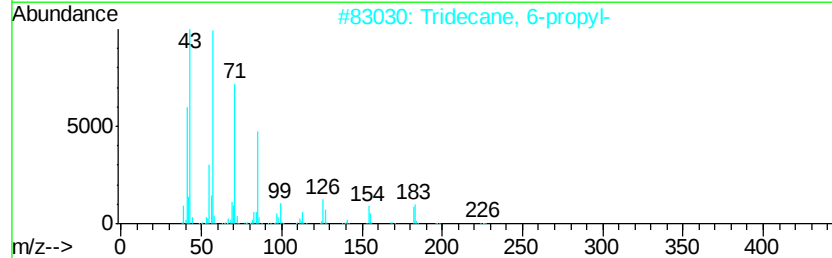
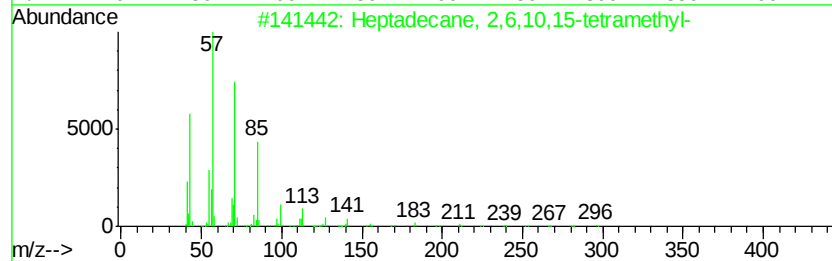
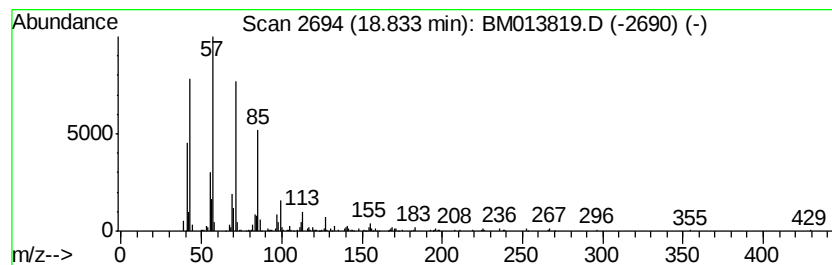
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.83 | 5.52 ng/ul | 437387 | Phenanthrene-d10 | 17.00 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 93   |
| 2    |    |   | Tridecane, 6-propyl-                | 226 | C16H34  | 055045-10-8 | 93   |
| 3    |    |   | Heneicosane                         | 296 | C21H44  | 000629-94-7 | 93   |
| 4    |    |   | Hexacosane                          | 366 | C26H54  | 000630-01-3 | 91   |
| 5    |    |   | Nonadecane                          | 268 | C19H40  | 000629-92-5 | 90   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

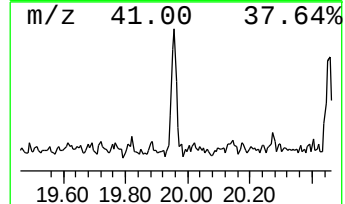
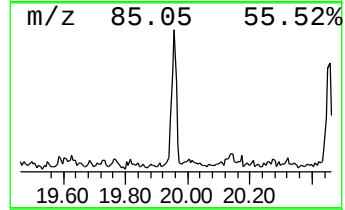
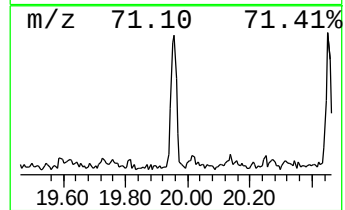
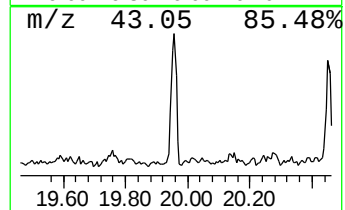
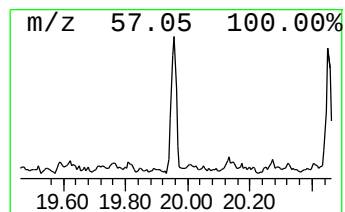
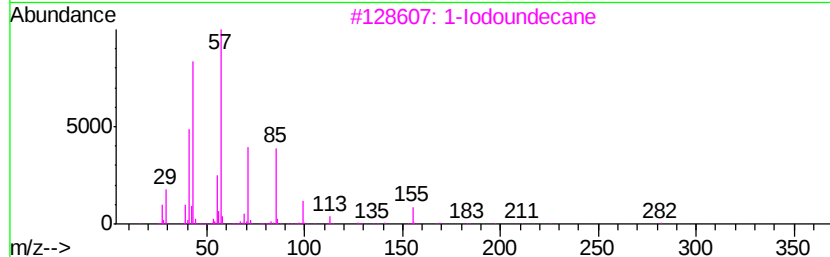
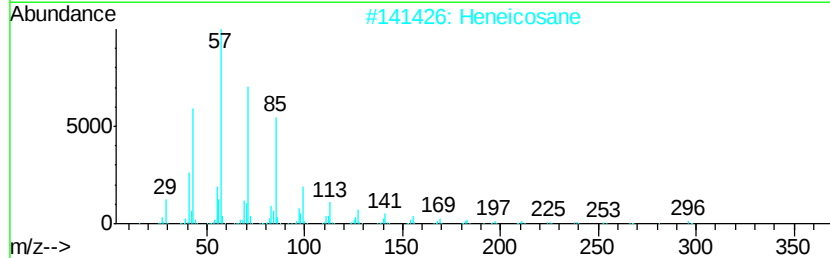
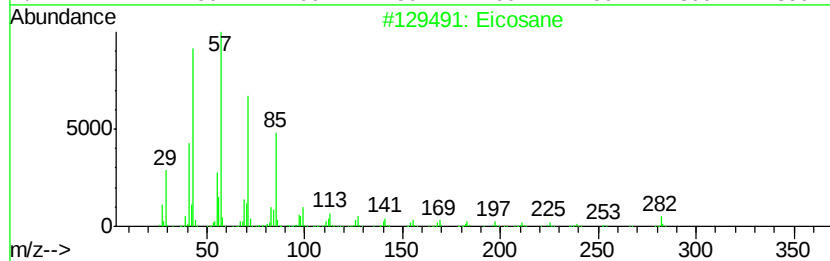
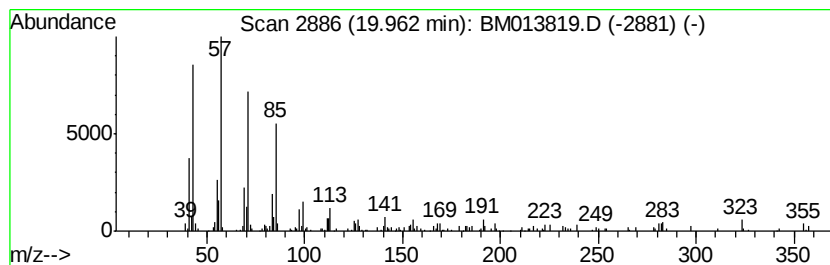
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.96 | 3.71 ng/ul | 341692 | Chrysene-d12     | 21.20 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Eicosane             | 282 | C20H42  | 000112-95-8 | 91   |
| 2    |    |   | Heneicosane          | 296 | C21H44  | 000629-94-7 | 80   |
| 3    |    |   | 1-Iodoundecane       | 282 | C11H23I | 004282-44-4 | 76   |
| 4    |    |   | Hexacosane           | 366 | C26H54  | 000630-01-3 | 74   |
| 5    |    |   | Eicosane, 10-methyl- | 296 | C21H44  | 054833-23-7 | 74   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

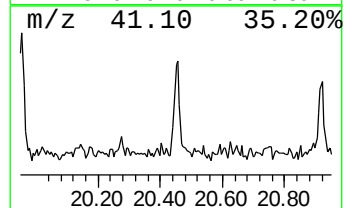
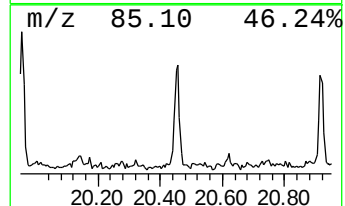
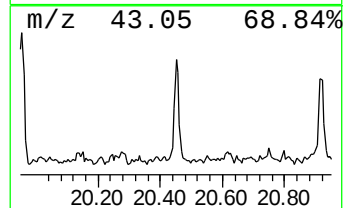
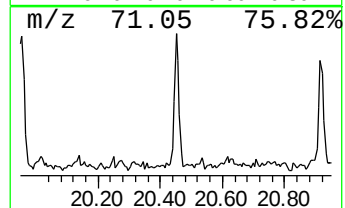
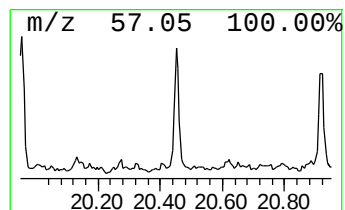
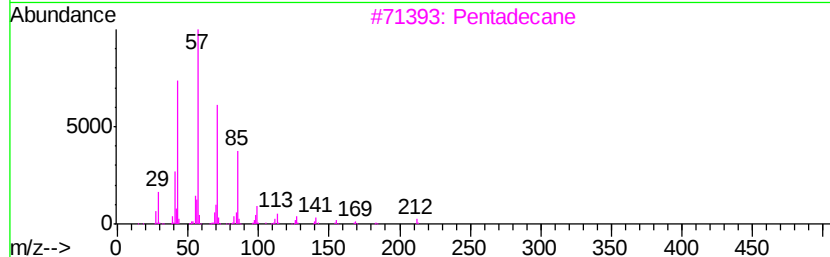
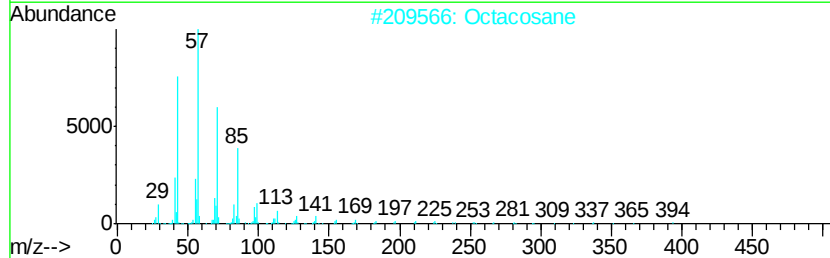
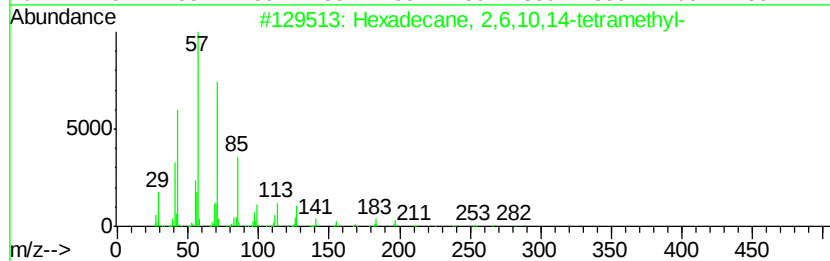
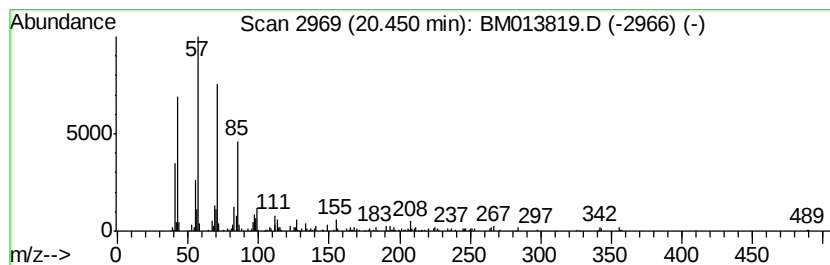
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 24

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.45 | 2.61 ng/ul | 239912 | Chrysene-d12     | 21.20 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8 | 90   |
| 2    |    |   | Octacosane                         | 394 | C28H58  | 000630-02-4 | 87   |
| 3    |    |   | Pentadecane                        | 212 | C15H32  | 000629-62-9 | 87   |
| 4    |    |   | Hexadecane                         | 226 | C16H34  | 000544-76-3 | 87   |
| 5    |    |   | Nonadecane                         | 268 | C19H40  | 000629-92-5 | 87   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

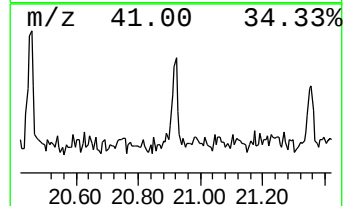
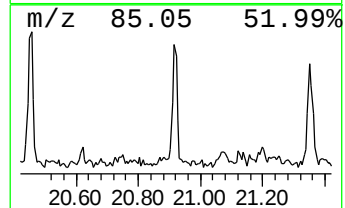
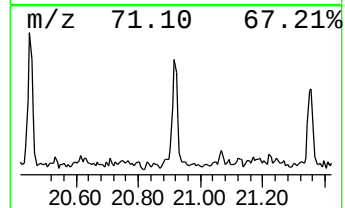
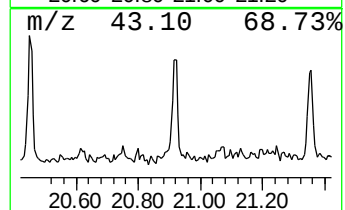
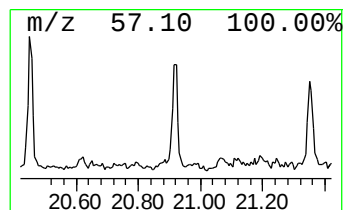
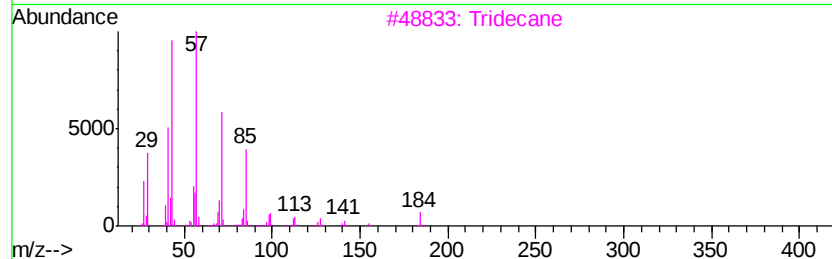
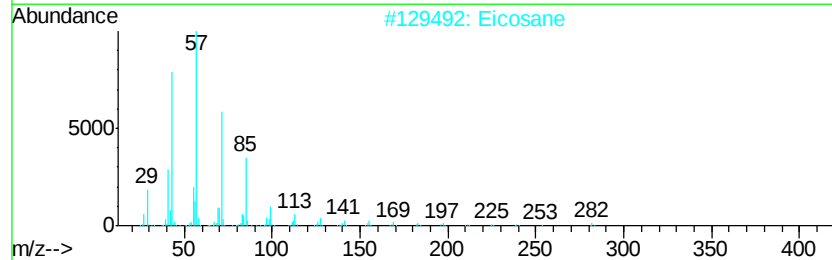
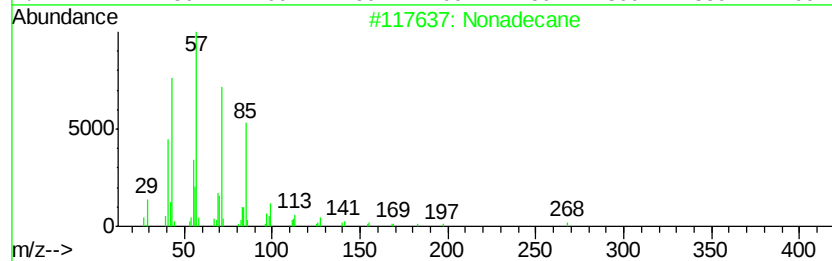
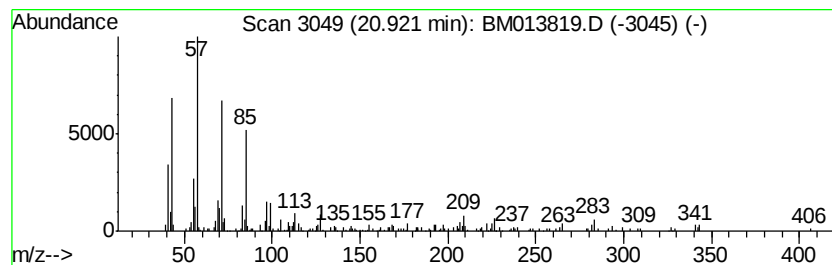
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 31 (DEL) Alkane: Straight-Chai... Concentration Rank 26

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.92 | 2.48 ng/ul | 228158 | Chrysene-d12     | 21.20 |

| Hit# | of | Tentative ID     | MW  | MolForm | CAS#        | Qual |
|------|----|------------------|-----|---------|-------------|------|
| 1    | 5  | Nonadecane       | 268 | C19H40  | 000629-92-5 | 91   |
| 2    |    | Eicosane         | 282 | C20H42  | 000112-95-8 | 83   |
| 3    |    | Tridecane        | 184 | C13H28  | 000629-50-5 | 81   |
| 4    |    | Nonane, 5-butyl- | 184 | C13H28  | 017312-63-9 | 81   |
| 5    |    | Hexadecane       | 226 | C16H34  | 000544-76-3 | 81   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013819.D  
Acq On : 25 Jan 2018 23:11  
Operator : SJ/JU  
Sample : J1222-08DL 20X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A81DL

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

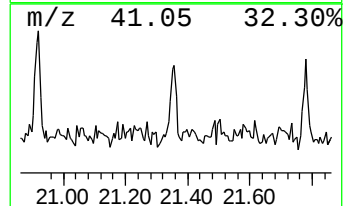
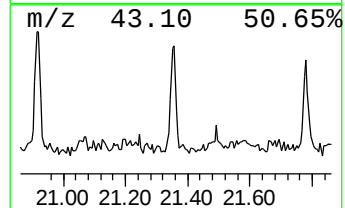
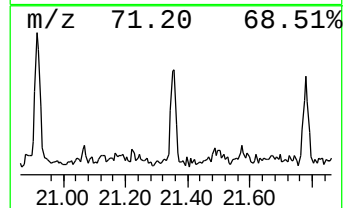
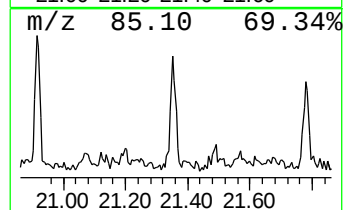
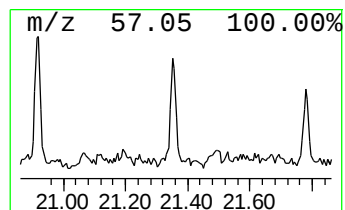
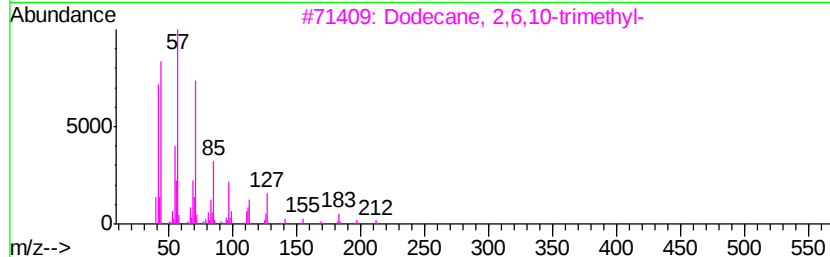
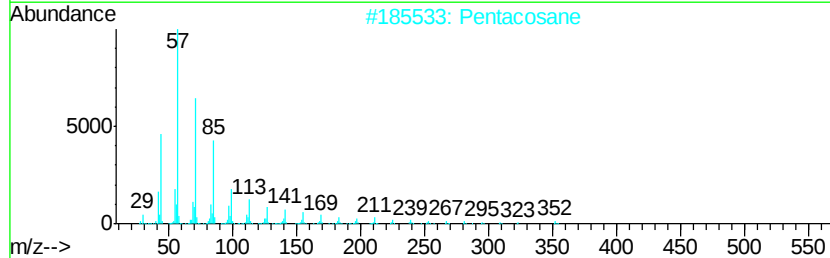
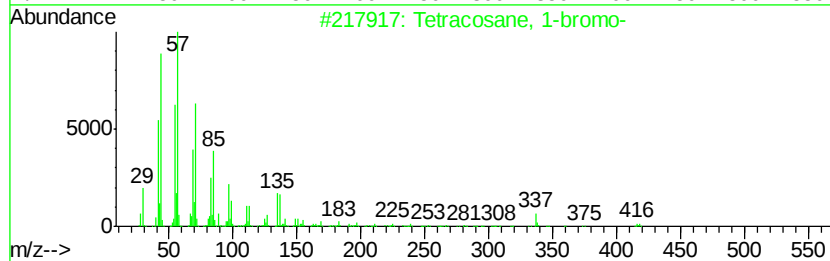
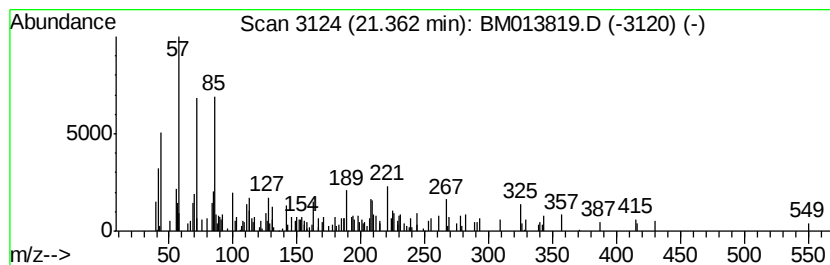
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 31

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 21.36 | 2.03 ng/ul | 187077 | Chrysene-d12     | 21.20 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-----------------------------|-----|----------|-------------|------|
| 1    |    |   | Tetracosane, 1-bromo-       | 416 | C24H49Br | 006946-24-3 | 81   |
| 2    |    |   | Pentacosane                 | 352 | C25H52   | 000629-99-2 | 80   |
| 3    |    |   | Dodecane, 2,6,10-trimethyl- | 212 | C15H32   | 003891-98-3 | 78   |
| 4    |    |   | Nonadecane                  | 268 | C19H40   | 000629-92-5 | 76   |
| 5    |    |   | Octadecane                  | 254 | C18H38   | 000593-45-3 | 76   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM012518\  
 Data File : BM013819.D  
 Acq On : 25 Jan 2018 23:11  
 Operator : SJ/JU  
 Sample : J1222-08DL 20X  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 F6A81DL

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011018.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 |
| unknown-01           | 4.76  | 3.3     | ng/ul | 143644   | 1                     | 7.59  | 879186  | 20.0 |
| Hexylene glycol      | 5.95  | 6.1     | ng/ul | 269222   | 1                     | 7.59  | 879186  | 20.0 |
| (DEL) Alkane: Str... | 7.21  | 3.1     | ng/ul | 137319   | 1                     | 7.59  | 879186  | 20.0 |
| unknown-02           | 8.23  | 2.2     | ng/ul | 95193    | 1                     | 7.59  | 879186  | 20.0 |
| Oxalic acid, 2-et... | 11.39 | 3.0     | ng/ul | 233686   | 2                     | 10.36 | 1546940 | 20.0 |
| (DEL) Alkane: Str... | 11.79 | 8.9     | ng/ul | 690949   | 2                     | 10.36 | 1546940 | 20.0 |
| Naphthalene, 1-me... | 12.26 | 4.7     | ng/ul | 365074   | 2                     | 10.36 | 1546940 | 20.0 |
| (DEL) Alkane: Str... | 12.76 | 2.1     | ng/ul | 234738   | 3                     | 14.24 | 2273760 | 20.0 |
| Naphthalene, 1,6-... | 13.41 | 3.2     | ng/ul | 368424   | 3                     | 14.24 | 2273760 | 20.0 |
| Naphthalene, 1,7-... | 13.56 | 3.5     | ng/ul | 403850   | 3                     | 14.24 | 2273760 | 20.0 |
| Sulfurous acid, 2... | 13.73 | 3.1     | ng/ul | 357539   | 3                     | 14.24 | 2273760 | 20.0 |
| (DEL) Alkane: Str... | 14.14 | 9.3     | ng/ul | 1052050  | 3                     | 14.24 | 2273760 | 20.0 |
| Naphthalene, 2-(1... | 14.46 | 2.1     | ng/ul | 239006   | 3                     | 14.24 | 2273760 | 20.0 |
| (DEL) Alkane: Str... | 14.77 | 2.0     | ng/ul | 230303   | 3                     | 14.24 | 2273760 | 20.0 |
| (DEL) Alkane: Str... | 15.10 | 8.6     | ng/ul | 972869   | 3                     | 14.24 | 2273760 | 20.0 |
| (DEL) Alkane: Str... | 15.52 | 2.9     | ng/ul | 332630   | 3                     | 14.24 | 2273760 | 20.0 |
| (4-Acetylphenyl)p... | 15.72 | 12.5    | ng/ul | 993482   | 4                     | 17.00 | 1586160 | 20.0 |
| (DEL) Alkane: Str... | 15.97 | 19.5    | ng/ul | 1545560  | 4                     | 17.00 | 1586160 | 20.0 |
| (DEL) Alkane: Str... | 16.76 | 11.3    | ng/ul | 893980   | 4                     | 17.00 | 1586160 | 20.0 |
| (DEL) Alkane: Str... | 16.82 | 5.4     | ng/ul | 430088   | 4                     | 17.00 | 1586160 | 20.0 |
| Benzene, 1,1'-(2,... | 16.90 | 2.1     | ng/ul | 164145   | 4                     | 17.00 | 1586160 | 20.0 |
| Benzene, 1,1'-eth... | 17.23 | 2.6     | ng/ul | 208982   | 4                     | 17.00 | 1586160 | 20.0 |
| (DEL) Alkane: Str... | 17.42 | 2.6     | ng/ul | 203583   | 4                     | 17.00 | 1586160 | 20.0 |
| (DEL) Alkane: Str... | 17.50 | 8.7     | ng/ul | 693466   | 4                     | 17.00 | 1586160 | 20.0 |
| Anthracene, 9-met... | 17.93 | 2.9     | ng/ul | 228929   | 4                     | 17.00 | 1586160 | 20.0 |
| (DEL) Alkane: Str... | 18.19 | 6.7     | ng/ul | 529704   | 4                     | 17.00 | 1586160 | 20.0 |
| (DEL) Alkane: Str... | 18.83 | 5.5     | ng/ul | 437387   | 4                     | 17.00 | 1586160 | 20.0 |
| (DEL) Alkane: Str... | 19.96 | 3.7     | ng/ul | 341692   | 5                     | 21.20 | 1840920 | 20.0 |
| (DEL) Alkane: Str... | 20.45 | 2.6     | ng/ul | 239912   | 5                     | 21.20 | 1840920 | 20.0 |
| (DEL) Alkane: Str... | 20.92 | 2.5     | ng/ul | 228158   | 5                     | 21.20 | 1840920 | 20.0 |
| (DEL) Alkane: Str... | 21.36 | 2.0     | ng/ul | 187077   | 5                     | 21.20 | 1840920 | 20.0 |