

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
 Data File : BM018190.D  
 Acq On : 15 Dec 2018 05:11  
 Operator : JU/SJ  
 Sample : J6238-01  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BE7Z4

## Integration Parameters: LSCINT.P

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

Filtering: 5  
 Min Area: 0.1 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
 Title : SVOA CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 3.593       | 115           | 120         | 129          | rBV      | 301684         | 391296        | 15.78%          | 0.788%        |
| 2         | 4.669       | 299           | 303         | 312          | rVV2     | 49776          | 95213         | 3.84%           | 0.192%        |
| 3         | 4.775       | 312           | 321         | 332          | rVV      | 80379          | 138404        | 5.58%           | 0.279%        |
| 4         | 5.187       | 386           | 391         | 403          | rVB      | 51896          | 102084        | 4.12%           | 0.206%        |
| 5         | 6.704       | 643           | 649         | 664          | rVB2     | 364911         | 766034        | 30.90%          | 1.542%        |
| 6         | 6.857       | 669           | 675         | 687          | rBV      | 315313         | 490164        | 19.77%          | 0.987%        |
| 7         | 7.045       | 700           | 707         | 715          | rVV      | 421583         | 678051        | 27.35%          | 1.365%        |
| 8         | 7.122       | 715           | 720         | 723          | rVV      | 52863          | 80554         | 3.25%           | 0.162%        |
| 9         | 7.504       | 779           | 785         | 799          | rVB      | 692780         | 1111490       | 44.84%          | 2.238%        |
| 10        | 8.228       | 901           | 908         | 917          | rBV      | 347602         | 631970        | 25.49%          | 1.273%        |
| 11        | 8.586       | 964           | 969         | 974          | rVB3     | 43817          | 73100         | 2.95%           | 0.147%        |
| 12        | 8.663       | 974           | 982         | 986          | rBV      | 401146         | 675794        | 27.26%          | 1.361%        |
| 13        | 8.710       | 986           | 990         | 998          | rVB      | 163953         | 259642        | 10.47%          | 0.523%        |
| 14        | 8.822       | 999           | 1009        | 1017         | rVB10    | 32798          | 107498        | 4.34%           | 0.216%        |
| 15        | 9.186       | 1065          | 1071        | 1077         | rBV6     | 55251          | 102023        | 4.12%           | 0.205%        |
| 16        | 9.375       | 1093          | 1103        | 1112         | rBV2     | 356510         | 708012        | 28.56%          | 1.426%        |
| 17        | 9.628       | 1140          | 1146        | 1150         | rVV6     | 38167          | 82687         | 3.34%           | 0.166%        |
| 18        | 9.692       | 1150          | 1157        | 1163         | rVV6     | 44963          | 107762        | 4.35%           | 0.217%        |
| 19        | 9.910       | 1188          | 1194        | 1208         | rBV      | 406428         | 822830        | 33.19%          | 1.657%        |
| 20        | 10.269      | 1244          | 1255        | 1261         | rBV      | 1001102        | 2079215       | 83.87%          | 4.187%        |
| 21        | 10.322      | 1261          | 1264        | 1270         | rVV      | 77294          | 126800        | 5.11%           | 0.255%        |
| 22        | 10.427      | 1277          | 1282        | 1288         | rVV2     | 100165         | 174990        | 7.06%           | 0.352%        |
| 23        | 11.098      | 1391          | 1396        | 1403         | rBV7     | 72116          | 136453        | 5.50%           | 0.275%        |
| 24        | 11.180      | 1403          | 1410        | 1416         | rVB5     | 76255          | 138594        | 5.59%           | 0.279%        |
| 25        | 11.286      | 1423          | 1428        | 1432         | rBV      | 159495         | 242901        | 9.80%           | 0.489%        |
| 26        | 11.445      | 1448          | 1455        | 1460         | rBV8     | 24189          | 68357         | 2.76%           | 0.138%        |
| 27        | 11.698      | 1489          | 1498        | 1504         | rBV      | 444542         | 703404        | 28.37%          | 1.416%        |
| 28        | 11.945      | 1534          | 1540        | 1545         | rVB      | 261357         | 471986        | 19.04%          | 0.950%        |
| 29        | 12.169      | 1570          | 1578        | 1588         | rVV      | 167460         | 353445        | 14.26%          | 0.712%        |
| 30        | 12.351      | 1602          | 1609        | 1612         | rBV5     | 43988          | 82041         | 3.31%           | 0.165%        |
| 31        | 12.398      | 1613          | 1617        | 1622         | rVV3     | 55296          | 105058        | 4.24%           | 0.212%        |
| 32        | 12.533      | 1636          | 1640        | 1650         | rVB4     | 92823          | 211099        | 8.52%           | 0.425%        |
| 33        | 12.621      | 1650          | 1655        | 1658         | rBV2     | 87662          | 145519        | 5.87%           | 0.293%        |
| 34        | 12.674      | 1659          | 1664        | 1669         | rVV2     | 222018         | 348495        | 14.06%          | 0.702%        |

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 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 12.721 | 1669 | 1672 | 1679 | rVB9 | 43095   | 69503   | 2.80%  | 0.140% |
| 36 | 12.904 | 1697 | 1703 | 1706 | rBV8 | 32318   | 82891   | 3.34%  | 0.167% |
| 37 | 12.974 | 1710 | 1715 | 1721 | rVV  | 631845  | 886021  | 35.74% | 1.784% |
| 38 | 13.063 | 1727 | 1730 | 1737 | rBV7 | 34146   | 77242   | 3.12%  | 0.156% |
| 39 | 13.133 | 1739 | 1742 | 1746 | rVB4 | 54957   | 68593   | 2.77%  | 0.138% |
| 40 | 13.180 | 1746 | 1750 | 1752 | rBV  | 73707   | 100977  | 4.07%  | 0.203% |
| 41 | 13.316 | 1766 | 1773 | 1774 | rBV2 | 182138  | 280876  | 11.33% | 0.566% |
| 42 | 13.474 | 1794 | 1800 | 1805 | rBV  | 270232  | 516715  | 20.84% | 1.040% |
| 43 | 13.527 | 1805 | 1809 | 1812 | rVV  | 180056  | 276297  | 11.15% | 0.556% |
| 44 | 13.586 | 1815 | 1819 | 1823 | rVV  | 847016  | 1157124 | 46.68% | 2.330% |
| 45 | 13.633 | 1823 | 1827 | 1832 | rVB2 | 402616  | 635035  | 25.62% | 1.279% |
| 46 | 13.686 | 1832 | 1836 | 1838 | rBV3 | 92097   | 123853  | 5.00%  | 0.249% |
| 47 | 13.851 | 1858 | 1864 | 1873 | rVB  | 972236  | 1568894 | 63.29% | 3.159% |
| 48 | 13.974 | 1881 | 1885 | 1889 | rBV4 | 80429   | 125964  | 5.08%  | 0.254% |
| 49 | 14.068 | 1894 | 1901 | 1905 | rBV  | 764054  | 1041879 | 42.03% | 2.098% |
| 50 | 14.163 | 1910 | 1917 | 1923 | rVB2 | 1379735 | 2155939 | 86.97% | 4.341% |
| 51 | 14.221 | 1923 | 1927 | 1930 | rBV2 | 62281   | 98798   | 3.99%  | 0.199% |
| 52 | 14.374 | 1949 | 1953 | 1956 | rBV  | 86614   | 150608  | 6.08%  | 0.303% |
| 53 | 14.410 | 1956 | 1959 | 1965 | rVB4 | 110015  | 198956  | 8.03%  | 0.401% |
| 54 | 14.568 | 1982 | 1986 | 1990 | rVB2 | 235967  | 317769  | 12.82% | 0.640% |
| 55 | 14.610 | 1990 | 1993 | 1995 | rVV4 | 54911   | 75679   | 3.05%  | 0.152% |
| 56 | 14.645 | 1995 | 1999 | 2003 | rVB2 | 155445  | 250241  | 10.09% | 0.504% |
| 57 | 14.692 | 2003 | 2007 | 2015 | rVB3 | 224525  | 341530  | 13.78% | 0.688% |
| 58 | 14.786 | 2016 | 2023 | 2028 | rBV3 | 132372  | 313080  | 12.63% | 0.630% |
| 59 | 14.839 | 2029 | 2032 | 2036 | rVB  | 99810   | 118248  | 4.77%  | 0.238% |
| 60 | 14.939 | 2045 | 2049 | 2055 | rBV6 | 126083  | 266917  | 10.77% | 0.537% |
| 61 | 14.992 | 2056 | 2058 | 2060 | rVV  | 74438   | 74740   | 3.01%  | 0.150% |
| 62 | 15.027 | 2060 | 2064 | 2069 | rVB  | 807898  | 950009  | 38.32% | 1.913% |
| 63 | 15.162 | 2082 | 2087 | 2092 | rVB  | 1208769 | 1647488 | 66.46% | 3.317% |
| 64 | 15.215 | 2092 | 2096 | 2100 | rVB4 | 64378   | 88458   | 3.57%  | 0.178% |
| 65 | 15.315 | 2109 | 2113 | 2121 | rBV7 | 133072  | 296806  | 11.97% | 0.598% |
| 66 | 15.433 | 2128 | 2133 | 2138 | rVB2 | 324104  | 486494  | 19.62% | 0.980% |
| 67 | 15.527 | 2144 | 2149 | 2155 | rVB6 | 134894  | 305654  | 12.33% | 0.615% |
| 68 | 15.586 | 2155 | 2159 | 2164 | rBV4 | 122834  | 222725  | 8.98%  | 0.448% |
| 69 | 15.645 | 2165 | 2169 | 2178 | rVB6 | 98235   | 219781  | 8.87%  | 0.443% |
| 70 | 15.739 | 2181 | 2185 | 2187 | rBV4 | 46918   | 78178   | 3.15%  | 0.157% |
| 71 | 15.898 | 2207 | 2212 | 2220 | rBV  | 850490  | 1853807 | 74.78% | 3.733% |

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Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |        |
|-----|--------|------|------|------|------|---------|---------|---------|--------|
| 72  | 16.580 | 2325 | 2328 | 2335 | rVB5 | 88782   | 129376  | 5.22%   | 0.261% |
| 73  | 16.686 | 2342 | 2346 | 2350 | rBV  | 556861  | 701434  | 28.29%  | 1.412% |
| 74  | 16.739 | 2351 | 2355 | 2365 | rVB  | 303733  | 547547  | 22.09%  | 1.103% |
| 75  | 16.921 | 2381 | 2386 | 2390 | rBV  | 1444882 | 1990270 | 80.28%  | 4.008% |
| 76  | 16.962 | 2390 | 2393 | 2397 | rVB  | 481228  | 597729  | 24.11%  | 1.204% |
| 77  | 17.021 | 2398 | 2403 | 2407 | rBV  | 1165474 | 1594137 | 64.30%  | 3.210% |
| 78  | 17.109 | 2416 | 2418 | 2424 | rVV7 | 58289   | 87795   | 3.54%   | 0.177% |
| 79  | 17.156 | 2424 | 2426 | 2432 | rVB5 | 56475   | 66342   | 2.68%   | 0.134% |
| 80  | 17.345 | 2455 | 2458 | 2463 | rVB5 | 92317   | 131561  | 5.31%   | 0.265% |
| 81  | 17.427 | 2468 | 2472 | 2478 | rVV  | 466917  | 576521  | 23.26%  | 1.161% |
| 82  | 17.503 | 2482 | 2485 | 2494 | rVB3 | 90170   | 196992  | 7.95%   | 0.397% |
| 83  | 17.703 | 2515 | 2519 | 2527 | rBV4 | 163681  | 410712  | 16.57%  | 0.827% |
| 84  | 17.798 | 2533 | 2535 | 2538 | rBV  | 79817   | 92978   | 3.75%   | 0.187% |
| 85  | 17.839 | 2538 | 2542 | 2551 | rVB2 | 691542  | 1045450 | 42.17%  | 2.105% |
| 86  | 17.986 | 2564 | 2567 | 2571 | rVB3 | 117128  | 151162  | 6.10%   | 0.304% |
| 87  | 18.121 | 2586 | 2590 | 2595 | rBV  | 361103  | 470840  | 18.99%  | 0.948% |
| 88  | 18.309 | 2619 | 2622 | 2625 | rVB  | 58437   | 66054   | 2.66%   | 0.133% |
| 89  | 18.727 | 2690 | 2693 | 2696 | rBV  | 127814  | 154793  | 6.24%   | 0.312% |
| 90  | 18.768 | 2696 | 2700 | 2706 | rVB2 | 289986  | 426048  | 17.19%  | 0.858% |
| 91  | 18.992 | 2734 | 2738 | 2747 | rBV  | 765148  | 1000121 | 40.34%  | 2.014% |
| 92  | 19.127 | 2758 | 2761 | 2768 | rBV4 | 244433  | 328626  | 13.26%  | 0.662% |
| 93  | 19.327 | 2791 | 2795 | 2798 | rBV2 | 1156136 | 1711691 | 69.05%  | 3.447% |
| 94  | 19.356 | 2798 | 2800 | 2810 | rVB  | 859076  | 1106557 | 44.64%  | 2.228% |
| 95  | 19.892 | 2889 | 2891 | 2895 | rVB  | 173271  | 176006  | 7.10%   | 0.354% |
| 96  | 20.392 | 2974 | 2976 | 2979 | rVB  | 155765  | 146906  | 5.93%   | 0.296% |
| 97  | 20.862 | 3053 | 3056 | 3063 | rBV3 | 576562  | 657608  | 26.53%  | 1.324% |
| 98  | 21.139 | 3097 | 3103 | 3107 | rBV2 | 1565424 | 2479049 | 100.00% | 4.992% |
| 99  | 22.656 | 3358 | 3361 | 3365 | rBV2 | 601849  | 896566  | 36.17%  | 1.805% |
| 100 | 23.168 | 3445 | 3448 | 3452 | rBV  | 835742  | 1386002 | 55.91%  | 2.791% |

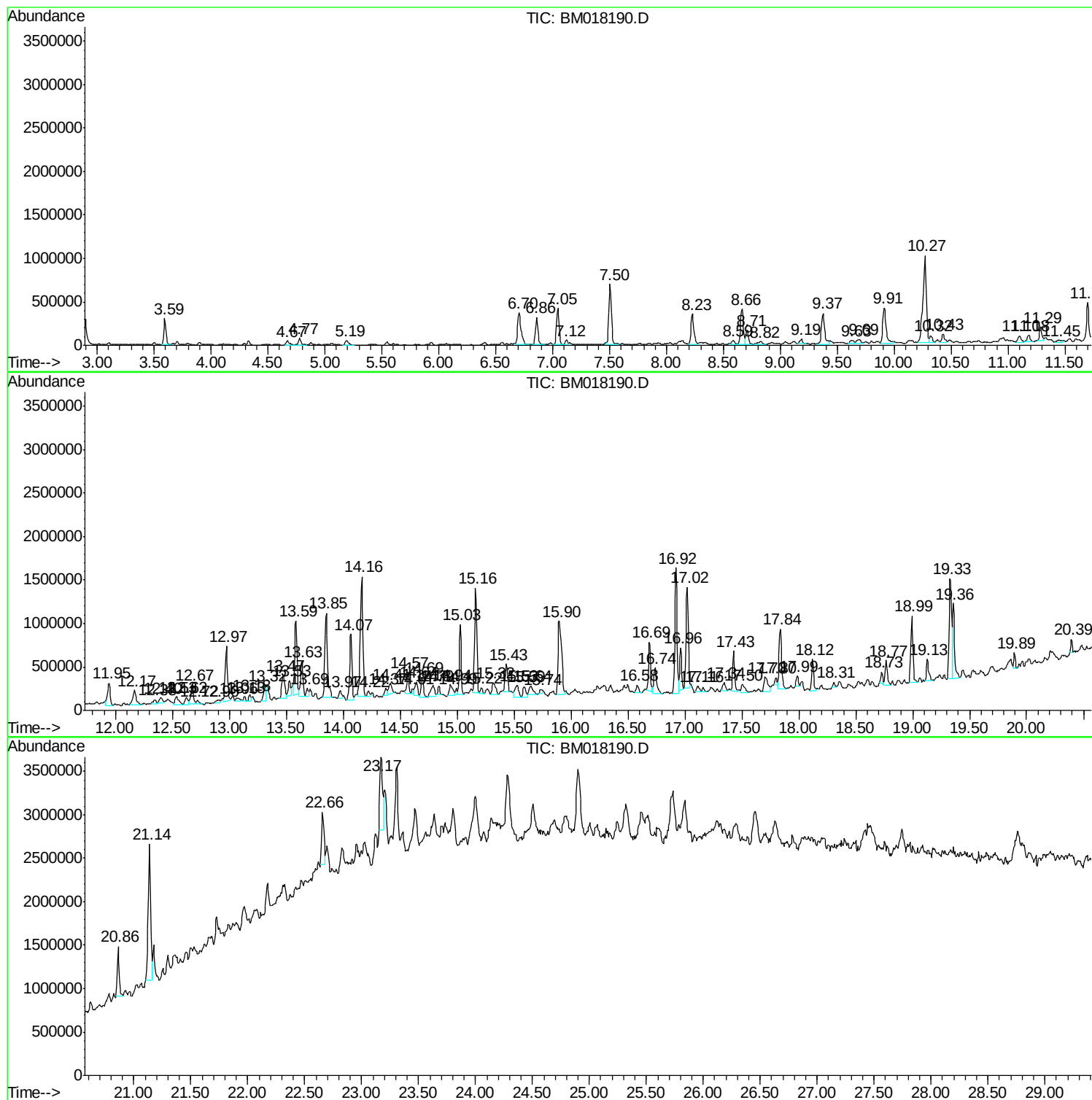
Sum of corrected areas: 49663607

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P



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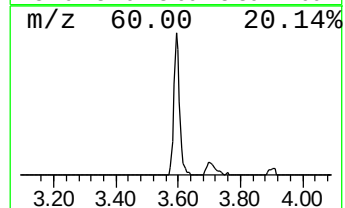
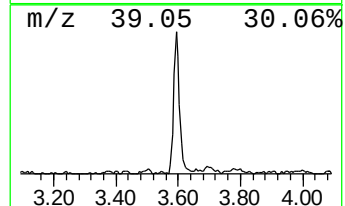
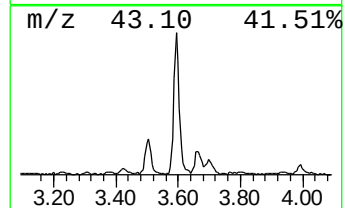
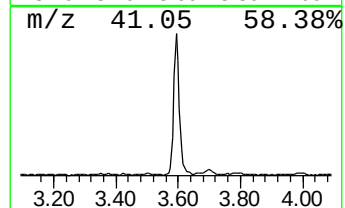
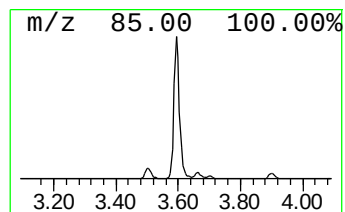
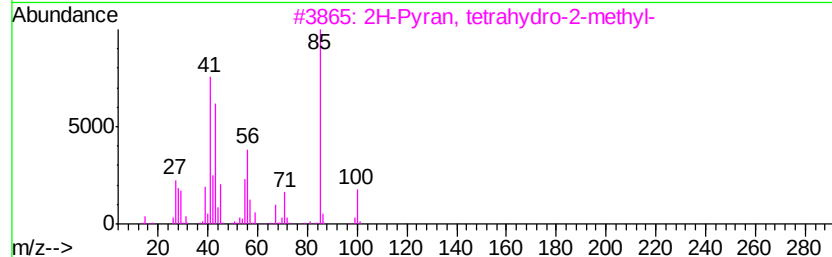
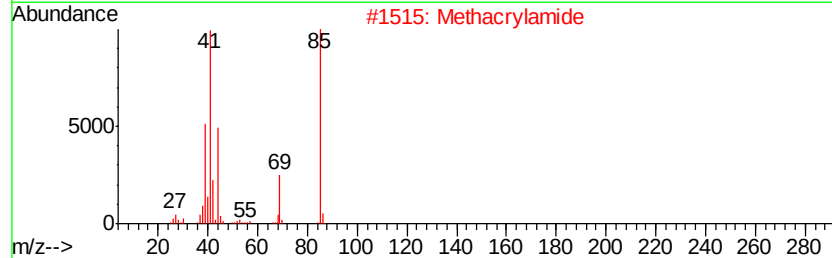
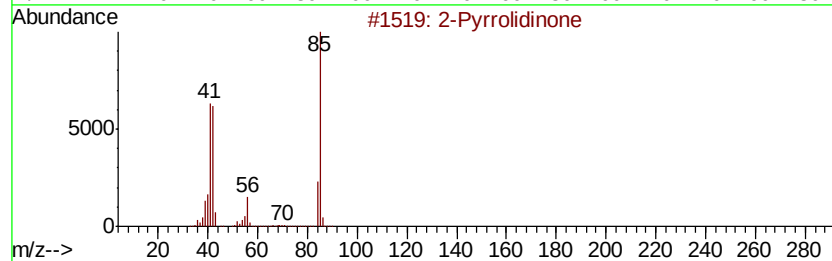
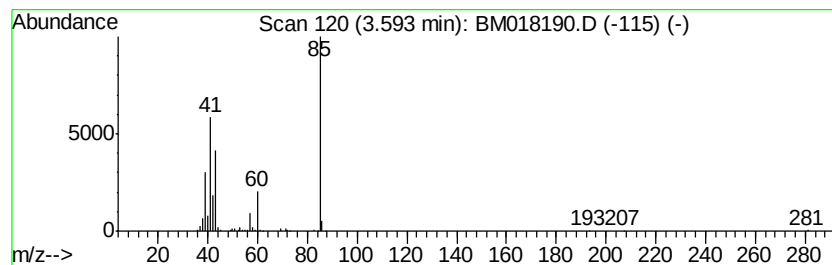
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 2-Pyrrolidinone Concentration Rank 7

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 3.59 | 7.04 ng/ul | 391296 | 1,4-Dichlorobenzene-d4 | 7.50 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pyrrolidinone                | 85  | C4H7NO   | 000616-45-5 | 53   |
| 2    |    |   | Methacrylamide                 | 85  | C4H7NO   | 000079-39-0 | 40   |
| 3    |    |   | 2H-Pyran, tetrahydro-2-methyl- | 100 | C6H12O   | 010141-72-7 | 38   |
| 4    |    |   | 2-Pyrrolidinone                | 85  | C4H7NO   | 000616-45-5 | 9    |
| 5    |    |   | 2-(Bromomethyl)acrylic acid    | 164 | C4H5BrO2 | 072707-66-5 | 9    |



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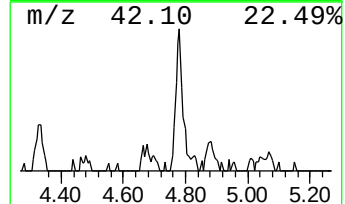
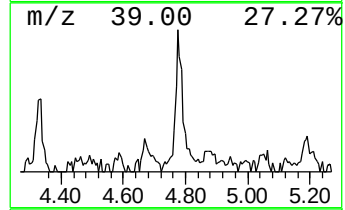
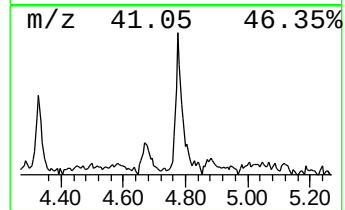
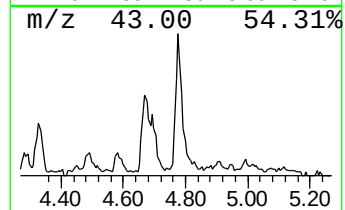
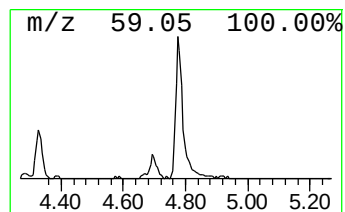
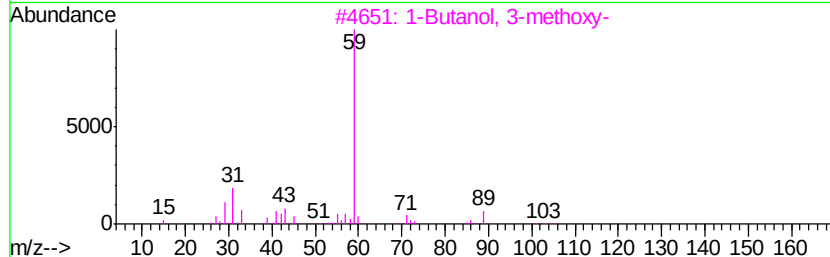
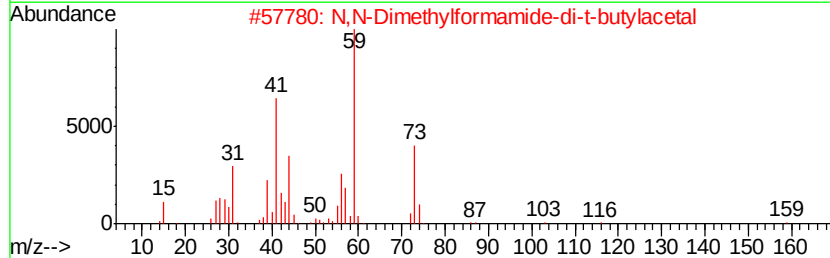
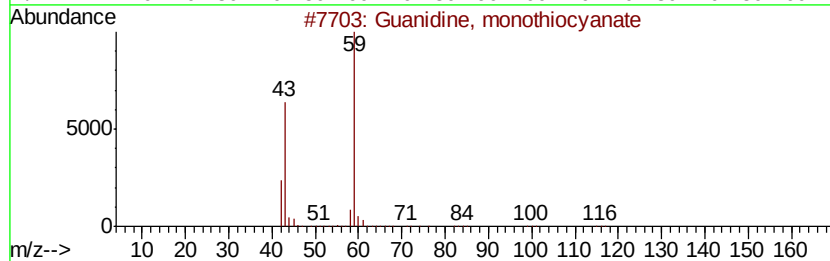
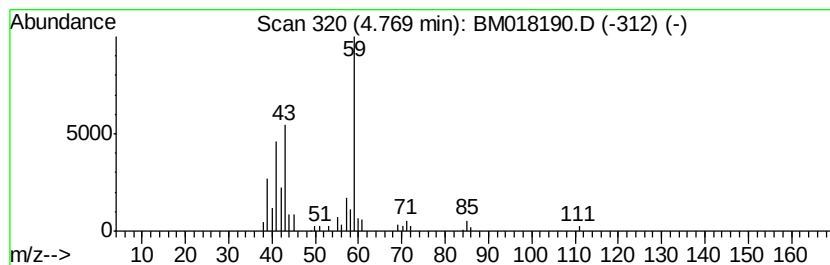
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Guanidine, monothiocyanate Concentration Rank 26

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 4.77 | 2.49 ng/ul | 138404 | 1,4-Dichlorobenzene-d4 | 7.50 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Guanidine, monothiocyanate          | 116 | C2H4N4S   | 056960-89-5  | 59   |
| 2    |    | N,N-Dimethylformamide-di-t-butyl... | 203 | C11H25NO2 | 036805-97-7  | 50   |
| 3    |    | 1-Butanol, 3-methoxy-               | 104 | C5H12O2   | 002517-43-3  | 45   |
| 4    |    | 4-O-Acetyl-2,5-di-O-methyl-3,6-d... | 215 | C10H17NO4 | 1000101-80-3 | 42   |
| 5    |    | Oxirane, tetramethyl-               | 100 | C6H12O    | 005076-20-0  | 40   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BE7Z4

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Quant Title : SVOA CALIBRATION

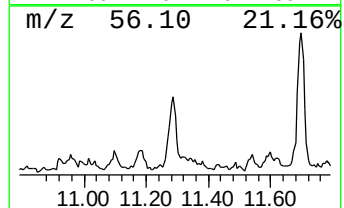
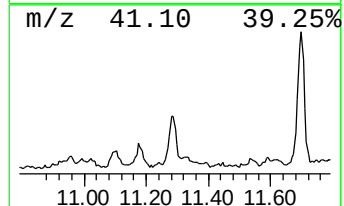
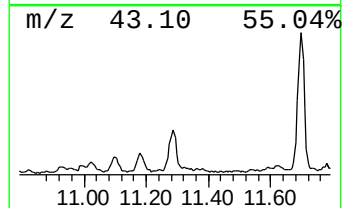
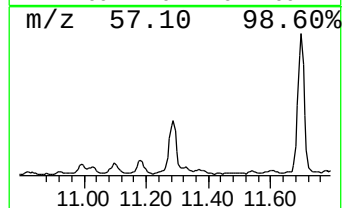
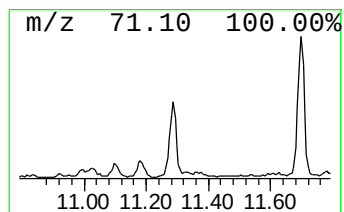
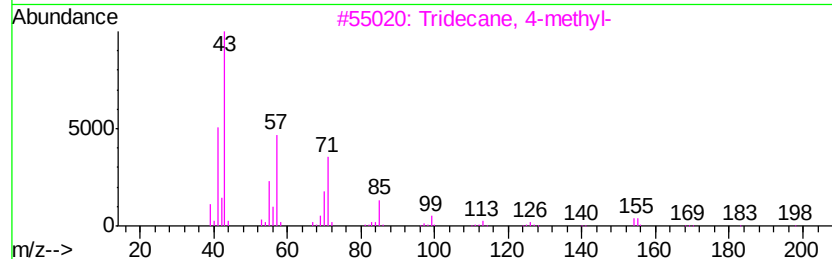
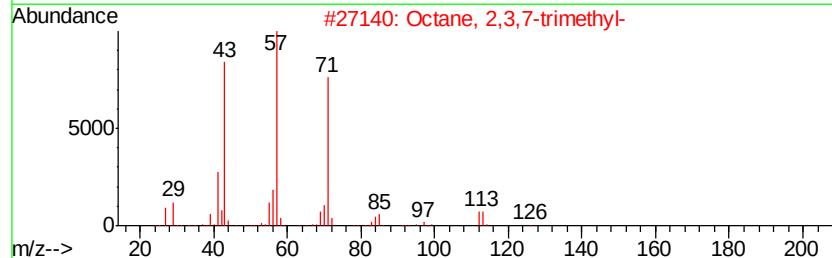
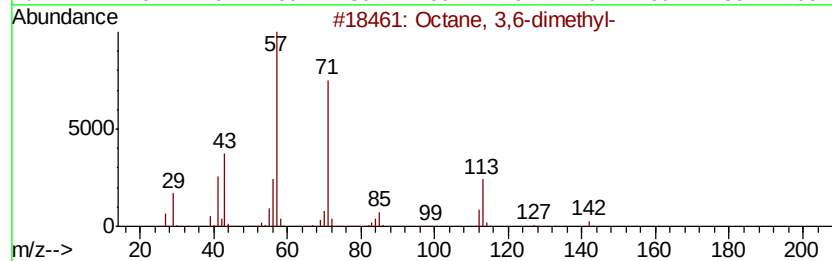
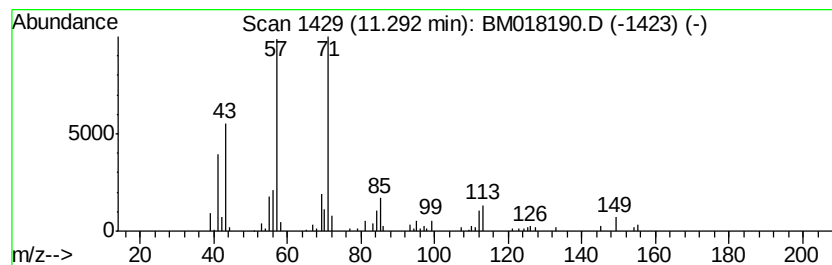
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TIC Integration Parameters: LSCINT.P

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Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 28

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.29 | 2.34 ng/ul | 242901 | Naphthalene-d8   | 10.27 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Octane, 3,6-dimethyl-       | 142 | C10H22  | 015869-94-0 | 64   |
| 2    |    | Octane, 2,3,7-trimethyl-    | 156 | C11H24  | 062016-34-6 | 64   |
| 3    |    | Tridecane, 4-methyl-        | 198 | C14H30  | 026730-12-1 | 59   |
| 4    |    | Dodecane, 2,7,10-trimethyl- | 212 | C15H32  | 074645-98-0 | 59   |
| 5    |    | Heptane, 3-ethyl-5-methyl-  | 142 | C10H22  | 052896-90-9 | 53   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BE7Z4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
Quant Title : SVOA CALIBRATION

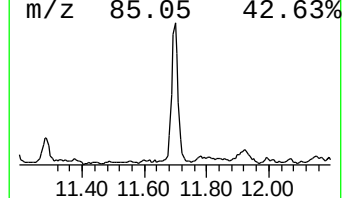
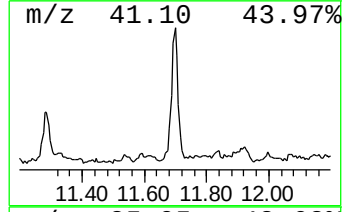
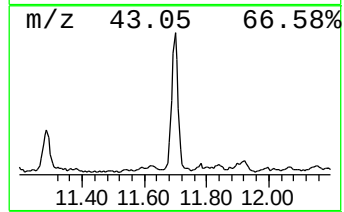
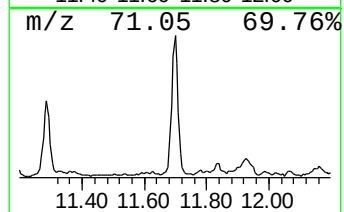
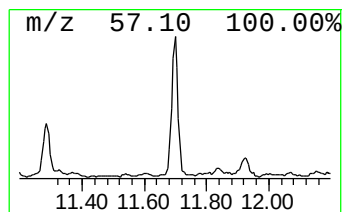
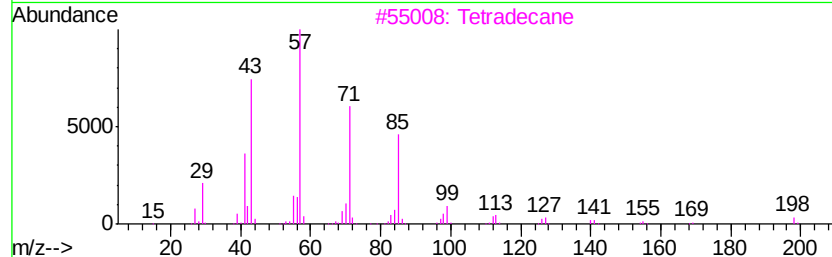
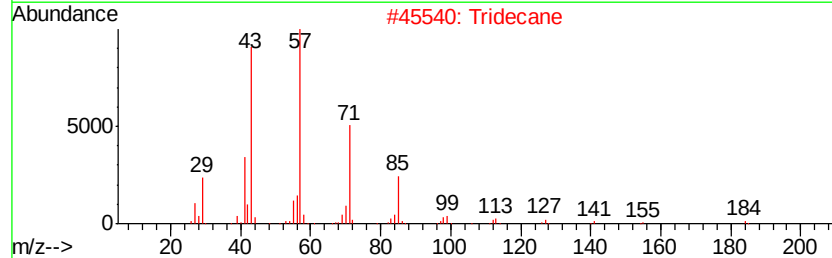
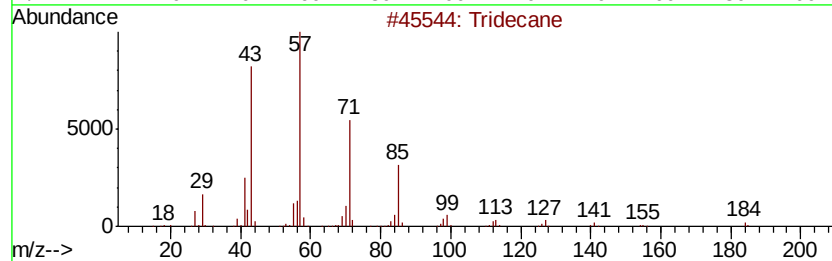
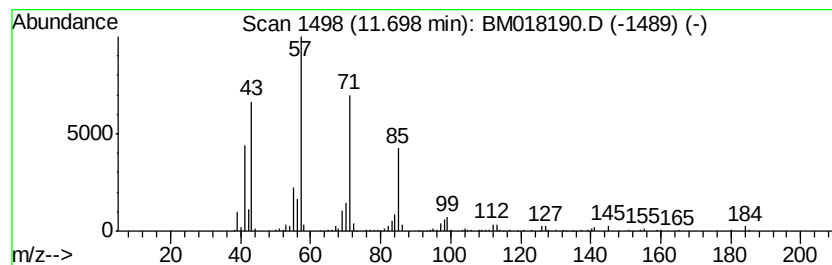
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.70 | 6.77 ng/ul | 703404 | Naphthalene-d8   | 10.27 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Tridecane                | 184 | C13H28  | 000629-50-5 | 94   |
| 2    |    | Tridecane                | 184 | C13H28  | 000629-50-5 | 81   |
| 3    |    | Tetradecane              | 198 | C14H30  | 000629-59-4 | 80   |
| 4    |    | Octane, 2,4,6-trimethyl- | 156 | C11H24  | 062016-37-9 | 80   |
| 5    |    | Decane, 2,3,5-trimethyl- | 184 | C13H28  | 062238-11-3 | 78   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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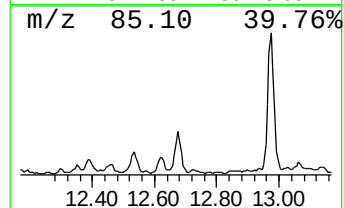
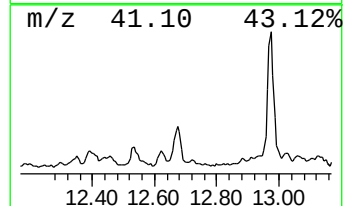
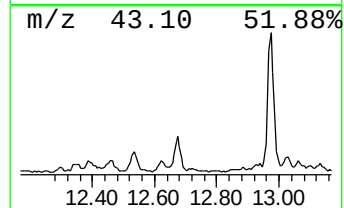
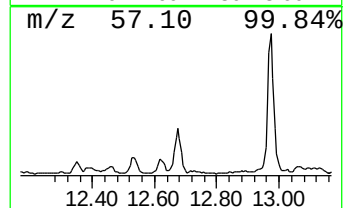
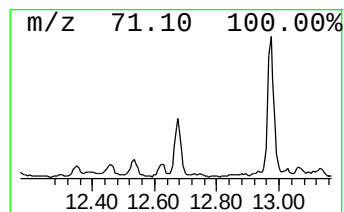
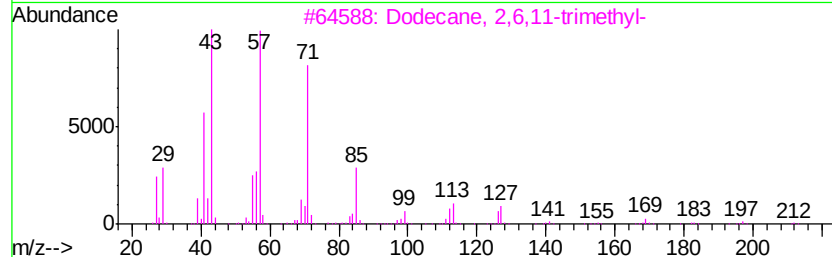
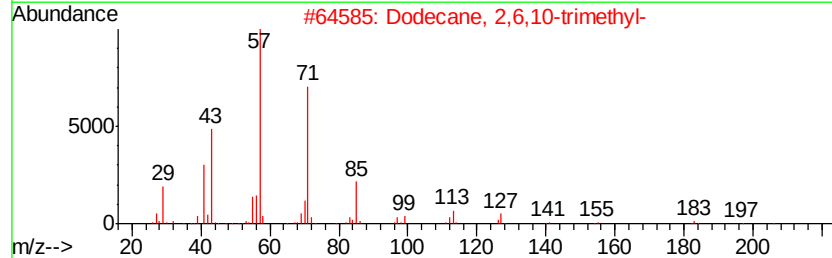
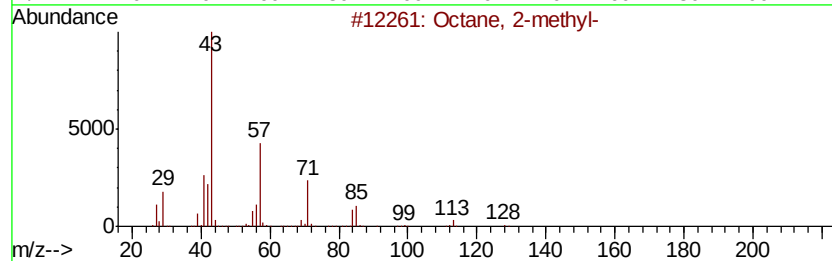
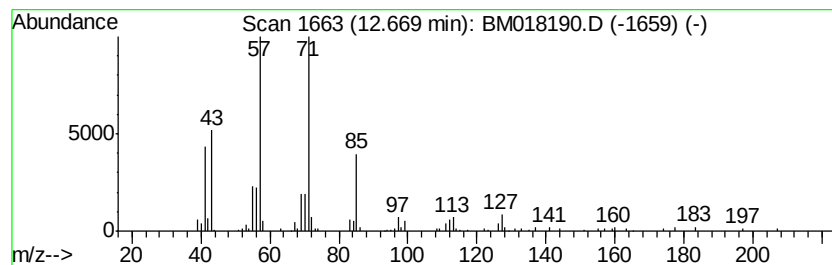
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 18

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.67 | 3.23 ng/ul | 348495 | Acenaphthene-d10 | 14.16 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Octane, 2-methyl-           | 128 | C9H20   | 003221-61-2 | 87   |
| 2    |    | Dodecane, 2,6,10-trimethyl- | 212 | C15H32  | 003891-98-3 | 86   |
| 3    |    | Dodecane, 2,6,11-trimethyl- | 212 | C15H32  | 031295-56-4 | 83   |
| 4    |    | Dodecane, 2,7,10-trimethyl- | 212 | C15H32  | 074645-98-0 | 80   |
| 5    |    | Dodecane, 2,6,10-trimethyl- | 212 | C15H32  | 003891-98-3 | 68   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
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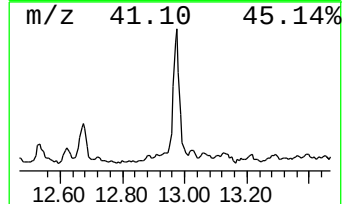
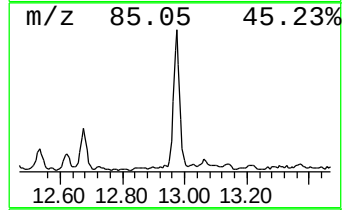
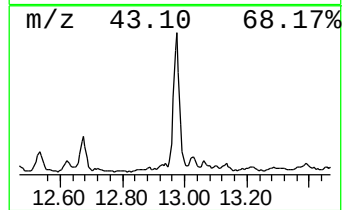
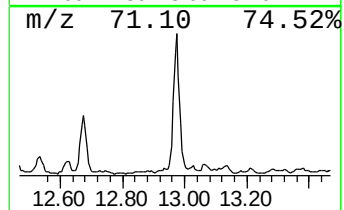
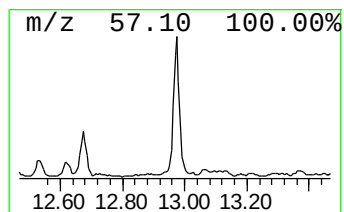
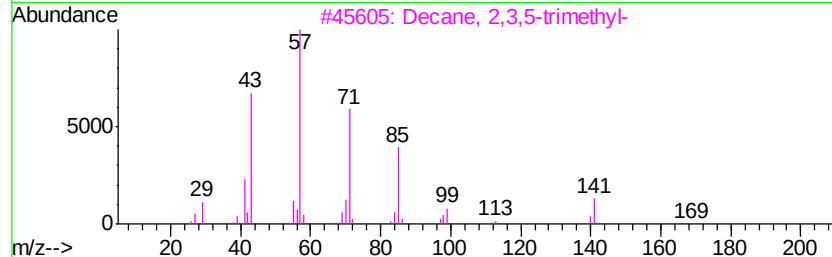
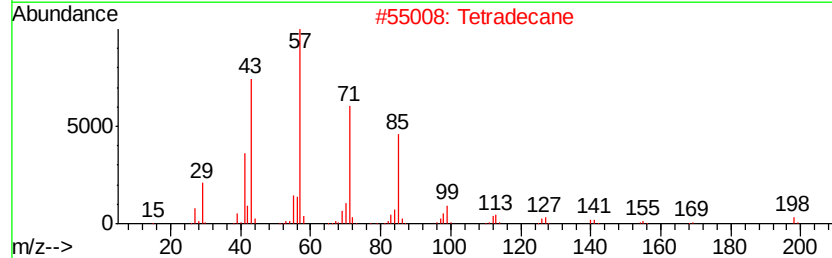
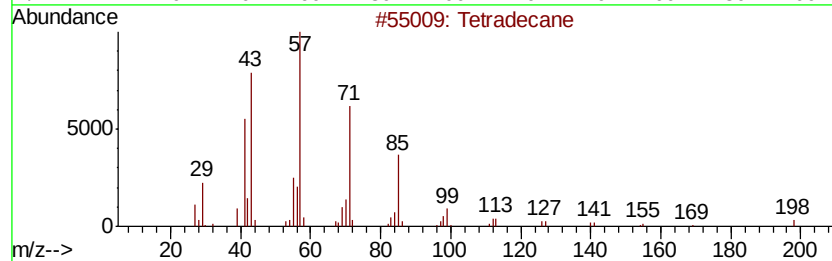
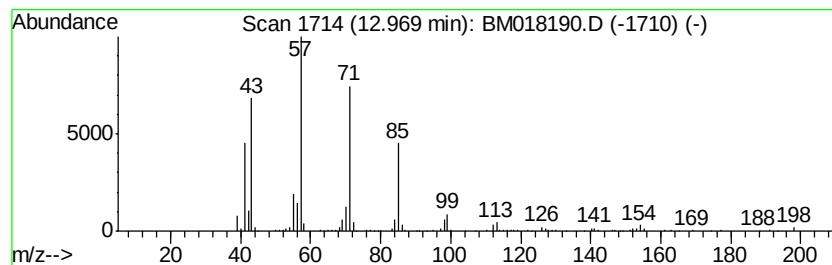
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 5

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.97 | 8.22 ng/ul | 886021 | Acenaphthene-d10 | 14.16 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Tetradecane               | 198 | C14H30  | 000629-59-4 | 87   |
| 2    |    | Tetradecane               | 198 | C14H30  | 000629-59-4 | 83   |
| 3    |    | Decane, 2,3,5-trimethyl-  | 184 | C13H28  | 062238-11-3 | 83   |
| 4    |    | 1,3-Dimethylcyclopentanol | 114 | C7H14O  | 019550-46-0 | 72   |
| 5    |    | Hexadecane                | 226 | C16H34  | 000544-76-3 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
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Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

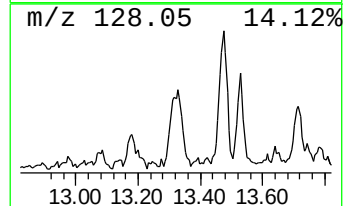
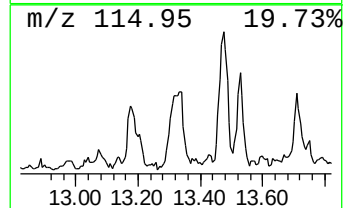
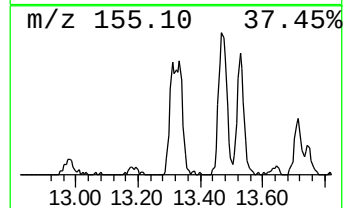
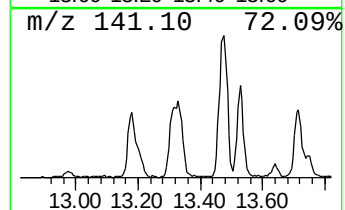
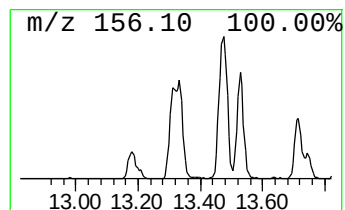
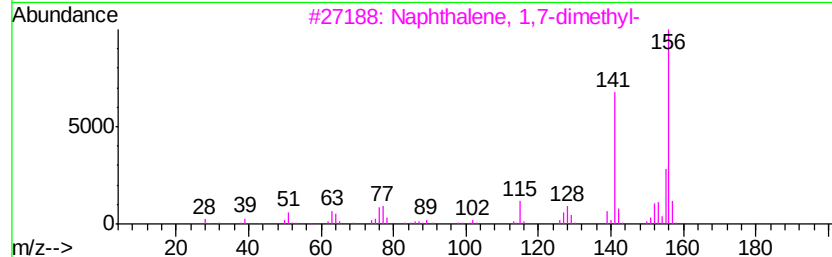
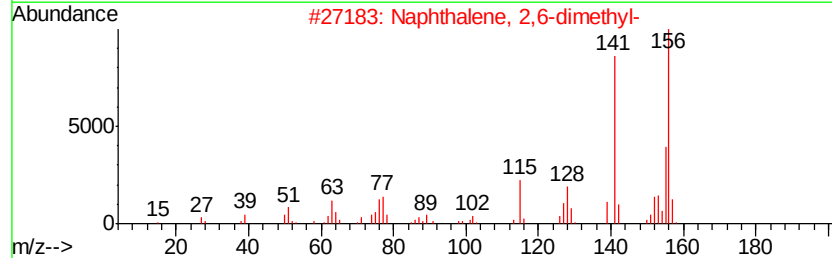
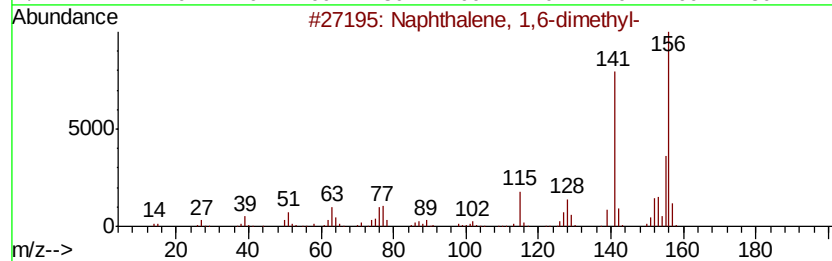
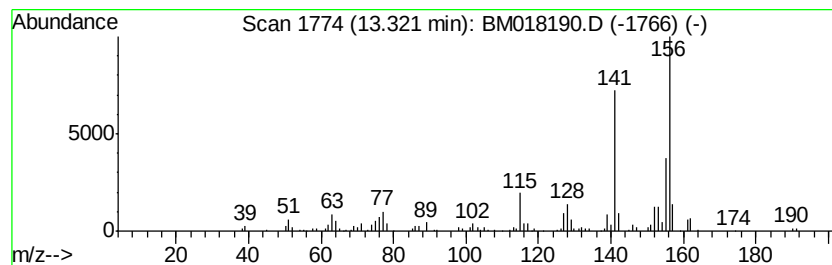
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ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Naphthalene, 1,6-dimethyl- Concentration Rank 24

| R.T.    | EstConc    | Area                       | Relative to ISTD | R.T.           |
|---------|------------|----------------------------|------------------|----------------|
| 13.32   | 2.61 ng/ul | 280876                     | Acenaphthene-d10 | 14.16          |
| Hit# of | 5          | Tentative ID               | MW MolForm       | CAS# Qual      |
| 1       |            | Naphthalene, 1,6-dimethyl- | 156 C12H12       | 000575-43-9 97 |
| 2       |            | Naphthalene, 2,6-dimethyl- | 156 C12H12       | 000581-42-0 96 |
| 3       |            | Naphthalene, 1,7-dimethyl- | 156 C12H12       | 000575-37-1 96 |
| 4       |            | Naphthalene, 2,7-dimethyl- | 156 C12H12       | 000582-16-1 96 |
| 5       |            | Naphthalene, 2,7-dimethyl- | 156 C12H12       | 000582-16-1 96 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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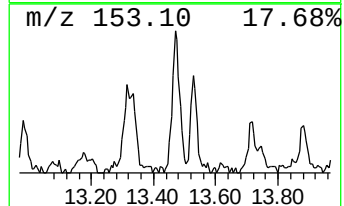
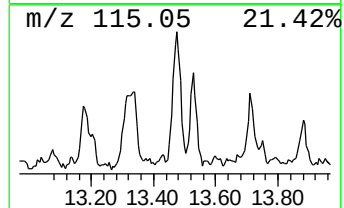
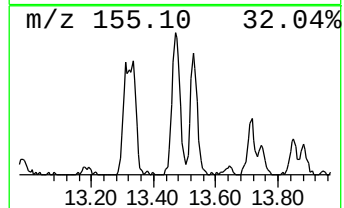
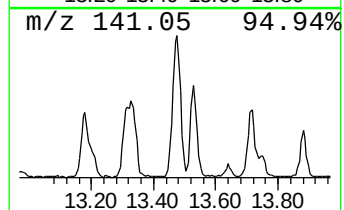
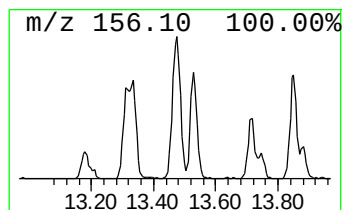
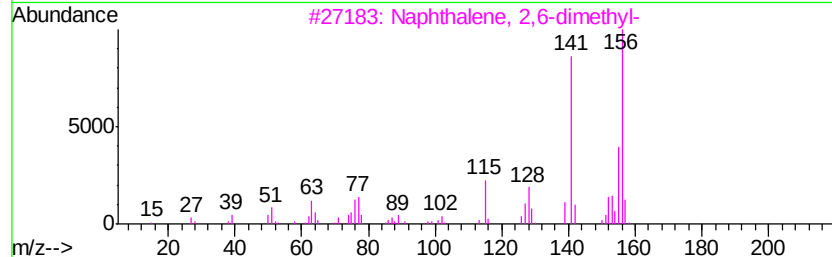
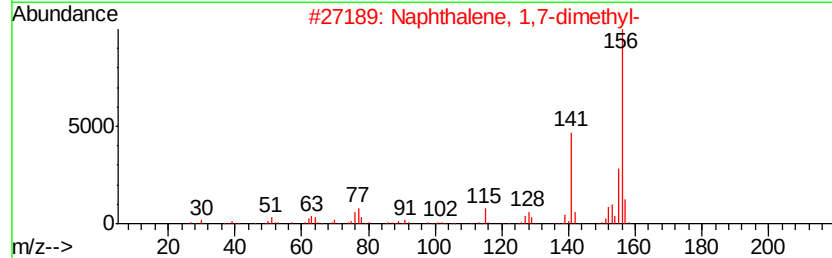
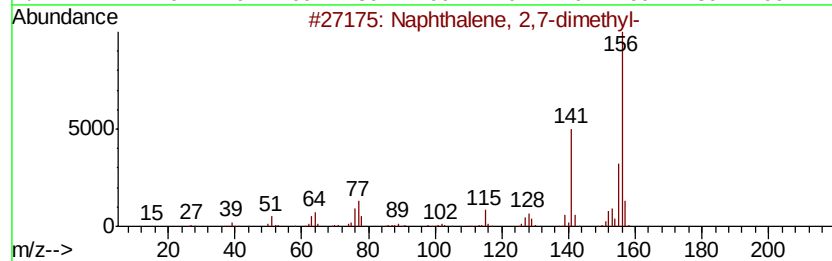
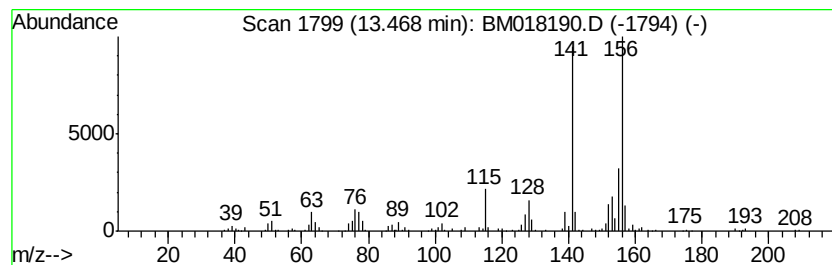
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Naphthalene, 2,7-dimethyl- Concentration Rank 12

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.47 | 4.79 ng/ul | 516715 | Acenaphthene-d10 | 14.16 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 97   |
| 2    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 97   |
| 3    |    | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 96   |
| 4    |    | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 96   |
| 5    |    | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 96   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
Quant Title : SVOA CALIBRATION

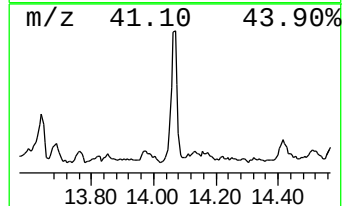
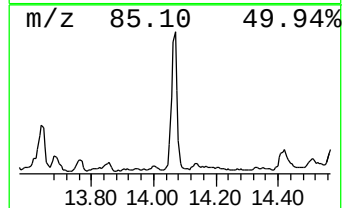
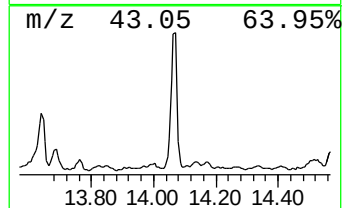
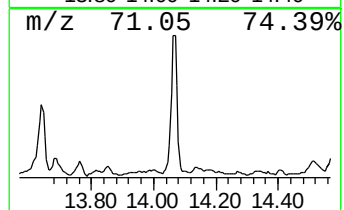
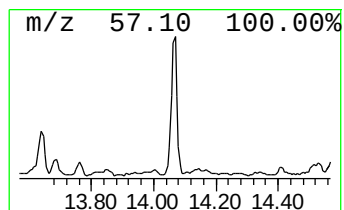
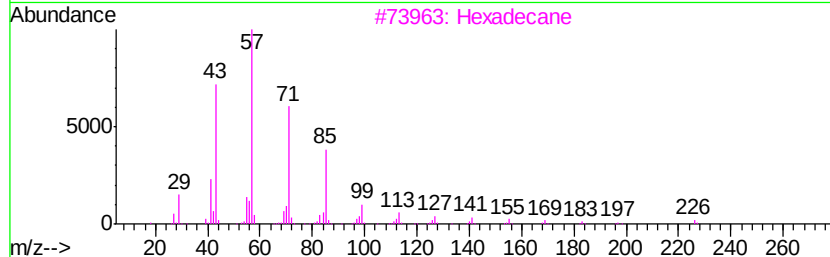
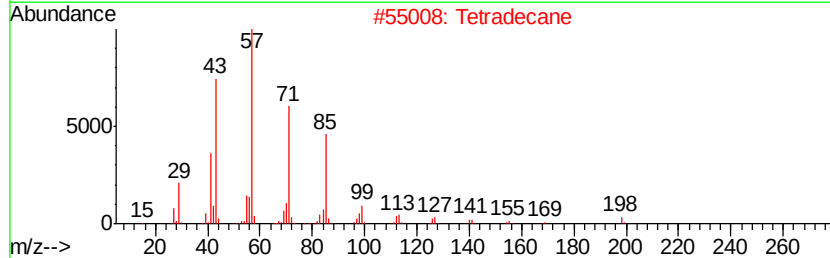
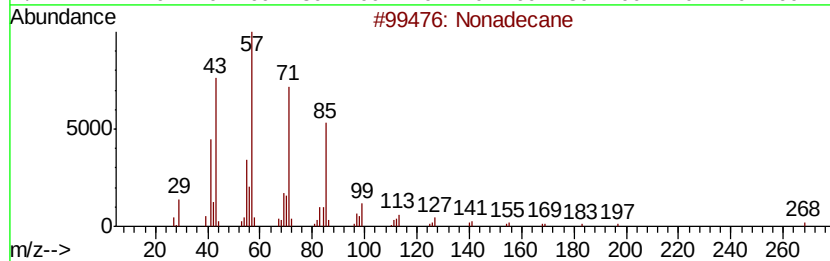
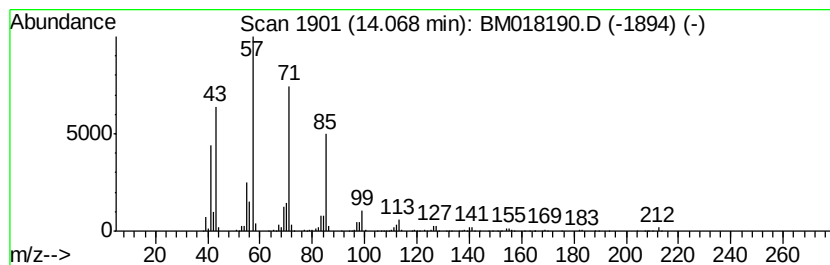
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 3

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 14.07 | 9.67 ng/ul | 1041880 | Acenaphthene-d10 | 14.16 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Nonadecane              | 268 | C19H40  | 000629-92-5 | 91   |
| 2    |    | Tetradecane             | 198 | C14H30  | 000629-59-4 | 91   |
| 3    |    | Hexadecane              | 226 | C16H34  | 000544-76-3 | 90   |
| 4    |    | Undecane, 5,7-dimethyl- | 184 | C13H28  | 017312-83-3 | 80   |
| 5    |    | Heptadecane             | 240 | C17H36  | 000629-78-7 | 78   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BE7Z4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
Quant Title : SVOA CALIBRATION

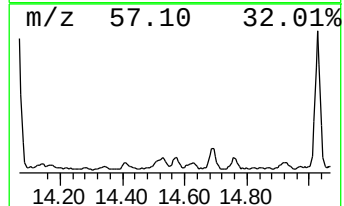
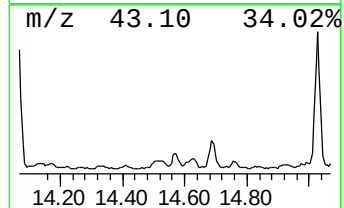
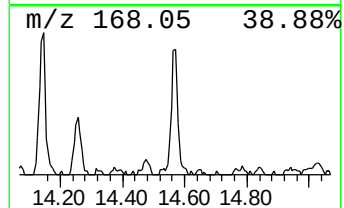
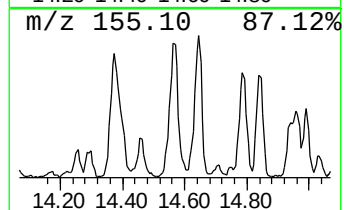
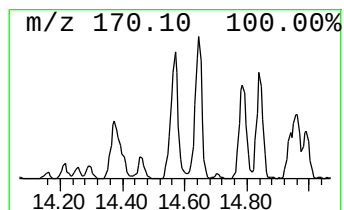
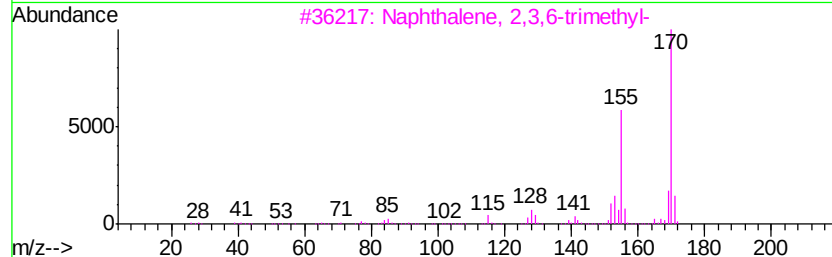
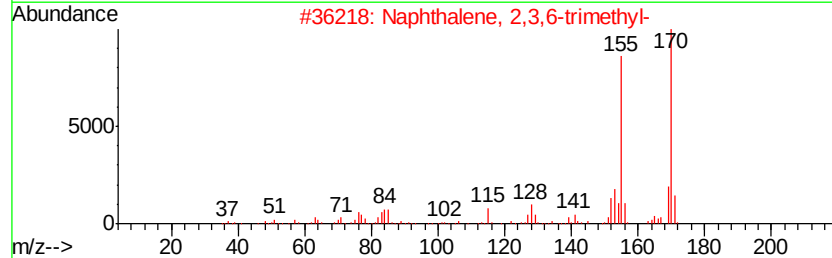
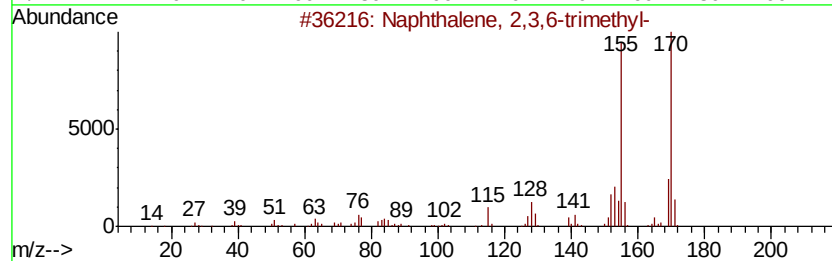
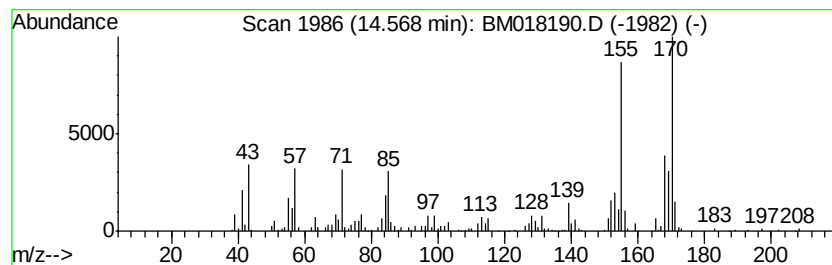
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Naphthalene, 2,3,6-trimethyl- Concentration Rank 20

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.57 | 2.95 ng/ul | 317769 | Acenaphthene-d10 | 14.16 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 95   |
| 2    |    | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 94   |
| 3    |    | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 93   |
| 4    |    | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 90   |
| 5    |    | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 76   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BE7Z4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
Quant Title : SVOA CALIBRATION

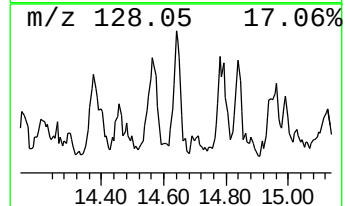
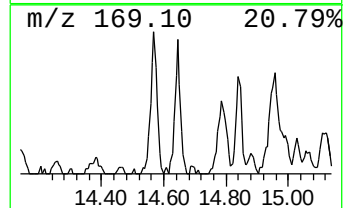
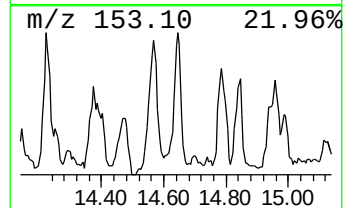
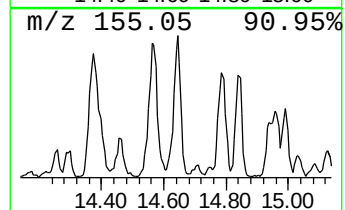
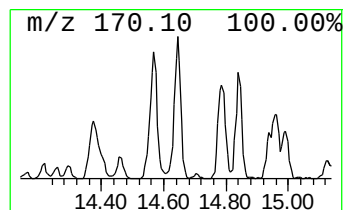
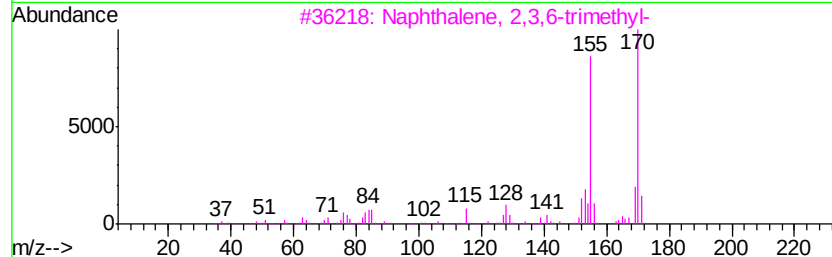
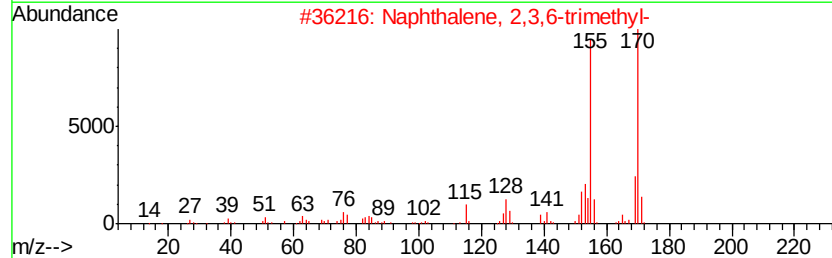
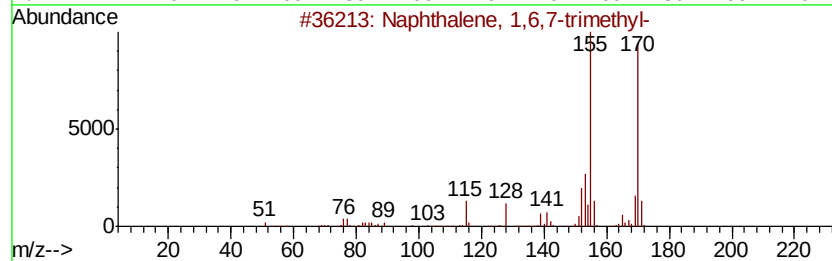
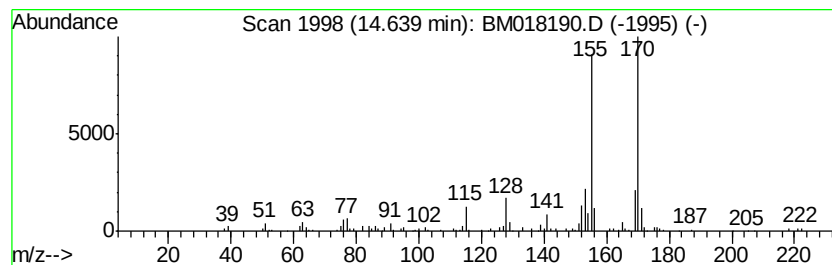
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 Naphthalene, 1,6,7-trimethyl- Concentration Rank 29

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.64 | 2.32 ng/ul | 250241 | Acenaphthene-d10 | 14.16 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 96   |
| 2    |    | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 95   |
| 3    |    | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 94   |
| 4    |    | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 94   |
| 5    |    | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 94   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BE7Z4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
Quant Title : SVOA CALIBRATION

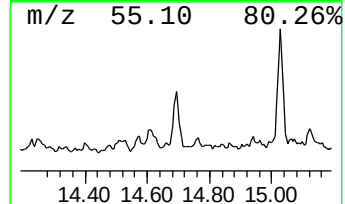
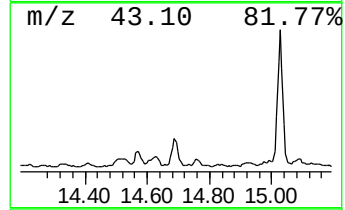
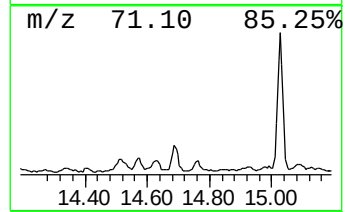
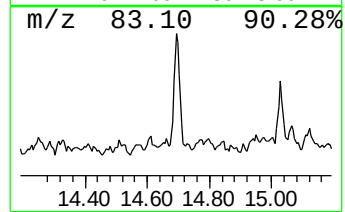
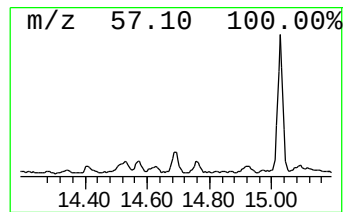
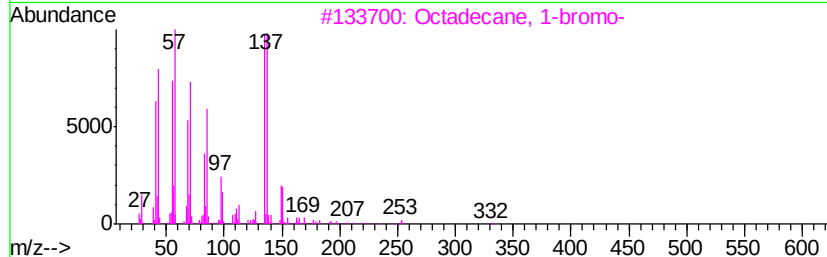
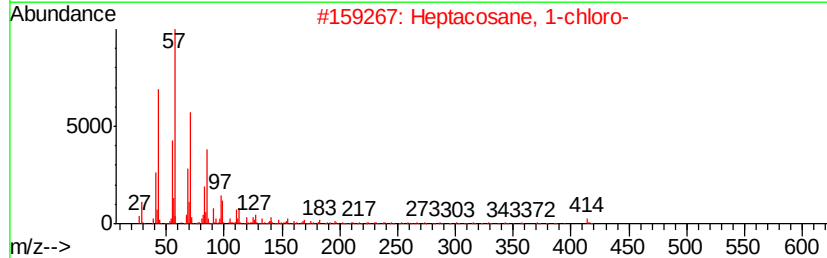
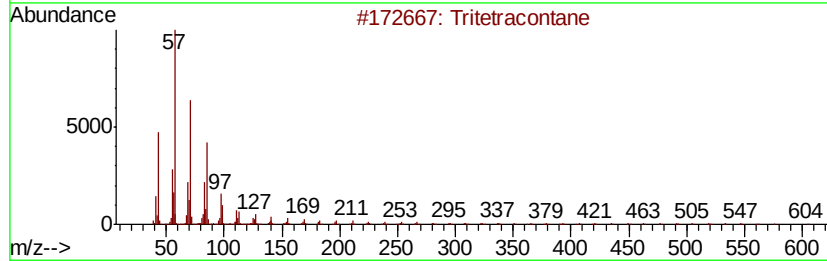
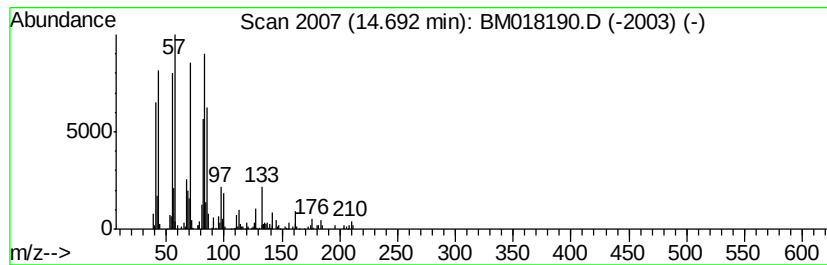
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 19

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.69 | 3.17 ng/ul | 341530 | Acenaphthene-d10 | 14.16 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm  | CAS#        | Qual |
|------|----|---|------------------------|-----|----------|-------------|------|
| 1    |    |   | Tritetracontane        | 605 | C43H88   | 007098-21-7 | 49   |
| 2    |    |   | Heptacosane, 1-chloro- | 414 | C27H55Cl | 062016-79-9 | 46   |
| 3    |    |   | Octadecane, 1-bromo-   | 332 | C18H37Br | 000112-89-0 | 46   |
| 4    |    |   | Octadecane, 1-bromo-   | 332 | C18H37Br | 000112-89-0 | 46   |
| 5    |    |   | Tetratetracontane      | 619 | C44H90   | 007098-22-8 | 46   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BE7Z4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
Quant Title : SVOA CALIBRATION

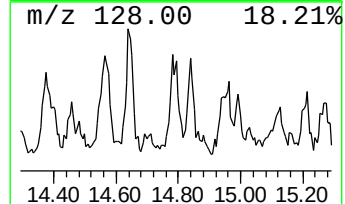
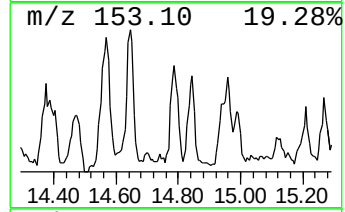
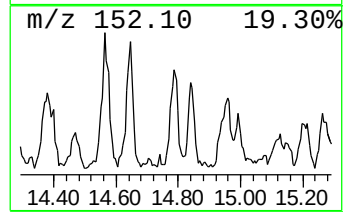
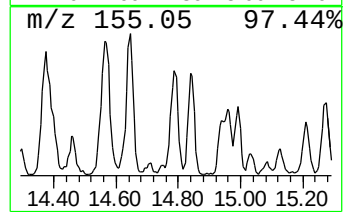
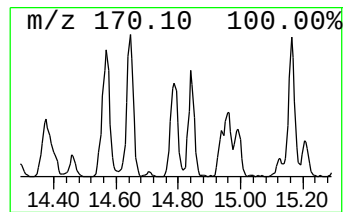
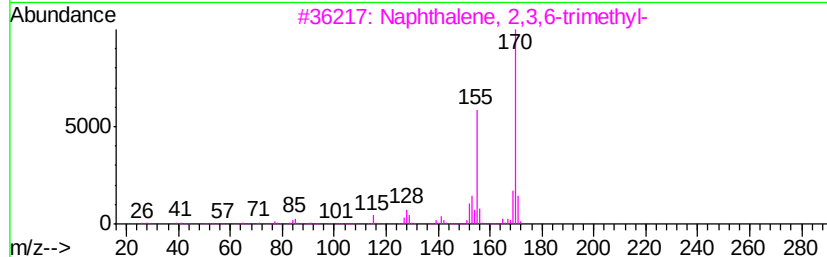
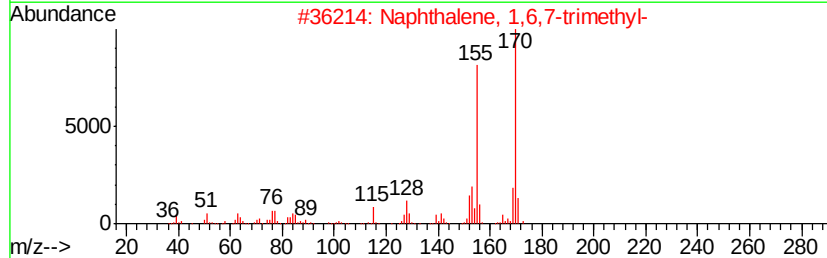
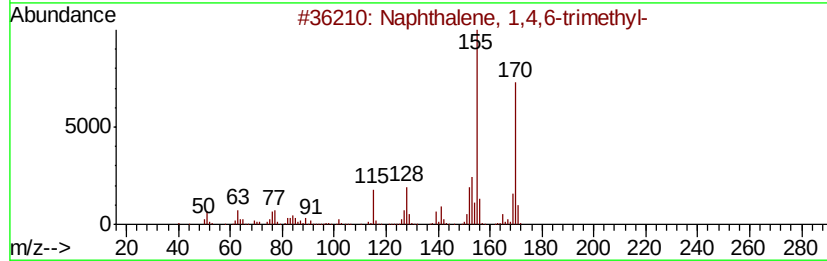
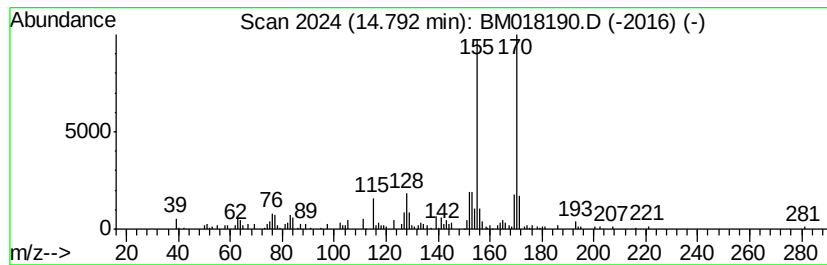
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Naphthalene, 1,4,6-trimethyl- Concentration Rank 21

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.79 | 2.90 ng/ul | 313080 | Acenaphthene-d10 | 14.16 |

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BE7Z4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
Quant Title : SVOA CALIBRATION

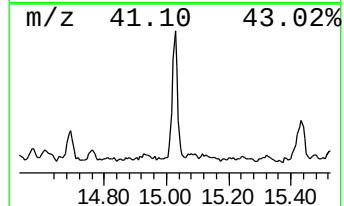
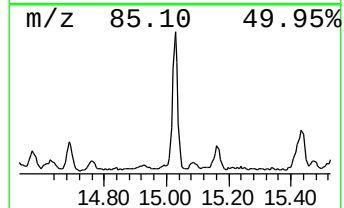
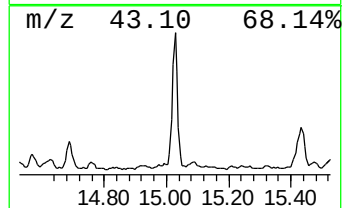
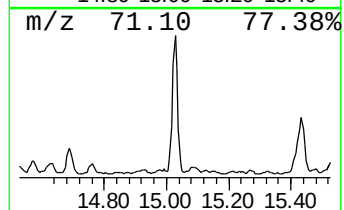
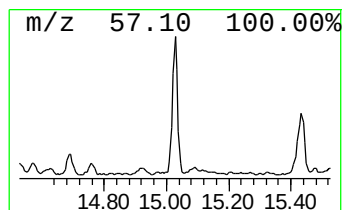
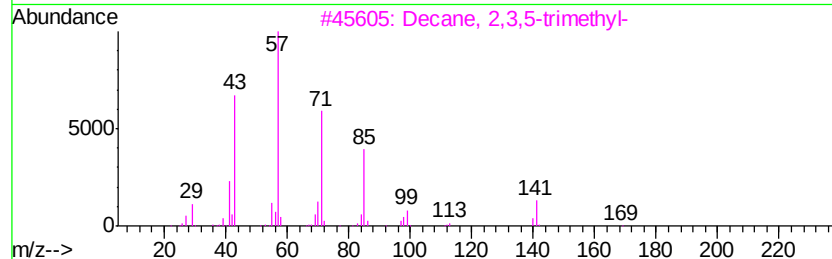
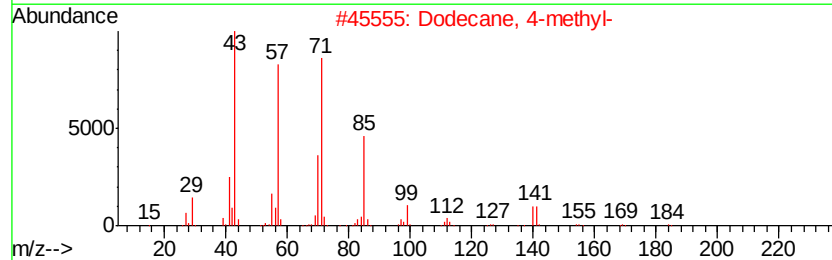
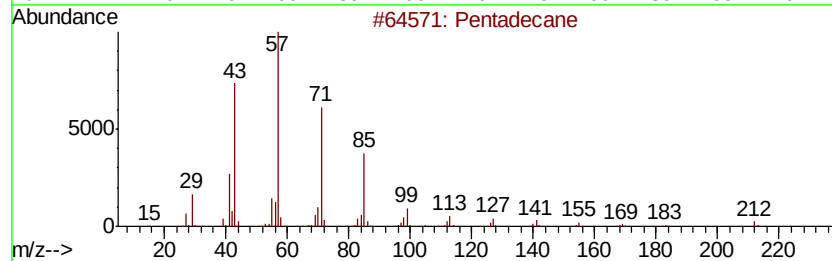
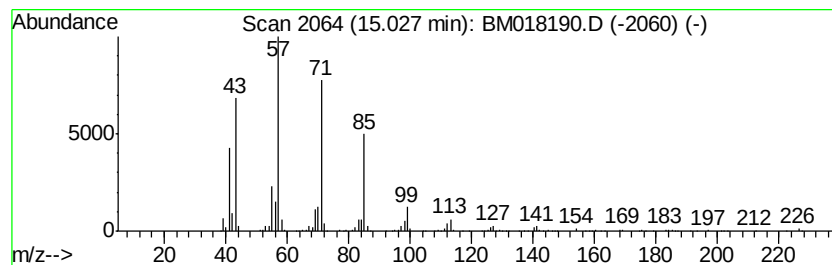
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.03 | 8.81 ng/ul | 950009 | Acenaphthene-d10 | 14.16 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentadecane              | 212 | C15H32  | 000629-62-9 | 91   |
| 2    |    |   | Dodecane, 4-methyl-      | 184 | C13H28  | 006117-97-1 | 90   |
| 3    |    |   | Decane, 2,3,5-trimethyl- | 184 | C13H28  | 062238-11-3 | 90   |
| 4    |    |   | Pentadecane              | 212 | C15H32  | 000629-62-9 | 90   |
| 5    |    |   | Hexadecane               | 226 | C16H34  | 000544-76-3 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

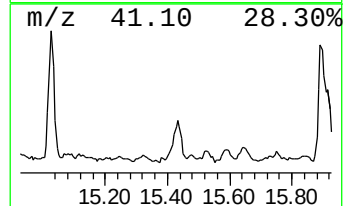
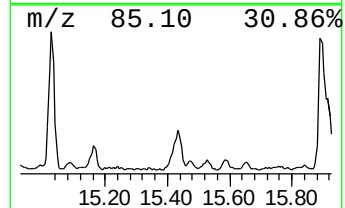
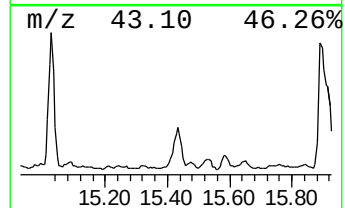
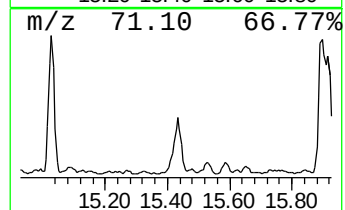
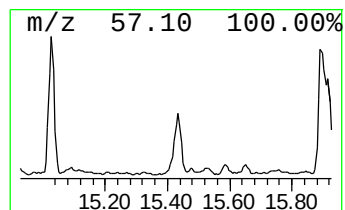
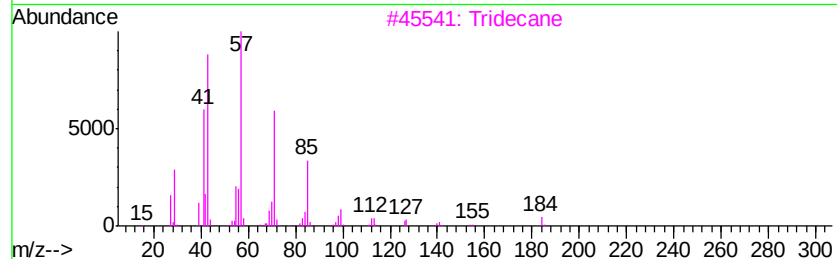
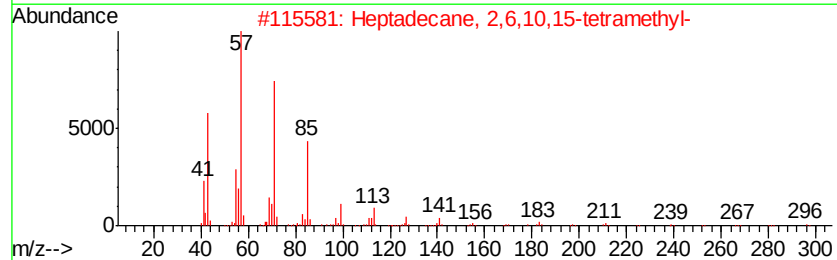
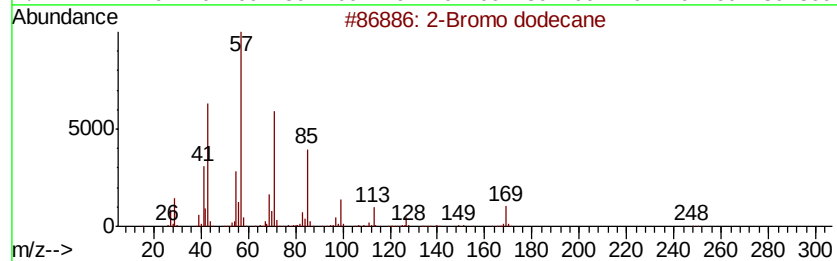
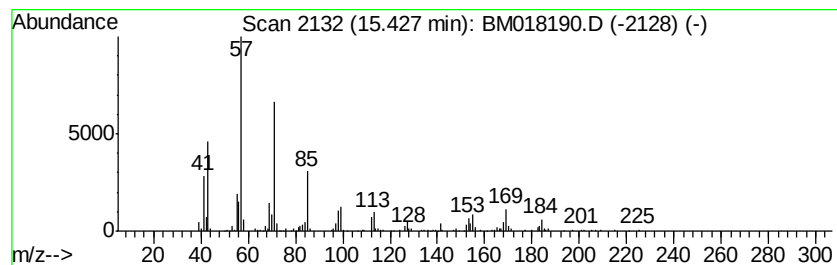
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BNA\_M  
ClientSampleId :  
BE7Z4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 2-Bromo dodecane Concentration Rank 14

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.           |
|---------|-------------------------------------|--------------|------------------|----------------|
| 15.43   | 4.51 ng/ul                          | 486494       | Acenaphthene-d10 | 14.16          |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual      |
| 1       | 2-Bromo dodecane                    | 248          | C12H25Br         | 013187-99-0 86 |
| 2       | Heptadecane, 2,6,10,15-tetramethyl- | 296          | C21H44           | 054833-48-6 74 |
| 3       | Tridecane                           | 184          | C13H28           | 000629-50-5 74 |
| 4       | Heptacosane                         | 380          | C27H56           | 000593-49-7 72 |
| 5       | Hexadecane, 2,6,10,14-tetramethyl-  | 282          | C20H42           | 000638-36-8 72 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BE7Z4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
Quant Title : SVOA CALIBRATION

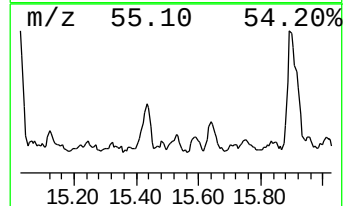
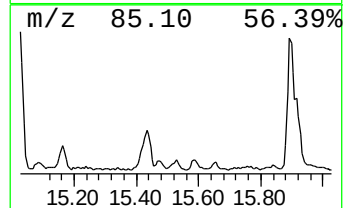
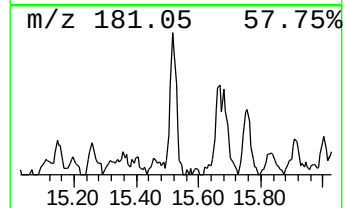
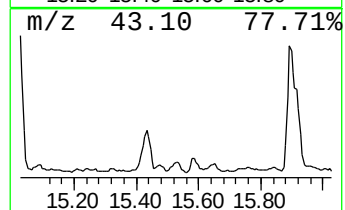
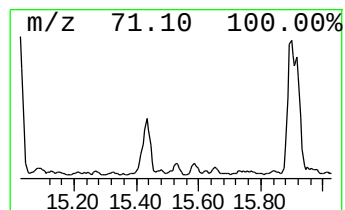
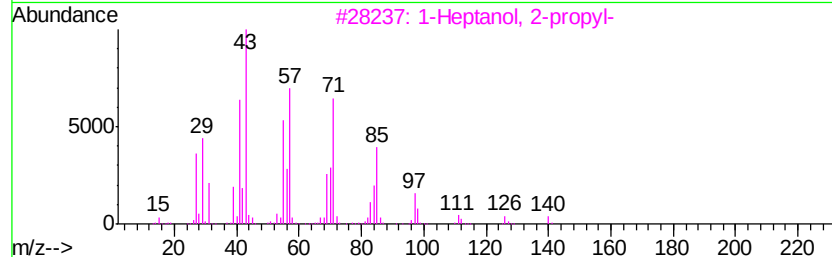
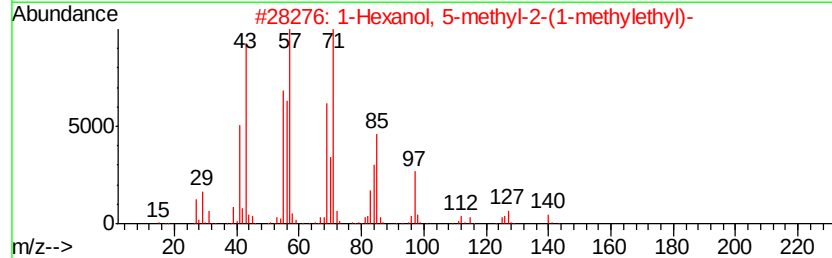
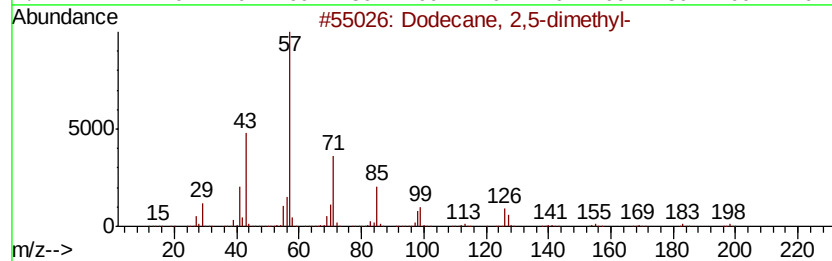
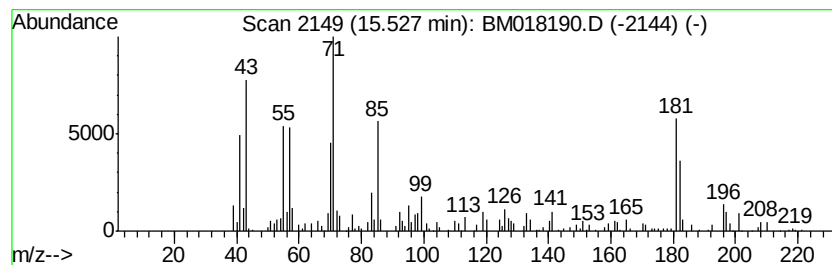
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 22

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.53 | 2.84 ng/ul | 305654 | Acenaphthene-d10 | 14.16 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Dodecane, 2,5-dimethyl-             | 198 | C14H30  | 056292-65-0 | 52   |
| 2    |    |   | 1-Hexanol, 5-methyl-2-(1-methyle... | 158 | C10H22O | 002051-33-4 | 47   |
| 3    |    |   | 1-Heptanol, 2-propyl-               | 158 | C10H22O | 010042-59-8 | 47   |
| 4    |    |   | Tridecane, 7-hexyl-                 | 268 | C19H40  | 007225-66-3 | 43   |
| 5    |    |   | Decane, 3,3,6-trimethyl-            | 184 | C13H28  | 062338-14-1 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BE7Z4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
Quant Title : SVOA CALIBRATION

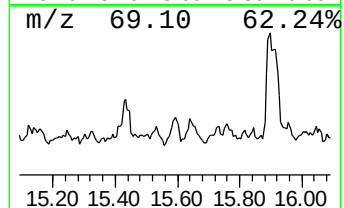
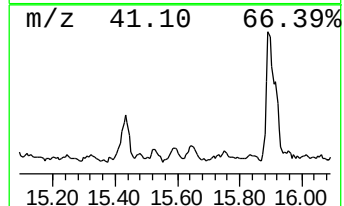
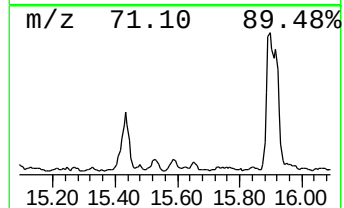
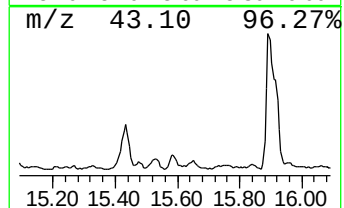
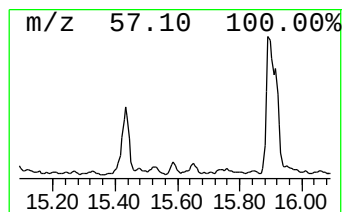
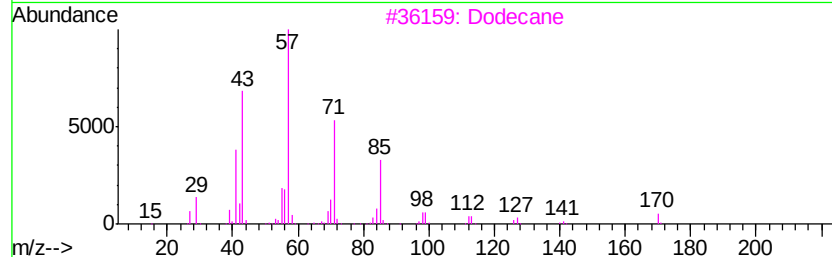
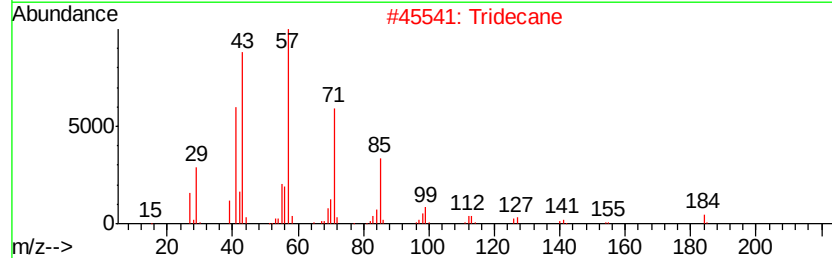
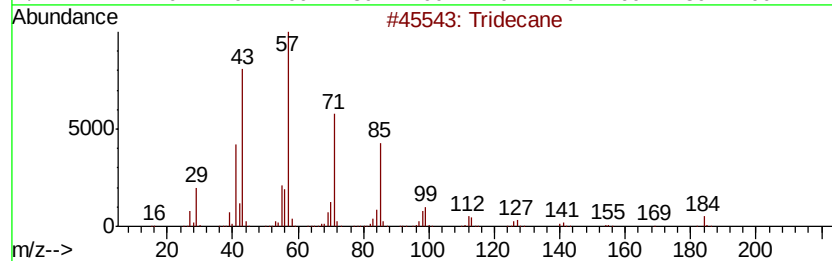
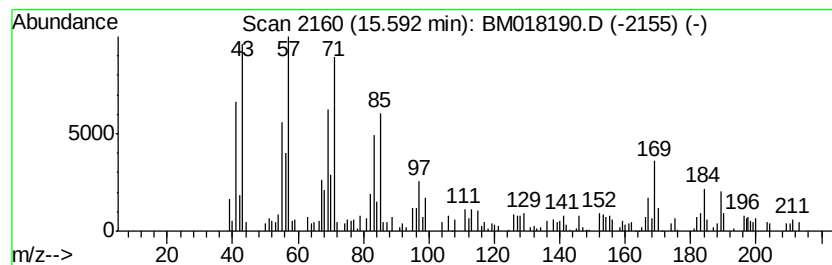
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 30

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.59 | 2.24 ng/ul | 222725 | Phenanthrene-d10 | 16.92 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Tridecane    | 184 | C13H28  | 000629-50-5 | 86   |
| 2    |    | Tridecane    | 184 | C13H28  | 000629-50-5 | 70   |
| 3    |    | Dodecane     | 170 | C12H26  | 000112-40-3 | 64   |
| 4    |    | Tridecane    | 184 | C13H28  | 000629-50-5 | 64   |
| 5    |    | Tridecane    | 184 | C13H28  | 000629-50-5 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BE7Z4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
Quant Title : SVOA CALIBRATION

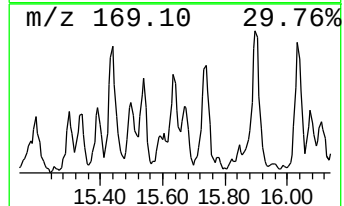
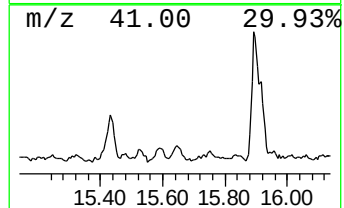
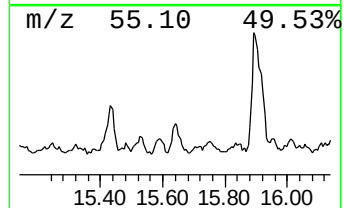
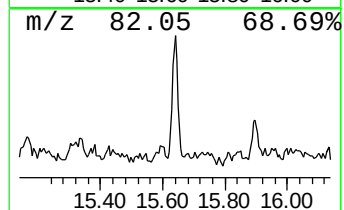
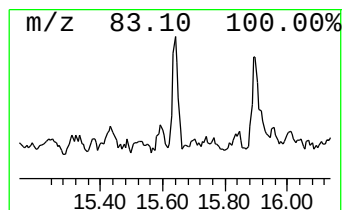
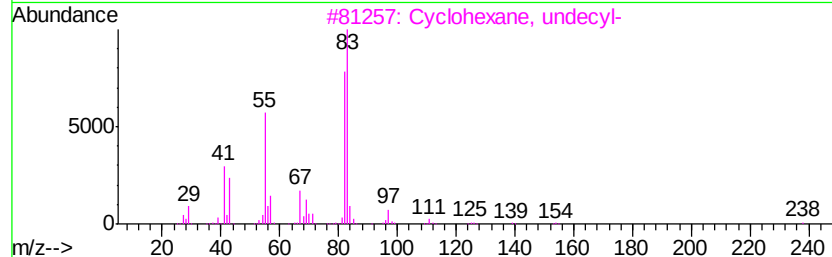
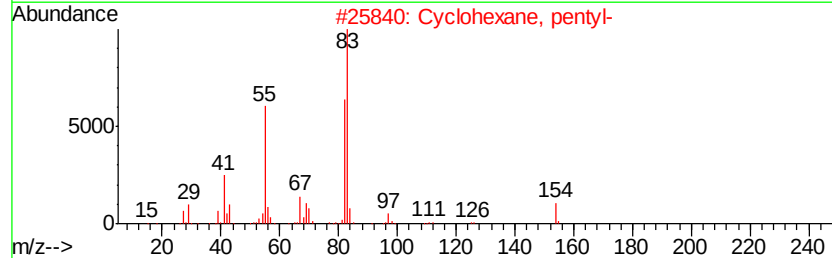
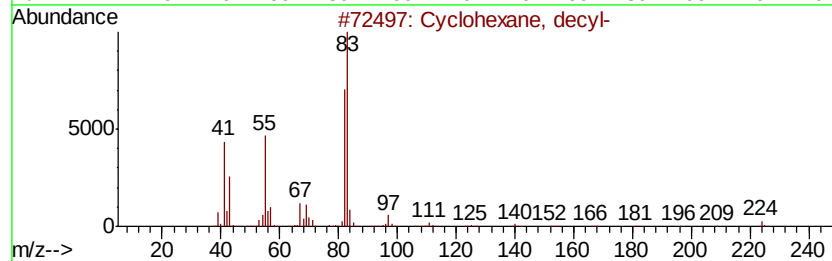
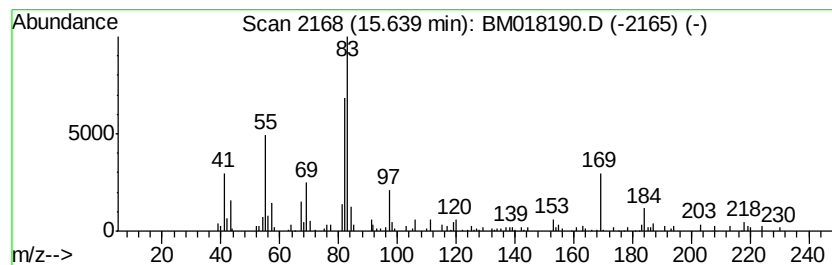
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 21 (DEL) Alkane: Cyclic15.64 Concentration Rank 31

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.64 | 2.21 ng/ul | 219781 | Phenanthrene-d10 | 16.92 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, decyl-            | 224 | C16H32  | 001795-16-0 | 58   |
| 2    |    | Cyclohexane, pentyl-           | 154 | C11H22  | 004292-92-6 | 52   |
| 3    |    | Cyclohexane, undecyl-          | 238 | C17H34  | 054105-66-7 | 50   |
| 4    |    | Cyclohexane, (4-methylpentyl)- | 168 | C12H24  | 061142-20-9 | 50   |
| 5    |    | Cyclohexane, pentyl-           | 154 | C11H22  | 004292-92-6 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BE7Z4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
Quant Title : SVOA CALIBRATION

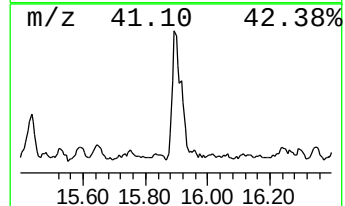
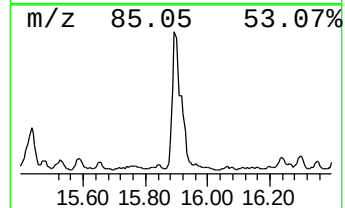
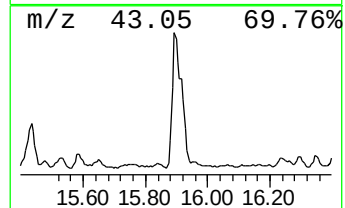
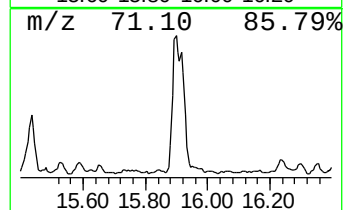
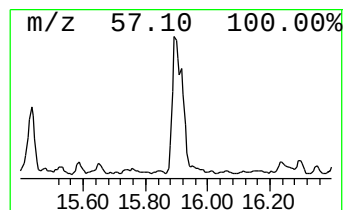
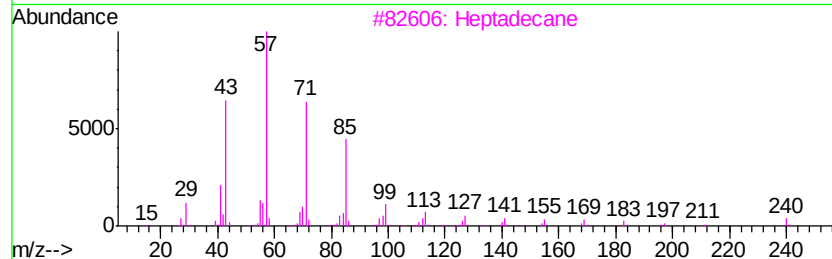
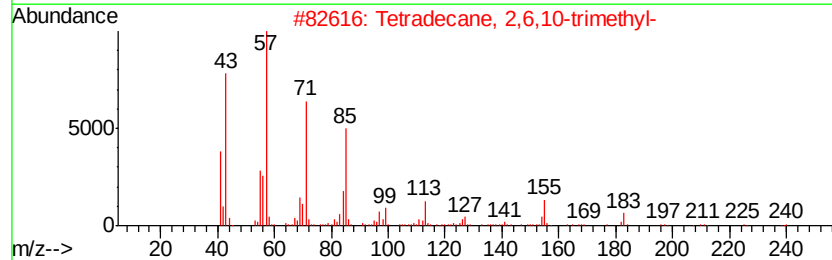
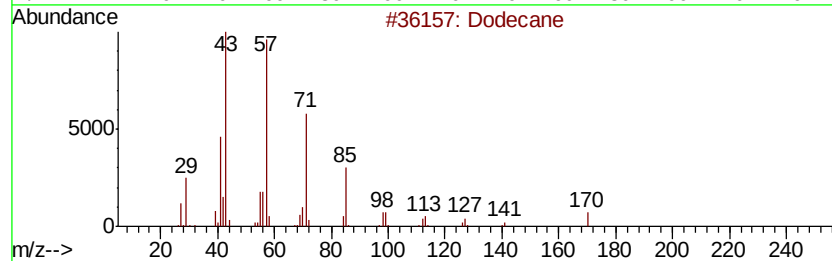
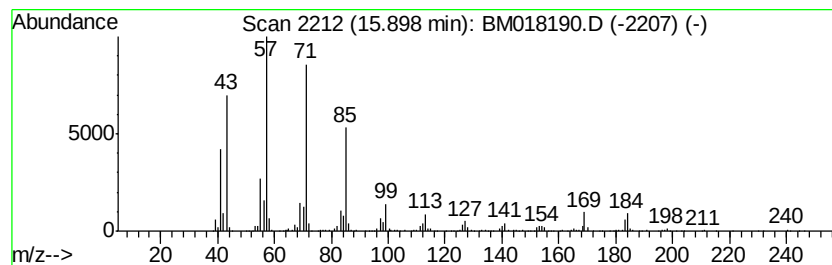
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 1

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.90 | 18.63 ng/ul | 1853810 | Phenanthrene-d10 | 16.92 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Dodecane                       | 170 | C12H26  | 000112-40-3 | 90   |
| 2    |    | Tetradecane, 2,6,10-trimethyl- | 240 | C17H36  | 014905-56-7 | 90   |
| 3    |    | Heptadecane                    | 240 | C17H36  | 000629-78-7 | 83   |
| 4    |    | Tetradecane                    | 198 | C14H30  | 000629-59-4 | 83   |
| 5    |    | Pentadecane                    | 212 | C15H32  | 000629-62-9 | 83   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BE7Z4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
Quant Title : SVOA CALIBRATION

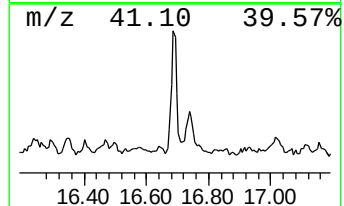
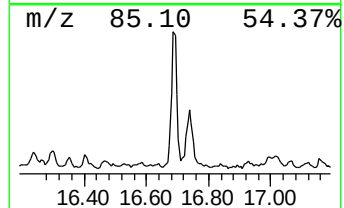
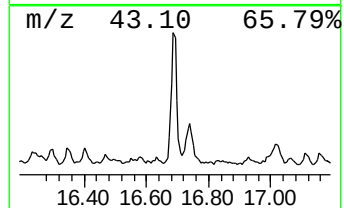
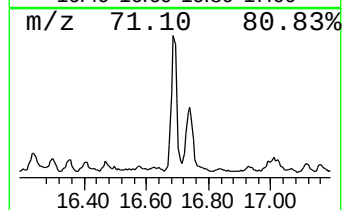
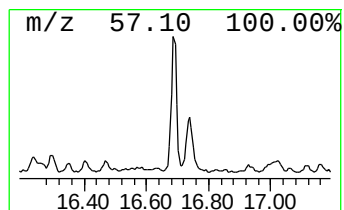
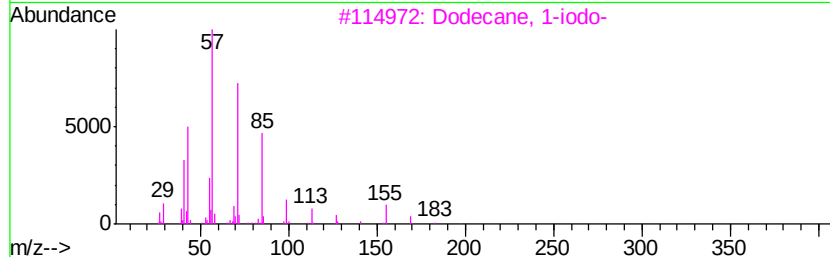
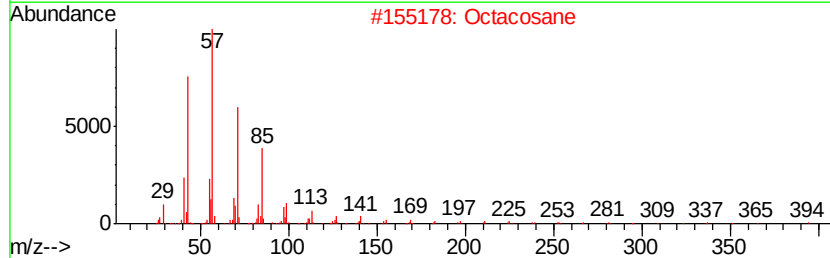
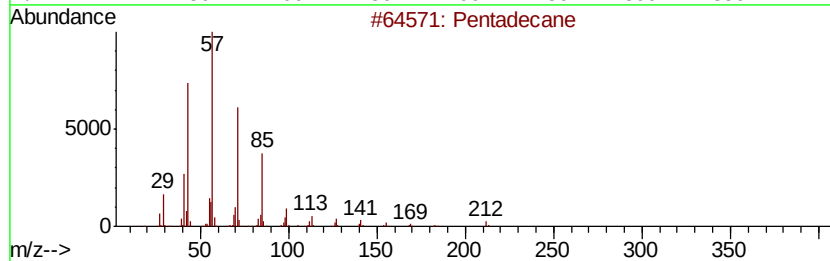
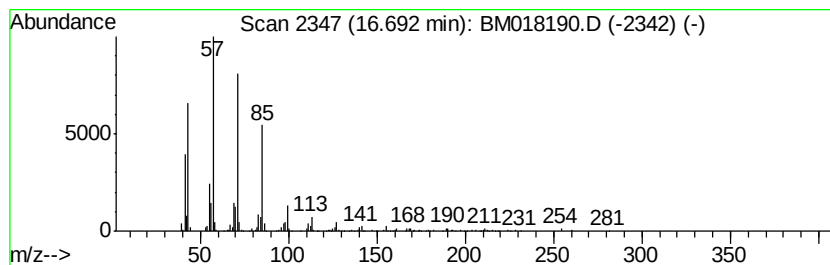
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.69 | 7.05 ng/ul | 701434 | Phenanthrene-d10 | 16.92 |

| Hit# | of | 5 | Tentative ID       | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------|-----|---------|-------------|------|
| 1    |    |   | Pentadecane        | 212 | C15H32  | 000629-62-9 | 94   |
| 2    |    |   | Octacosane         | 394 | C28H58  | 000630-02-4 | 90   |
| 3    |    |   | Dodecane, 1-iodo-  | 296 | C12H25I | 004292-19-7 | 86   |
| 4    |    |   | Tridecane, 1-iodo- | 310 | C13H27I | 035599-77-0 | 83   |
| 5    |    |   | Hexadecane         | 226 | C16H34  | 000544-76-3 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
 Data File : BM018190.D  
 Acq On : 15 Dec 2018 05:11  
 Operator : JU/SJ  
 Sample : J6238-01  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BE7Z4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
 Quant Title : SVOA CALIBRATION

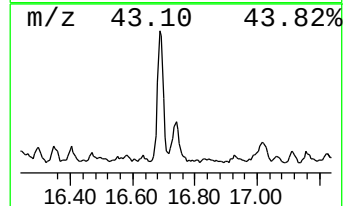
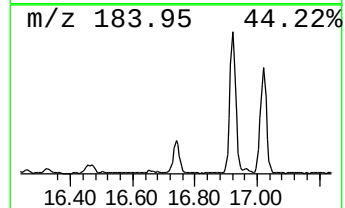
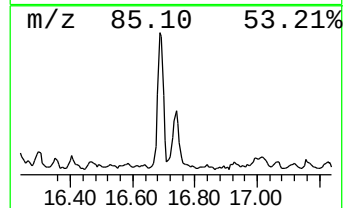
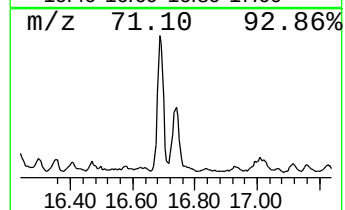
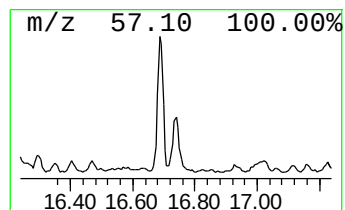
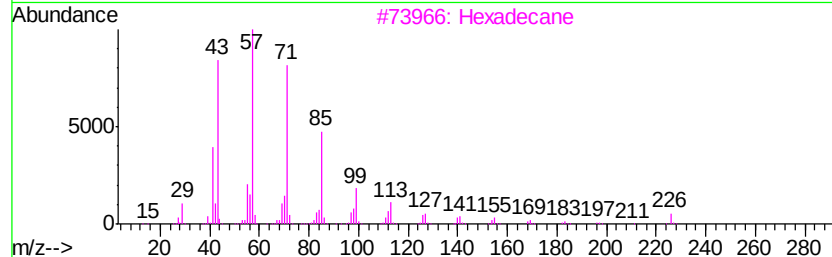
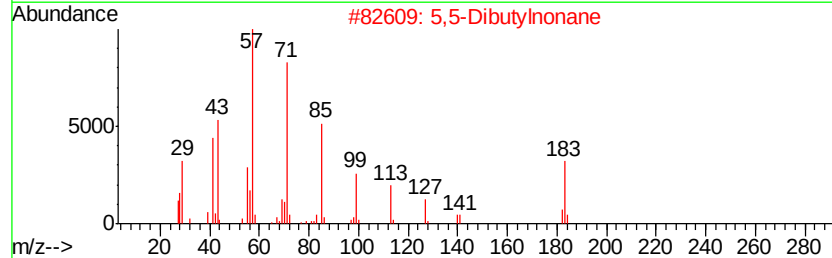
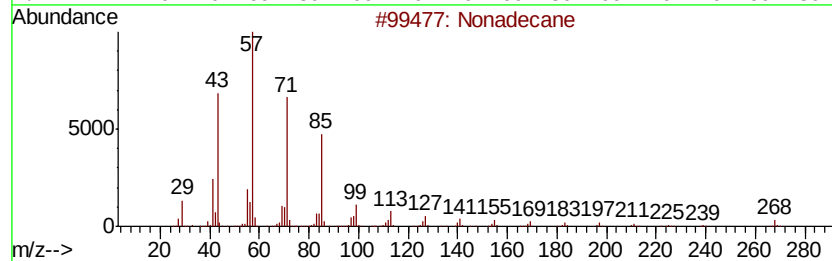
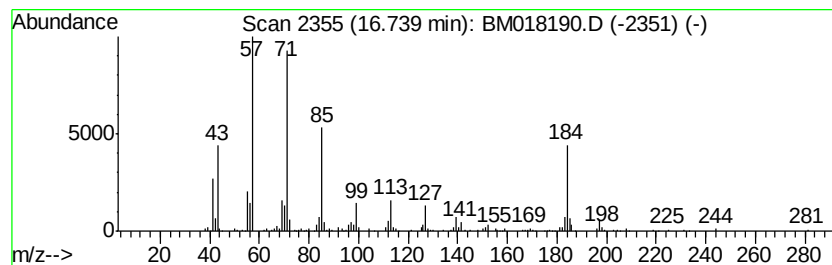
TIC Library : C:\DATABASE\NIST02.L  
 TIC Integration Parameters: LSCINT.P

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 Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.74 | 5.50 ng/ul | 547547 | Phenanthrene-d10 | 16.92 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Nonadecane                          | 268 | C19H40  | 000629-92-5 | 64   |
| 2    |    |   | 5,5-Dibutylnonane                   | 240 | C17H36  | 006008-17-9 | 64   |
| 3    |    |   | Hexadecane                          | 226 | C16H34  | 000544-76-3 | 64   |
| 4    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 64   |
| 5    |    |   | Pentadecane, 8-heptyl-              | 310 | C22H46  | 071005-15-7 | 59   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

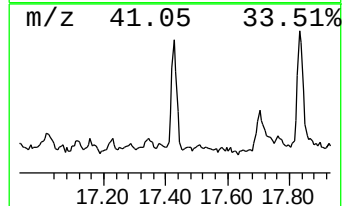
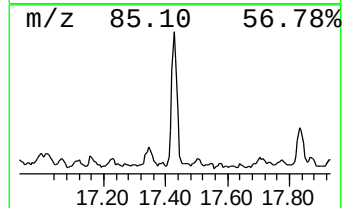
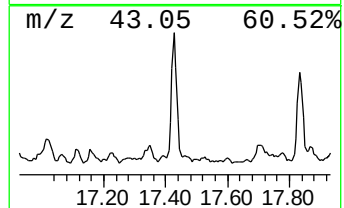
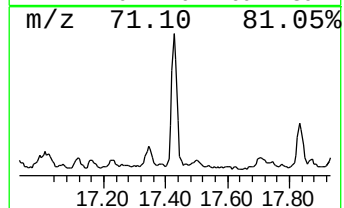
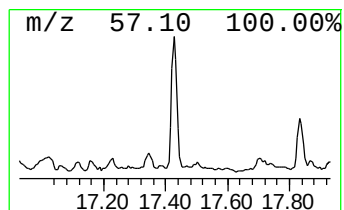
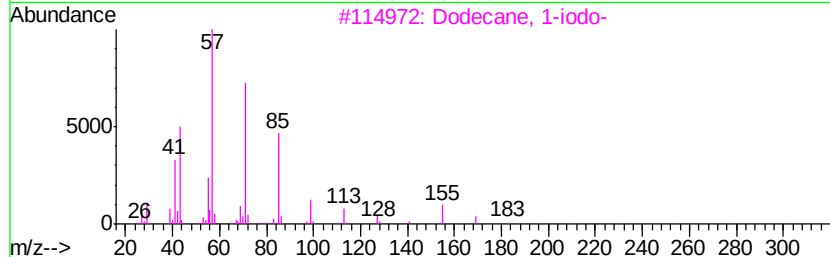
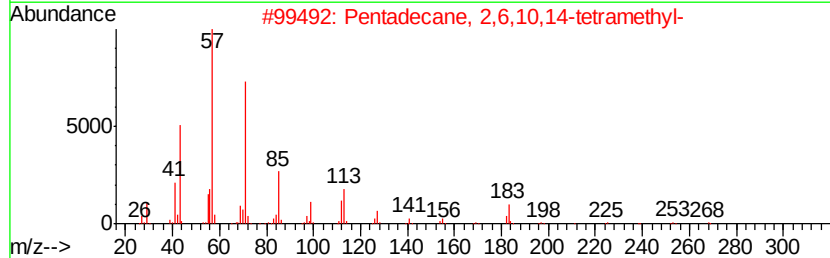
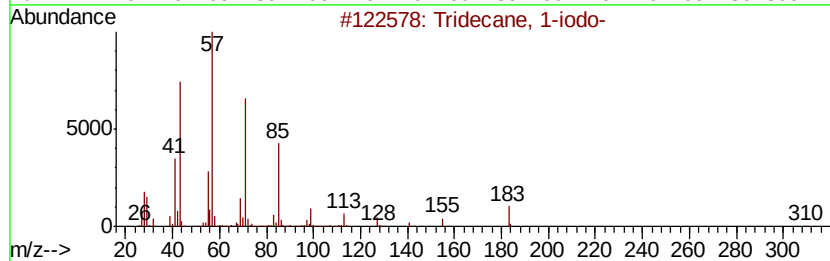
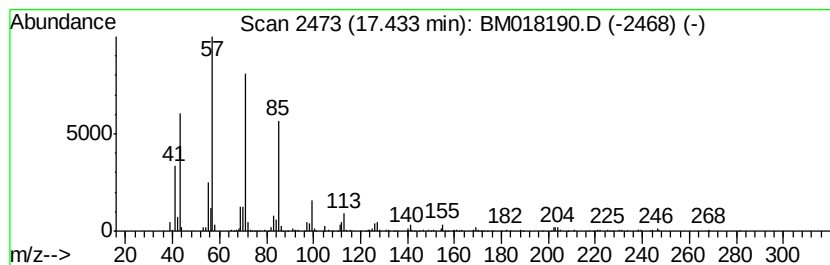
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ClientSampleId :  
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 25 Tridecane, 1-iodo- Concentration Rank 9

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.      |
|---------|-------------------------------------|--------------|------------------|-----------|
| 17.43   | 5.79 ng/ul                          | 576521       | Phenanthrene-d10 | 16.92     |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Tridecane, 1-iodo-                  | 310 C13H27I  | 035599-77-0      | 91        |
| 2       | Pentadecane, 2,6,10,14-tetramethyl- | 268 C19H40   | 001921-70-6      | 90        |
| 3       | Dodecane, 1-iodo-                   | 296 C12H25I  | 004292-19-7      | 90        |
| 4       | Dodecane, 2-methyl-6-propyl-        | 226 C16H34   | 055045-08-4      | 87        |
| 5       | Tridecane                           | 184 C13H28   | 000629-50-5      | 87        |



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Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
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Quant Title : SVOA CALIBRATION

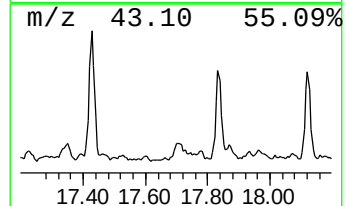
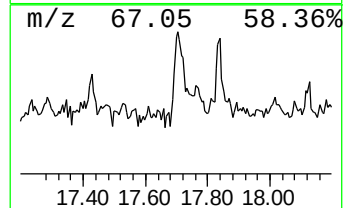
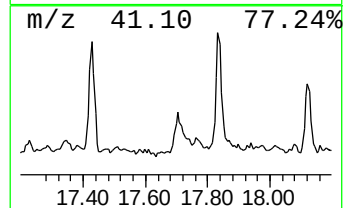
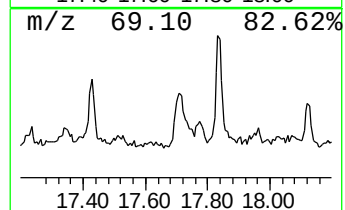
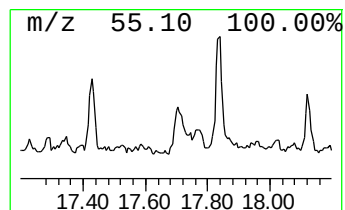
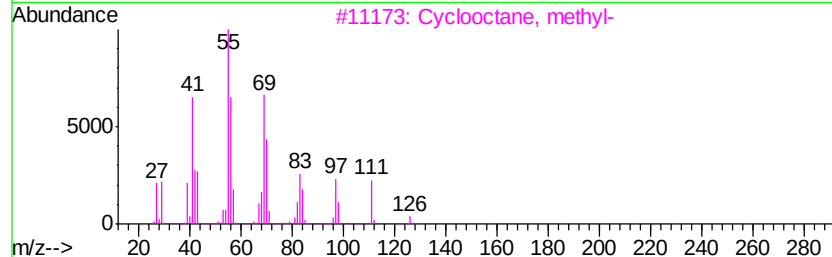
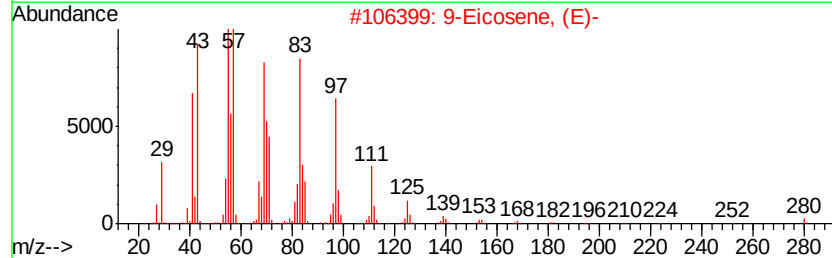
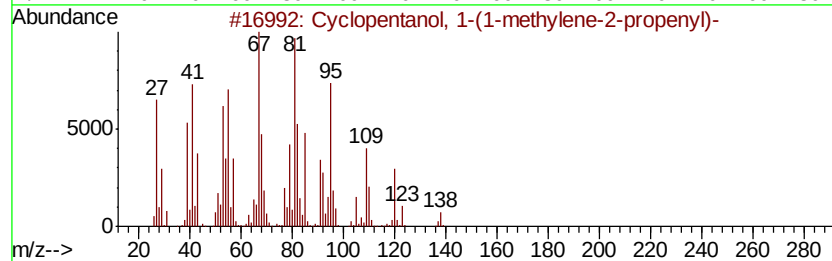
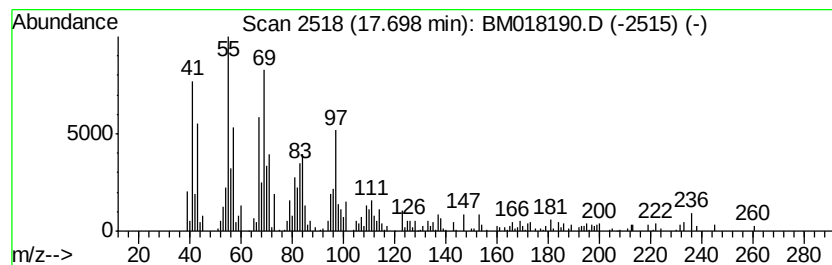
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TIC Integration Parameters: LSCINT.P

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Peak Number 26 Cyclopentanol, 1-(1-methyle... Concentration Rank 16

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.70 | 4.13 ng/ul | 410712 | Phenanthrene-d10 | 16.92 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Cyclopentanol, 1-(1-methylene-2-... | 138 | C9H14O   | 078158-11-9 | 53   |
| 2    |    |   | 9-Eicosene, (E)-                    | 280 | C20H40   | 074685-29-3 | 49   |
| 3    |    |   | Cyclooctane, methyl-                | 126 | C9H18    | 001502-38-1 | 49   |
| 4    |    |   | 6-Octenal, 3,7-dimethyl-            | 154 | C10H18O  | 000106-23-0 | 46   |
| 5    |    |   | 1,2-Cyclohexanediol, cyclic sulf... | 162 | C6H10O3S | 019456-18-9 | 46   |



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Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
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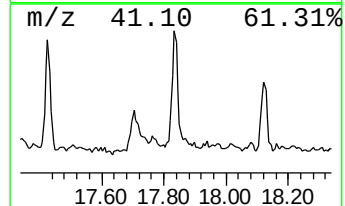
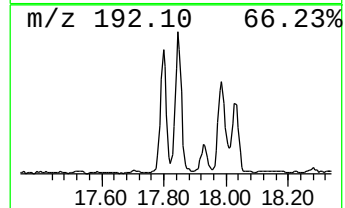
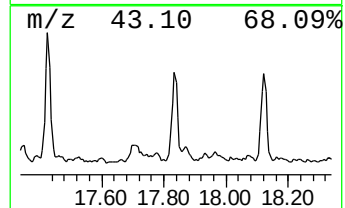
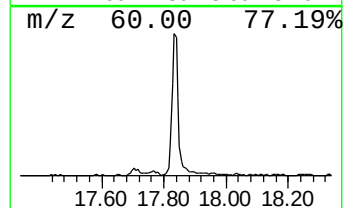
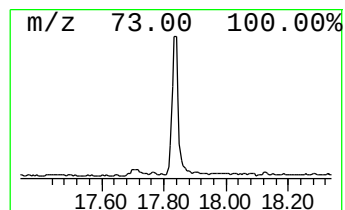
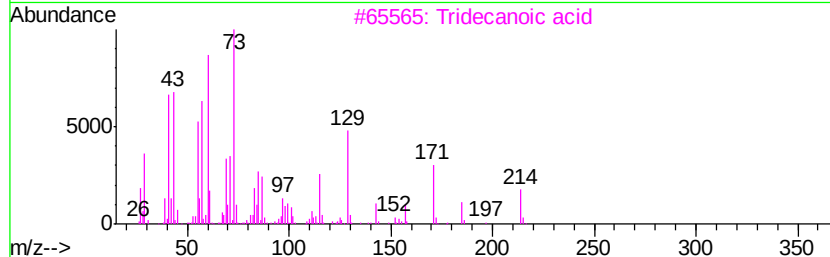
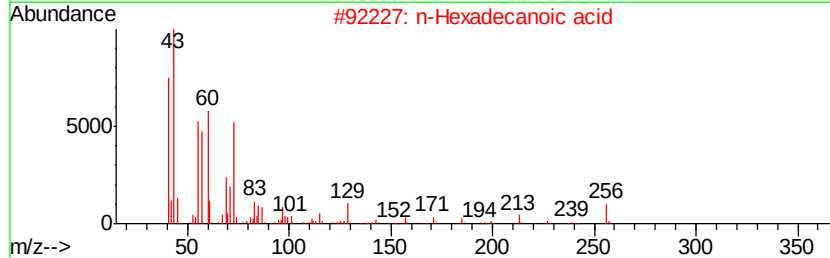
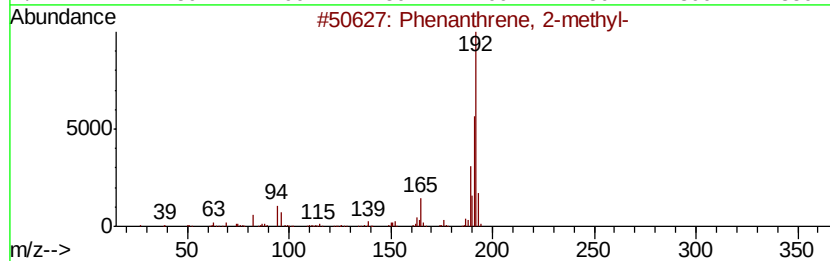
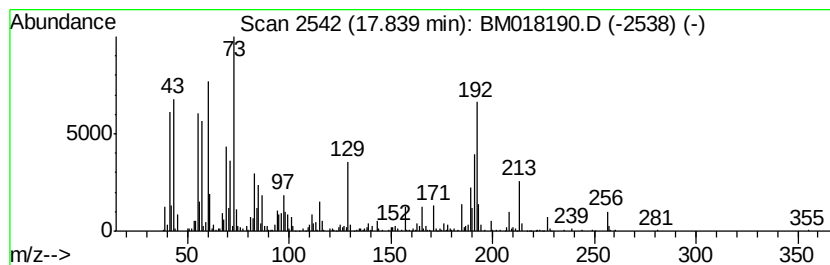
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TIC Integration Parameters: LSCINT.P

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Peak Number 27 Phenanthrene, 2-methyl- Concentration Rank 2

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 17.84 | 10.51 ng/ul | 1045450 | Phenanthrene-d10 | 16.92 |

| Hit# | of | Tentative ID            | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------|-----|----------|-------------|------|
| 1    | 5  | Phenanthrene, 2-methyl- | 192 | C15H12   | 002531-84-2 | 90   |
| 2    |    | n-Hexadecanoic acid     | 256 | C16H32O2 | 000057-10-3 | 89   |
| 3    |    | Tridecanoic acid        | 214 | C13H26O2 | 000638-53-9 | 86   |
| 4    |    | Anthracene, 2-methyl-   | 192 | C15H12   | 000613-12-7 | 83   |
| 5    |    | Tetradecanoic acid      | 228 | C14H28O2 | 000544-63-8 | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

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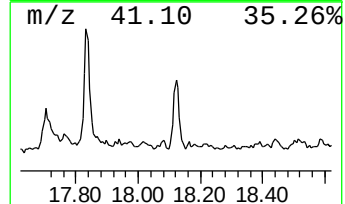
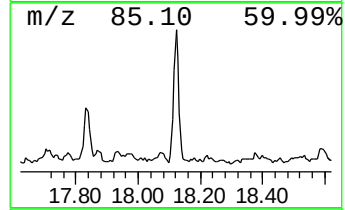
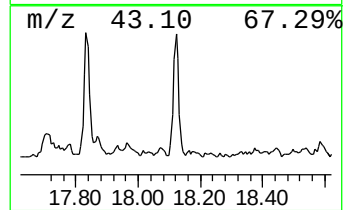
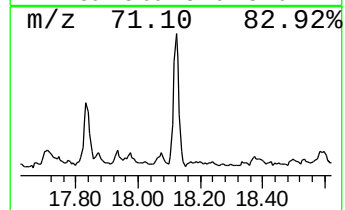
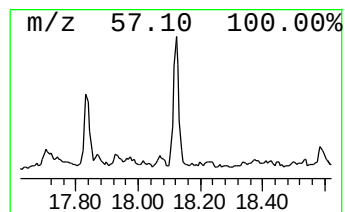
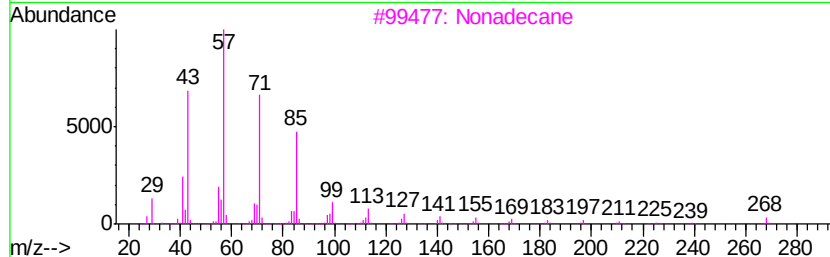
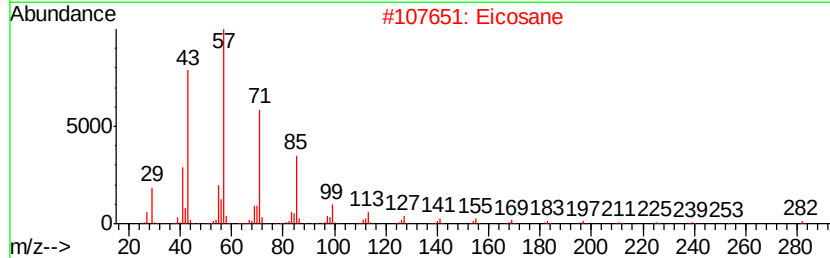
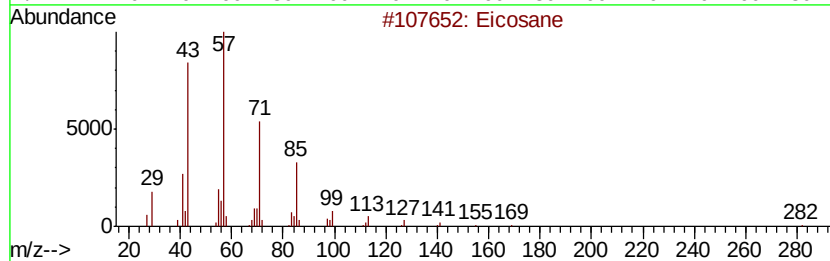
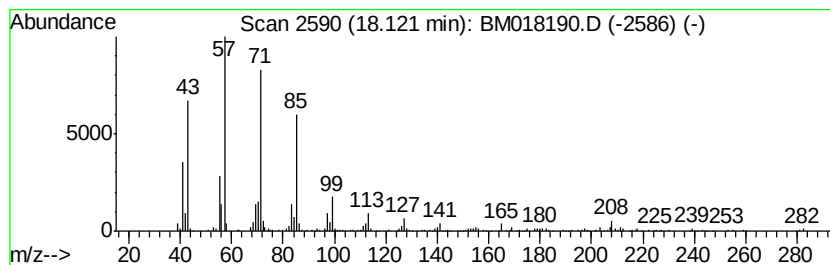
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TIC Integration Parameters: LSCINT.P

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Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 13

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.12 | 4.73 ng/ul | 470840 | Phenanthrene-d10 | 16.92 |

| Hit# | of                     | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------------------------|---|--------------|-----|---------|-------------|------|
| 1    | Eicosane               |   |              | 282 | C20H42  | 000112-95-8 | 93   |
| 2    | Eicosane               |   |              | 282 | C20H42  | 000112-95-8 | 87   |
| 3    | Nonadecane             |   |              | 268 | C19H40  | 000629-92-5 | 87   |
| 4    | Tridecane, 7-propyl-   |   |              | 226 | C16H34  | 055045-09-5 | 87   |
| 5    | Pentadecane, 8-heptyl- |   |              | 310 | C22H46  | 071005-15-7 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
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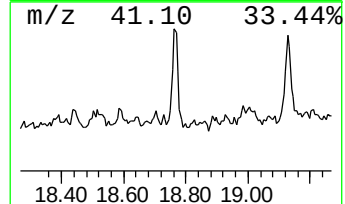
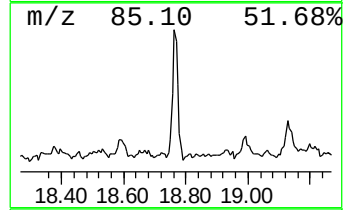
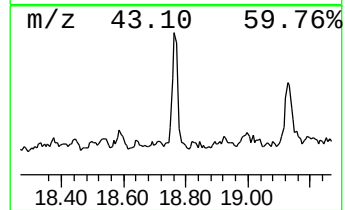
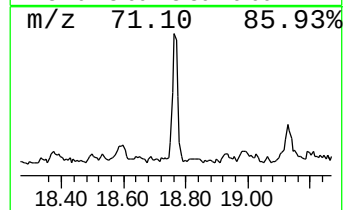
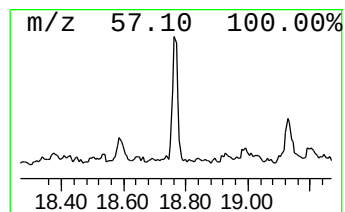
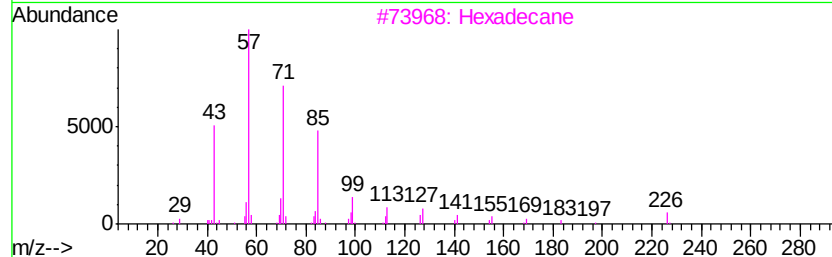
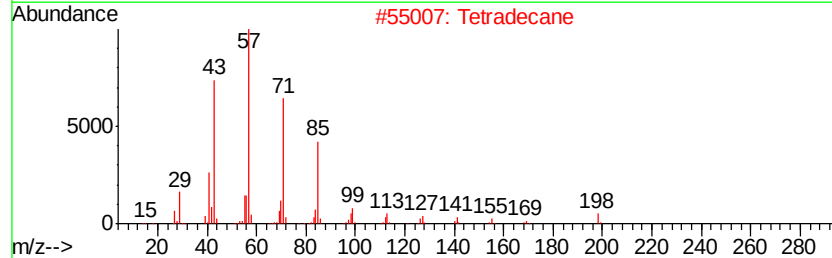
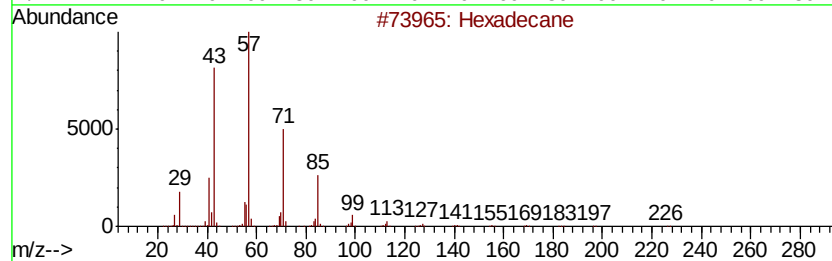
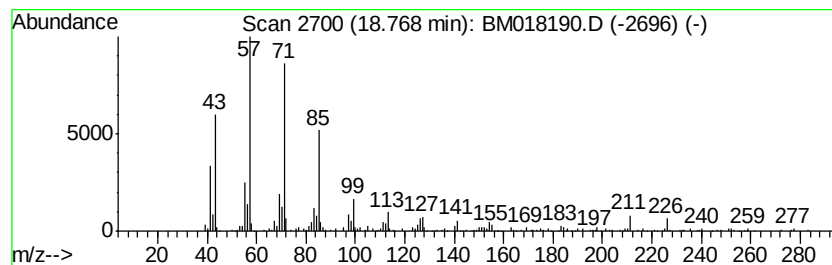
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TIC Integration Parameters: LSCINT.P

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Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.77 | 4.28 ng/ul | 426048 | Phenanthrene-d10 | 16.92 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 96   |
| 2    |    |   | Tetradecane  | 198 | C14H30  | 000629-59-4 | 94   |
| 3    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 93   |
| 4    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 93   |
| 5    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BE7Z4

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Quant Title : SVOA CALIBRATION

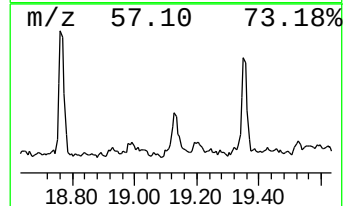
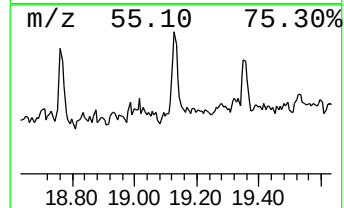
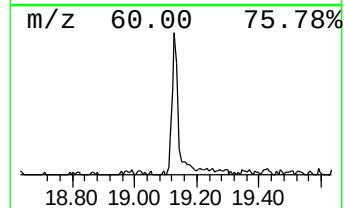
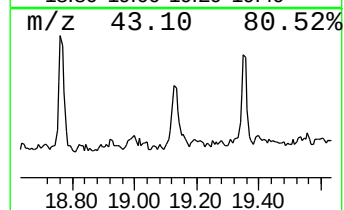
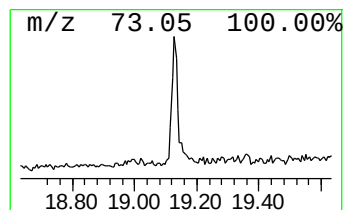
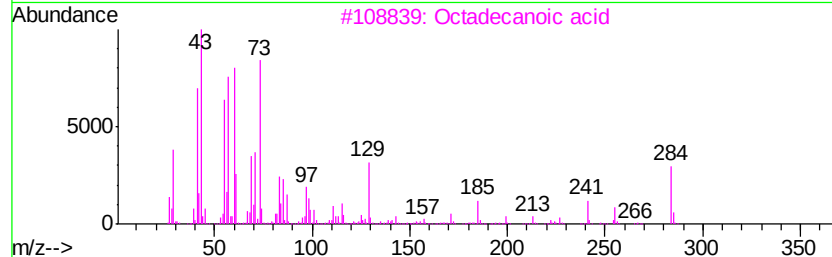
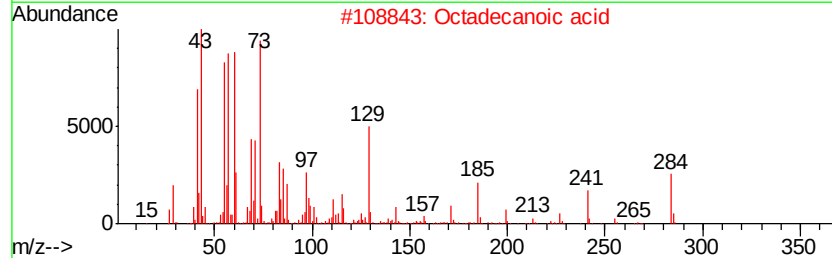
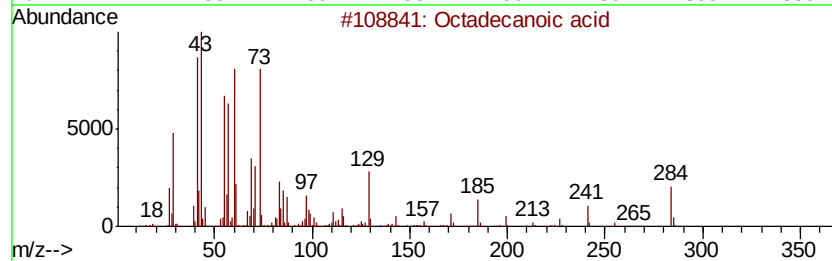
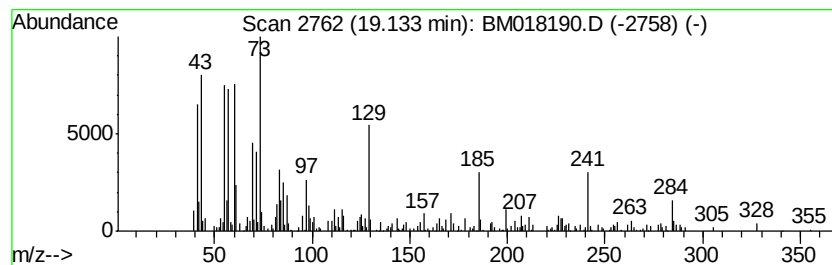
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TIC Integration Parameters: LSCINT.P

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Peak Number 30 Octadecanoic acid Concentration Rank 23

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.13 | 2.65 ng/ul | 328626 | Chrysene-d12     | 21.14 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Octadecanoic acid                   | 284 | C18H36O2 | 000057-11-4 | 98   |
| 2    |    |   | Octadecanoic acid                   | 284 | C18H36O2 | 000057-11-4 | 96   |
| 3    |    |   | Octadecanoic acid                   | 284 | C18H36O2 | 000057-11-4 | 93   |
| 4    |    |   | Octadecanoic acid, 2-(2-hydroxye... | 372 | C22H44O4 | 000106-11-6 | 86   |
| 5    |    |   | Tetradecanoic acid                  | 228 | C14H28O2 | 000544-63-8 | 74   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121418\  
Data File : BM018190.D  
Acq On : 15 Dec 2018 05:11  
Operator : JU/SJ  
Sample : J6238-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
Quant Title : SVOA CALIBRATION

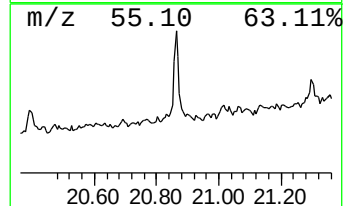
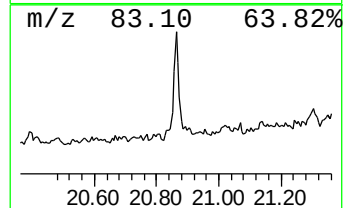
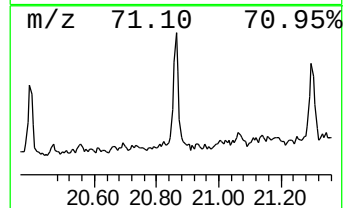
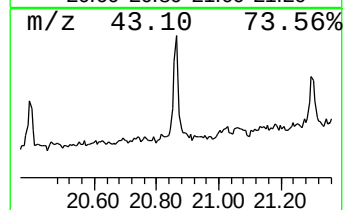
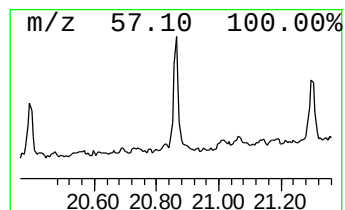
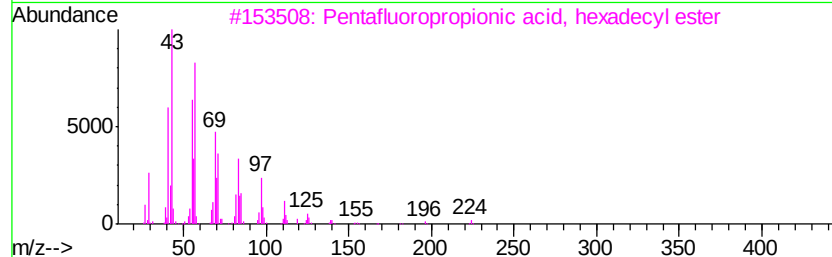
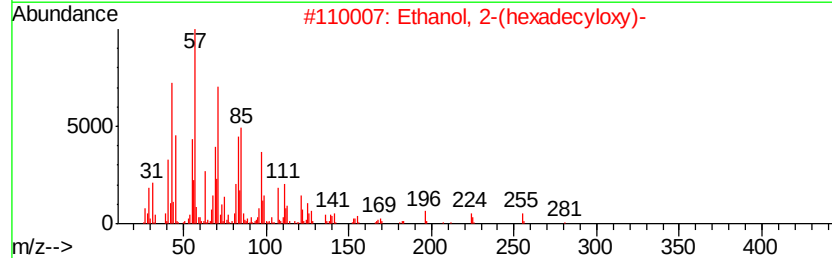
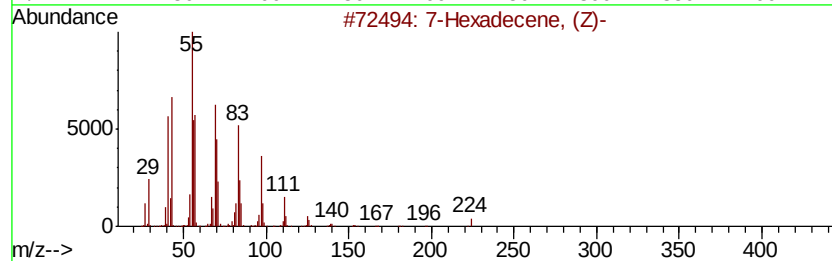
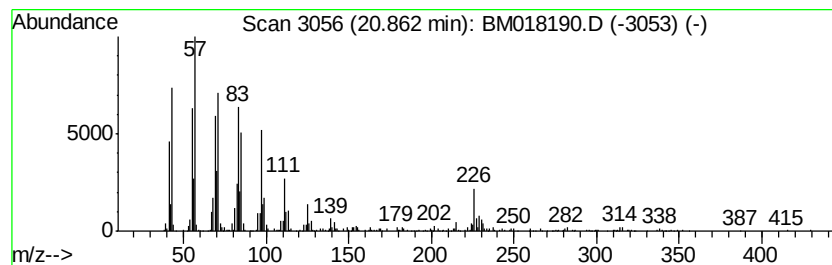
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 31 (DEL) Alkane: Cyclic20.86 Concentration Rank 11

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.86 | 5.31 ng/ul | 657608 | Chrysene-d12     | 21.14 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | 7-Hexadecene, (Z)-                  | 224 | C16H32     | 035507-09-6 | 89   |
| 2    |    | Ethanol, 2-(hexadecyloxy)-          | 286 | C18H38O2   | 002136-71-2 | 83   |
| 3    |    | Pentafluoropropionic acid, hexad... | 388 | C19H33F5O2 | 006222-07-7 | 81   |
| 4    |    | Ethanol, 2-(tetradecyloxy)-         | 258 | C16H34O2   | 002136-70-1 | 81   |
| 5    |    | Trifluoroacetic acid, n-octadecy... | 366 | C20H37F3O2 | 079392-43-1 | 74   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM121418\  
 Data File : BM018190.D  
 Acq On : 15 Dec 2018 05:11  
 Operator : JU/SJ  
 Sample : J6238-01  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BE7Z4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121318MA.M  
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| 2-Pyrrolidinone      | 3.59  | 7.0     | ng/ul | 391296   | 1                     | 7.50  | 1111490 | 20.0 |
| Guanidine, monoth... | 4.77  | 2.5     | ng/ul | 138404   | 1                     | 7.50  | 1111490 | 20.0 |
| (DEL) Alkane: Str... | 11.29 | 2.3     | ng/ul | 242901   | 2                     | 10.27 | 2079220 | 20.0 |
| (DEL) Alkane: Str... | 11.70 | 6.8     | ng/ul | 703404   | 2                     | 10.27 | 2079220 | 20.0 |
| (DEL) Alkane: Str... | 12.67 | 3.2     | ng/ul | 348495   | 3                     | 14.16 | 2155940 | 20.0 |
| (DEL) Alkane: Str... | 12.97 | 8.2     | ng/ul | 886021   | 3                     | 14.16 | 2155940 | 20.0 |
| Naphthalene, 1,6-... | 13.32 | 2.6     | ng/ul | 280876   | 3                     | 14.16 | 2155940 | 20.0 |
| Naphthalene, 2,7-... | 13.47 | 4.8     | ng/ul | 516715   | 3                     | 14.16 | 2155940 | 20.0 |
| (DEL) Alkane: Str... | 14.07 | 9.7     | ng/ul | 1041880  | 3                     | 14.16 | 2155940 | 20.0 |
| Naphthalene, 2,3,... | 14.57 | 3.0     | ng/ul | 317769   | 3                     | 14.16 | 2155940 | 20.0 |
| Naphthalene, 1,6,... | 14.64 | 2.3     | ng/ul | 250241   | 3                     | 14.16 | 2155940 | 20.0 |
| (DEL) Alkane: Str... | 14.69 | 3.2     | ng/ul | 341530   | 3                     | 14.16 | 2155940 | 20.0 |
| Naphthalene, 1,4,... | 14.79 | 2.9     | ng/ul | 313080   | 3                     | 14.16 | 2155940 | 20.0 |
| (DEL) Alkane: Str... | 15.03 | 8.8     | ng/ul | 950009   | 3                     | 14.16 | 2155940 | 20.0 |
| 2-Bromo dodecane     | 15.43 | 4.5     | ng/ul | 486494   | 3                     | 14.16 | 2155940 | 20.0 |
| (DEL) Alkane: Str... | 15.53 | 2.8     | ng/ul | 305654   | 3                     | 14.16 | 2155940 | 20.0 |
| (DEL) Alkane: Str... | 15.59 | 2.2     | ng/ul | 222725   | 4                     | 16.92 | 1990270 | 20.0 |
| (DEL) Alkane: Cyc... | 15.64 | 2.2     | ng/ul | 219781   | 4                     | 16.92 | 1990270 | 20.0 |
| (DEL) Alkane: Str... | 15.90 | 18.6    | ng/ul | 1853810  | 4                     | 16.92 | 1990270 | 20.0 |
| (DEL) Alkane: Str... | 16.69 | 7.0     | ng/ul | 701434   | 4                     | 16.92 | 1990270 | 20.0 |
| (DEL) Alkane: Str... | 16.74 | 5.5     | ng/ul | 547547   | 4                     | 16.92 | 1990270 | 20.0 |
| Tridecane, 1-iodo-   | 17.43 | 5.8     | ng/ul | 576521   | 4                     | 16.92 | 1990270 | 20.0 |
| Cyclopentanol, 1-... | 17.70 | 4.1     | ng/ul | 410712   | 4                     | 16.92 | 1990270 | 20.0 |
| Phenanthrene, 2-m... | 17.84 | 10.5    | ng/ul | 1045450  | 4                     | 16.92 | 1990270 | 20.0 |
| (DEL) Alkane: Str... | 18.12 | 4.7     | ng/ul | 470840   | 4                     | 16.92 | 1990270 | 20.0 |
| (DEL) Alkane: Str... | 18.77 | 4.3     | ng/ul | 426048   | 4                     | 16.92 | 1990270 | 20.0 |
| Octadecanoic acid    | 19.13 | 2.6     | ng/ul | 328626   | 5                     | 21.14 | 2479050 | 20.0 |
| (DEL) Alkane: Cyc... | 20.86 | 5.3     | ng/ul | 657608   | 5                     | 21.14 | 2479050 | 20.0 |