

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
 Data File : BM013809.D  
 Acq On : 25 Jan 2018 17:13  
 Operator : SJ/JU  
 Sample : J1243-04  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 F6B27

## Integration Parameters: LSCINT.P

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

Filtering: 5  
 Min Area: 1 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

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

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

Signal : TIC

| peak<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         | 4.763       | 296           | 302         | 308          | rVB      | 140217         | 204315        | 4.83%           | 0.481%        |
| 2         | 5.198       | 372           | 376         | 382          | rVB4     | 40266          | 59058         | 1.40%           | 0.139%        |
| 3         | 5.263       | 382           | 387         | 396          | rBV3     | 27050          | 62319         | 1.47%           | 0.147%        |
| 4         | 5.622       | 443           | 448         | 456          | rVB      | 202815         | 285571        | 6.75%           | 0.672%        |
| 5         | 5.869       | 486           | 490         | 498          | rBV8     | 26721          | 62011         | 1.47%           | 0.146%        |
| 6         | 5.957       | 498           | 505         | 509          | rBV      | 94327          | 155126        | 3.67%           | 0.365%        |
| 7         | 5.992       | 509           | 511         | 519          | rVB6     | 32185          | 50716         | 1.20%           | 0.119%        |
| 8         | 6.081       | 523           | 526         | 534          | rVB7     | 28797          | 56929         | 1.35%           | 0.134%        |
| 9         | 6.157       | 534           | 539         | 544          | rBV      | 47087          | 66666         | 1.58%           | 0.157%        |
| 10        | 6.234       | 548           | 552         | 554          | rBV3     | 34028          | 42900         | 1.01%           | 0.101%        |
| 11        | 6.492       | 588           | 596         | 601          | rBV5     | 29745          | 78163         | 1.85%           | 0.184%        |
| 12        | 6.581       | 608           | 611         | 617          | rVB4     | 60965          | 91591         | 2.16%           | 0.216%        |
| 13        | 6.640       | 617           | 621         | 625          | rBV2     | 62977          | 91703         | 2.17%           | 0.216%        |
| 14        | 6.687       | 625           | 629         | 633          | rVV2     | 51026          | 61743         | 1.46%           | 0.145%        |
| 15        | 6.751       | 633           | 640         | 641          | rVV5     | 71617          | 108465        | 2.56%           | 0.255%        |
| 16        | 6.781       | 641           | 645         | 658          | rVB2     | 463423         | 924952        | 21.86%          | 2.177%        |
| 17        | 6.939       | 663           | 672         | 684          | rBV2     | 367115         | 636159        | 15.04%          | 1.497%        |
| 18        | 7.128       | 699           | 704         | 713          | rVB      | 509795         | 836573        | 19.77%          | 1.969%        |
| 19        | 7.210       | 713           | 718         | 722          | rBV      | 315647         | 453873        | 10.73%          | 1.068%        |
| 20        | 7.498       | 761           | 767         | 772          | rBV7     | 22917          | 42568         | 1.01%           | 0.100%        |
| 21        | 7.592       | 772           | 783         | 789          | rVV      | 496450         | 878359        | 20.76%          | 2.067%        |
| 22        | 7.663       | 790           | 795         | 799          | rVV      | 104111         | 169798        | 4.01%           | 0.400%        |
| 23        | 7.698       | 799           | 801         | 809          | rVV5     | 29220          | 50909         | 1.20%           | 0.120%        |
| 24        | 7.863       | 821           | 829         | 836          | rVB4     | 32053          | 68916         | 1.63%           | 0.162%        |
| 25        | 8.022       | 851           | 856         | 861          | rBV5     | 32637          | 58981         | 1.39%           | 0.139%        |
| 26        | 8.169       | 878           | 881         | 885          | rVB2     | 34346          | 47683         | 1.13%           | 0.112%        |
| 27        | 8.234       | 885           | 892         | 899          | rBV4     | 87646          | 189379        | 4.48%           | 0.446%        |
| 28        | 8.310       | 899           | 905         | 915          | rBV2     | 567776         | 993819        | 23.49%          | 2.339%        |
| 29        | 8.686       | 964           | 969         | 973          | rBV2     | 37855          | 55176         | 1.30%           | 0.130%        |
| 30        | 8.751       | 973           | 980         | 985          | rVV      | 467107         | 816432        | 19.30%          | 1.922%        |
| 31        | 8.804       | 985           | 989         | 999          | rVB      | 361868         | 543178        | 12.84%          | 1.278%        |
| 32        | 8.928       | 1005          | 1010        | 1015         | rVB6     | 41960          | 79932         | 1.89%           | 0.188%        |
| 33        | 9.057       | 1030          | 1032        | 1041         | rVB6     | 29664          | 52363         | 1.24%           | 0.123%        |
| 34        | 9.145       | 1041          | 1047        | 1053         | rBV7     | 25246          | 59353         | 1.40%           | 0.140%        |

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
 Data File : BM013809.D  
 Acq On : 25 Jan 2018 17:13  
 Operator : SJ/JU  
 Sample : J1243-04  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 F6B27

## Integration Parameters: LSCINT.P

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

Filtering: 5  
 Min Area: 1 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

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

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

|    |        |      |      |      |      |         |         |         |        |
|----|--------|------|------|------|------|---------|---------|---------|--------|
| 35 | 9.216  | 1053 | 1059 | 1064 | rBV7 | 29736   | 68199   | 1.61%   | 0.161% |
| 36 | 9.469  | 1093 | 1102 | 1111 | rBV  | 411432  | 824812  | 19.49%  | 1.941% |
| 37 | 9.722  | 1140 | 1145 | 1149 | rVB5 | 29986   | 49283   | 1.16%   | 0.116% |
| 38 | 9.792  | 1149 | 1157 | 1163 | rBV6 | 57264   | 127694  | 3.02%   | 0.301% |
| 39 | 9.898  | 1169 | 1175 | 1179 | rBV3 | 48890   | 75525   | 1.79%   | 0.178% |
| 40 | 10.004 | 1187 | 1193 | 1202 | rBV  | 568353  | 979946  | 23.16%  | 2.307% |
| 41 | 10.369 | 1244 | 1255 | 1259 | rBV2 | 594072  | 1516495 | 35.84%  | 3.569% |
| 42 | 10.416 | 1259 | 1263 | 1270 | rVB  | 300790  | 502897  | 11.89%  | 1.184% |
| 43 | 10.522 | 1275 | 1281 | 1289 | rVV2 | 516335  | 980084  | 23.16%  | 2.307% |
| 44 | 10.592 | 1290 | 1293 | 1299 | rVB7 | 28211   | 46628   | 1.10%   | 0.110% |
| 45 | 11.063 | 1365 | 1373 | 1377 | rBV6 | 27206   | 62863   | 1.49%   | 0.148% |
| 46 | 11.204 | 1390 | 1397 | 1402 | rBV7 | 26817   | 60817   | 1.44%   | 0.143% |
| 47 | 11.280 | 1405 | 1410 | 1418 | rVB3 | 51858   | 101406  | 2.40%   | 0.239% |
| 48 | 11.386 | 1423 | 1428 | 1433 | rBV2 | 104970  | 173421  | 4.10%   | 0.408% |
| 49 | 11.704 | 1475 | 1482 | 1489 | rBV  | 108284  | 182028  | 4.30%   | 0.428% |
| 50 | 11.798 | 1489 | 1498 | 1503 | rBV  | 276580  | 426167  | 10.07%  | 1.003% |
| 51 | 12.039 | 1530 | 1539 | 1545 | rVB2 | 62551   | 131992  | 3.12%   | 0.311% |
| 52 | 12.263 | 1569 | 1577 | 1584 | rBV2 | 42980   | 85040   | 2.01%   | 0.200% |
| 53 | 12.445 | 1601 | 1608 | 1612 | rBV3 | 56069   | 90492   | 2.14%   | 0.213% |
| 54 | 12.769 | 1658 | 1663 | 1669 | rVB4 | 53228   | 88296   | 2.09%   | 0.208% |
| 55 | 12.986 | 1694 | 1700 | 1707 | rBV4 | 41245   | 68936   | 1.63%   | 0.162% |
| 56 | 13.063 | 1707 | 1713 | 1720 | rVV  | 187963  | 261974  | 6.19%   | 0.617% |
| 57 | 13.304 | 1749 | 1754 | 1760 | rVB6 | 29825   | 43448   | 1.03%   | 0.102% |
| 58 | 13.663 | 1809 | 1815 | 1820 | rBV  | 971738  | 1375698 | 32.51%  | 3.238% |
| 59 | 13.710 | 1820 | 1823 | 1831 | rVV  | 401362  | 622281  | 14.71%  | 1.465% |
| 60 | 13.933 | 1854 | 1861 | 1870 | rVB  | 1101894 | 1661350 | 39.27%  | 3.910% |
| 61 | 14.151 | 1892 | 1898 | 1902 | rBV  | 119087  | 154988  | 3.66%   | 0.365% |
| 62 | 14.245 | 1905 | 1914 | 1921 | rBV2 | 774472  | 1233571 | 29.16%  | 2.903% |
| 63 | 14.486 | 1949 | 1955 | 1972 | rBV2 | 253103  | 599322  | 14.16%  | 1.411% |
| 64 | 14.798 | 1999 | 2008 | 2014 | rBV  | 578989  | 891476  | 21.07%  | 2.098% |
| 65 | 14.886 | 2014 | 2023 | 2031 | rVB  | 230597  | 397988  | 9.41%   | 0.937% |
| 66 | 15.110 | 2055 | 2061 | 2065 | rVB3 | 77357   | 105907  | 2.50%   | 0.249% |
| 67 | 15.245 | 2077 | 2084 | 2090 | rVB  | 1354979 | 1906049 | 45.05%  | 4.486% |
| 68 | 15.386 | 2102 | 2108 | 2121 | rBV  | 329747  | 565216  | 13.36%  | 1.330% |
| 69 | 15.615 | 2142 | 2147 | 2161 | rVB  | 819799  | 1176272 | 27.80%  | 2.769% |
| 70 | 15.974 | 2201 | 2208 | 2217 | rBV6 | 71889   | 172964  | 4.09%   | 0.407% |
| 71 | 16.657 | 2315 | 2324 | 2338 | rBV  | 2920141 | 4231045 | 100.00% | 9.959% |

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
 Data File : BM013809.D  
 Acq On : 25 Jan 2018 17:13  
 Operator : SJ/JU  
 Sample : J1243-04  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 F6B27

## Integration Parameters: LSCINT.P

Integrator: RTE  
 Smoothing : OFF Filtering: 5  
 Sampling : 1 Min Area: 1 % of largest Peak  
 Start Thrs: 0.2 Max Peaks: 100  
 Stop Thrs : 0 Peak Location: TOP

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

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

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 72 | 16.762 | 2338 | 2342 | 2347 | rVB3 | 35546   | 48993   | 1.16%  | 0.115% |
| 73 | 16.998 | 2377 | 2382 | 2388 | rBV2 | 990323  | 1386446 | 32.77% | 3.263% |
| 74 | 17.098 | 2393 | 2399 | 2407 | rVB  | 1497613 | 2053405 | 48.53% | 4.833% |
| 75 | 17.903 | 2531 | 2536 | 2547 | rBV2 | 178956  | 302012  | 7.14%  | 0.711% |
| 76 | 19.198 | 2752 | 2756 | 2762 | rBV6 | 66831   | 100128  | 2.37%  | 0.236% |
| 77 | 19.403 | 2785 | 2791 | 2798 | rBV2 | 1649540 | 2363834 | 55.87% | 5.564% |
| 78 | 19.462 | 2798 | 2801 | 2806 | rVB5 | 40815   | 65401   | 1.55%  | 0.154% |
| 79 | 20.915 | 3045 | 3048 | 3054 | rBV3 | 206680  | 269009  | 6.36%  | 0.633% |
| 80 | 21.197 | 3091 | 3096 | 3106 | rBV  | 1206237 | 1723926 | 40.74% | 4.058% |
| 81 | 23.256 | 3440 | 3446 | 3455 | rVB  | 1203066 | 2208731 | 52.20% | 5.199% |
| 82 | 23.391 | 3464 | 3469 | 3479 | rVB  | 735185  | 1485803 | 35.12% | 3.497% |

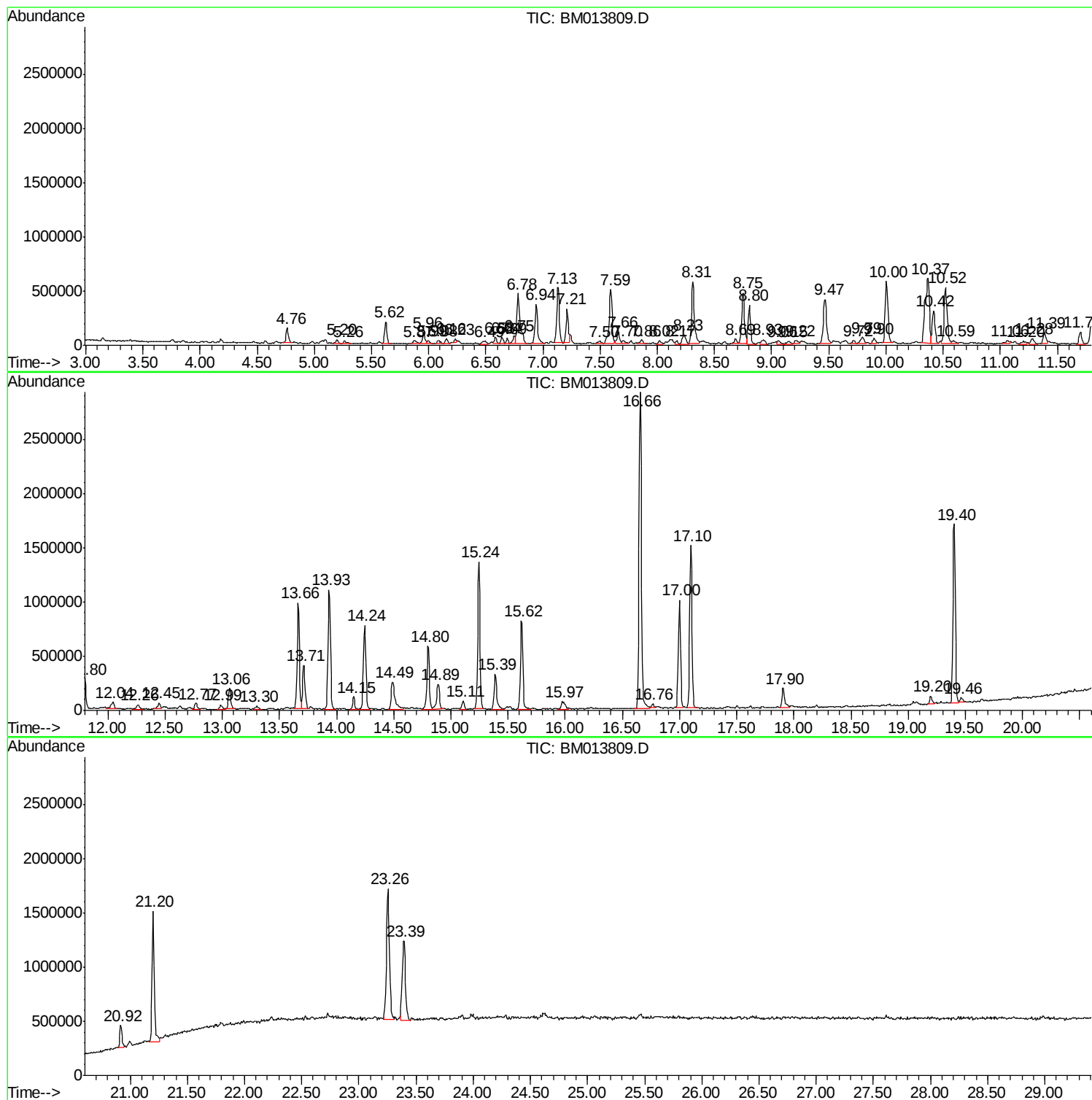
Sum of corrected areas: 42485937

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

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



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

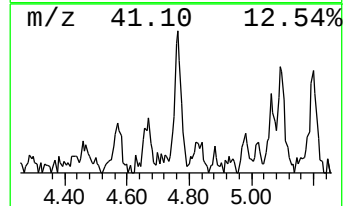
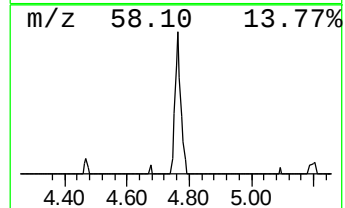
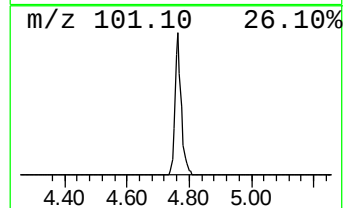
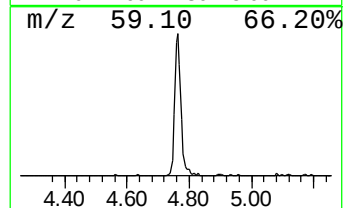
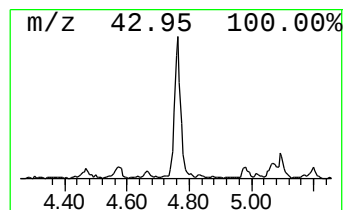
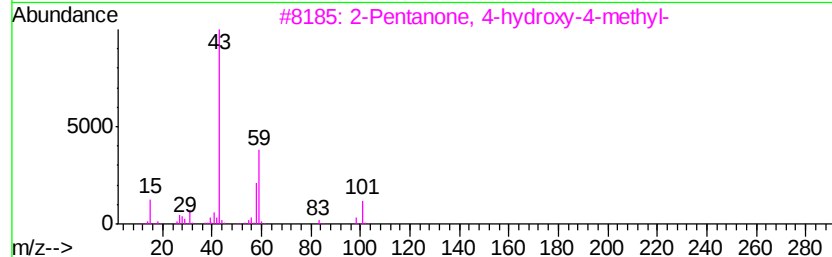
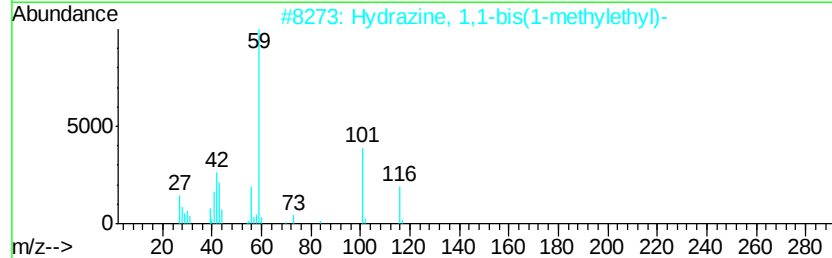
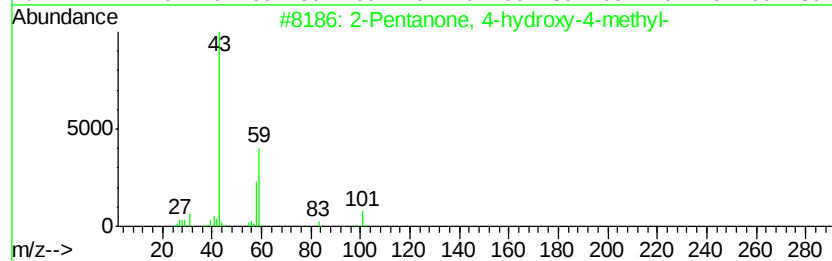
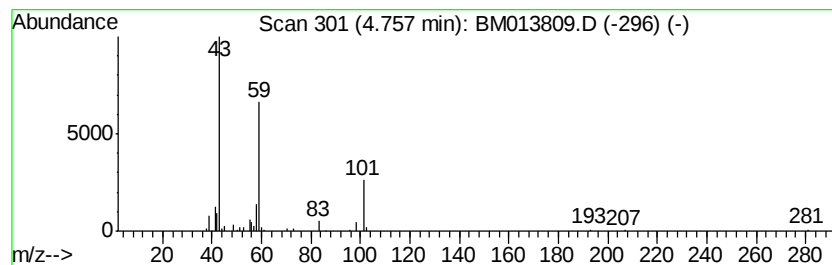
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 4.76 | 4.65 ng/ul | 204315 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-   | 116 | C6H12O2 | 000123-42-2 | 50   |
| 2    |    | Hydrazine, 1,1-bis(1-methylethyl)- | 116 | C6H16N2 | 000921-14-2 | 47   |
| 3    |    | 2-Pentanone, 4-hydroxy-4-methyl-   | 116 | C6H12O2 | 000123-42-2 | 39   |
| 4    |    | Dimethadione                       | 129 | C5H7NO3 | 000695-53-4 | 33   |
| 5    |    | 3-Hydroxy-3-methyl-2-butanone      | 102 | C5H10O2 | 000115-22-0 | 27   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

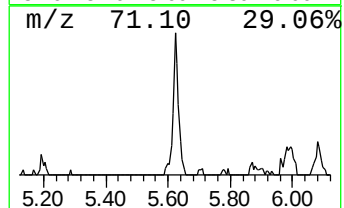
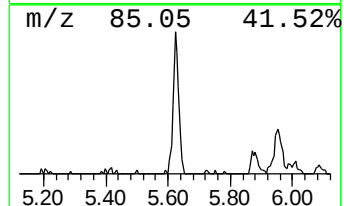
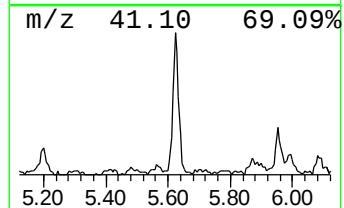
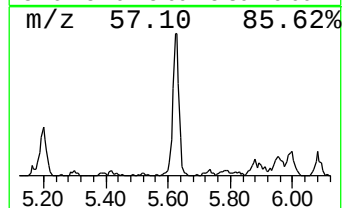
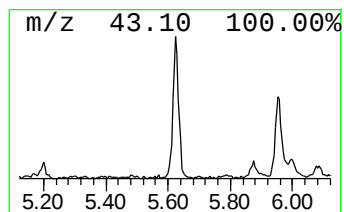
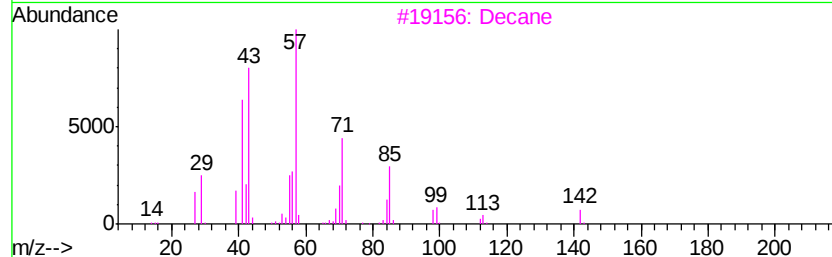
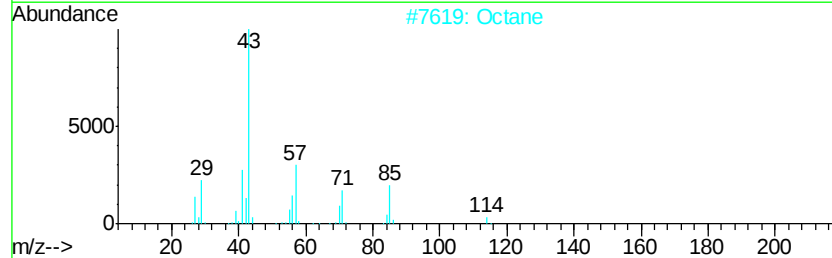
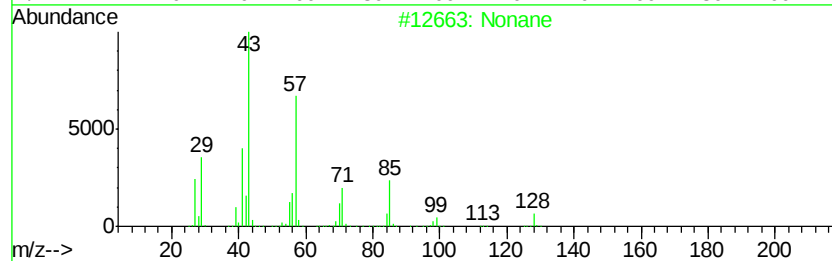
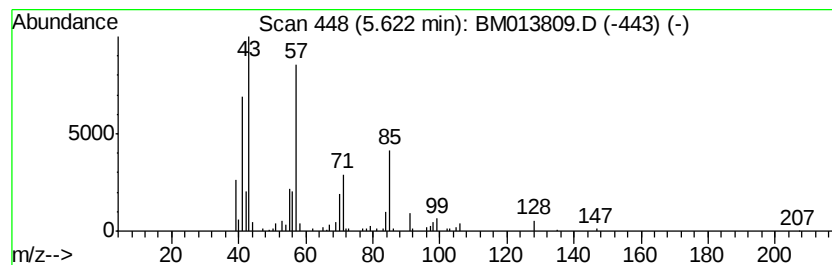
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 5

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 5.62 | 6.50 ng/ul | 285571 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Nonane       | 128 | C9H20   | 000111-84-2 | 91   |
| 2    |    | Octane       | 114 | C8H18   | 000111-65-9 | 59   |
| 3    |    | Decane       | 142 | C10H22  | 000124-18-5 | 59   |
| 4    |    | Nonane       | 128 | C9H20   | 000111-84-2 | 58   |
| 5    |    | Decane       | 142 | C10H22  | 000124-18-5 | 53   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

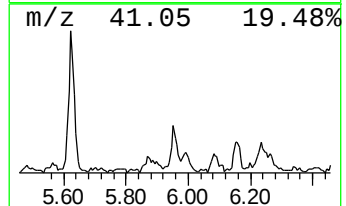
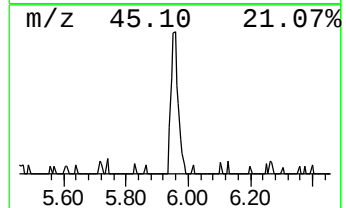
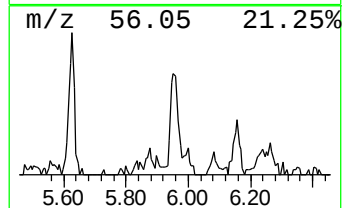
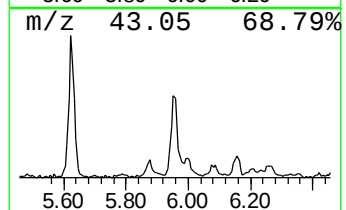
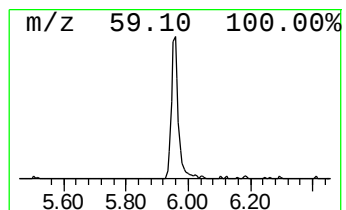
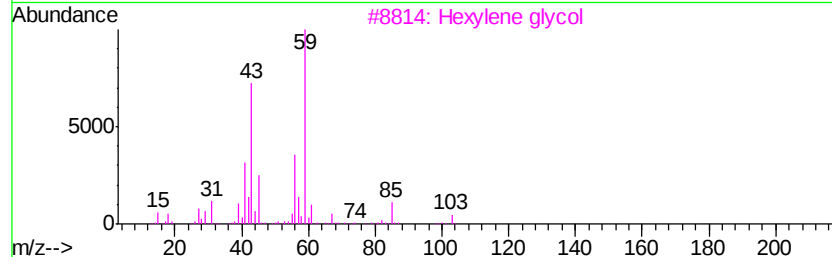
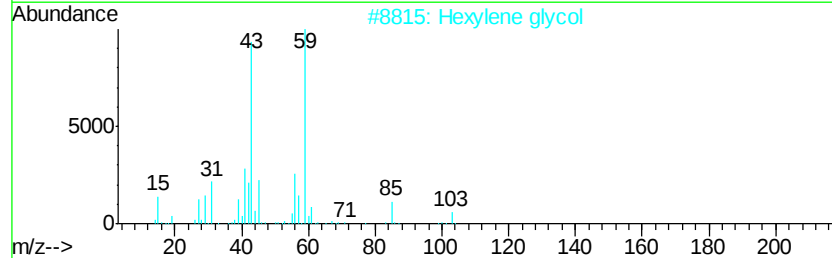
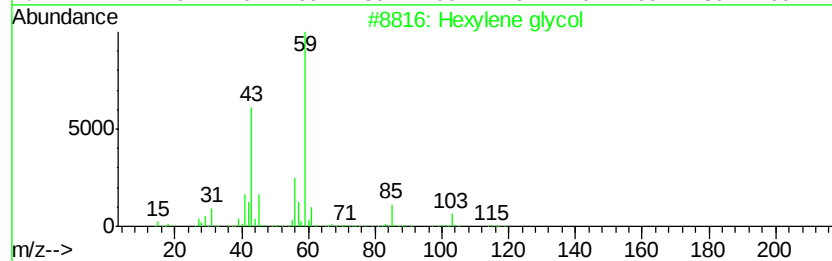
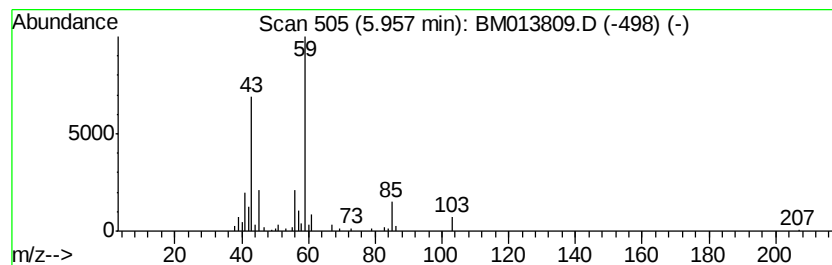
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 Hexylene glycol Concentration Rank 12

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 5.96 | 3.53 ng/ul | 155126 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexylene glycol           | 118 | C6H14O2 | 000107-41-5 | 86   |
| 2    |    |   | Hexylene glycol           | 118 | C6H14O2 | 000107-41-5 | 47   |
| 3    |    |   | Hexylene glycol           | 118 | C6H14O2 | 000107-41-5 | 47   |
| 4    |    |   | Hexylene glycol           | 118 | C6H14O2 | 000107-41-5 | 40   |
| 5    |    |   | Oxetane, 2,2,4-trimethyl- | 100 | C6H12O  | 023120-44-7 | 40   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

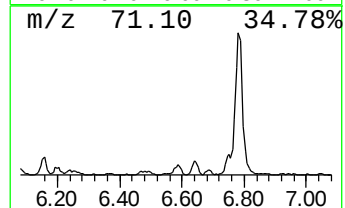
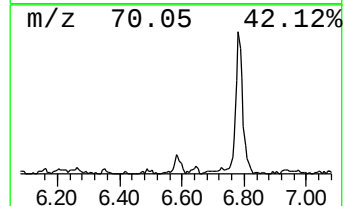
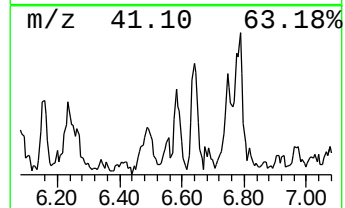
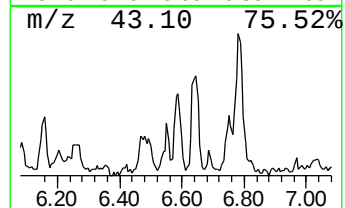
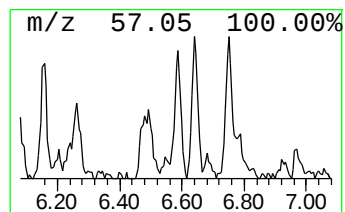
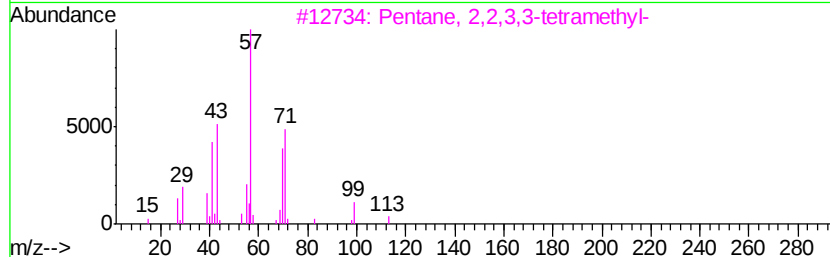
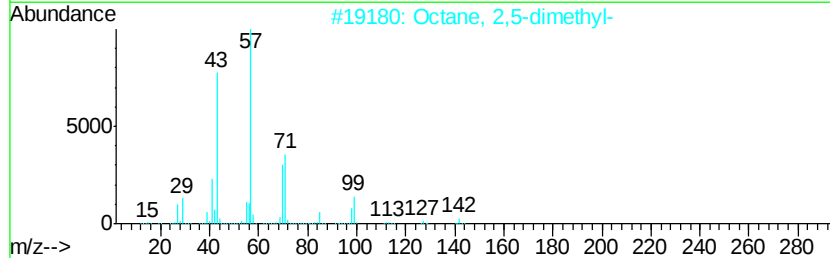
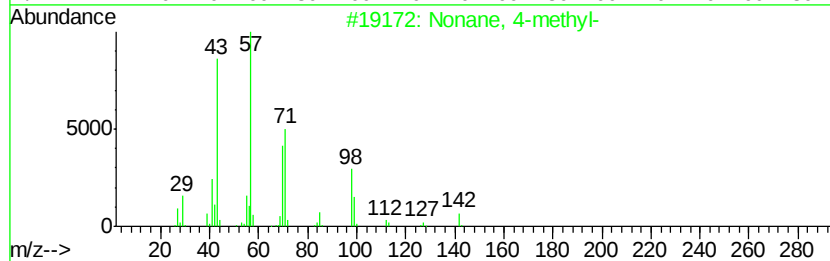
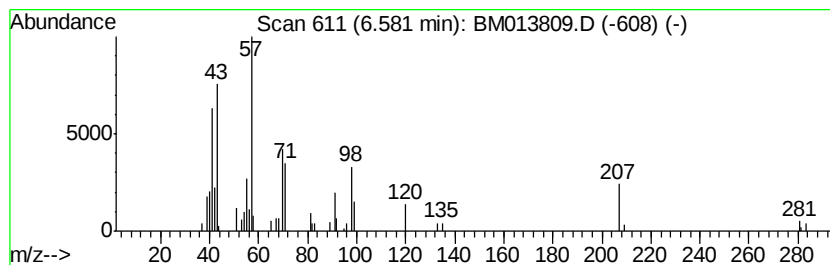
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 19

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 6.58 | 2.09 ng/ul | 91591 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Nonane, 4-methyl-             | 142 | C10H22  | 017301-94-9 | 50   |
| 2    |    | Octane, 2,5-dimethyl-         | 142 | C10H22  | 015869-89-3 | 37   |
| 3    |    | Pentane, 2,2,3,3-tetramethyl- | 128 | C9H20   | 007154-79-2 | 37   |
| 4    |    | Heptane, 3-ethyl-             | 128 | C9H20   | 015869-80-4 | 32   |
| 5    |    | Octane, 2,3-dimethyl-         | 142 | C10H22  | 007146-60-3 | 32   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

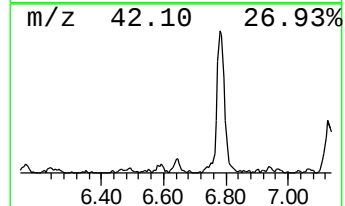
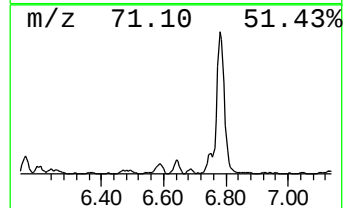
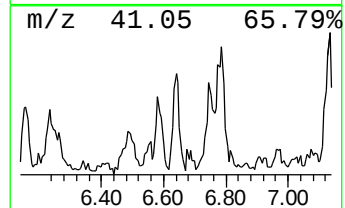
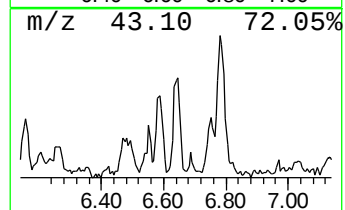
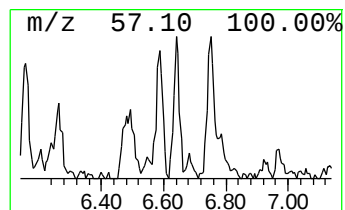
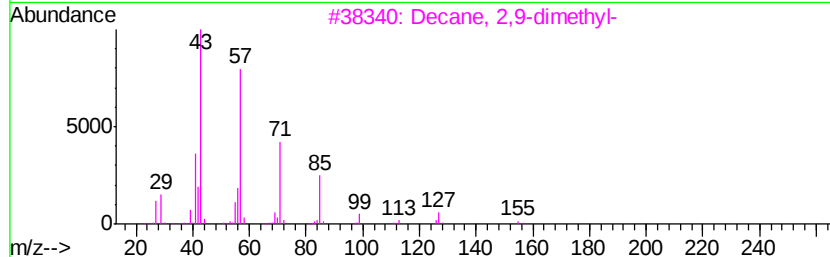
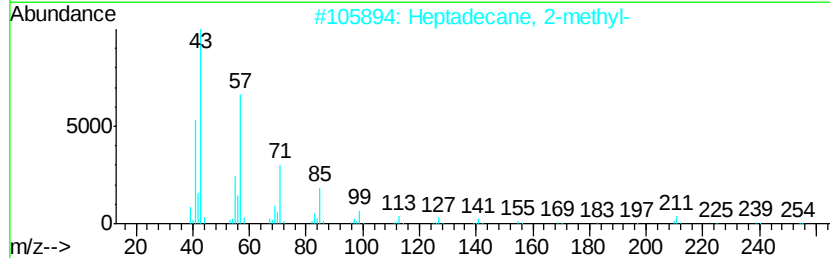
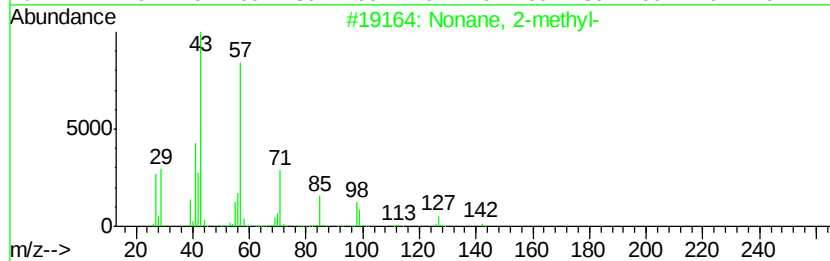
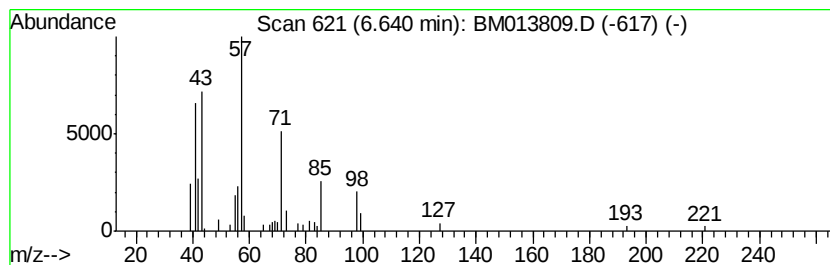
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 18

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 6.64 | 2.09 ng/ul | 91703 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|---------|---|-------------------------|-----|---------|-------------|------|
| 1       |   | Nonane, 2-methyl-       | 142 | C10H22  | 000871-83-0 | 72   |
| 2       |   | Heptadecane, 2-methyl-  | 254 | C18H38  | 001560-89-0 | 64   |
| 3       |   | Decane, 2,9-dimethyl-   | 170 | C12H26  | 001002-17-1 | 50   |
| 4       |   | Undecane, 2,9-dimethyl- | 184 | C13H28  | 017301-26-7 | 50   |
| 5       |   | Tridecane, 2-methyl-    | 198 | C14H30  | 001560-96-9 | 50   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

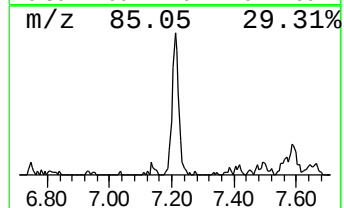
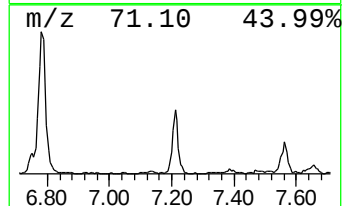
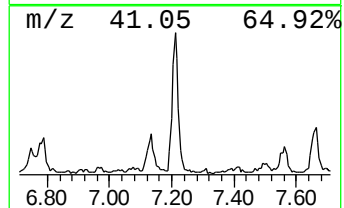
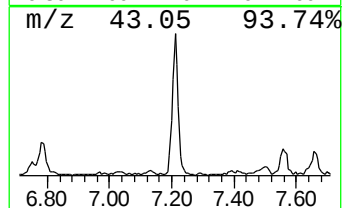
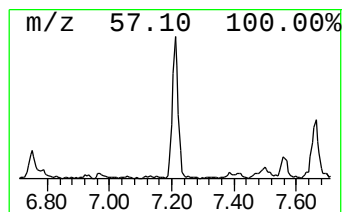
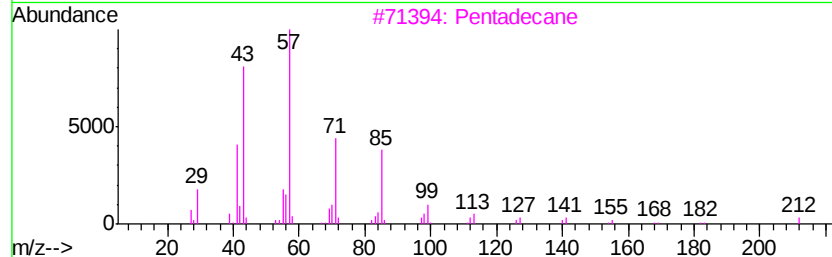
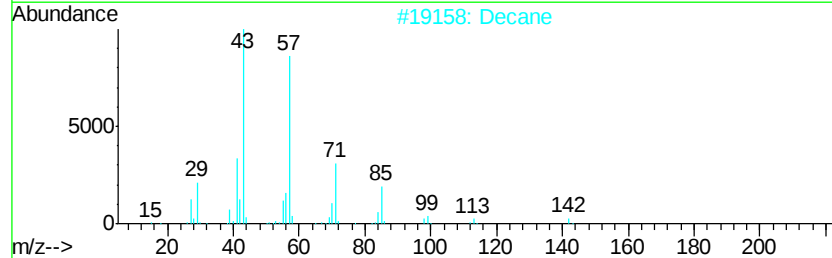
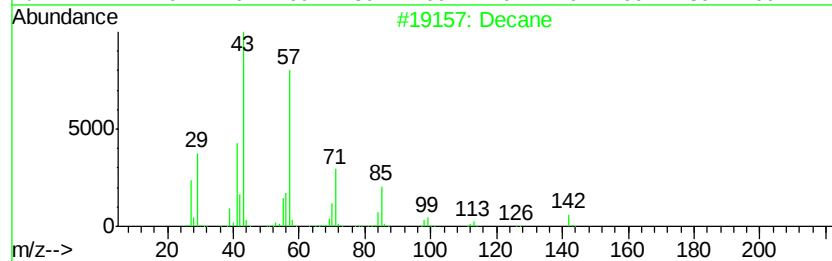
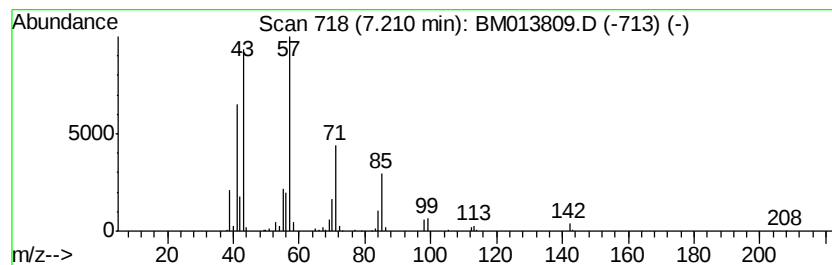
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 7.21 | 10.33 ng/ul | 453873 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Decane                   | 142 | C10H22  | 000124-18-5 | 90   |
| 2    |    | Decane                   | 142 | C10H22  | 000124-18-5 | 87   |
| 3    |    | Pentadecane              | 212 | C15H32  | 000629-62-9 | 78   |
| 4    |    | Decane, 2,3,6-trimethyl- | 184 | C13H28  | 062238-12-4 | 72   |
| 5    |    | Decane                   | 142 | C10H22  | 000124-18-5 | 68   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

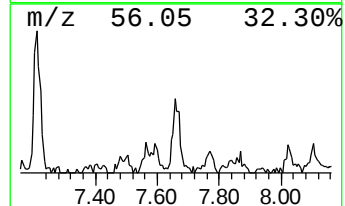
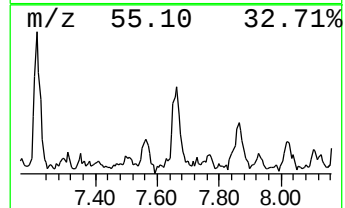
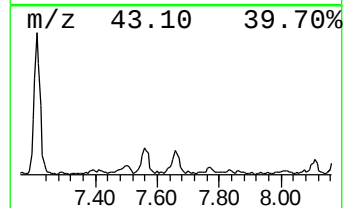
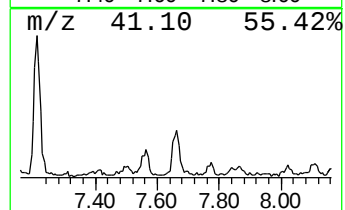
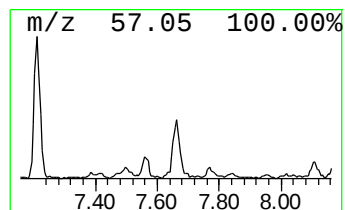
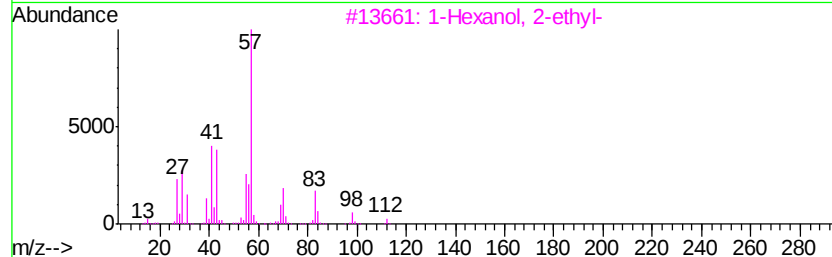
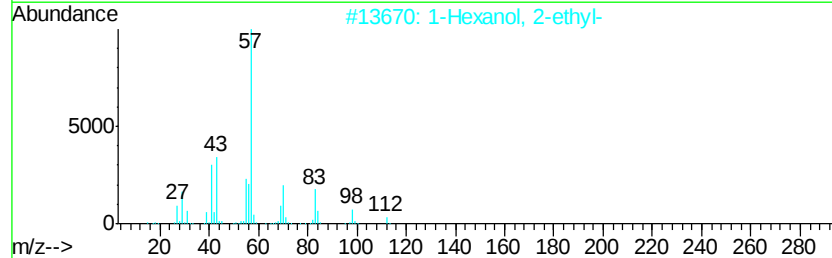
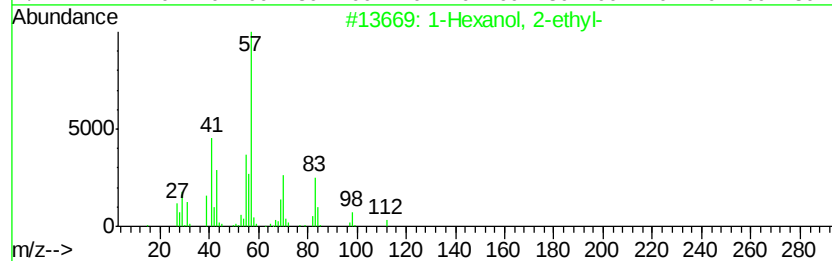
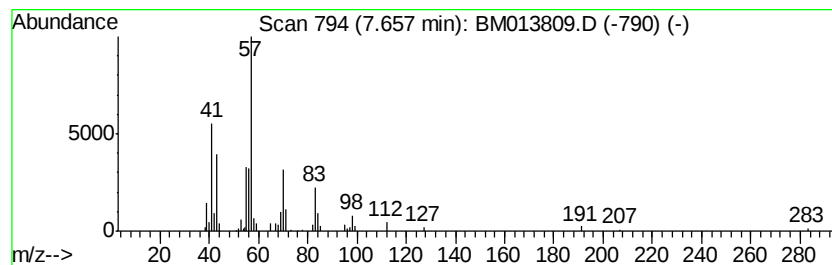
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 1-Hexanol, 2-ethyl- Concentration Rank 11

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.66 | 3.87 ng/ul | 169798 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------|-----|---------|-------------|------|
| 1    |    |   | 1-Hexanol, 2-ethyl- | 130 | C8H18O  | 000104-76-7 | 78   |
| 2    |    |   | 1-Hexanol, 2-ethyl- | 130 | C8H18O  | 000104-76-7 | 78   |
| 3    |    |   | 1-Hexanol, 2-ethyl- | 130 | C8H18O  | 000104-76-7 | 78   |
| 4    |    |   | 1-Dodecene          | 168 | C12H24  | 000112-41-4 | 53   |
| 5    |    |   | 2-Propyl-1-pentanol | 130 | C8H18O  | 058175-57-8 | 53   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

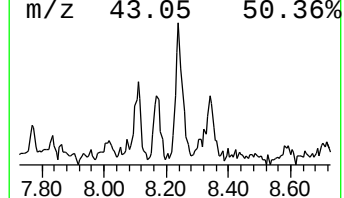
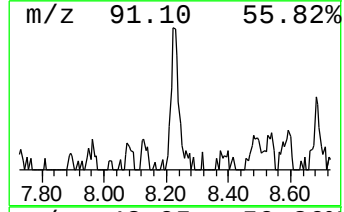
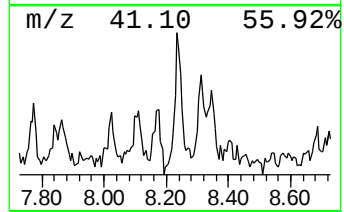
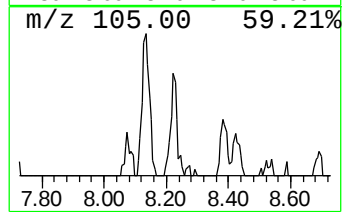
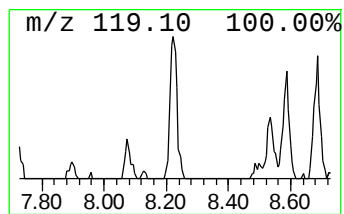
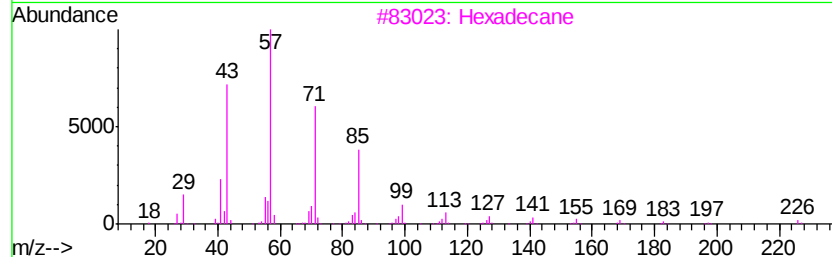
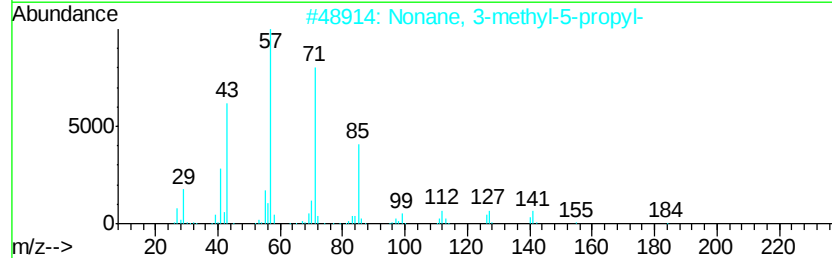
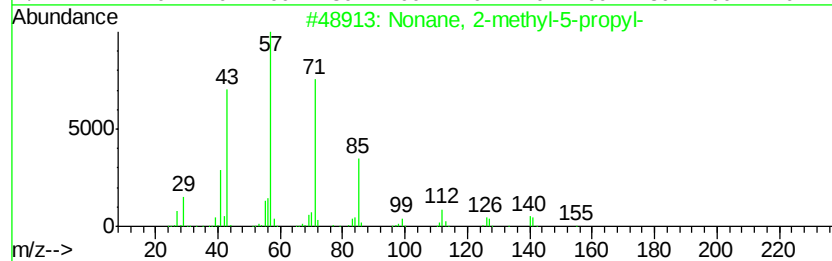
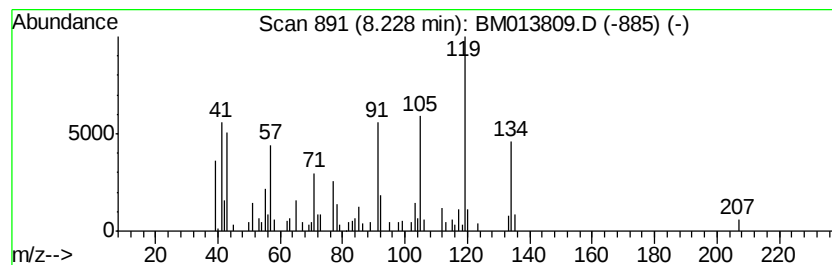
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.23 | 4.31 ng/ul | 189379 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|---------|---|----------------------------|-----|---------|-------------|------|
| 1       |   | Nonane, 2-methyl-5-propyl- | 184 | C13H28  | 031081-17-1 | 47   |
| 2       |   | Nonane, 3-methyl-5-propyl- | 184 | C13H28  | 031081-18-2 | 47   |
| 3       |   | Hexadecane                 | 226 | C16H34  | 000544-76-3 | 43   |
| 4       |   | Pentadecane                | 212 | C15H32  | 000629-62-9 | 43   |
| 5       |   | Decane, 1-iodo-            | 268 | C10H21I | 002050-77-3 | 27   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

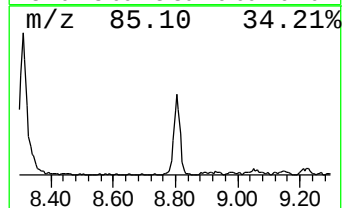
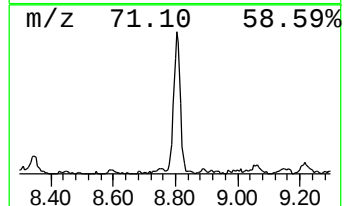
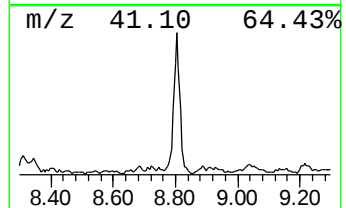
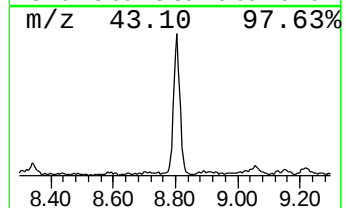
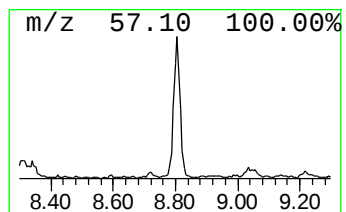
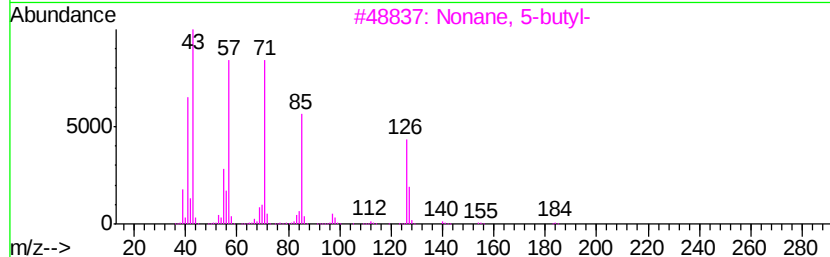
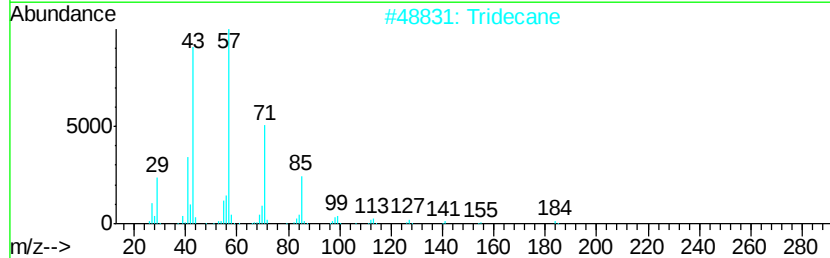
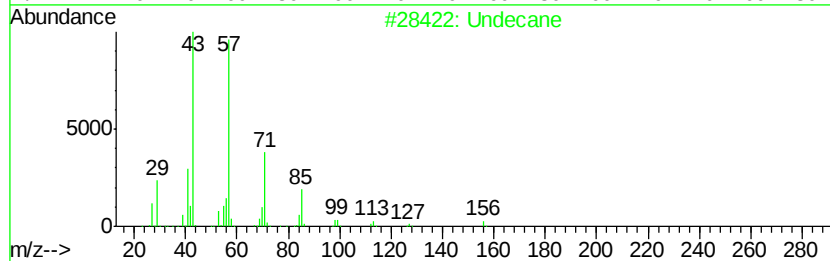
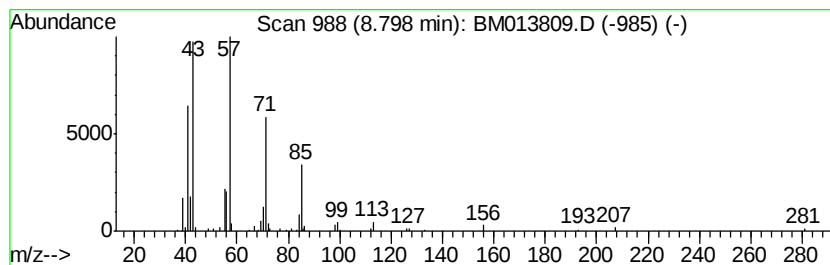
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 3

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 8.80 | 12.37 ng/ul | 543178 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# | of | Tentative ID                    | MW  | MolForm  | CAS#         | Qual |
|------|----|---------------------------------|-----|----------|--------------|------|
| 1    | 5  | Undecane                        | 156 | C11H24   | 001120-21-4  | 76   |
| 2    |    | Tridecane                       | 184 | C13H28   | 000629-50-5  | 72   |
| 3    |    | Nonane, 5-butyl-                | 184 | C13H28   | 017312-63-9  | 64   |
| 4    |    | Oxalic acid, octyl propyl ester | 244 | C13H24O4 | 1000309-26-1 | 59   |
| 5    |    | Bacchotricuneatin c             | 342 | C20H22O5 | 066563-30-2  | 58   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

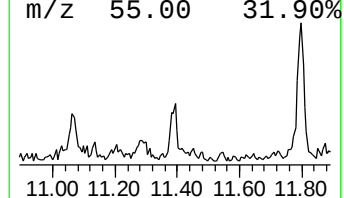
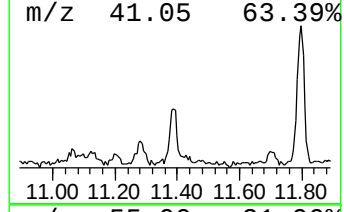
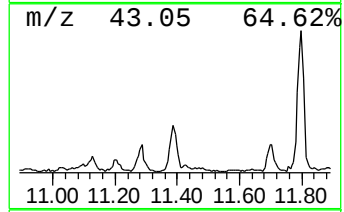
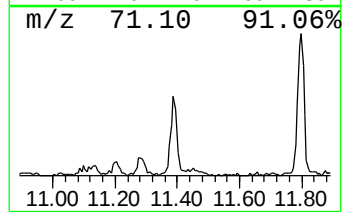
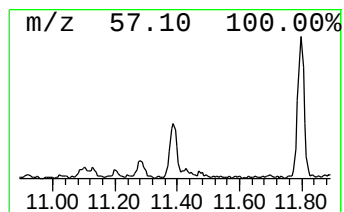
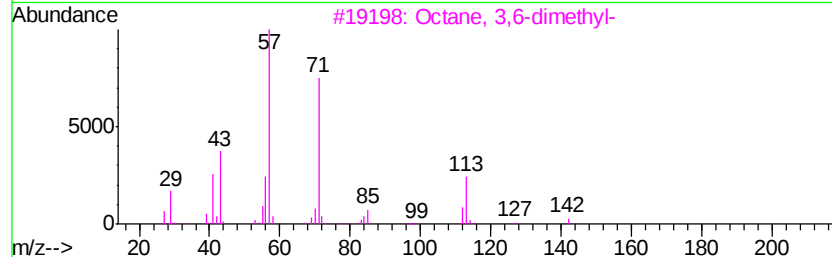
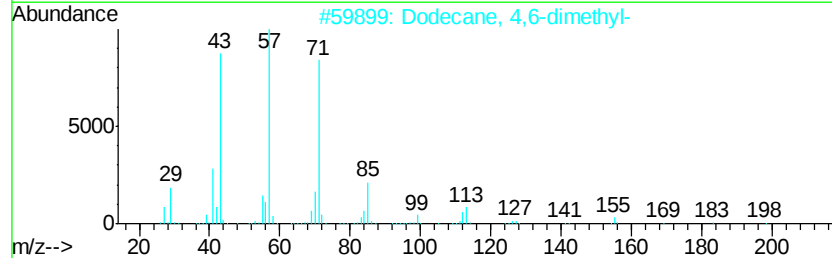
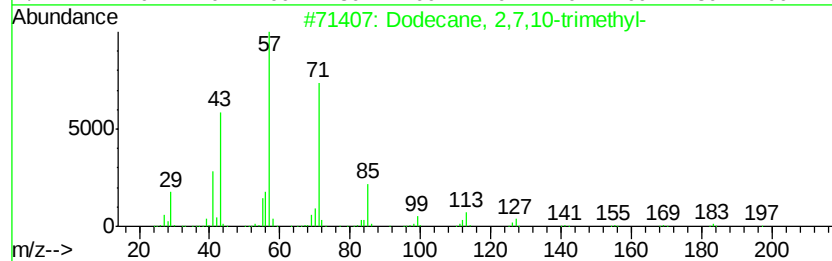
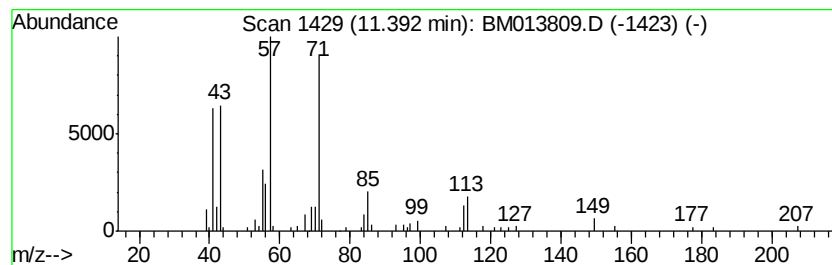
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 17

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.39 | 2.29 ng/ul | 173421 | Naphthalene-d8   | 10.37 |

| Hit# | of | Tentative ID                       | MW  | MolForm   | CAS#         | Qual |
|------|----|------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Dodecane, 2,7,10-trimethyl-        | 212 | C15H32    | 074645-98-0  | 86   |
| 2    |    | Dodecane, 4,6-dimethyl-            | 198 | C14H30    | 061141-72-8  | 64   |
| 3    |    | Octane, 3,6-dimethyl-              | 142 | C10H22    | 015869-94-0  | 64   |
| 4    |    | Sulfurous acid, hexyl pentyl ester | 236 | C11H24O3S | 1000309-14-1 | 58   |
| 5    |    | Dodecane, 2,6,11-trimethyl-        | 212 | C15H32    | 031295-56-4  | 53   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

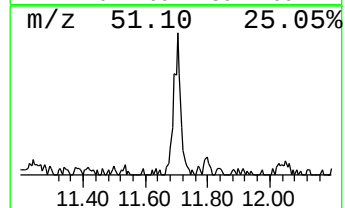
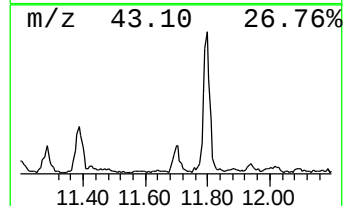
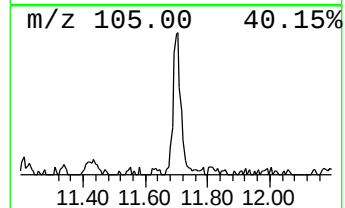
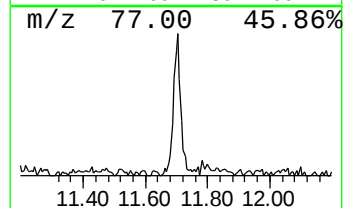
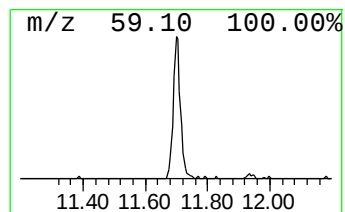
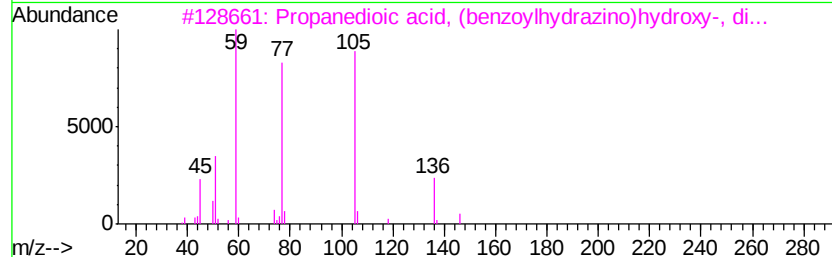
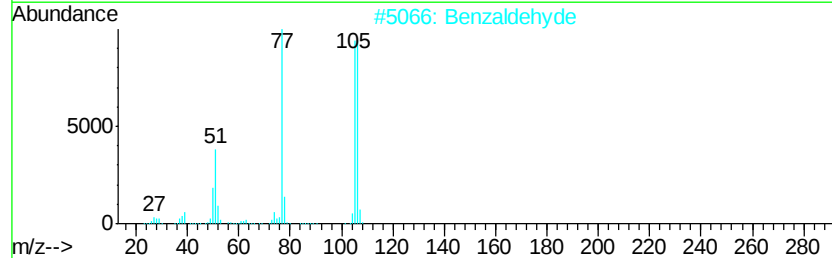
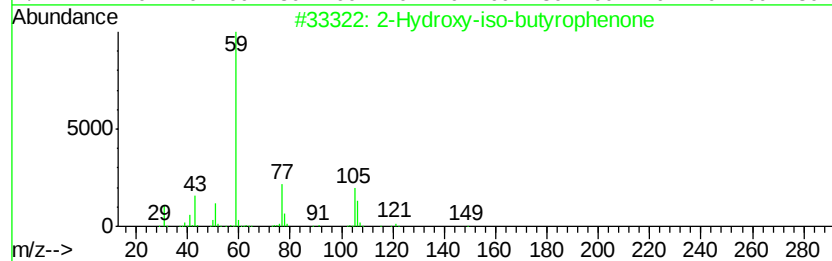
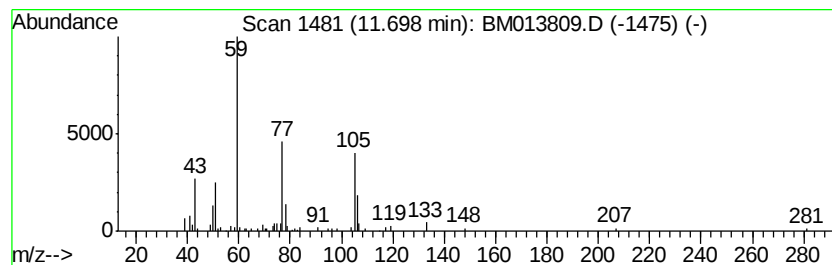
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 unknown-01 Concentration Rank 16

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.70 | 2.40 ng/ul | 182028 | Napthalene-d8    | 10.37 |

| Hit# | of | 5 | Tentative ID                                         | MW  | MolForm    | CAS#         | Qual |
|------|----|---|------------------------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 2-Hydroxy-iso-butyrophenone                          | 164 | C10H12O2   | 007473-98-5  | 43   |
| 2    |    |   | Benzaldehyde                                         | 106 | C7H6O      | 000100-52-7  | 43   |
| 3    |    |   | Propanedioic acid, (benzoylhydrazino)hydroxy-, di... | 282 | C12H14N2O6 | 1000142-99-9 | 40   |
| 4    |    |   | N-Benzoyl-D-threo-2-methylthreonine                  | 237 | C12H15NO4  | 131644-54-7  | 39   |
| 5    |    |   | .alpha.-Benzamido-2-hydroxycinnam...                 | 283 | C16H13NO4  | 1000223-61-7 | 27   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

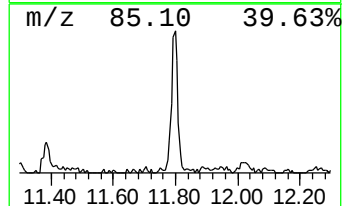
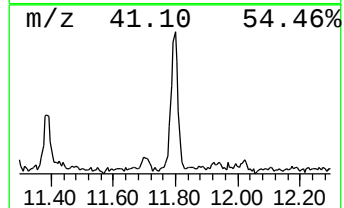
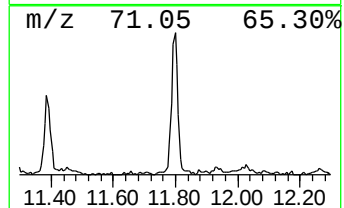
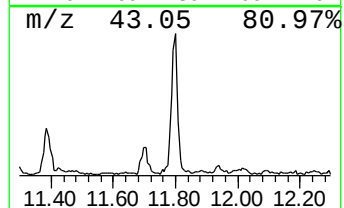
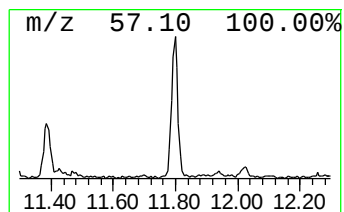
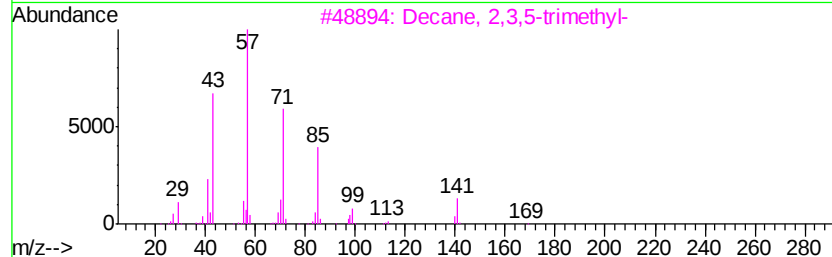
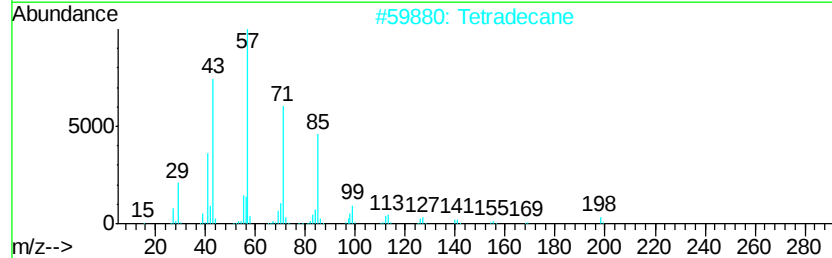
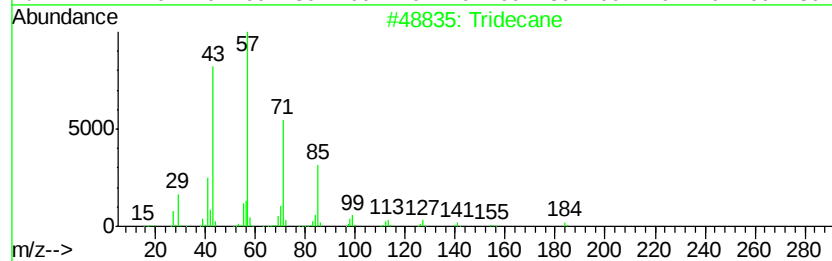
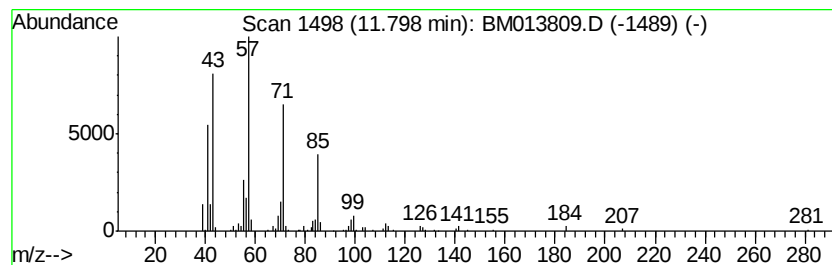
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.80 | 5.62 ng/ul | 426167 | Napthalene-d8    | 10.37 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Tridecane                | 184 | C13H28  | 000629-50-5 | 94   |
| 2    |    |   | Tetradecane              | 198 | C14H30  | 000629-59-4 | 90   |
| 3    |    |   | Decane, 2,3,5-trimethyl- | 184 | C13H28  | 062238-11-3 | 90   |
| 4    |    |   | Tetradecane              | 198 | C14H30  | 000629-59-4 | 83   |
| 5    |    |   | Pentadecane              | 212 | C15H32  | 000629-62-9 | 83   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

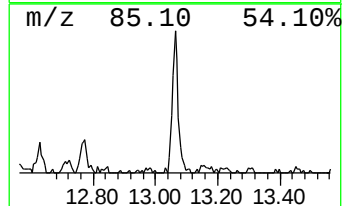
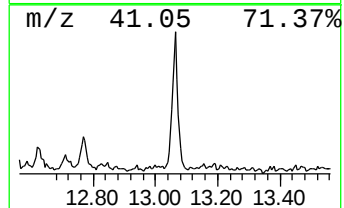
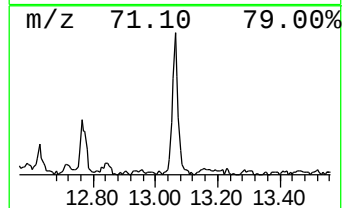
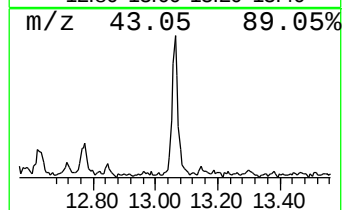
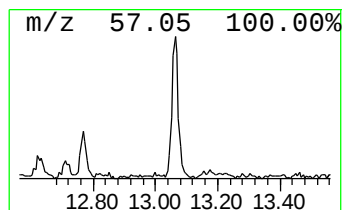
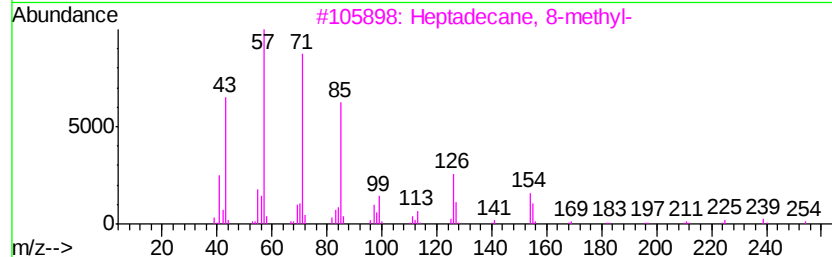
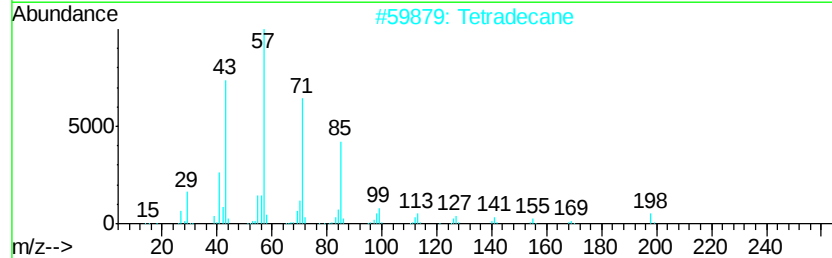
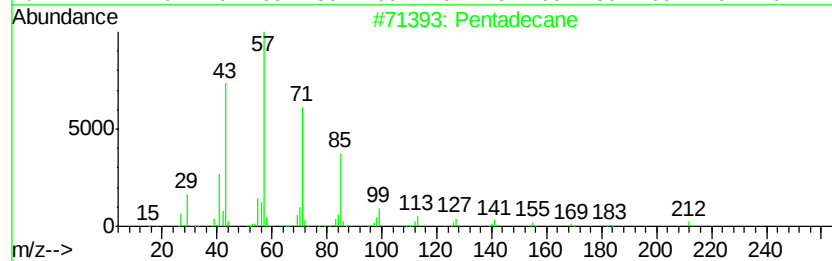
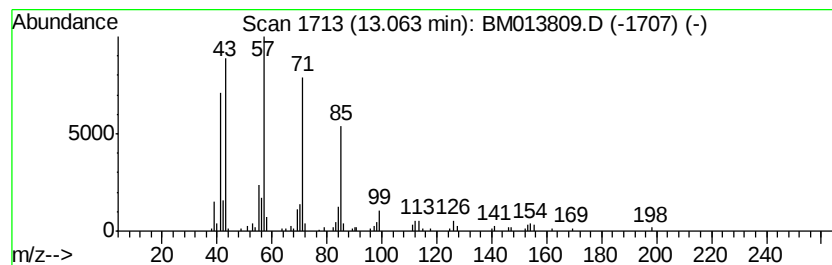
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.06 | 4.25 ng/ul | 261974 | Acenaphthene-d10 | 14.24 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentadecane            | 212 | C15H32  | 000629-62-9 | 87   |
| 2    |    |   | Tetradecane            | 198 | C14H30  | 000629-59-4 | 86   |
| 3    |    |   | Heptadecane, 8-methyl- | 254 | C18H38  | 013287-23-5 | 86   |
| 4    |    |   | Tridecane              | 184 | C13H28  | 000629-50-5 | 86   |
| 5    |    |   | Tetradecane            | 198 | C14H30  | 000629-59-4 | 86   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

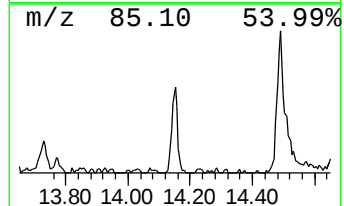
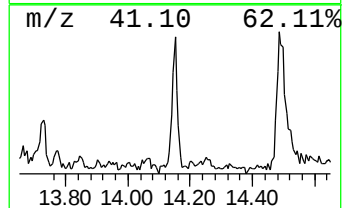
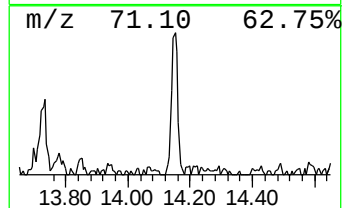
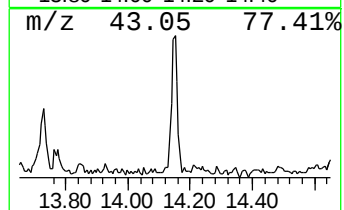
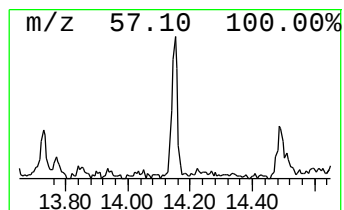
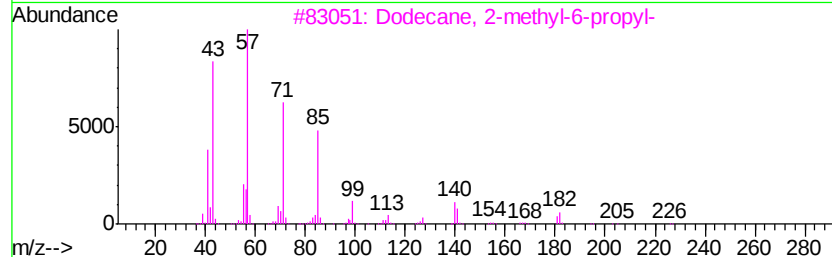
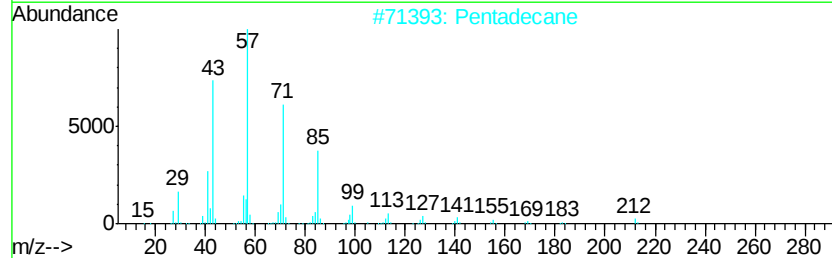
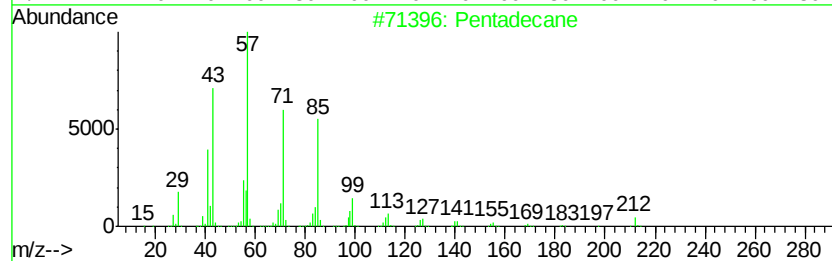
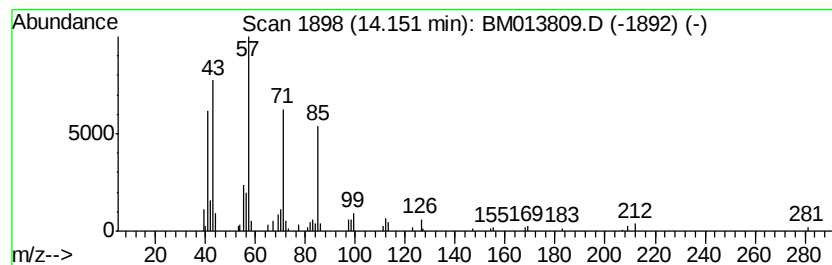
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.15 | 2.51 ng/ul | 154988 | Acenaphthene-d10 | 14.24 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentadecane                  | 212 | C15H32  | 000629-62-9 | 91   |
| 2    |    |   | Pentadecane                  | 212 | C15H32  | 000629-62-9 | 81   |
| 3    |    |   | Dodecane, 2-methyl-6-propyl- | 226 | C16H34  | 055045-08-4 | 78   |
| 4    |    |   | Decane, 2,4-dimethyl-        | 170 | C12H26  | 002801-84-5 | 72   |
| 5    |    |   | Tetradecane                  | 198 | C14H30  | 000629-59-4 | 72   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

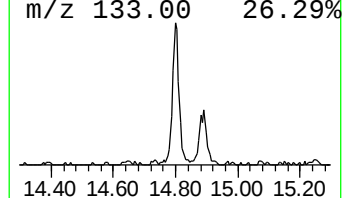
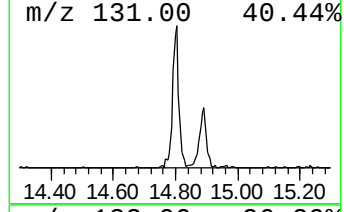
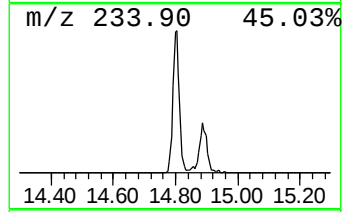
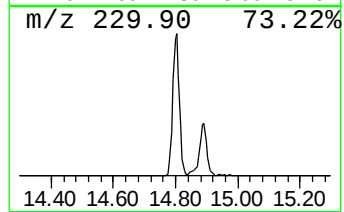
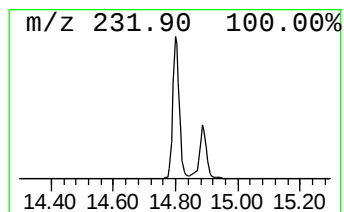
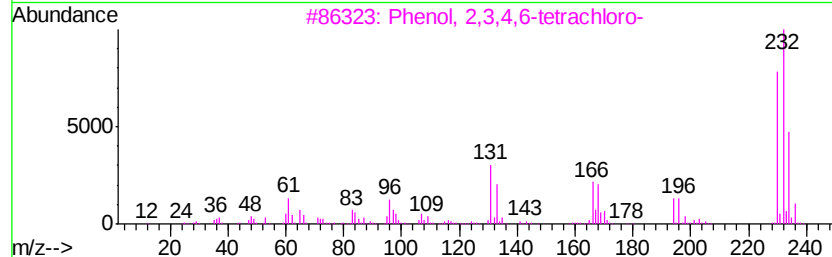
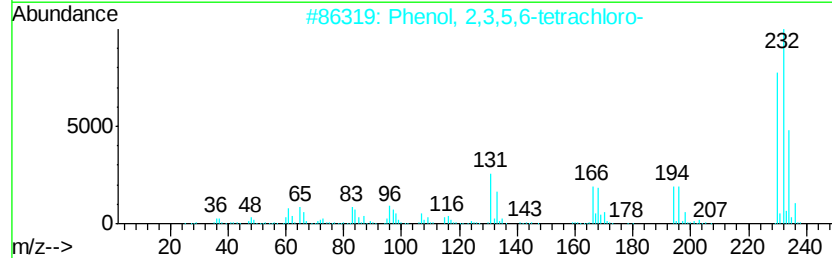
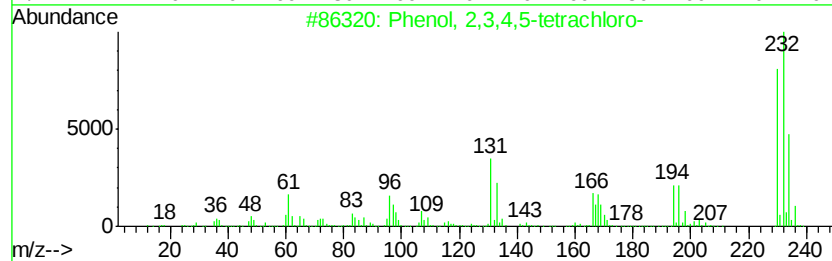
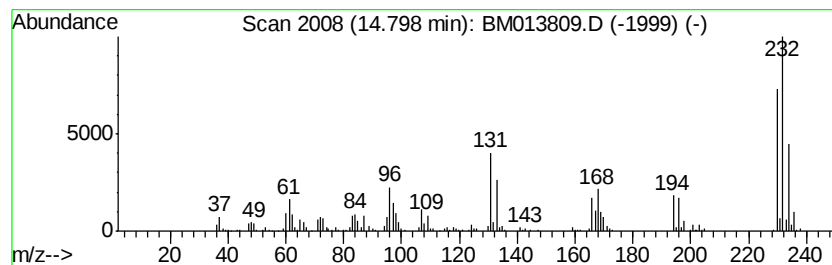
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Phenol, 2,3,4,5-tetrachloro- Concentration Rank 2

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 14.80 | 14.45 ng/ul | 891476 | Acenaphthene-d10 | 14.24 |

| Hit# | of | Tentative ID                 | MW  | MolForm  | CAS#        | Qual |
|------|----|------------------------------|-----|----------|-------------|------|
| 1    | 5  | Phenol, 2,3,4,5-tetrachloro- | 230 | C6H2Cl4O | 004901-51-3 | 99   |
| 2    |    | Phenol, 2,3,5,6-tetrachloro- | 230 | C6H2Cl4O | 000935-95-5 | 98   |
| 3    |    | Phenol, 2,3,4,6-tetrachloro- | 230 | C6H2Cl4O | 000058-90-2 | 96   |
| 4    |    | Phenol, 2,3,4,6-tetrachloro- | 230 | C6H2Cl4O | 000058-90-2 | 95   |
| 5    |    | Phenol, 2,3,4,6-tetrachloro- | 230 | C6H2Cl4O | 000058-90-2 | 86   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

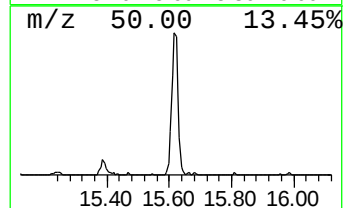
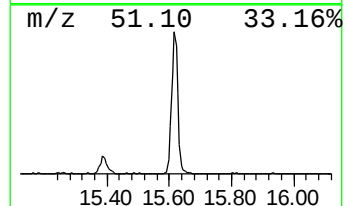
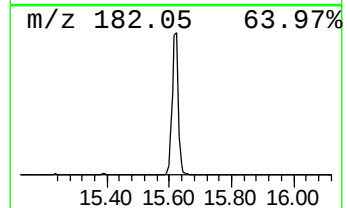
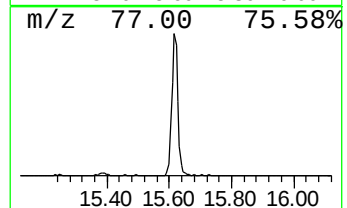
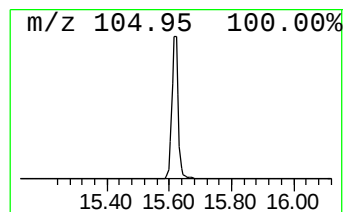
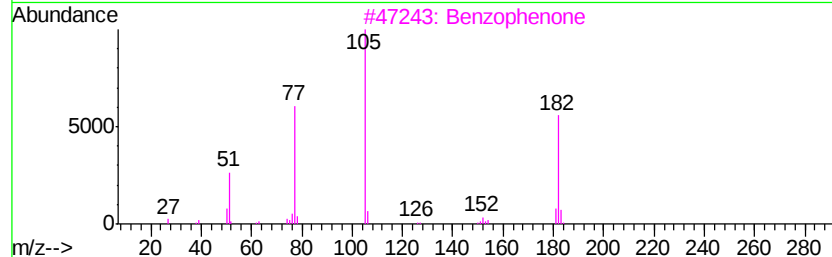
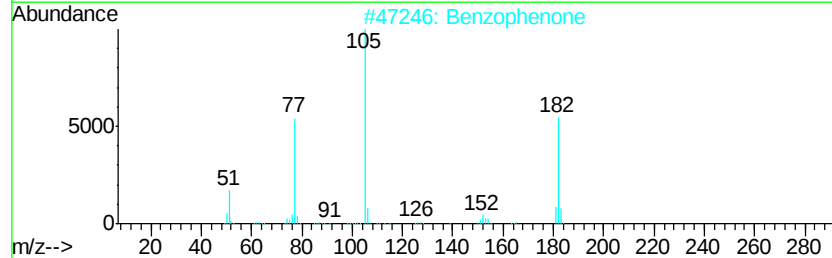
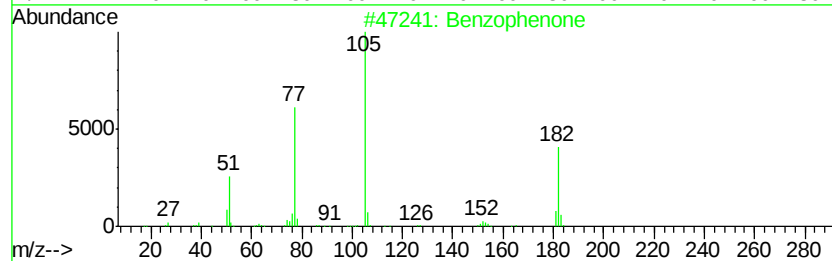
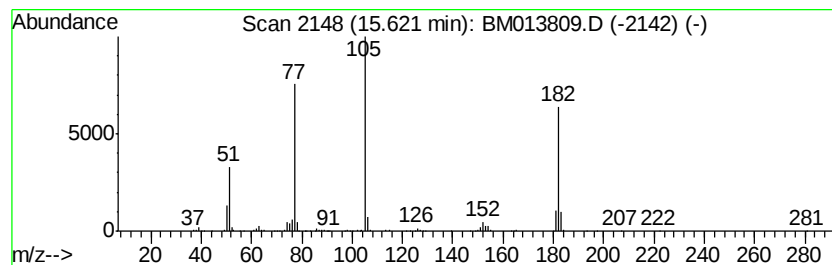
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Benzophenone Concentration Rank 1

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.62 | 19.07 ng/ul | 1176270 | Acenaphthene-d10 | 14.24 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Benzophenone | 182 | C13H10O | 000119-61-9 | 97   |
| 2    |    | Benzophenone | 182 | C13H10O | 000119-61-9 | 97   |
| 3    |    | Benzophenone | 182 | C13H10O | 000119-61-9 | 96   |
| 4    |    | Benzophenone | 182 | C13H10O | 000119-61-9 | 93   |
| 5    |    | Benzophenone | 182 | C13H10O | 000119-61-9 | 91   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

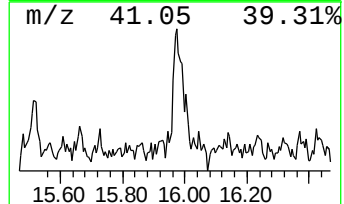
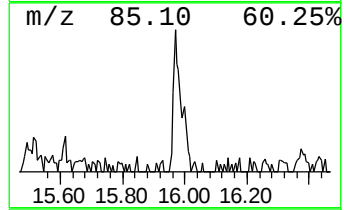
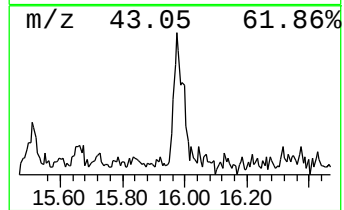
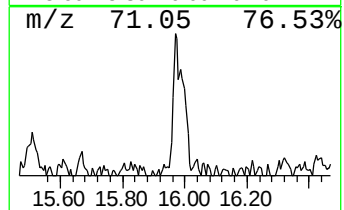
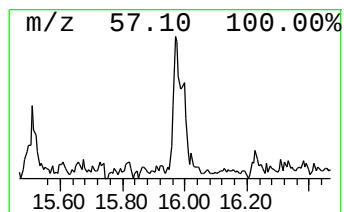
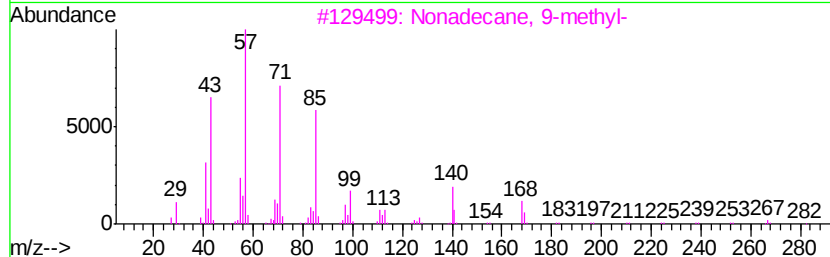
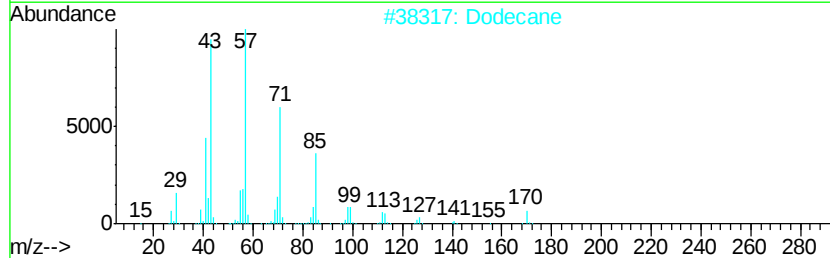
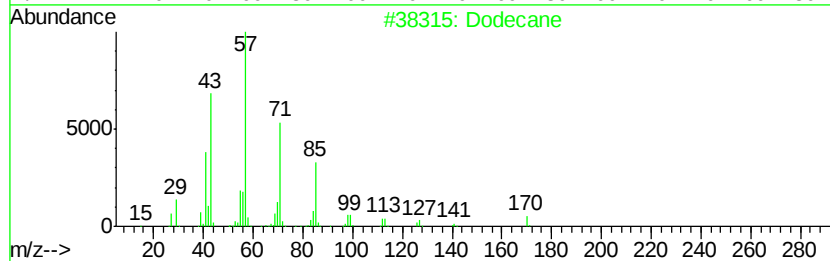
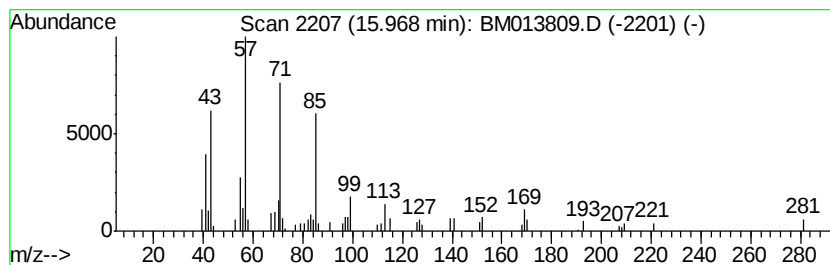
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.97 | 2.50 ng/ul | 172964 | Phenanthrene-d10 | 17.00 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Dodecane              | 170 | C12H26  | 000112-40-3 | 76   |
| 2    |    |   | Dodecane              | 170 | C12H26  | 000112-40-3 | 64   |
| 3    |    |   | Nonadecane, 9-methyl- | 282 | C20H42  | 013287-24-6 | 62   |
| 4    |    |   | Nonadecane            | 268 | C19H40  | 000629-92-5 | 58   |
| 5    |    |   | Tetracosane           | 338 | C24H50  | 000646-31-1 | 58   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

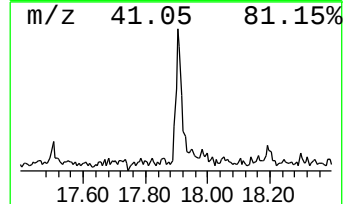
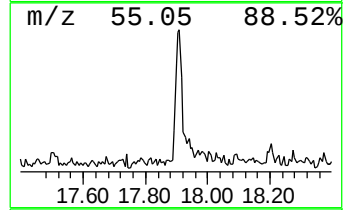
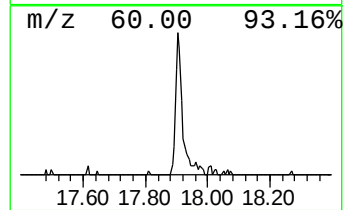
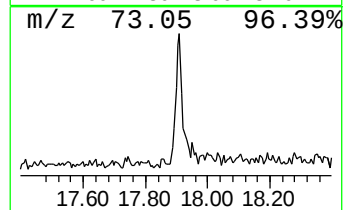
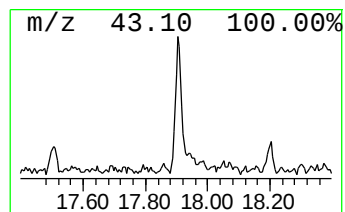
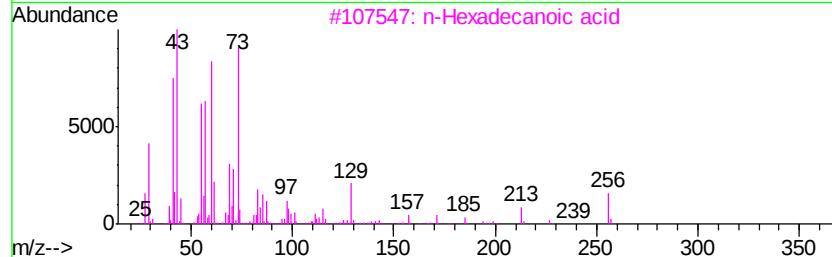
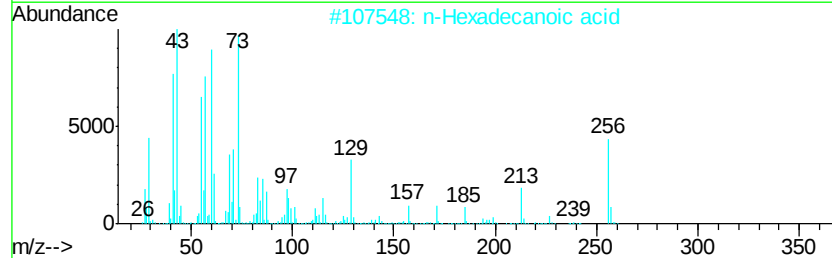
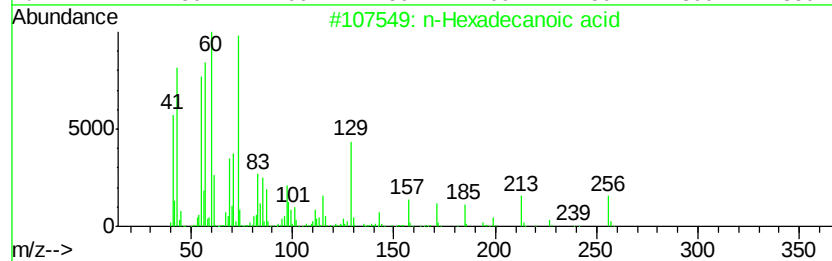
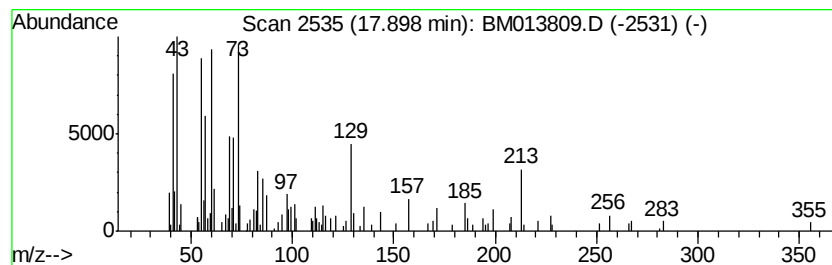
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 n-Hexadecanoic acid Concentration Rank 8

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.90 | 4.36 ng/ul | 302012 | Phenanthrene-d10 | 17.00 |

| Hit# | of | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------|-----|----------|-------------|------|
| 1    | 5  | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 90   |
| 2    |    | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 68   |
| 3    |    | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 60   |
| 4    |    | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 58   |
| 5    |    | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 58   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM012518\  
Data File : BM013809.D  
Acq On : 25 Jan 2018 17:13  
Operator : SJ/JU  
Sample : J1243-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

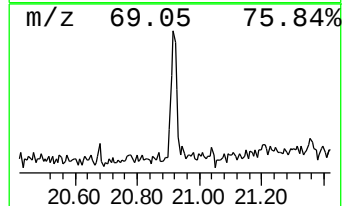
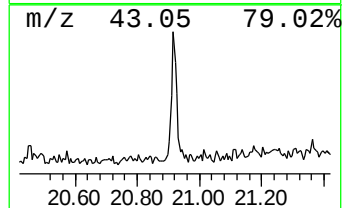
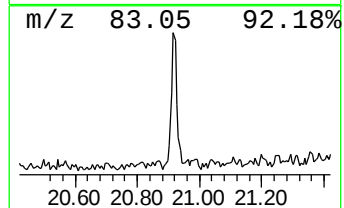
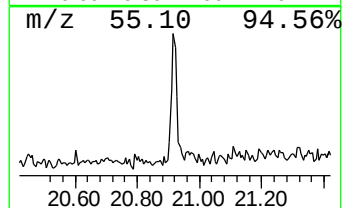
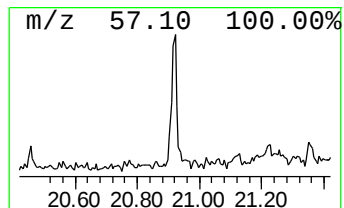
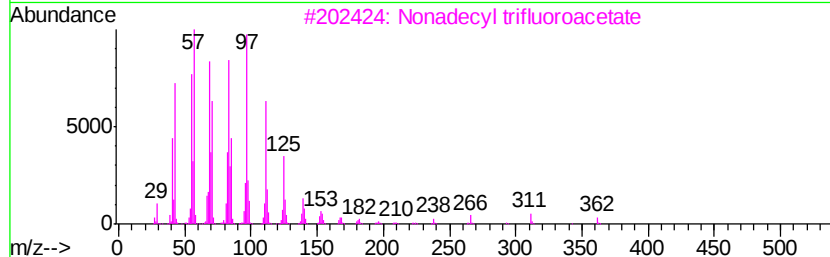
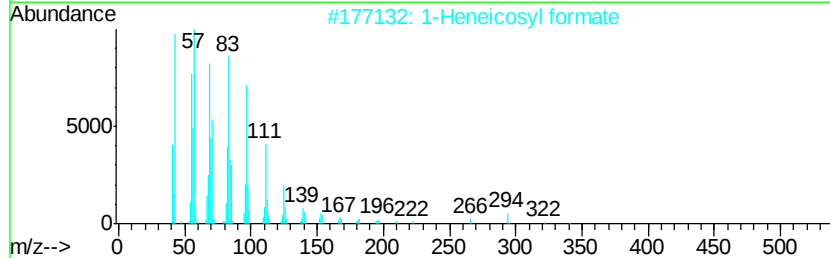
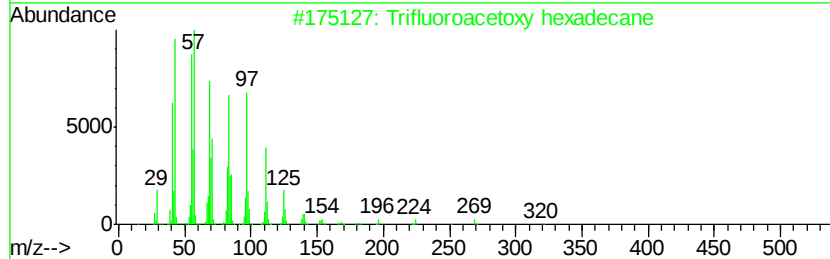
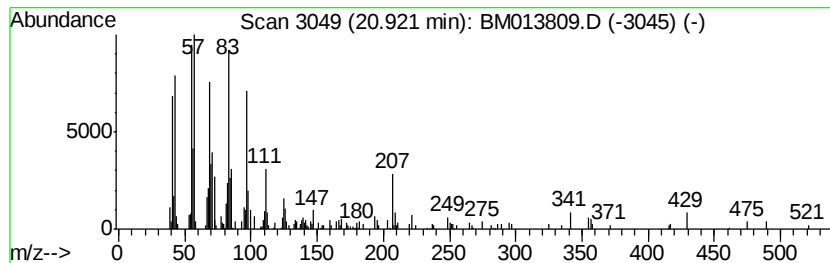
Instrument :  
BNA\_M  
ClientSampleId :  
F6B27

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

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

\*\*\*\*\*  
Peak Number 19 Trifluoroacetoxy hexadecane Concentration Rank 13

| R.T.    | EstConc    | Area                              | Relative to ISTD | R.T.       |              |      |
|---------|------------|-----------------------------------|------------------|------------|--------------|------|
| 20.92   | 3.12 ng/ul | 269009                            | Chrysene-d12     | 21.20      |              |      |
| Hit# of | 5          | Tentative ID                      | MW               | MolForm    | CAS#         | Qual |
| 1       |            | Trifluoroacetoxy hexadecane       | 338              | C18H33F3O2 | 006222-03-3  | 87   |
| 2       |            | 1-Heneicosyl formate              | 340              | C22H44O2   | 077899-03-7  | 76   |
| 3       |            | Nonadecyl trifluoroacetate        | 380              | C21H39F3O2 | 1000351-76-3 | 76   |
| 4       |            | Bromoacetic acid, hexadecyl ester | 362              | C18H35BrO2 | 005454-48-8  | 70   |
| 5       |            | 1-Docosene                        | 308              | C22H44     | 001599-67-3  | 64   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM012518\  
 Data File : BM013809.D  
 Acq On : 25 Jan 2018 17:13  
 Operator : SJ/JU  
 Sample : J1243-04  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 F6B27

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

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| 2-Pentanone, 4-hy... | 4.76  | 4.7     | ng/ul | 204315   | 1                     | 7.59  | 878359  | 20.0 |
| (DEL) Alkane: Str... | 5.62  | 6.5     | ng/ul | 285571   | 1                     | 7.59  | 878359  | 20.0 |
| Hexylene glycol      | 5.96  | 3.5     | ng/ul | 155126   | 1                     | 7.59  | 878359  | 20.0 |
| (DEL) Alkane: Str... | 6.58  | 2.1     | ng/ul | 91591    | 1                     | 7.59  | 878359  | 20.0 |
| (DEL) Alkane: Str... | 6.64  | 2.1     | ng/ul | 91703    | 1                     | 7.59  | 878359  | 20.0 |
| (DEL) Alkane: Str... | 7.21  | 10.3    | ng/ul | 453873   | 1                     | 7.59  | 878359  | 20.0 |
| 1-Hexanol, 2-ethyl-  | 7.66  | 3.9     | ng/ul | 169798   | 1                     | 7.59  | 878359  | 20.0 |
| (DEL) Alkane: Str... | 8.23  | 4.3     | ng/ul | 189379   | 1                     | 7.59  | 878359  | 20.0 |
| (DEL) Alkane: Str... | 8.80  | 12.4    | ng/ul | 543178   | 1                     | 7.59  | 878359  | 20.0 |
| (DEL) Alkane: Str... | 11.39 | 2.3     | ng/ul | 173421   | 2                     | 10.37 | 1516500 | 20.0 |
| unknown-01           | 11.70 | 2.4     | ng/ul | 182028   | 2                     | 10.37 | 1516500 | 20.0 |
| (DEL) Alkane: Str... | 11.80 | 5.6     | ng/ul | 426167   | 2                     | 10.37 | 1516500 | 20.0 |
| (DEL) Alkane: Str... | 13.06 | 4.3     | ng/ul | 261974   | 3                     | 14.24 | 1233570 | 20.0 |
| (DEL) Alkane: Str... | 14.15 | 2.5     | ng/ul | 154988   | 3                     | 14.24 | 1233570 | 20.0 |
| Phenol, 2,3,4,5-t... | 14.80 | 14.4    | ng/ul | 891476   | 3                     | 14.24 | 1233570 | 20.0 |
| Benzophenone         | 15.62 | 19.1    | ng/ul | 1176270  | 3                     | 14.24 | 1233570 | 20.0 |
| (DEL) Alkane: Str... | 15.97 | 2.5     | ng/ul | 172964   | 4                     | 17.00 | 1386450 | 20.0 |
| n-Hexadecanoic acid  | 17.90 | 4.4     | ng/ul | 302012   | 4                     | 17.00 | 1386450 | 20.0 |
| Trifluoroacetoxy ... | 20.92 | 3.1     | ng/ul | 269009   | 5                     | 21.20 | 1723930 | 20.0 |