

Data Path : U:\HPCHEM1\BNA M\DATA\BM012818\  
 Data File : BM013946.D  
 Acq On : 29 Jan 2018 05:24  
 Operator : SJ/JU  
 Sample : J1244-04  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 F6B45

## 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         | 3.140       | 23            | 26          | 30           | rVB3     | 29073          | 34653         | 1.17%           | 0.083%        |
| 2         | 3.751       | 126           | 130         | 135          | rVB      | 41040          | 59570         | 2.00%           | 0.143%        |
| 3         | 4.163       | 194           | 200         | 205          | rBV      | 117782         | 171283        | 5.76%           | 0.410%        |
| 4         | 4.734       | 291           | 297         | 309          | rVB2     | 202566         | 479536        | 16.13%          | 1.149%        |
| 5         | 5.345       | 394           | 401         | 409          | rVB      | 96204          | 147833        | 4.97%           | 0.354%        |
| 6         | 6.763       | 635           | 642         | 654          | rBV2     | 588253         | 1079408       | 36.30%          | 2.586%        |
| 7         | 6.922       | 662           | 669         | 676          | rBV      | 473667         | 715495        | 24.06%          | 1.714%        |
| 8         | 7.110       | 695           | 701         | 710          | rBV      | 642725         | 1050219       | 35.32%          | 2.516%        |
| 9         | 7.575       | 773           | 780         | 787          | rBV      | 796110         | 1249952       | 42.03%          | 2.995%        |
| 10        | 8.292       | 894           | 902         | 911          | rBV      | 710636         | 1203896       | 40.48%          | 2.884%        |
| 11        | 8.404       | 916           | 921         | 931          | rVB2     | 33607          | 59803         | 2.01%           | 0.143%        |
| 12        | 8.733       | 970           | 977         | 987          | rBV      | 640722         | 1064632       | 35.80%          | 2.551%        |
| 13        | 8.863       | 995           | 999         | 1008         | rVB7     | 23778          | 47127         | 1.58%           | 0.113%        |
| 14        | 9.128       | 1039          | 1044        | 1052         | rVB3     | 31330          | 47775         | 1.61%           | 0.114%        |
| 15        | 9.445       | 1090          | 1098        | 1109         | rBV      | 551035         | 963857        | 32.41%          | 2.309%        |
| 16        | 9.986       | 1183          | 1190        | 1203         | rBV      | 777404         | 1406779       | 47.31%          | 3.370%        |
| 17        | 10.351      | 1241          | 1252        | 1258         | rBV      | 1037251        | 1744915       | 58.68%          | 4.180%        |
| 18        | 10.504      | 1270          | 1278        | 1292         | rBV      | 485398         | 982961        | 33.05%          | 2.355%        |
| 19        | 11.898      | 1508          | 1515        | 1520         | rBV5     | 36397          | 71486         | 2.40%           | 0.171%        |
| 20        | 13.651      | 1805          | 1813        | 1817         | rBV      | 1415827        | 2033447       | 68.38%          | 4.872%        |
| 21        | 13.698      | 1817          | 1821        | 1827         | rVB      | 406237         | 585879        | 19.70%          | 1.404%        |
| 22        | 13.915      | 1850          | 1858        | 1869         | rBV      | 1553844        | 2453542       | 82.50%          | 5.878%        |
| 23        | 14.133      | 1888          | 1895        | 1903         | rBV2     | 53999          | 78113         | 2.63%           | 0.187%        |
| 24        | 14.227      | 1903          | 1911        | 1919         | rVV2     | 1409586        | 2211117       | 74.35%          | 5.297%        |
| 25        | 14.474      | 1947          | 1953        | 1969         | rBV      | 302440         | 704684        | 23.70%          | 1.688%        |
| 26        | 14.639      | 1977          | 1981        | 1984         | rBV5     | 21728          | 35237         | 1.18%           | 0.084%        |
| 27        | 15.039      | 2045          | 2049        | 2052         | rBV      | 39372          | 48232         | 1.62%           | 0.116%        |
| 28        | 15.092      | 2052          | 2058        | 2064         | rVV3     | 69289          | 91162         | 3.07%           | 0.218%        |
| 29        | 15.227      | 2075          | 2081        | 2087         | rBV      | 1962435        | 2780402       | 93.50%          | 6.661%        |
| 30        | 15.374      | 2100          | 2106        | 2116         | rBV      | 425070         | 668876        | 22.49%          | 1.602%        |
| 31        | 15.474      | 2119          | 2123        | 2125         | rBV2     | 21530          | 33403         | 1.12%           | 0.080%        |
| 32        | 15.604      | 2139          | 2145        | 2150         | rBV      | 242832         | 347792        | 11.70%          | 0.833%        |
| 33        | 15.957      | 2201          | 2205        | 2213         | rBV4     | 74376          | 151420        | 5.09%           | 0.363%        |
| 34        | 16.645      | 2316          | 2322        | 2329         | rBV      | 387146         | 634492        | 21.34%          | 1.520%        |

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ALS Vial : 28 Sample Multiplier: 1

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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

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Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |         |        |
|----|--------|------|------|------|------|---------|---------|---------|--------|
| 35 | 16.751 | 2337 | 2340 | 2345 | rBV2 | 57821   | 72498   | 2.44%   | 0.174% |
| 36 | 16.986 | 2374 | 2380 | 2385 | rBV2 | 1545286 | 2269899 | 76.33%  | 5.438% |
| 37 | 17.021 | 2385 | 2386 | 2391 | rVB  | 76282   | 85569   | 2.88%   | 0.205% |
| 38 | 17.086 | 2391 | 2397 | 2402 | rBV  | 1853672 | 2697157 | 90.70%  | 6.462% |
| 39 | 17.486 | 2462 | 2465 | 2470 | rVB6 | 51608   | 66333   | 2.23%   | 0.159% |
| 40 | 17.892 | 2529 | 2534 | 2544 | rBV2 | 363630  | 580116  | 19.51%  | 1.390% |
| 41 | 19.050 | 2727 | 2731 | 2738 | rVB  | 316723  | 456347  | 15.35%  | 1.093% |
| 42 | 19.180 | 2750 | 2753 | 2761 | rBV2 | 284346  | 409568  | 13.77%  | 0.981% |
| 43 | 19.386 | 2783 | 2788 | 2797 | rBV2 | 2016466 | 2973848 | 100.00% | 7.125% |
| 44 | 20.915 | 3045 | 3048 | 3054 | rVB4 | 199698  | 278168  | 9.35%   | 0.666% |
| 45 | 21.186 | 3090 | 3094 | 3099 | rBV  | 1614396 | 1999682 | 67.24%  | 4.791% |
| 46 | 23.238 | 3438 | 3443 | 3459 | rVB2 | 1294683 | 2553578 | 85.87%  | 6.118% |
| 47 | 23.374 | 3461 | 3466 | 3473 | rVB  | 1047404 | 1858487 | 62.49%  | 4.453% |

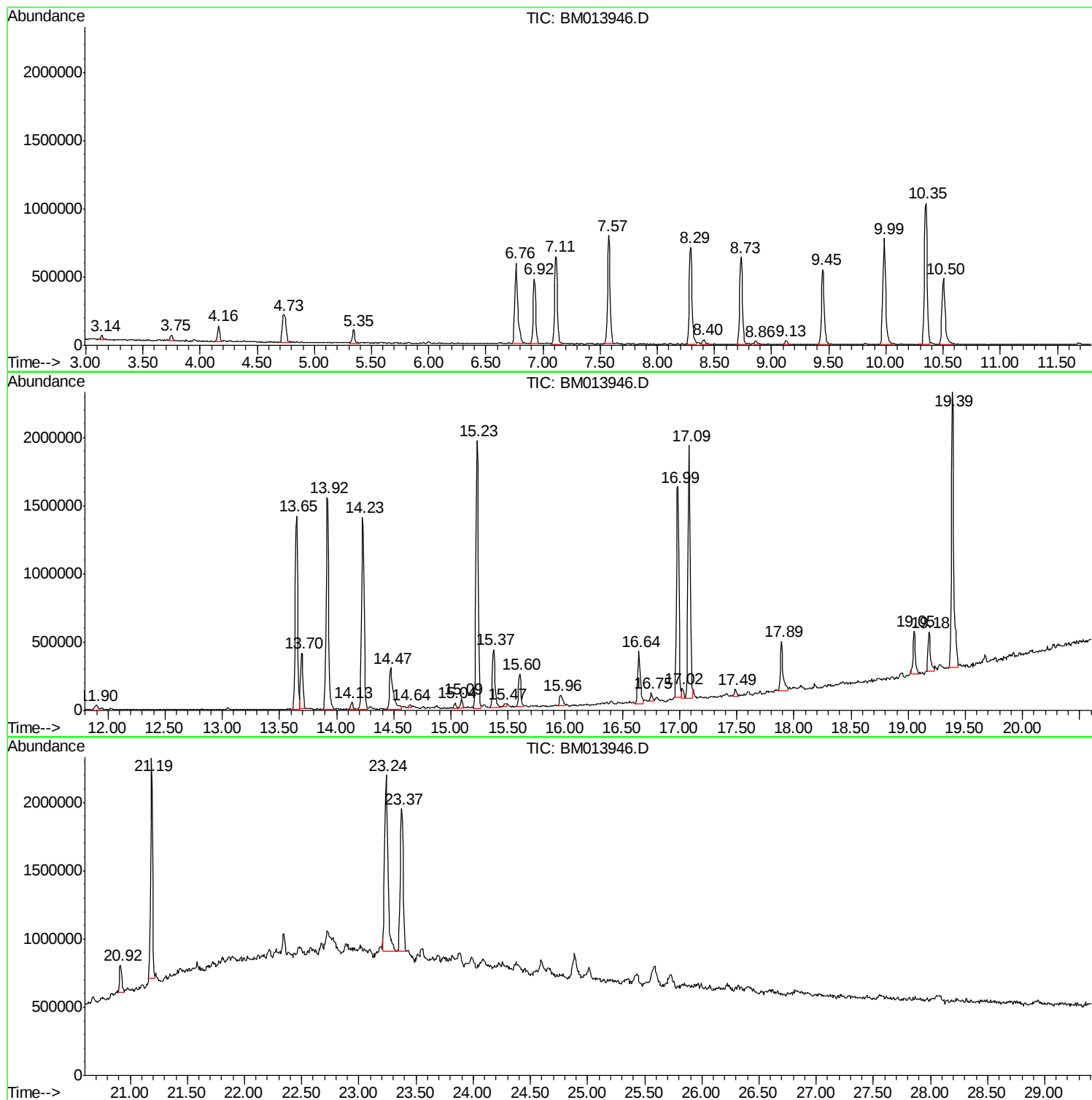
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

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



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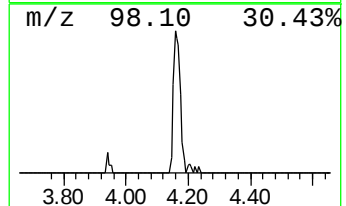
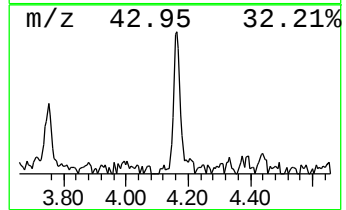
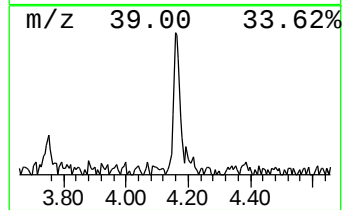
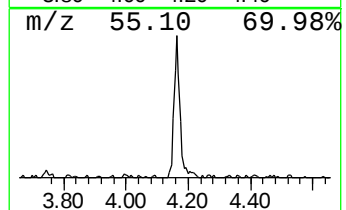
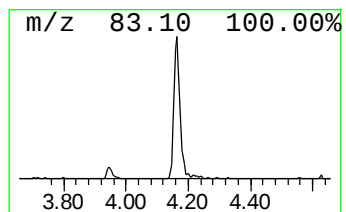
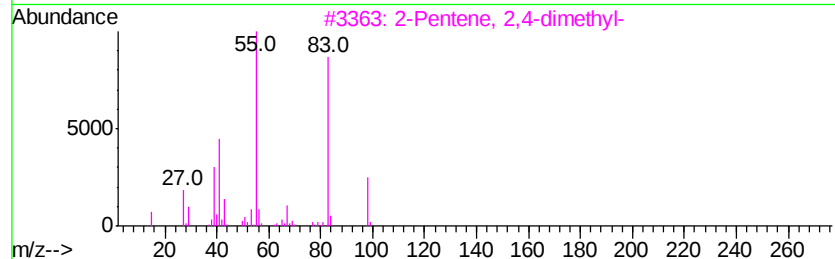
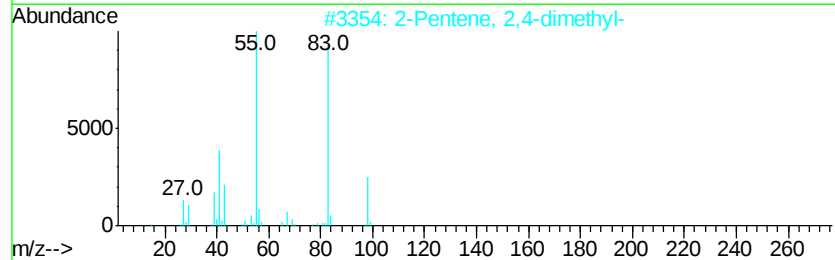
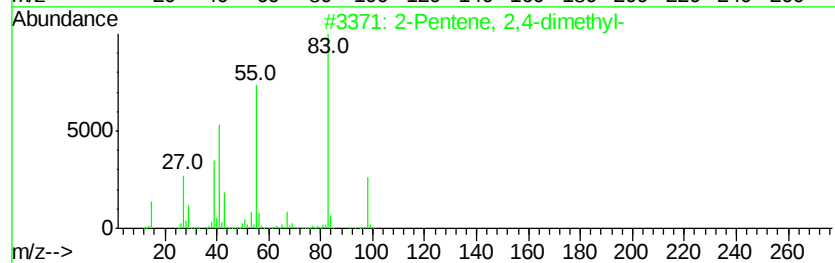
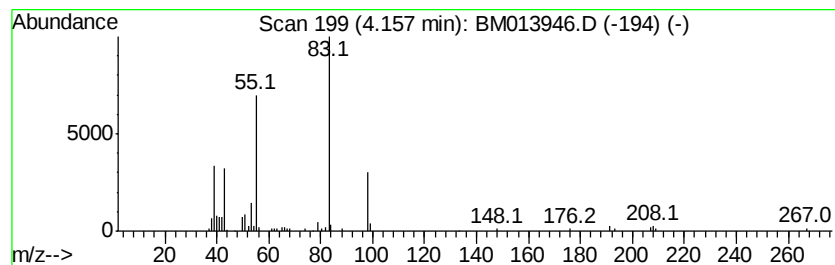
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 1 (DEL) Alkane: Cyclic4.16 Concentration Rank 6

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 4.16 | 2.74 ng/ul | 171283 | 1,4-Dichlorobenzene-d4 | 7.57 |

| Hit# | of                             | 5 | Tentative ID | MW | MolForm | CAS#        | Qual |
|------|--------------------------------|---|--------------|----|---------|-------------|------|
| 1    | 2-Pentene, 2,4-dimethyl-       |   |              | 98 | C7H14   | 000625-65-0 | 86   |
| 2    | 2-Pentene, 2,4-dimethyl-       |   |              | 98 | C7H14   | 000625-65-0 | 80   |
| 3    | 2-Pentene, 2,4-dimethyl-       |   |              | 98 | C7H14   | 000625-65-0 | 78   |
| 4    | 2-Pentene, 2,3-dimethyl-       |   |              | 98 | C7H14   | 010574-37-5 | 78   |
| 5    | 2-Pentene, 3,4-dimethyl-, (E)- |   |              | 98 | C7H14   | 004914-92-5 | 78   |



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

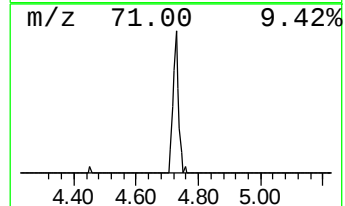
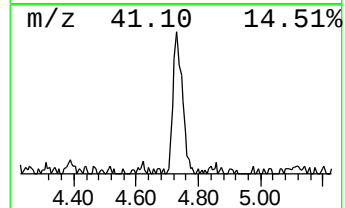
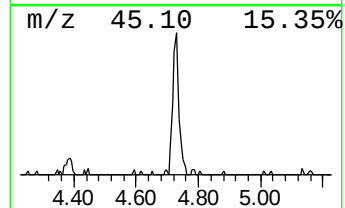
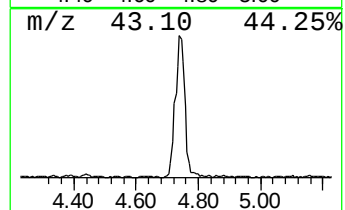
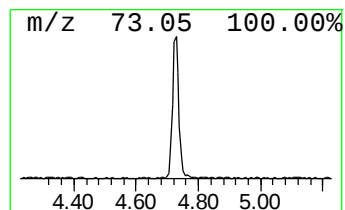
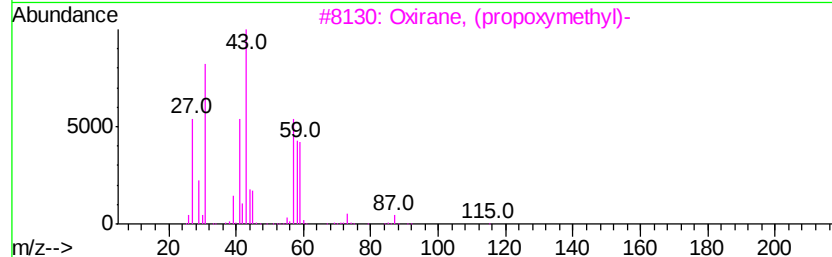
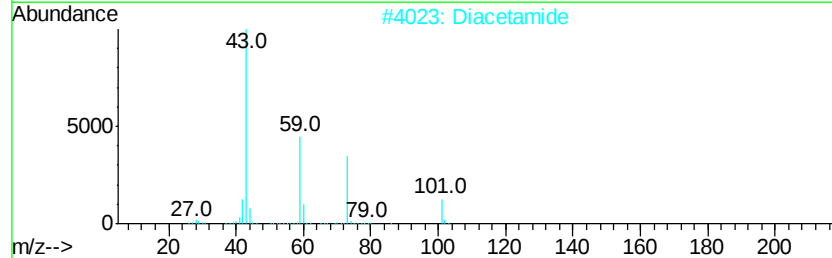
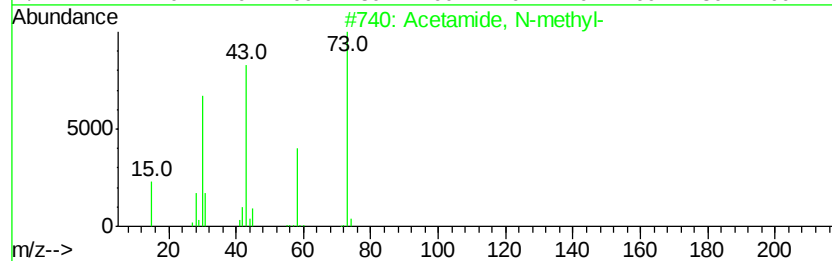
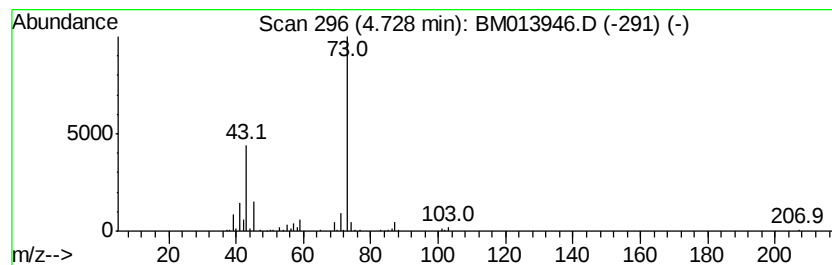
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 unknown-01 Concentration Rank 1

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 4.73 | 7.67 ng/ul | 479536 | 1,4-Dichlorobenzene-d4 | 7.57 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Acetamide, N-methyl-      | 73  | C3H7NO  | 000079-16-3 | 43   |
| 2    |    | Diacetamide               | 101 | C4H7NO2 | 000625-77-4 | 38   |
| 3    |    | Oxirane, (propoxymethyl)- | 116 | C6H12O2 | 003126-95-2 | 33   |
| 4    |    | Oxirane, (propoxymethyl)- | 116 | C6H12O2 | 003126-95-2 | 33   |
| 5    |    | Propanoic acid, 2-methyl- | 88  | C4H8O2  | 000079-31-2 | 25   |



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

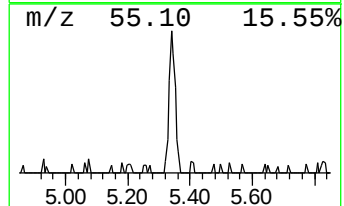
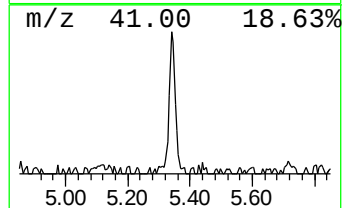
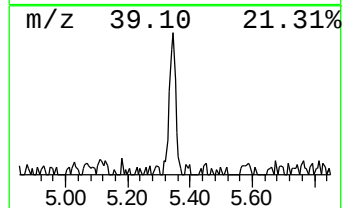
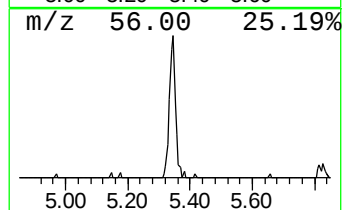
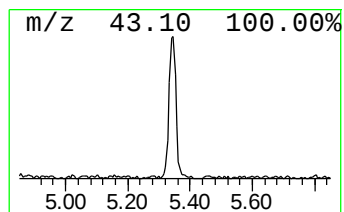
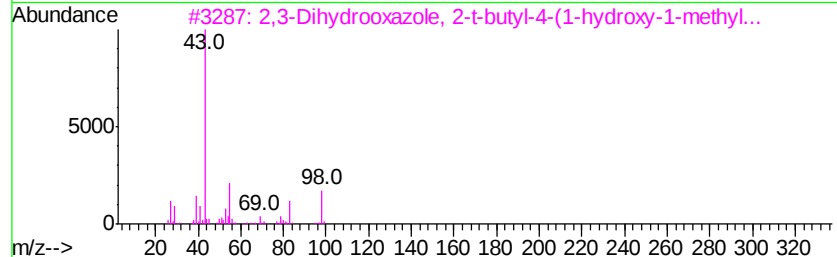
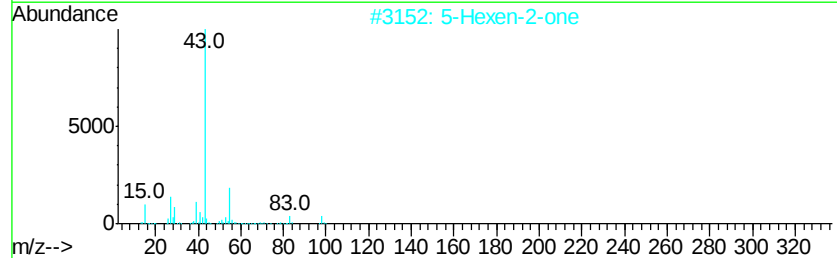
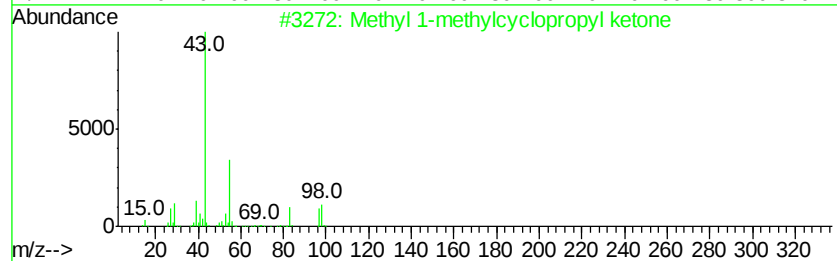
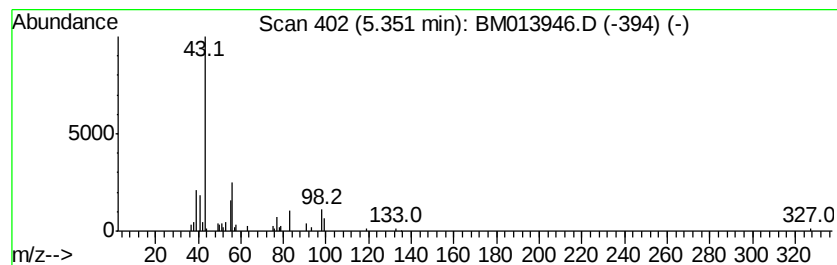
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 unknown-02 Concentration Rank 7

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 5.35 | 2.37 ng/ul | 147833 | 1,4-Dichlorobenzene-d4 | 7.57 |

| Hit# | of | Tentative ID                        | MW | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|----|---------|-------------|------|
| 1    | 5  | Methyl 1-methylcyclopropyl ketone   | 98 | C6H10O  | 001567-75-5 | 35   |
| 2    |    | 5-Hexen-2-one                       | 98 | C6H10O  | 000109-49-9 | 35   |
| 3    |    | 2,3-Dihydrooxazole, 2-t-butyl-4-... | 98 | C6H10O  | 036808-01-2 | 12   |
| 4    |    | 5-Hexen-2-one                       | 98 | C6H10O  | 000109-49-9 | 9    |
| 5    |    | 5-Hexen-2-one                       | 98 | C6H10O  | 000109-49-9 | 9    |



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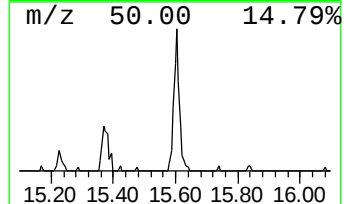
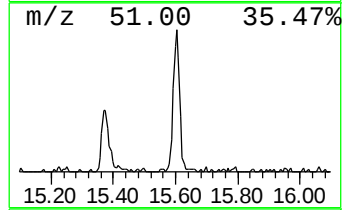
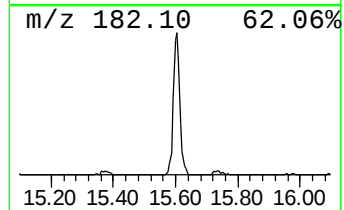
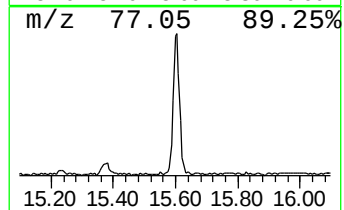
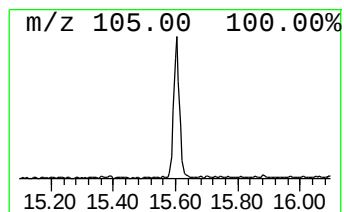
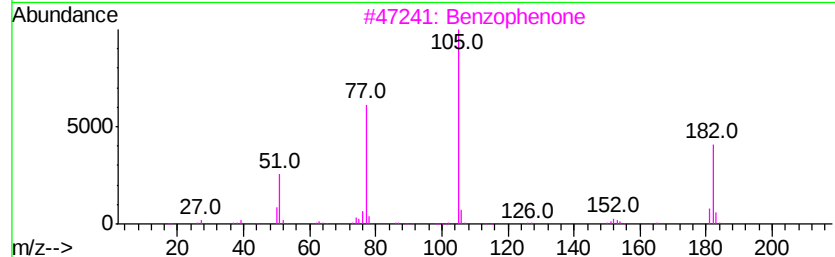
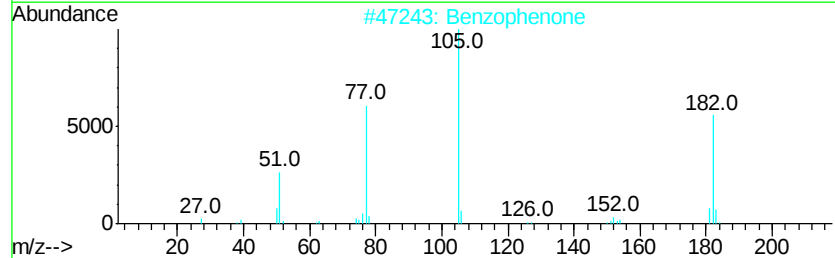
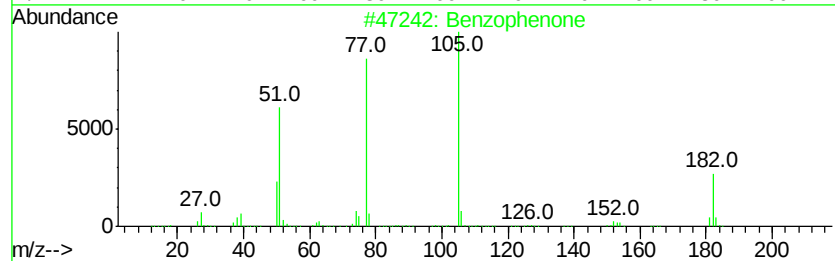
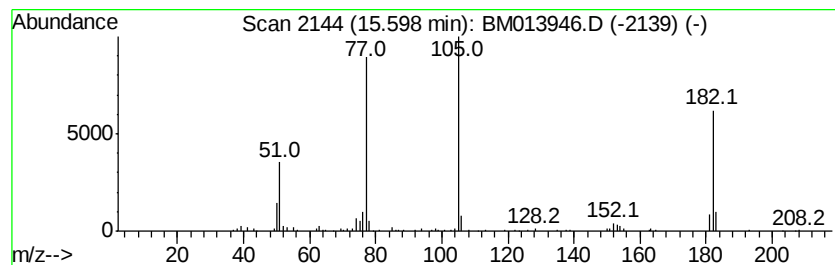
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Benzophenone Concentration Rank 4

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.60 | 3.06 ng/ul | 347792 | Phenanthrene-d10 | 16.98 |

| Hit# of | 5                                   | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|-------------------------------------|--------------|-----|---------|-------------|------|
| 1       | Benzophenone                        |              | 182 | C13H10O | 000119-61-9 | 95   |
| 2       | Benzophenone                        |              | 182 | C13H10O | 000119-61-9 | 95   |
| 3       | Benzophenone                        |              | 182 | C13H10O | 000119-61-9 | 94   |
| 4       | Benzophenone                        |              | 182 | C13H10O | 000119-61-9 | 90   |
| 5       | Benzenecarbothioic acid, S-methy... |              | 152 | C8H8OS  | 005925-68-8 | 58   |



Data Path : U:\HPCHEM1\BNA M\DATA\BM012818\  
Data File : BM013946.D  
Acq On : 29 Jan 2018 05:24  
Operator : SJ/JU  
Sample : J1244-04  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B45

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

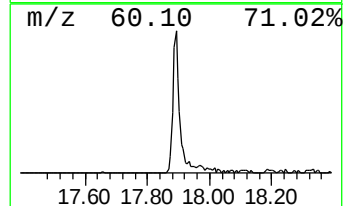
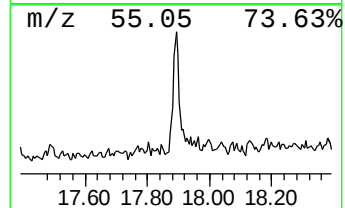
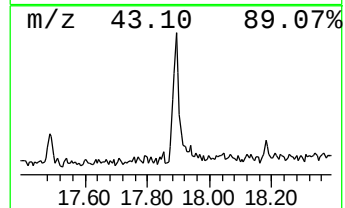
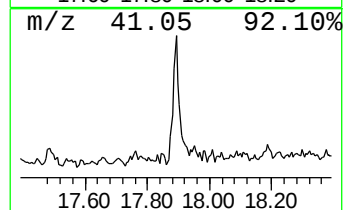
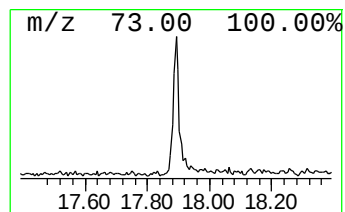
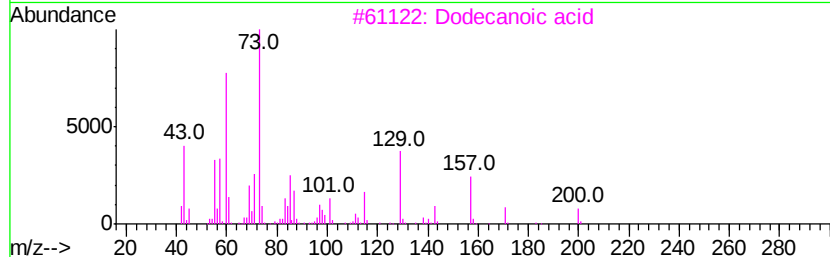
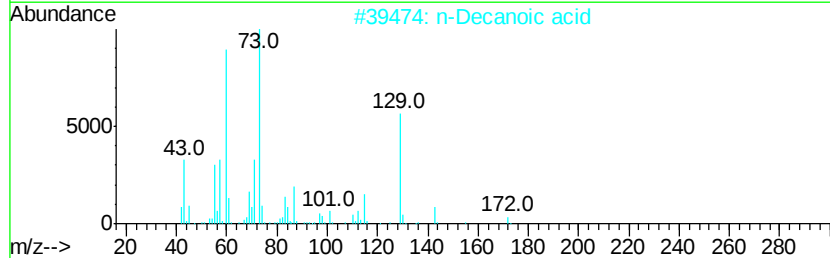
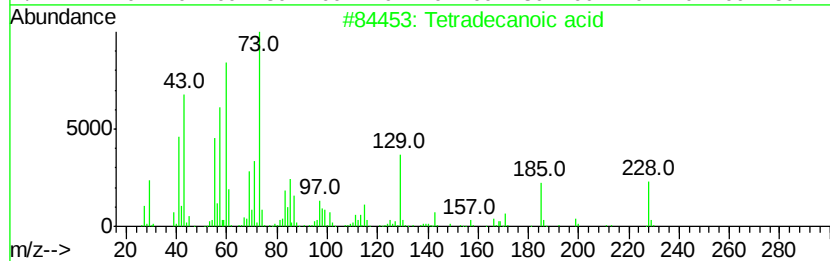
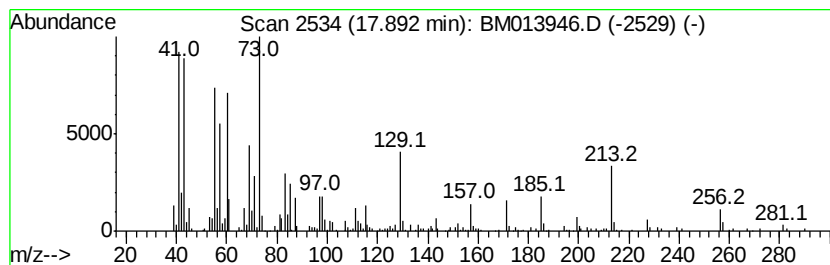
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Tetradecanoic acid Concentration Rank 2

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.89 | 5.11 ng/ul | 580116 | Phenanthrene-d10 | 16.98 |

| Hit# of | 5 | Tentative ID       | MW  | MolForm  | CAS#        | Qual |
|---------|---|--------------------|-----|----------|-------------|------|
| 1       |   | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 94   |
| 2       |   | n-Decanoic acid    | 172 | C10H20O2 | 000334-48-5 | 70   |
| 3       |   | Dodecanoic acid    | 200 | C12H24O2 | 000143-07-7 | 58   |
| 4       |   | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 58   |
| 5       |   | Tridecanoic acid   | 214 | C13H26O2 | 000638-53-9 | 58   |



Data Path : U:\HPCHEM1\BNA M\DATA\BM012818\  
Data File : BM013946.D  
Acq On : 29 Jan 2018 05:24  
Operator : SJ/JU  
Sample : J1244-04  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B45

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

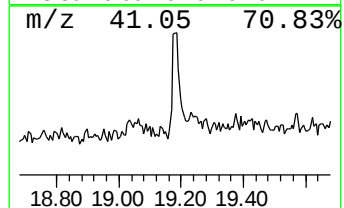
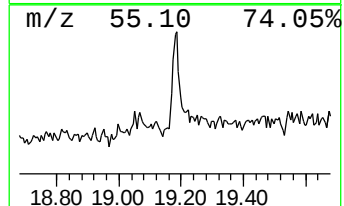
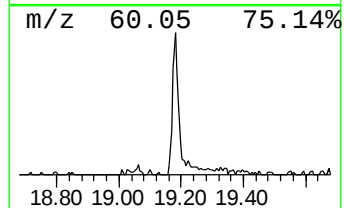
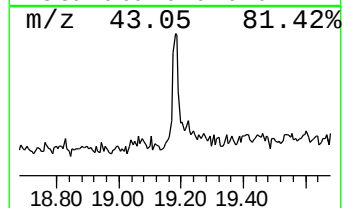
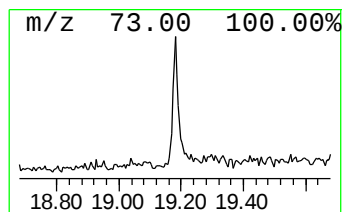
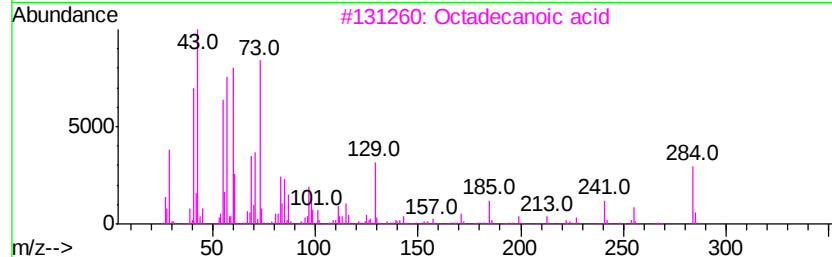
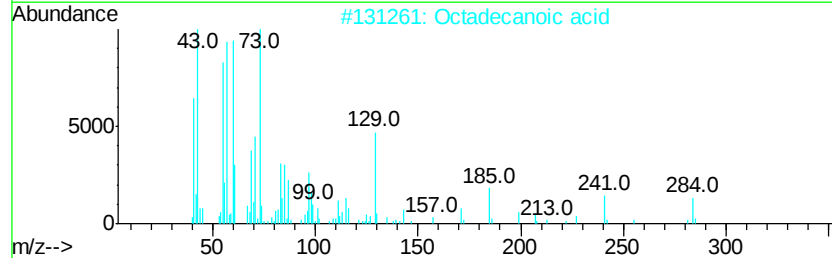
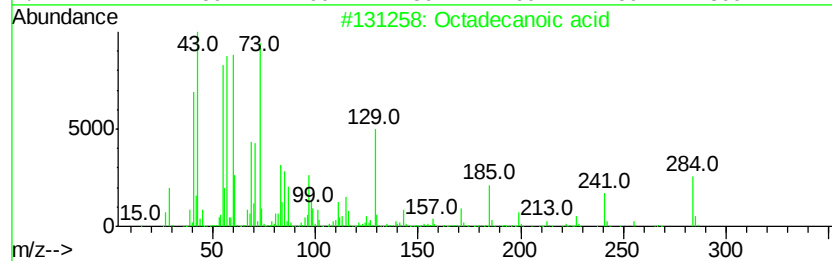
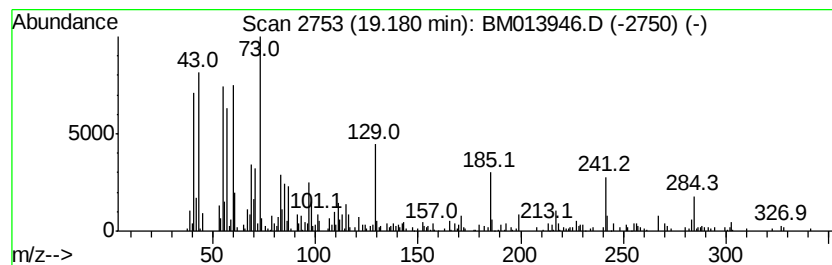
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Octadecanoic acid Concentration Rank 3

| R.T.  | EstConc    | Area   | Relative to ISTD |  | R.T.  |
|-------|------------|--------|------------------|--|-------|
| 19.18 | 4.10 ng/ul | 409568 | Chrysene-d12     |  | 21.19 |

| Hit# | of | 5 | Tentative ID       | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------|-----|----------|-------------|------|
| 1    |    |   | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 95   |
| 2    |    |   | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 94   |
| 3    |    |   | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 93   |
| 4    |    |   | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 93   |
| 5    |    |   | Heptadecanoic acid | 270 | C17H34O2 | 000506-12-7 | 91   |



Data Path : U:\HPCHEM1\BNA M\DATA\BM012818\  
Data File : BM013946.D  
Acq On : 29 Jan 2018 05:24  
Operator : SJ/JU  
Sample : J1244-04  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B45

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

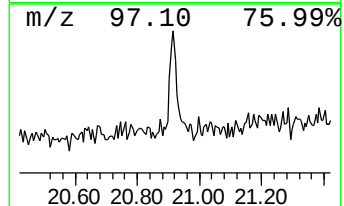
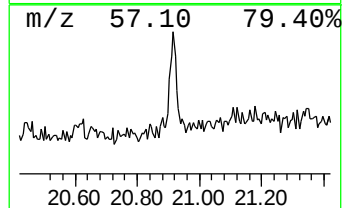
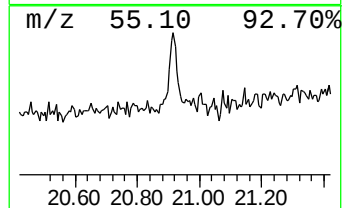
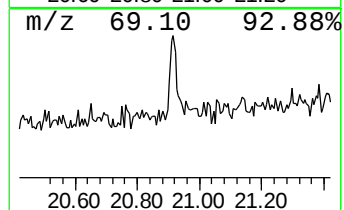
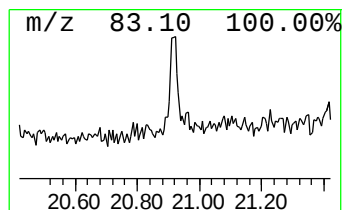
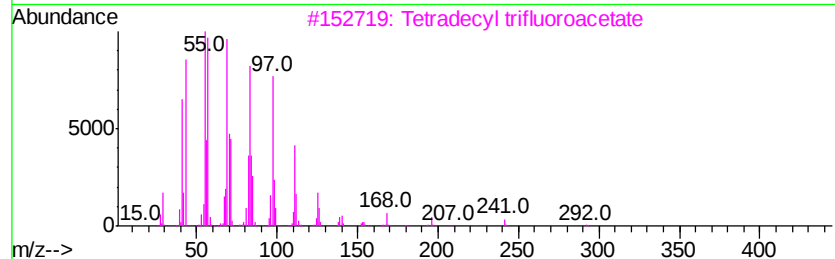
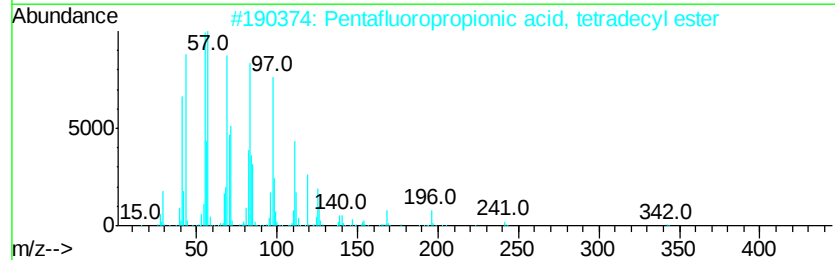
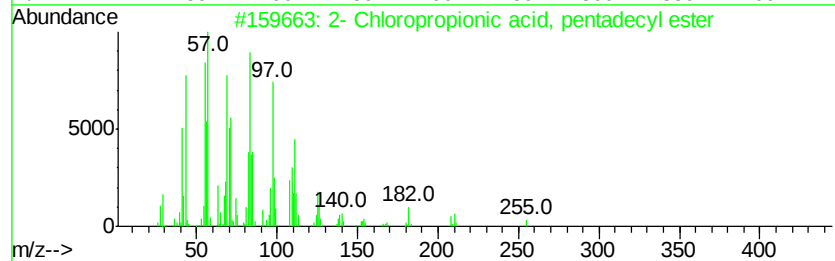
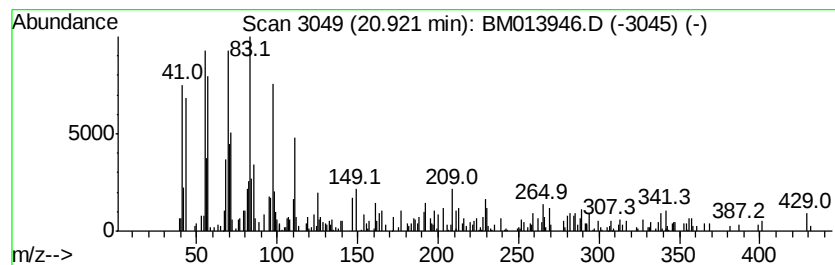
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 2- Chloropropionic acid, pe... Concentration Rank 5

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.92 | 2.78 ng/ul | 278168 | Chrysene-d12     | 21.19 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | 2- Chloropropionic acid, pentade... | 318 | C18H35ClO2 | 1000292-44-1 | 93   |
| 2    |    | Pentafluoropropionic acid, tetra... | 360 | C17H29F5O2 | 006222-06-6  | 93   |
| 3    |    | Tetradecyl trifluoroacetate         | 310 | C16H29F3O2 | 006222-02-2  | 93   |
| 4    |    | Trifluoroacetoxy hexadecane         | 338 | C18H33F3O2 | 006222-03-3  | 90   |
| 5    |    | Trifluoroacetic acid, pentadecyl... | 324 | C17H31F3O2 | 959010-23-2  | 90   |



Data Path : U:\HPCHEM1\BNA\_M\DATA\BM012818\  
Data File : BM013946.D  
Acq On : 29 Jan 2018 05:24  
Operator : SJ/JU  
Sample : J1244-04  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6B45

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 |
| (DEL) Alkane: Cyc... | 4.16  | 2.7     | ng/ul | 171283   | 1                     | 7.57  | 1249950 | 20.0 |
| unknown-01           | 4.73  | 7.7     | ng/ul | 479536   | 1                     | 7.57  | 1249950 | 20.0 |
| unknown-02           | 5.35  | 2.4     | ng/ul | 147833   | 1                     | 7.57  | 1249950 | 20.0 |
| Benzophenone         | 15.60 | 3.1     | ng/ul | 347792   | 4                     | 16.98 | 2269900 | 20.0 |
| Tetradecanoic acid   | 17.89 | 5.1     | ng/ul | 580116   | 4                     | 16.98 | 2269900 | 20.0 |
| Octadecanoic acid    | 19.18 | 4.1     | ng/ul | 409568   | 5                     | 21.19 | 1999680 | 20.0 |
| 2- Chloropropioni... | 20.92 | 2.8     | ng/ul | 278168   | 5                     | 21.19 | 1999680 | 20.0 |