

Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\  
 Data File : BM012482.D  
 Acq On : 27 Oct 2017 12:31  
 Operator : SJ/JU  
 Sample : I5873-07  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 B-18(5-6)\_101217

## 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-BM102417.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         | 3.146       | 6             | 10          | 21           | rVB      | 175995         | 239387        | 1.77%           | 0.133%        |
| 2         | 3.363       | 43            | 47          | 55           | rVB2     | 95406          | 137104        | 1.01%           | 0.076%        |
| 3         | 3.869       | 129           | 133         | 141          | rVB      | 314358         | 407649        | 3.01%           | 0.226%        |
| 4         | 5.005       | 322           | 326         | 332          | rBV      | 147219         | 195855        | 1.44%           | 0.109%        |
| 5         | 5.534       | 406           | 416         | 423          | rVB3     | 60116          | 139504        | 1.03%           | 0.077%        |
| 6         | 7.040       | 664           | 672         | 686          | rVB2     | 1487804        | 2846703       | 21.00%          | 1.578%        |
| 7         | 7.146       | 686           | 690         | 694          | rBV      | 136522         | 180243        | 1.33%           | 0.100%        |
| 8         | 7.204       | 694           | 700         | 709          | rVB      | 1280536        | 2007952       | 14.81%          | 1.113%        |
| 9         | 7.404       | 728           | 734         | 741          | rBV      | 1543256        | 2479831       | 18.30%          | 1.375%        |
| 10        | 7.798       | 795           | 801         | 805          | rBV      | 278451         | 417939        | 3.08%           | 0.232%        |
| 11        | 7.875       | 805           | 814         | 821          | rVB      | 1402430        | 2439983       | 18.00%          | 1.353%        |
| 12        | 8.057       | 841           | 845         | 850          | rBV      | 92275          | 139136        | 1.03%           | 0.077%        |
| 13        | 8.575       | 925           | 933         | 942          | rBV      | 1722147        | 2691475       | 19.86%          | 1.492%        |
| 14        | 9.028       | 1002          | 1010        | 1017         | rBV      | 1423179        | 2335430       | 17.23%          | 1.295%        |
| 15        | 9.457       | 1077          | 1083        | 1092         | rVB      | 551460         | 837248        | 6.18%           | 0.464%        |
| 16        | 9.745       | 1121          | 1132        | 1140         | rBV      | 1205002        | 1984038       | 14.64%          | 1.100%        |
| 17        | 10.287      | 1216          | 1224        | 1234         | rBV      | 1903888        | 3103385       | 22.90%          | 1.721%        |
| 18        | 10.663      | 1280          | 1288        | 1295         | rBV      | 1768638        | 2958255       | 21.83%          | 1.640%        |
| 19        | 10.798      | 1303          | 1311        | 1324         | rBV      | 1120361        | 2007939       | 14.81%          | 1.113%        |
| 20        | 11.951      | 1500          | 1507        | 1519         | rBV      | 909190         | 1341959       | 9.90%           | 0.744%        |
| 21        | 12.075      | 1519          | 1528        | 1536         | rBV      | 1997656        | 3191667       | 23.55%          | 1.769%        |
| 22        | 13.010      | 1681          | 1687        | 1693         | rBV      | 425334         | 667883        | 4.93%           | 0.370%        |
| 23        | 13.910      | 1830          | 1840        | 1845         | rBV      | 3155063        | 4261743       | 31.44%          | 2.363%        |
| 24        | 13.957      | 1845          | 1848        | 1852         | rVB      | 149076         | 203513        | 1.50%           | 0.113%        |
| 25        | 14.192      | 1877          | 1888        | 1900         | rBV      | 3550540        | 5707051       | 42.11%          | 3.164%        |
| 26        | 14.504      | 1932          | 1941        | 1947         | rBV2     | 2415065        | 3730304       | 27.52%          | 2.068%        |
| 27        | 14.563      | 1947          | 1951        | 1960         | rVB      | 103135         | 170821        | 1.26%           | 0.095%        |
| 28        | 14.698      | 1967          | 1974        | 1985         | rBV      | 771100         | 1225853       | 9.04%           | 0.680%        |
| 29        | 14.904      | 2004          | 2009        | 2017         | rVB      | 144212         | 260936        | 1.93%           | 0.145%        |
| 30        | 15.492      | 2102          | 2109        | 2115         | rBV      | 4519830        | 6313597       | 46.58%          | 3.500%        |
| 31        | 15.551      | 2115          | 2119        | 2124         | rVV      | 300898         | 428261        | 3.16%           | 0.237%        |
| 32        | 15.610      | 2124          | 2129        | 2136         | rVV      | 655612         | 925944        | 6.83%           | 0.513%        |
| 33        | 15.845      | 2164          | 2169        | 2179         | rVB2     | 140129         | 261836        | 1.93%           | 0.145%        |
| 34        | 15.992      | 2189          | 2194        | 2199         | rVB      | 108917         | 199882        | 1.47%           | 0.111%        |

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 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |          |         |        |
|----|--------|------|------|------|------|---------|----------|---------|--------|
| 35 | 16.221 | 2228 | 2233 | 2241 | rBV3 | 120220  | 244668   | 1.81%   | 0.136% |
| 36 | 16.551 | 2278 | 2289 | 2294 | rBV5 | 83163   | 231568   | 1.71%   | 0.128% |
| 37 | 16.780 | 2320 | 2328 | 2331 | rBV4 | 124121  | 228139   | 1.68%   | 0.126% |
| 38 | 16.898 | 2344 | 2348 | 2354 | rVB  | 185068  | 274496   | 2.03%   | 0.152% |
| 39 | 16.974 | 2357 | 2361 | 2363 | rBV  | 111974  | 163026   | 1.20%   | 0.090% |
| 40 | 17.063 | 2372 | 2376 | 2380 | rVB  | 182896  | 236292   | 1.74%   | 0.131% |
| 41 | 17.245 | 2401 | 2407 | 2410 | rBV  | 3022008 | 4306568  | 31.77%  | 2.388% |
| 42 | 17.286 | 2410 | 2414 | 2418 | rVV  | 4164466 | 5591570  | 41.25%  | 3.100% |
| 43 | 17.339 | 2418 | 2423 | 2427 | rVV2 | 4622407 | 6799239  | 50.16%  | 3.770% |
| 44 | 17.374 | 2427 | 2429 | 2434 | rVV  | 1109946 | 1320404  | 9.74%   | 0.732% |
| 45 | 17.427 | 2434 | 2438 | 2444 | rVB4 | 112929  | 204249   | 1.51%   | 0.113% |
| 46 | 17.639 | 2468 | 2474 | 2479 | rBV2 | 256533  | 454380   | 3.35%   | 0.252% |
| 47 | 18.139 | 2551 | 2559 | 2561 | rBV2 | 975863  | 2045153  | 15.09%  | 1.134% |
| 48 | 18.163 | 2561 | 2563 | 2569 | rVV  | 891437  | 1209896  | 8.93%   | 0.671% |
| 49 | 18.245 | 2573 | 2577 | 2582 | rVV  | 410933  | 567123   | 4.18%   | 0.314% |
| 50 | 18.310 | 2583 | 2588 | 2592 | rVV2 | 757146  | 1467228  | 10.82%  | 0.813% |
| 51 | 18.345 | 2592 | 2594 | 2600 | rVB  | 436891  | 566528   | 4.18%   | 0.314% |
| 52 | 18.615 | 2636 | 2640 | 2644 | rBV  | 622698  | 801997   | 5.92%   | 0.445% |
| 53 | 18.657 | 2644 | 2647 | 2652 | rVB2 | 267142  | 354811   | 2.62%   | 0.197% |
| 54 | 18.862 | 2679 | 2682 | 2686 | rVB2 | 181788  | 201679   | 1.49%   | 0.112% |
| 55 | 18.915 | 2688 | 2691 | 2694 | rVV  | 171017  | 221370   | 1.63%   | 0.123% |
| 56 | 18.951 | 2694 | 2697 | 2701 | rVB  | 218074  | 282911   | 2.09%   | 0.157% |
| 57 | 19.033 | 2706 | 2711 | 2715 | rBV  | 412535  | 752952   | 5.56%   | 0.417% |
| 58 | 19.174 | 2731 | 2735 | 2743 | rBV  | 587678  | 955499   | 7.05%   | 0.530% |
| 59 | 19.298 | 2750 | 2756 | 2762 | rVB  | 8966281 | 11975562 | 88.35%  | 6.639% |
| 60 | 19.409 | 2773 | 2775 | 2779 | rVB  | 557421  | 553512   | 4.08%   | 0.307% |
| 61 | 19.627 | 2807 | 2812 | 2815 | rBV2 | 6925806 | 10712403 | 79.03%  | 5.939% |
| 62 | 19.657 | 2815 | 2817 | 2825 | rVV  | 7846170 | 9867930  | 72.80%  | 5.471% |
| 63 | 19.727 | 2826 | 2829 | 2836 | rVB2 | 477321  | 809186   | 5.97%   | 0.449% |
| 64 | 19.833 | 2845 | 2847 | 2853 | rVB  | 564953  | 703294   | 5.19%   | 0.390% |
| 65 | 19.980 | 2867 | 2872 | 2879 | rBV2 | 652933  | 1764716  | 13.02%  | 0.978% |
| 66 | 20.451 | 2949 | 2952 | 2955 | rBV  | 438637  | 490908   | 3.62%   | 0.272% |
| 67 | 20.898 | 3024 | 3028 | 3033 | rBV  | 1268910 | 1725701  | 12.73%  | 0.957% |
| 68 | 21.127 | 3064 | 3067 | 3074 | rVB3 | 1307080 | 2433894  | 17.96%  | 1.349% |
| 69 | 21.192 | 3075 | 3078 | 3085 | rVB  | 1234177 | 1571666  | 11.60%  | 0.871% |
| 70 | 21.403 | 3109 | 3114 | 3119 | rBV3 | 6879700 | 13554192 | 100.00% | 7.515% |
| 71 | 21.451 | 3119 | 3122 | 3131 | rVB  | 4778173 | 6096063  | 44.98%  | 3.380% |

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Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 72 | 23.027 | 3385 | 3390 | 3395 | rBV  | 4261195 | 9570294 | 70.61% | 5.306% |
| 73 | 23.521 | 3470 | 3474 | 3478 | rBV  | 1942038 | 3236785 | 23.88% | 1.794% |
| 74 | 23.574 | 3479 | 3483 | 3487 | rVV2 | 4045210 | 7267800 | 53.62% | 4.029% |
| 75 | 23.621 | 3487 | 3491 | 3499 | rVB  | 4309836 | 7706108 | 56.85% | 4.272% |
| 76 | 23.721 | 3503 | 3508 | 3513 | rBV  | 2606993 | 4411538 | 32.55% | 2.446% |

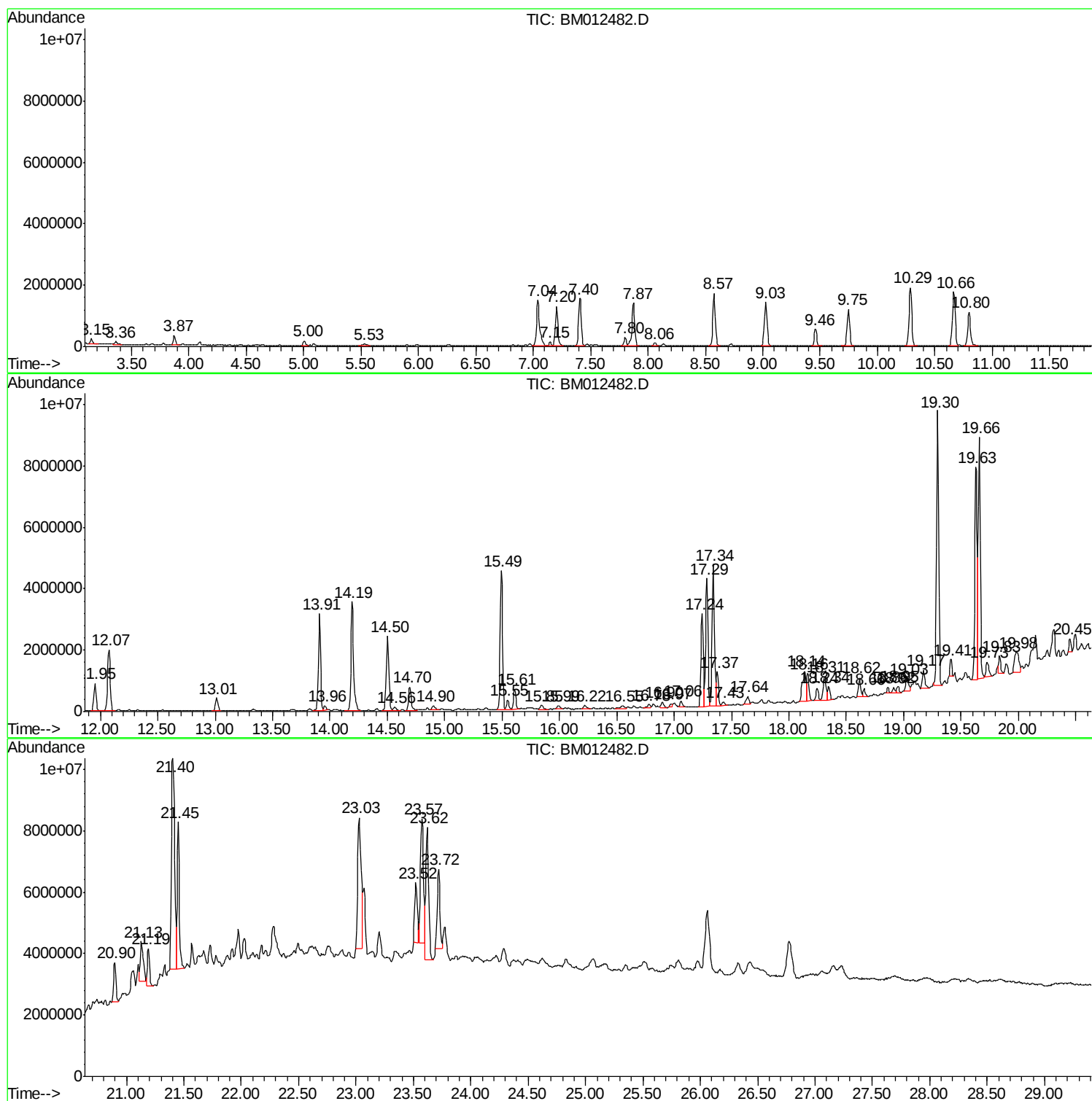
Sum of corrected areas: 180373604

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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.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-BM102417.M  
Quant Title : SVOA CALIBRATION

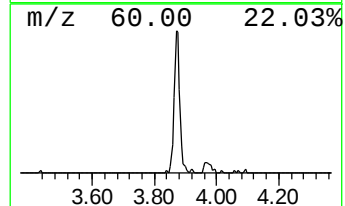
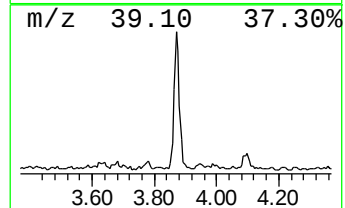
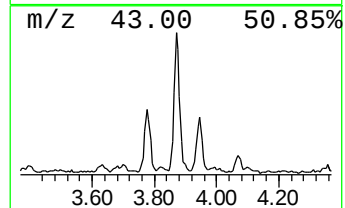
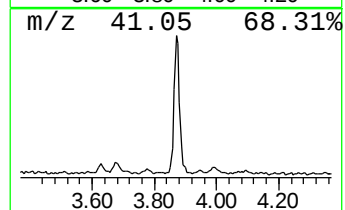
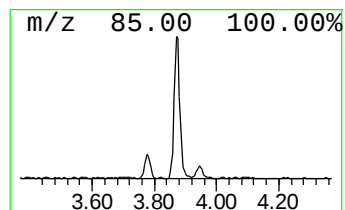
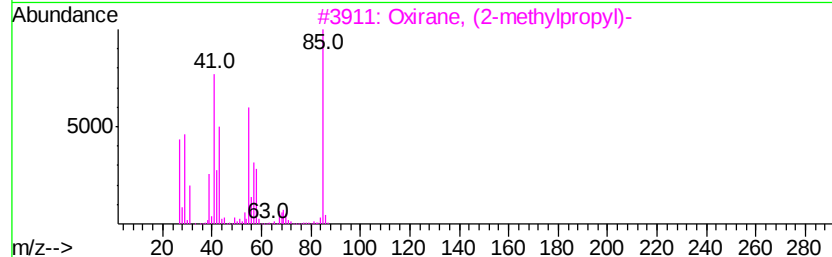
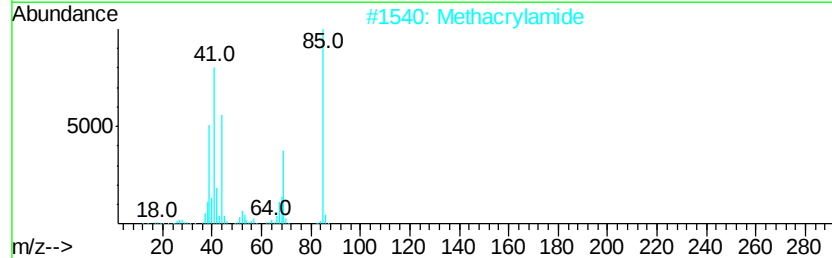
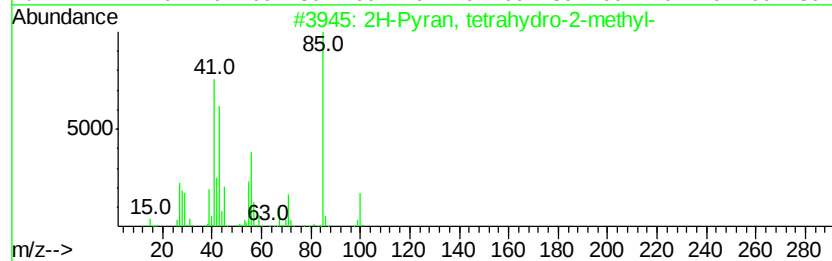
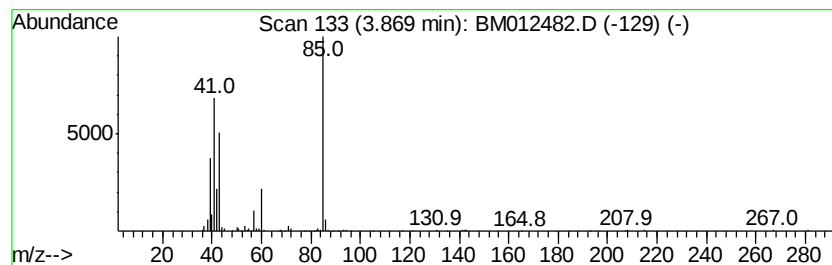
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 unknown-01 Concentration Rank 14

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 3.87 | 3.34 ng/ul | 407649 | 1,4-Dichlorobenzene-d4 | 7.87 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2H-Pyran, tetrahydro-2-methyl-       | 100 | C6H12O   | 010141-72-7 | 38   |
| 2    |    |   | Methacrylamide                       | 85  | C4H7NO   | 000079-39-0 | 38   |
| 3    |    |   | Oxirane, (2-methylpropyl)-           | 100 | C6H12O   | 023850-78-4 | 36   |
| 4    |    |   | Hexane, 1,1'-[methylenebis(oxym)]... | 216 | C13H28O2 | 054815-12-2 | 33   |
| 5    |    |   | Thiazole                             | 85  | C3H3NS   | 000288-47-1 | 28   |



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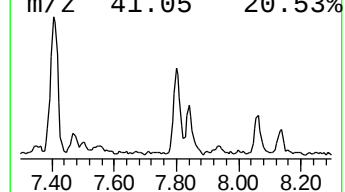
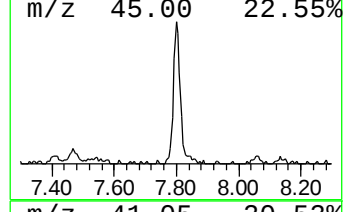
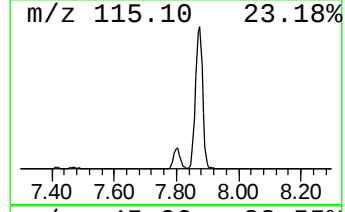
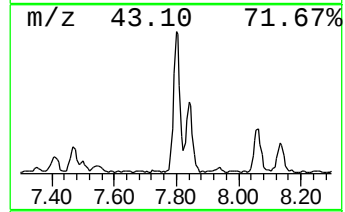
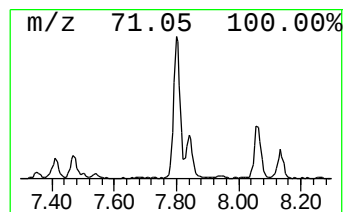
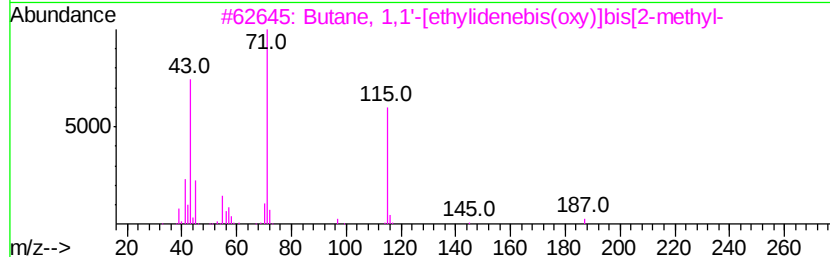
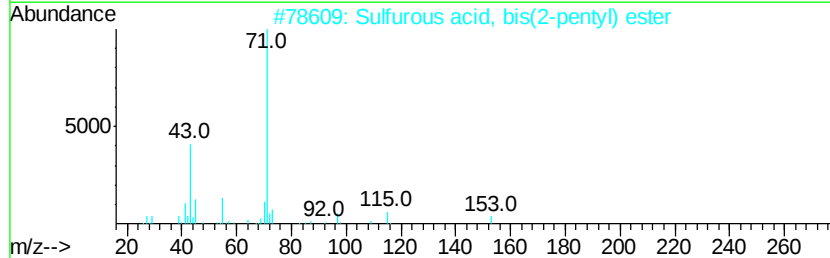
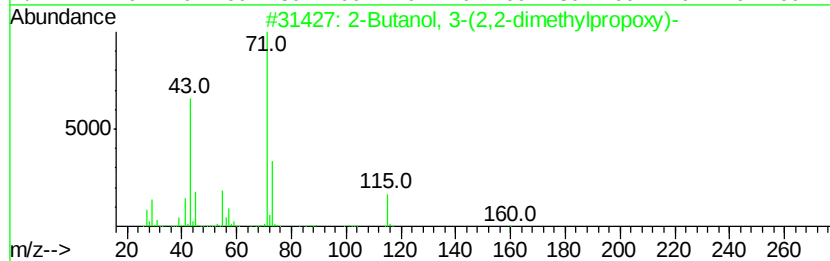
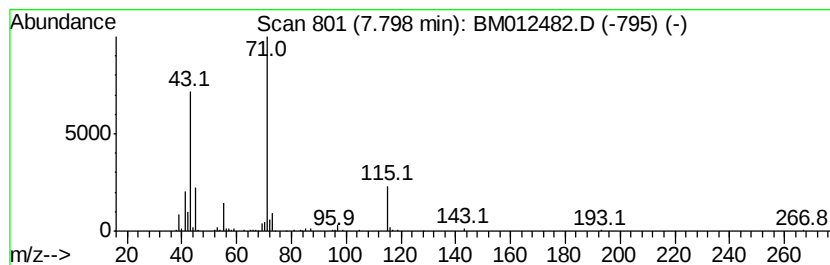
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 2-Butanol, 3-(2,2-dimethylp... Concentration Rank 13

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.80 | 3.43 ng/ul | 417939 | 1,4-Dichlorobenzene-d4 | 7.87 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm   | CAS#         | Qual |
|------|----|---|--------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 2-Butanol, 3-(2,2-dimethylpropoxy)-  | 160 | C9H20O2   | 074793-66-1  | 64   |
| 2    |    |   | Sulfurous acid, bis(2-pentyl) ester  | 222 | C10H22O3S | 1000309-15-3 | 59   |
| 3    |    |   | Butane, 1,1'-[ethylidenebis(oxo)]... | 202 | C12H26O2  | 013535-43-8  | 56   |
| 4    |    |   | Sulfurous acid, isobutyl 2-pentyl... | 208 | C9H20O3S  | 1000309-15-2 | 50   |
| 5    |    |   | Sulfurous acid, 2-pentyl pentyl ...  | 222 | C10H22O3S | 1000309-15-4 | 50   |



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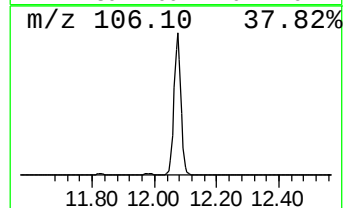
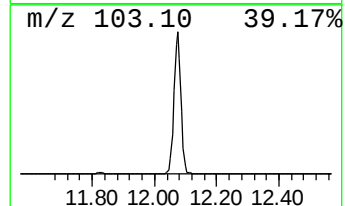
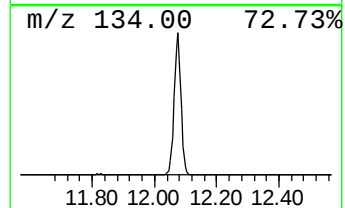
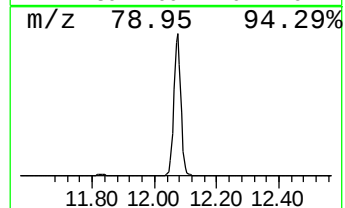
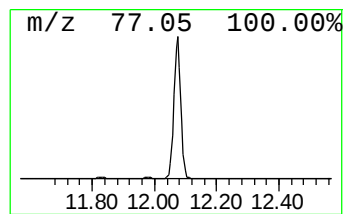
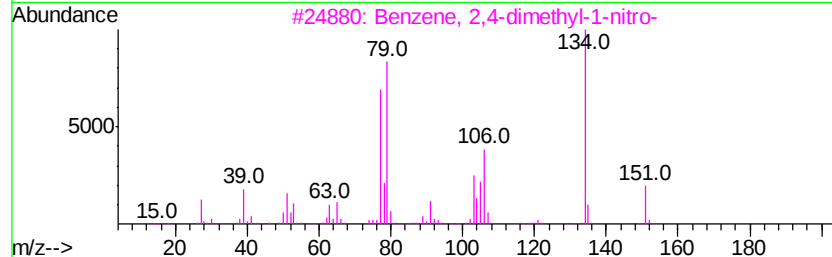
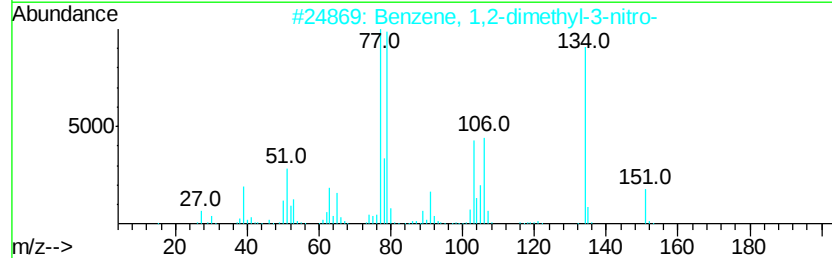
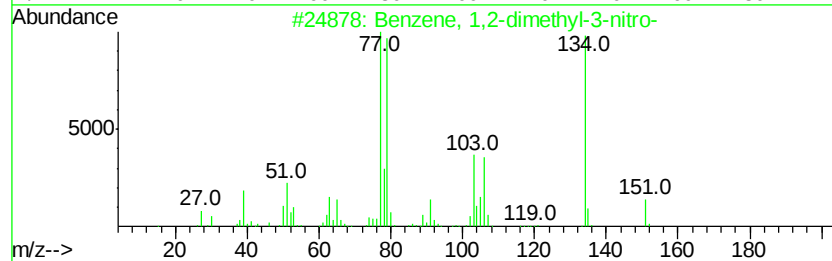
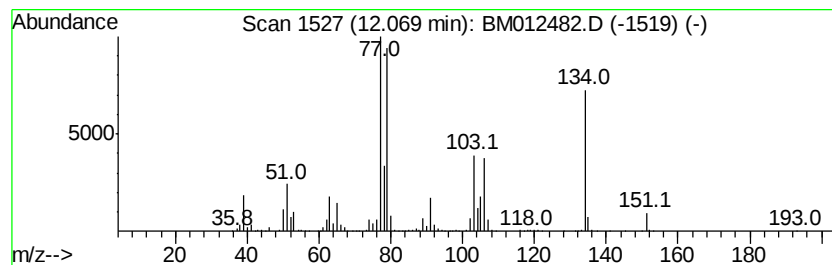
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Benzene, 1,2-dimethyl-3-nitro- Concentration Rank 1

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 12.07 | 21.58 ng/ul | 3191670 | Napthalene-d8    | 10.66 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2-dimethyl-3-nitro- | 151 | C8H9NO2 | 000083-41-0 | 97   |
| 2    |    | Benzene, 1,2-dimethyl-3-nitro- | 151 | C8H9NO2 | 000083-41-0 | 97   |
| 3    |    | Benzene, 2,4-dimethyl-1-nitro- | 151 | C8H9NO2 | 000089-87-2 | 95   |
| 4    |    | Benzene, 1,4-dimethyl-2-nitro- | 151 | C8H9NO2 | 000089-58-7 | 91   |
| 5    |    | Benzene, 1,3-dimethyl-2-nitro- | 151 | C8H9NO2 | 000081-20-9 | 91   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\  
Data File : BM012482.D  
Acq On : 27 Oct 2017 12:31  
Operator : SJ/JU  
Sample : I5873-07  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-18(5-6)\_101217

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M  
Quant Title : SVOA CALIBRATION

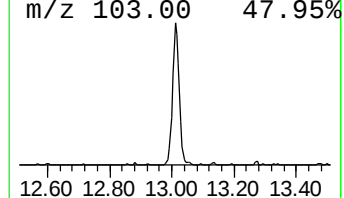
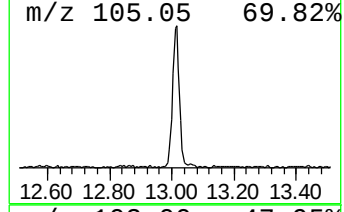
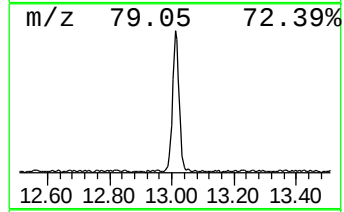
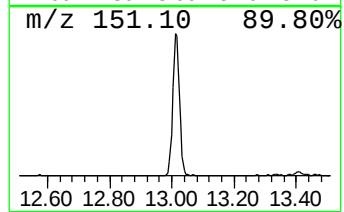
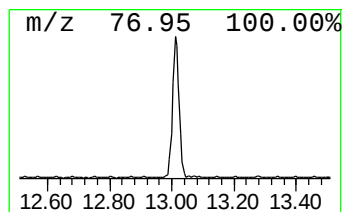
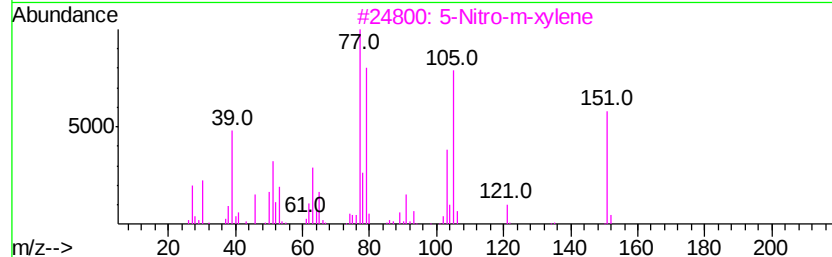
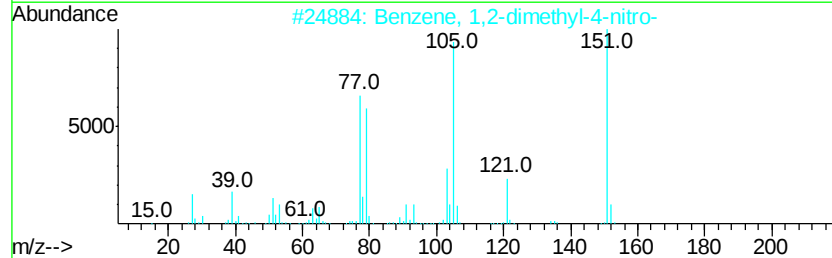
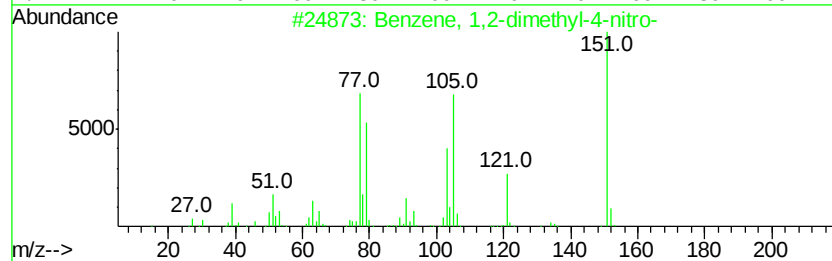
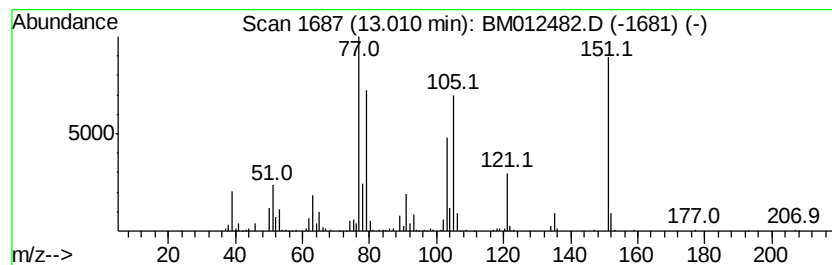
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Benzene, 1,2-dimethyl-4-nitro- Concentration Rank 11

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.01 | 3.58 ng/ul | 667883 | Acenaphthene-d10 | 14.50 |

| Hit# of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|---------|---|--------------------------------|-----|---------|-------------|------|
| 1       |   | Benzene, 1,2-dimethyl-4-nitro- | 151 | C8H9NO2 | 000099-51-4 | 96   |
| 2       |   | Benzene, 1,2-dimethyl-4-nitro- | 151 | C8H9NO2 | 000099-51-4 | 87   |
| 3       |   | 5-Nitro-m-xylene               | 151 | C8H9NO2 | 000099-12-7 | 87   |
| 4       |   | 5-Nitro-m-xylene               | 151 | C8H9NO2 | 000099-12-7 | 70   |
| 5       |   | 5-Nitro-m-xylene               | 151 | C8H9NO2 | 000099-12-7 | 58   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\  
Data File : BM012482.D  
Acq On : 27 Oct 2017 12:31  
Operator : SJ/JU  
Sample : I5873-07  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-18(5-6)\_101217

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M  
Quant Title : SVOA CALIBRATION

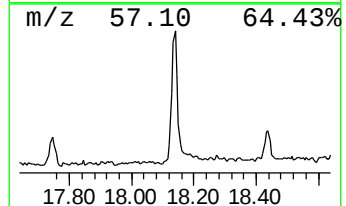
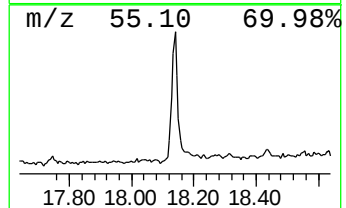
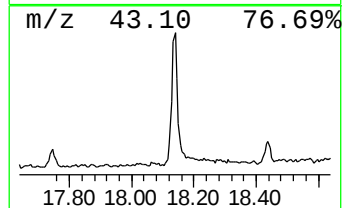
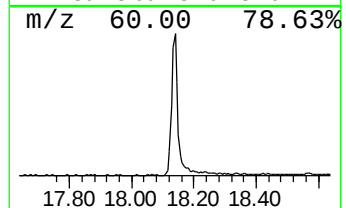
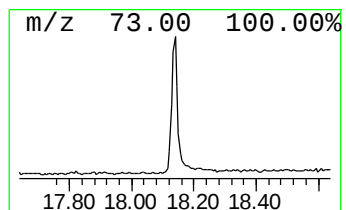
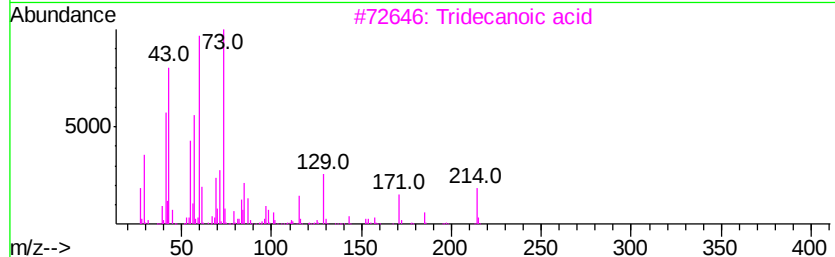
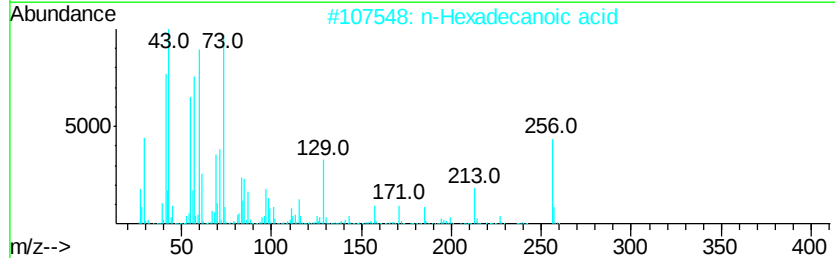
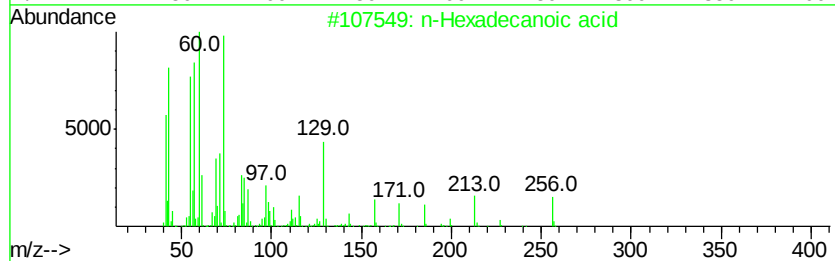
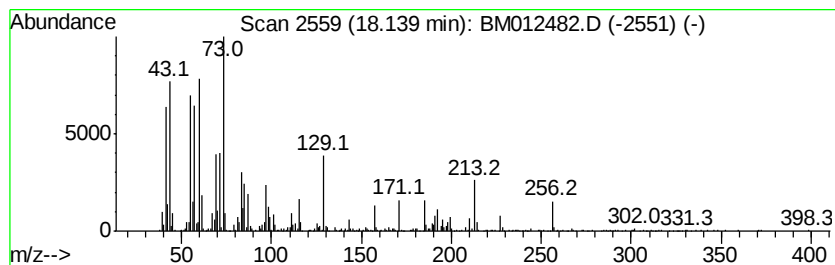
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 n-Hexadecanoic acid Concentration Rank 3

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 18.14 | 9.50 ng/ul | 2045150 | Phenanthrene-d10 | 17.24 |

| Hit# | of | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------|-----|----------|-------------|------|
| 1    | 5  | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 99   |
| 2    |    | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 93   |
| 3    |    | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 90   |
| 4    |    | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 90   |
| 5    |    | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 86   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\  
Data File : BM012482.D  
Acq On : 27 Oct 2017 12:31  
Operator : SJ/JU  
Sample : I5873-07  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-18(5-6)\_101217

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M  
Quant Title : SVOA CALIBRATION

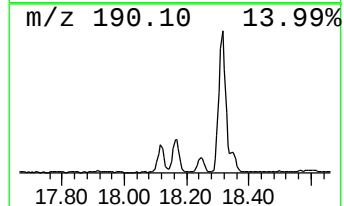
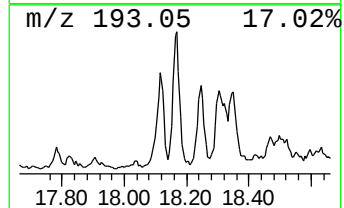
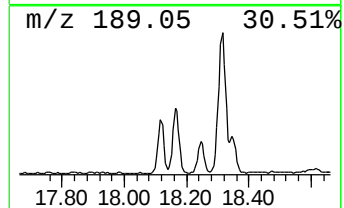
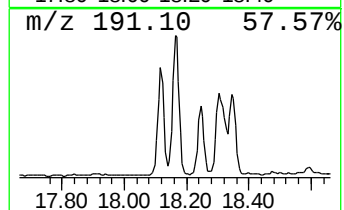
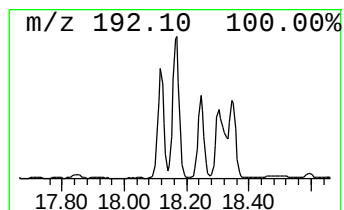
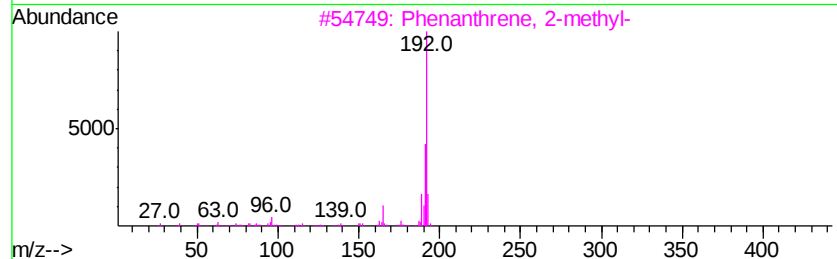
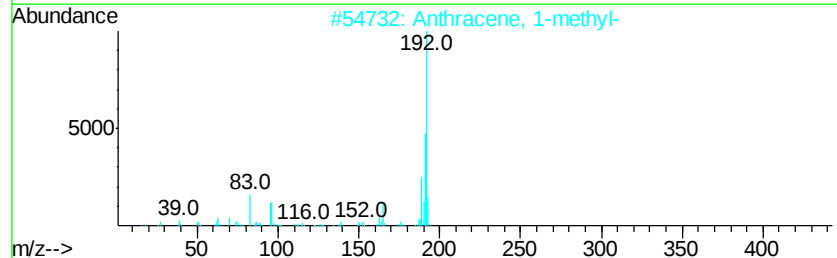
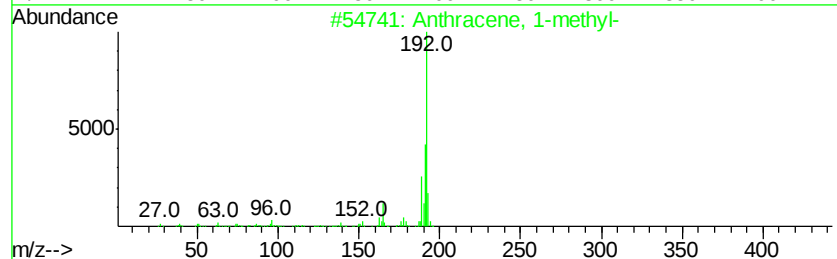
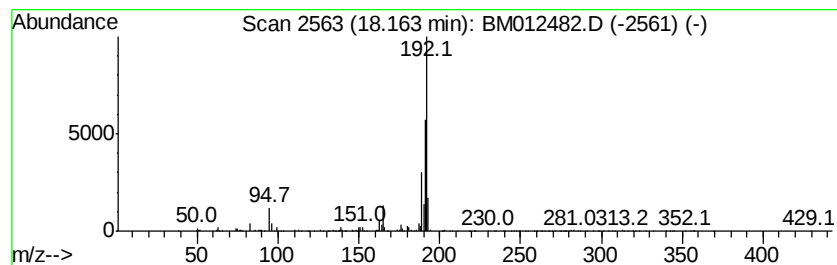
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Anthracene, 1-methyl- Concentration Rank 7

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 18.16 | 5.62 ng/ul | 1209900 | Phenanthrene-d10 | 17.24 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Anthracene, 1-methyl-   | 192 | C15H12  | 000610-48-0 | 95   |
| 2    |    |   | Anthracene, 1-methyl-   | 192 | C15H12  | 000610-48-0 | 93   |
| 3    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 93   |
| 4    |    |   | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 93   |
| 5    |    |   | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 93   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\  
Data File : BM012482.D  
Acq On : 27 Oct 2017 12:31  
Operator : SJ/JU  
Sample : I5873-07  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-18(5-6)\_101217

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M  
Quant Title : SVOA CALIBRATION

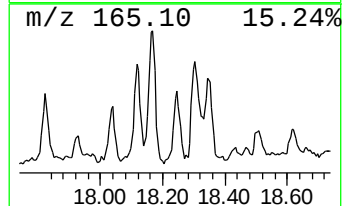
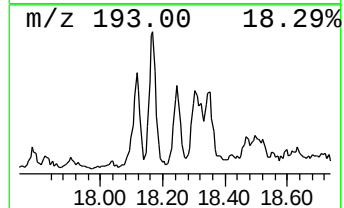
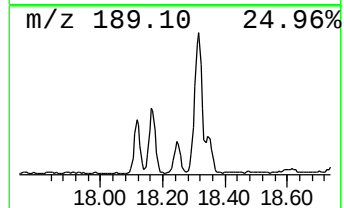
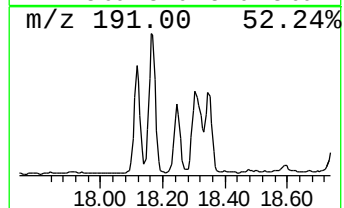
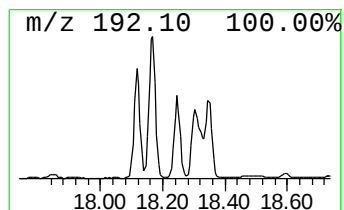
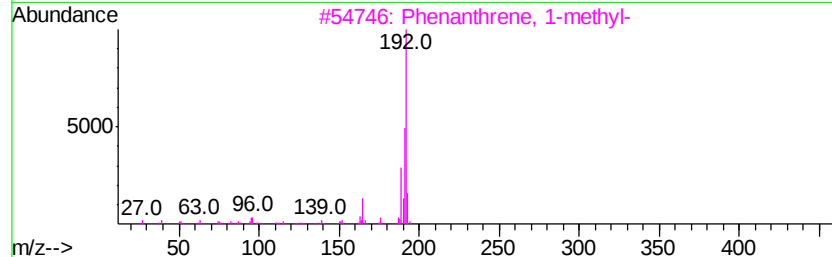
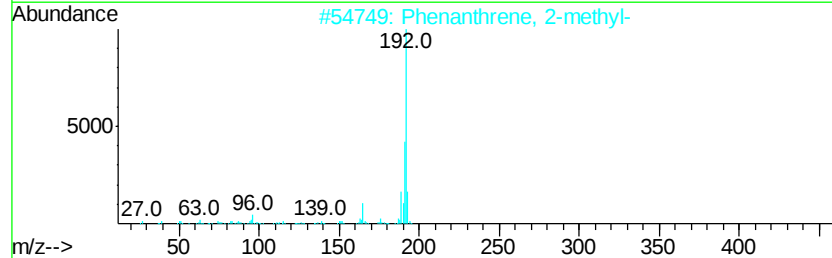
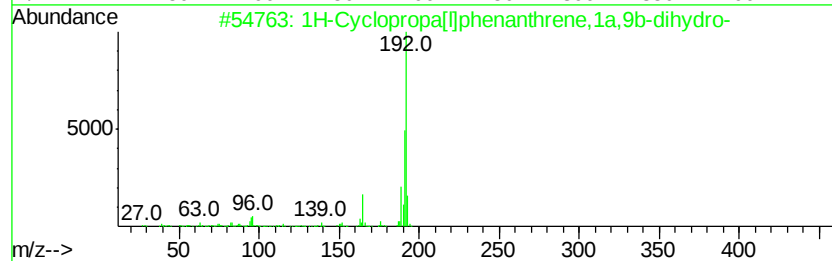
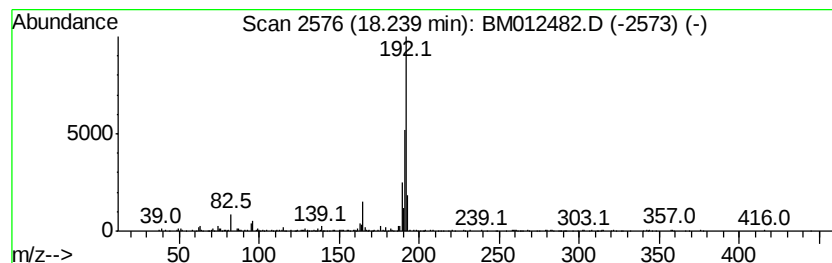
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 1H-Cyclopropa[l]phenanthren... Concentration Rank 15

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.24 | 2.63 ng/ul | 567123 | Phenanthrene-d10 | 17.24 |

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



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\  
Data File : BM012482.D  
Acq On : 27 Oct 2017 12:31  
Operator : SJ/JU  
Sample : I5873-07  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

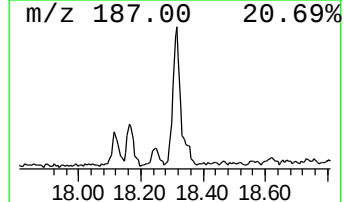
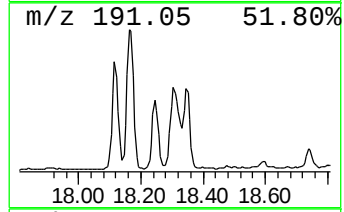
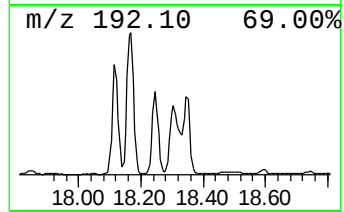
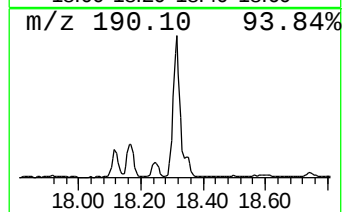
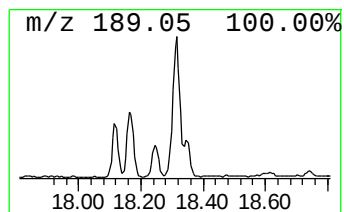
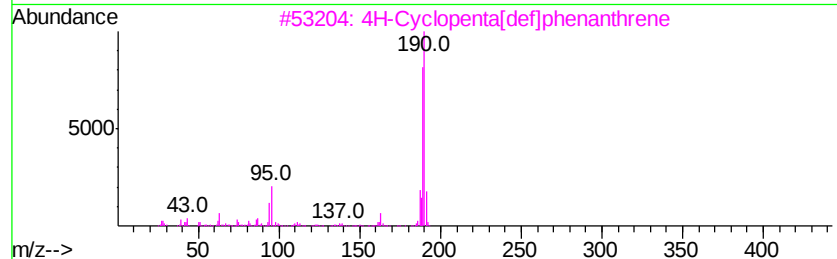
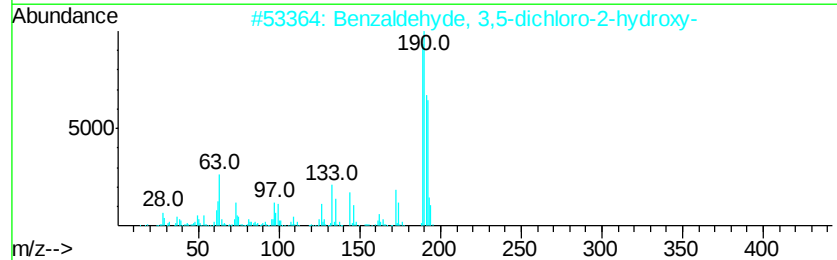
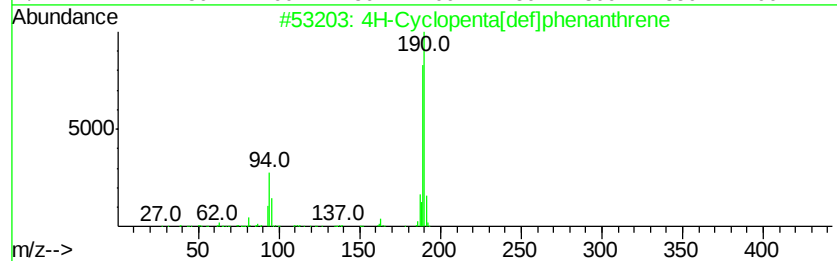
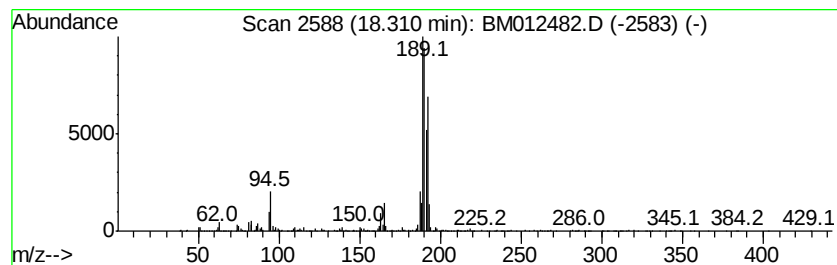
Instrument :  
BNA\_M  
ClientSampleId :  
B-18(5-6)\_101217

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 11 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 18.31     | 6.81 ng/ul                          | 1467230 | Phenanthrene-d10 | 17.24       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | 4H-Cyclopenta[def]phenanthrene      | 190     | C15H10           | 000203-64-5 | 70   |
| 2         | Benzaldehyde, 3,5-dichloro-2-hyd... | 190     | C7H4Cl2O2        | 000090-60-8 | 64   |
| 3         | 4H-Cyclopenta[def]phenanthrene      | 190     | C15H10           | 000203-64-5 | 60   |
| 4         | 6H-Cyclobuta[flk]phenanthrene       | 190     | C15H10           | 083469-43-6 | 58   |
| 5         | Methyl diselenide                   | 190     | C2H6Se2          | 007101-31-7 | 41   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\  
Data File : BM012482.D  
Acq On : 27 Oct 2017 12:31  
Operator : SJ/JU  
Sample : I5873-07  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-18(5-6)\_101217

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M  
Quant Title : SVOA CALIBRATION

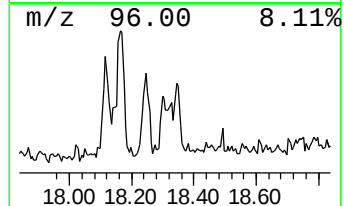
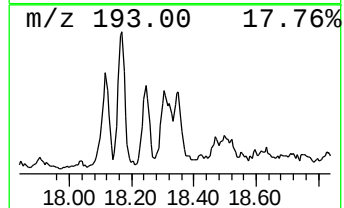
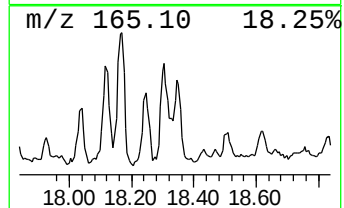
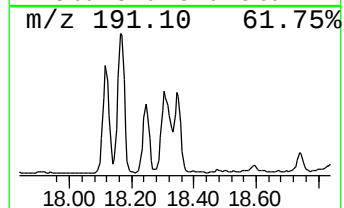
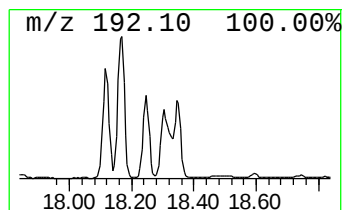
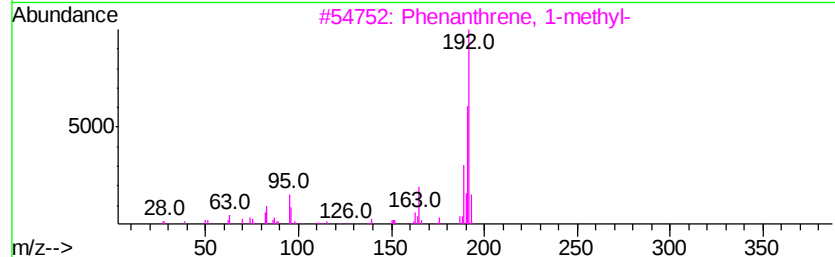
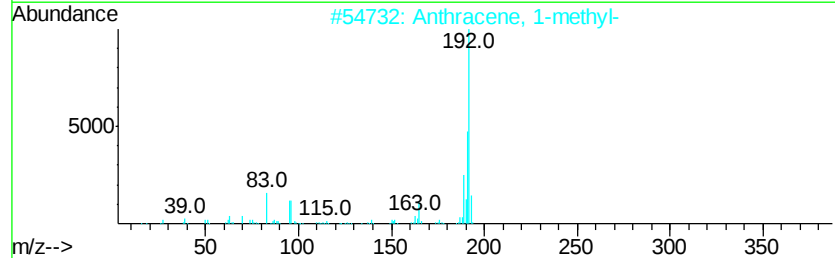
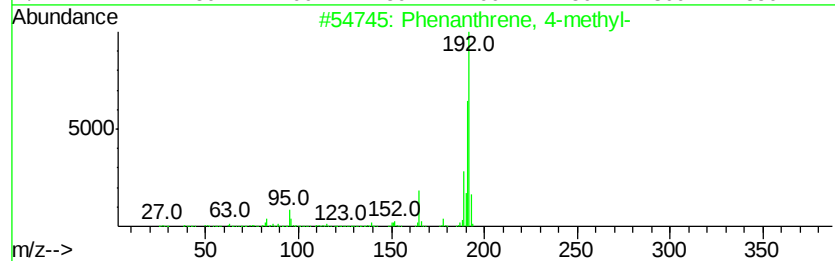
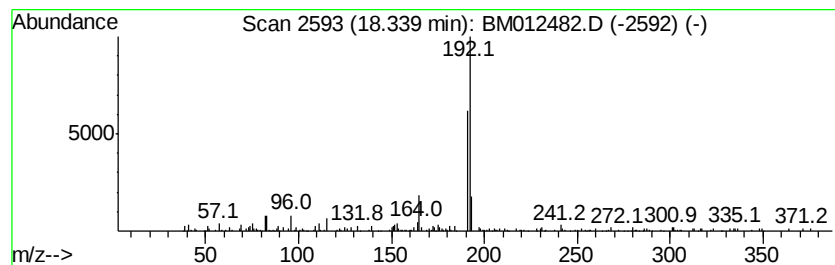
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Phenanthrene, 4-methyl- Concentration Rank 16

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.34 | 2.63 ng/ul | 566528 | Phenanthrene-d10 | 17.24 |

| Hit# | of | 5 | Tentative ID                            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 4-methyl-                 | 192 | C15H12  | 000832-64-4 | 87   |
| 2    |    |   | Anthracene, 1-methyl-                   | 192 | C15H12  | 000610-48-0 | 87   |
| 3    |    |   | Phenanthrene, 1-methyl-                 | 192 | C15H12  | 000832-69-9 | 83   |
| 4    |    |   | 1H-Cyclopropa[1,1]phenanthrene, 1a, ... | 192 | C15H12  | 000949-41-7 | 83   |
| 5    |    |   | Anthracene, 1-methyl-                   | 192 | C15H12  | 000610-48-0 | 83   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\  
Data File : BM012482.D  
Acq On : 27 Oct 2017 12:31  
Operator : SJ/JU  
Sample : I5873-07  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-18(5-6)\_101217

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M  
Quant Title : SVOA CALIBRATION

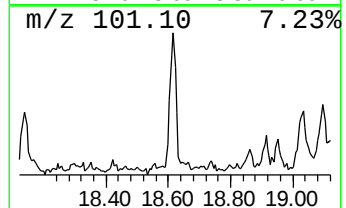
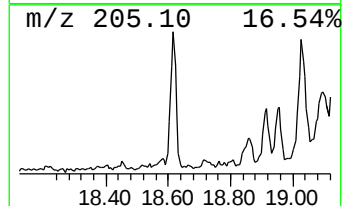
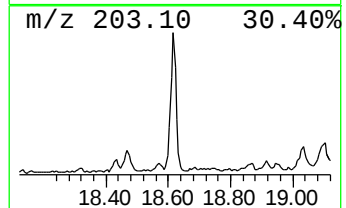
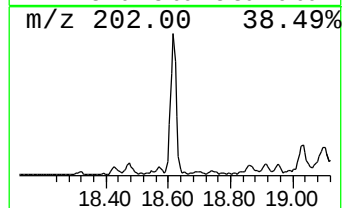
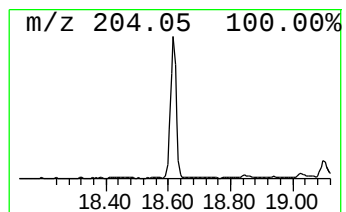
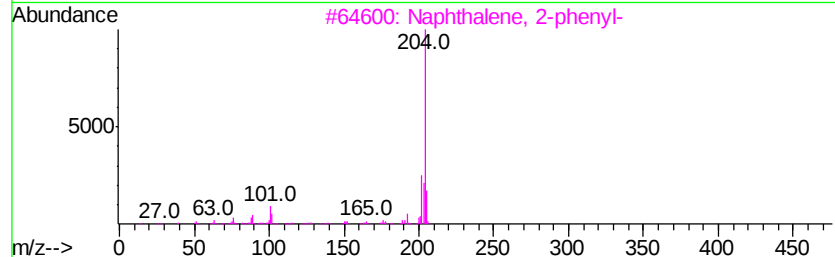
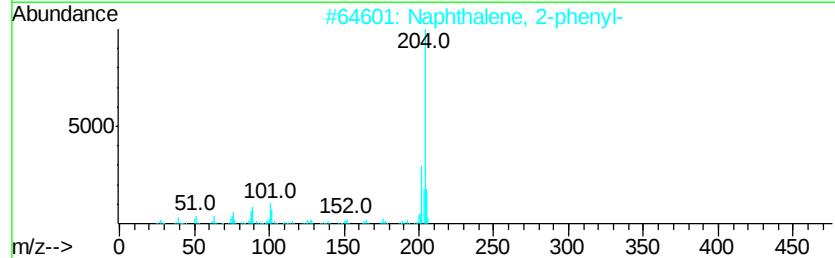
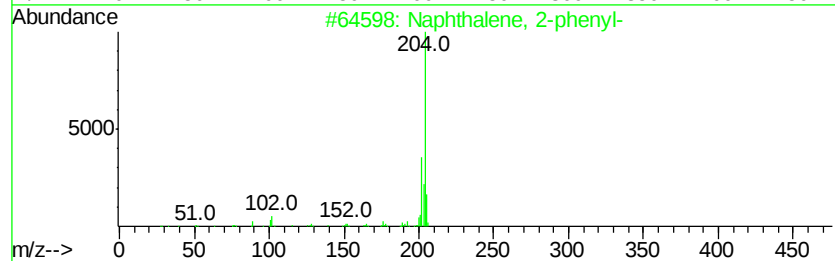
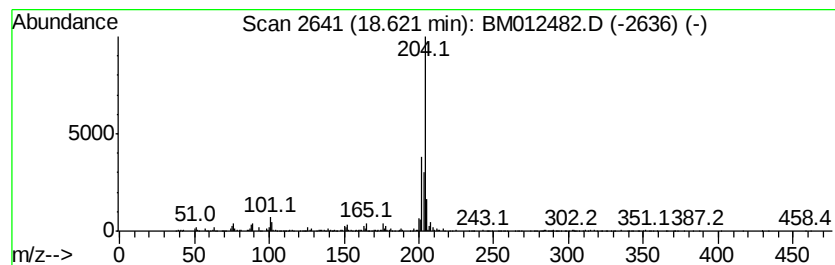
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TIC Integration Parameters: LSCINT.P

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Peak Number 13 Naphthalene, 2-phenyl- Concentration Rank 9

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.62 | 3.72 ng/ul | 801997 | Phenanthrene-d10 | 17.24 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2-phenyl-              | 204 | C16H12  | 000612-94-2 | 70   |
| 2    |    |   | Naphthalene, 2-phenyl-              | 204 | C16H12  | 000612-94-2 | 64   |
| 3    |    |   | Naphthalene, 2-phenyl-              | 204 | C16H12  | 000612-94-2 | 62   |
| 4    |    |   | Tricyclo[8.2.2.2(4,7)]hexadeca-2... | 204 | C16H12  | 006572-60-7 | 62   |
| 5    |    |   | Naphthalene, 2-phenyl-              | 204 | C16H12  | 000612-94-2 | 58   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\  
Data File : BM012482.D  
Acq On : 27 Oct 2017 12:31  
Operator : SJ/JU  
Sample : I5873-07  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-18(5-6)\_101217

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M  
Quant Title : SVOA CALIBRATION

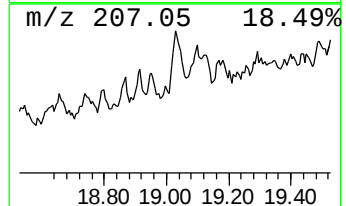
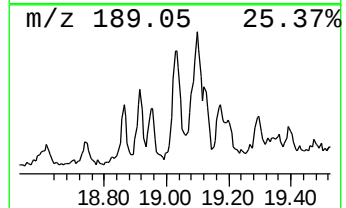
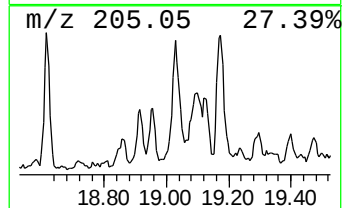
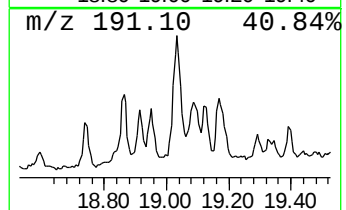
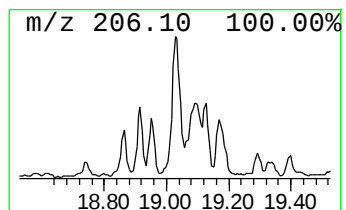
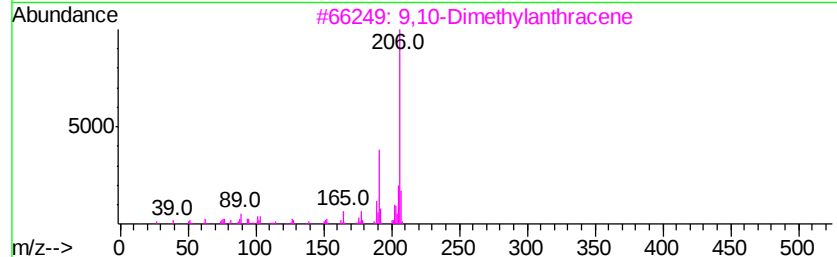
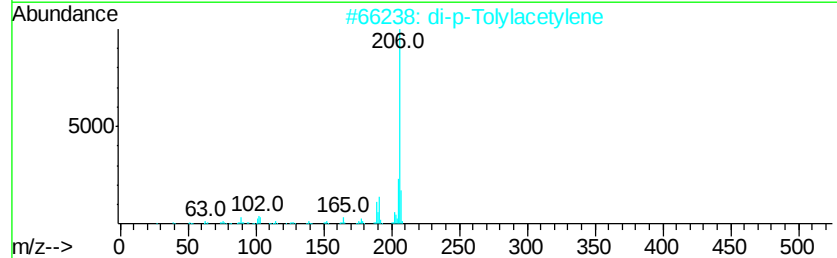
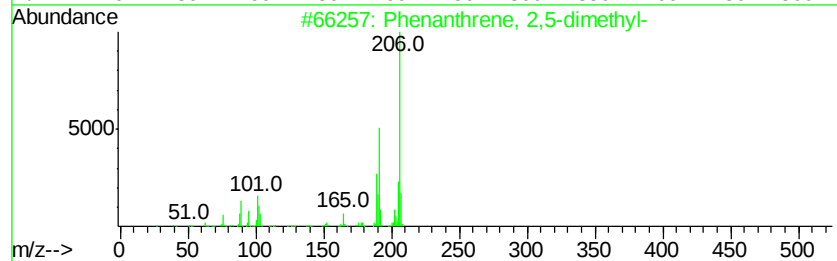
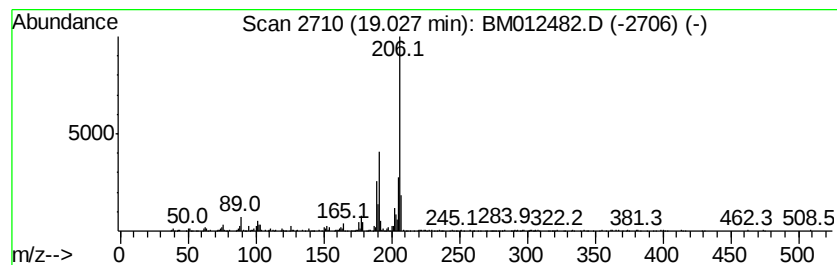
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TIC Integration Parameters: LSCINT.P

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Peak Number 14 Phenanthrene, 2,5-dimethyl- Concentration Rank 12

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.03 | 3.50 ng/ul | 752952 | Phenanthrene-d10 | 17.24 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 94   |
| 2    |    |   | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 93   |
| 3    |    |   | 9,10-Dimethylantracene      | 206 | C16H14  | 000781-43-1 | 91   |
| 4    |    |   | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 91   |
| 5    |    |   | 9,10-Dimethylantracene      | 206 | C16H14  | 000781-43-1 | 91   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\  
Data File : BM012482.D  
Acq On : 27 Oct 2017 12:31  
Operator : SJ/JU  
Sample : I5873-07  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-18(5-6)\_101217

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M  
Quant Title : SVOA CALIBRATION

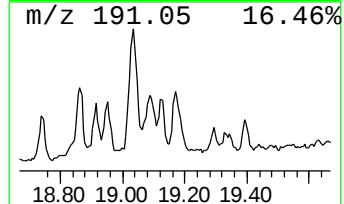
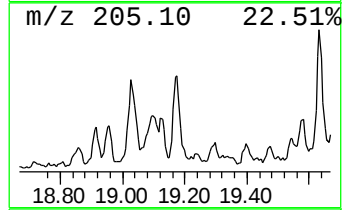
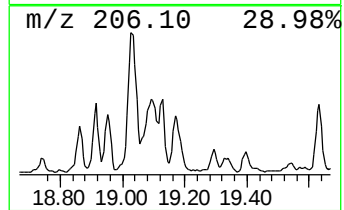
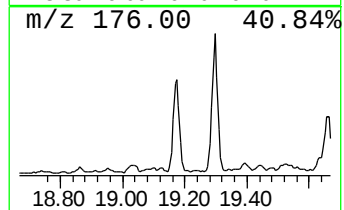
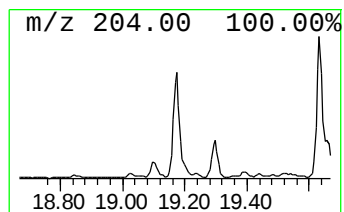
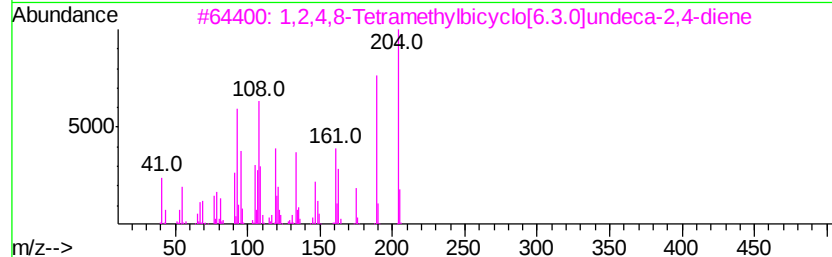
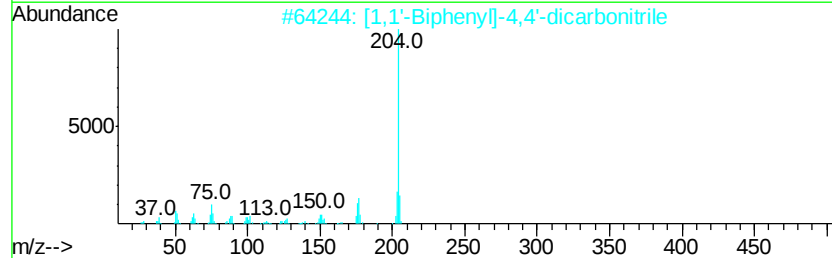
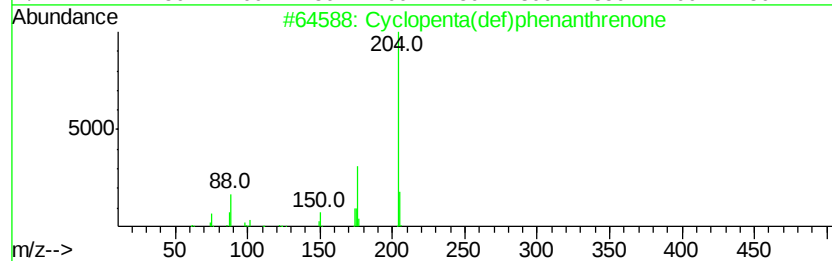
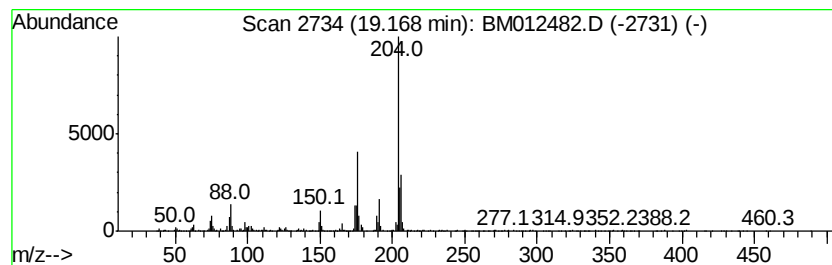
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 Cyclopenta(def)phenanthrenone Concentration Rank 8

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.17 | 4.44 ng/ul | 955499 | Phenanthrene-d10 | 17.24 |

| Hit# | of | Tentative ID                                      | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclopenta(def)phenanthrenone                     | 204 | C15H8O  | 005737-13-3 | 90   |
| 2    |    | [1,1'-Biphenyl]-4,4'-dicarbonitrile               | 204 | C14H8N2 | 001591-30-6 | 47   |
| 3    |    | 1,2,4,8-Tetramethylbicyclo[6.3.0]undeca-2,4-diene | 204 | C15H24  | 137235-51-9 | 45   |
| 4    |    | [1,1'-Biphenyl]-2,2'-dicarbonitrile               | 204 | C14H8N2 | 004341-02-0 | 43   |
| 5    |    | 9-Acridinecarbonitrile                            | 204 | C14H8N2 | 005326-19-2 | 43   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\  
Data File : BM012482.D  
Acq On : 27 Oct 2017 12:31  
Operator : SJ/JU  
Sample : I5873-07  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

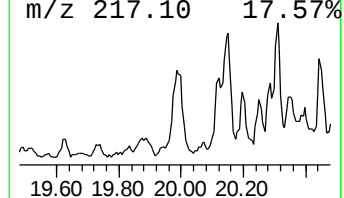
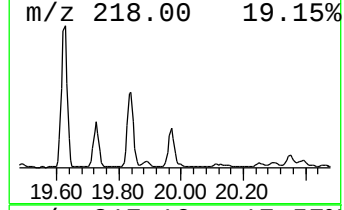
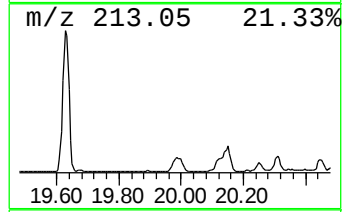
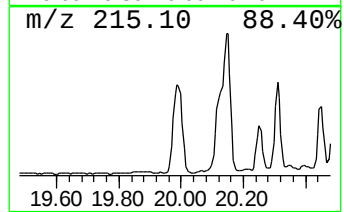
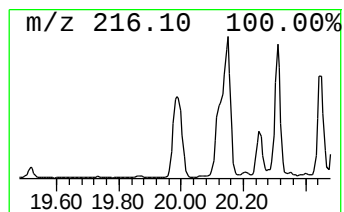
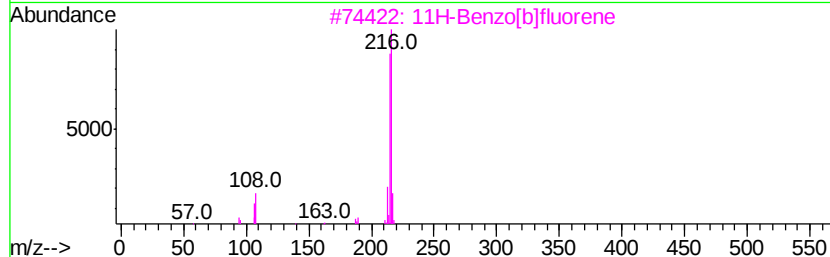
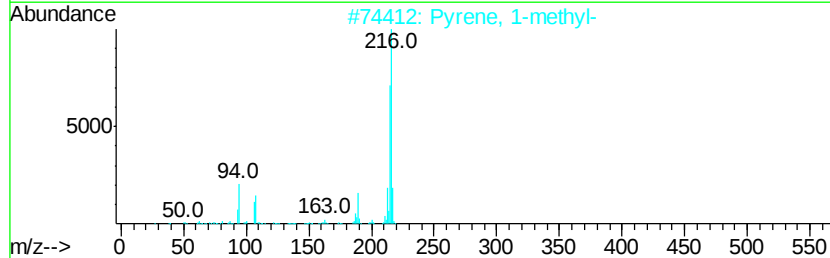
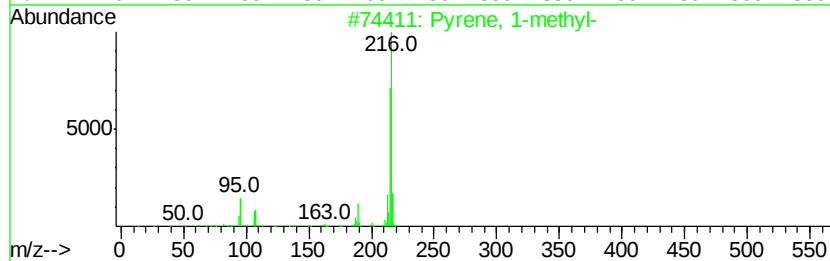
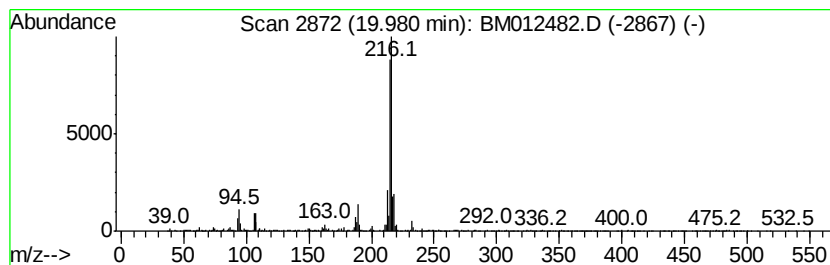
Instrument :  
BNA\_M  
ClientSampleId :  
B-18(5-6)\_101217

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 16 Pyrene, 1-methyl- Concentration Rank 17

| R.T.    | EstConc    | Area                 | Relative to ISTD | R.T.           |
|---------|------------|----------------------|------------------|----------------|
| 19.98   | 2.60 ng/ul | 1764720              | Chrysene-d12     | 21.42          |
| Hit# of | 5          | Tentative ID         | MW MolForm       | CAS# Qual      |
| 1       |            | Pyrene, 1-methyl-    | 216 C17H12       | 002381-21-7 93 |
| 2       |            | Pyrene, 1-methyl-    | 216 C17H12       | 002381-21-7 90 |
| 3       |            | 11H-Benzo[b]fluorene | 216 C17H12       | 000243-17-4 89 |
| 4       |            | 11H-Benzo[b]fluorene | 216 C17H12       | 000243-17-4 87 |
| 5       |            | 11H-Benzo[a]fluorene | 216 C17H12       | 000238-84-6 81 |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\  
Data File : BM012482.D  
Acq On : 27 Oct 2017 12:31  
Operator : SJ/JU  
Sample : I5873-07  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-18(5-6)\_101217

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M  
Quant Title : SVOA CALIBRATION

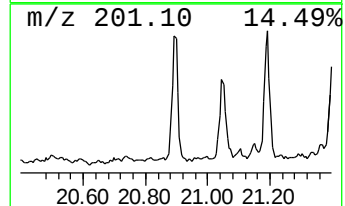
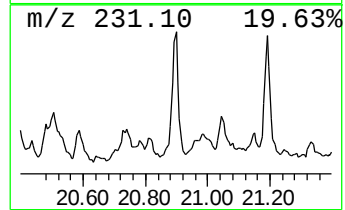
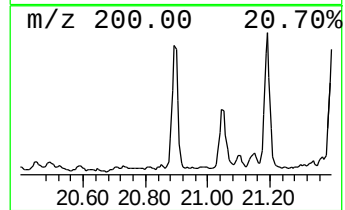
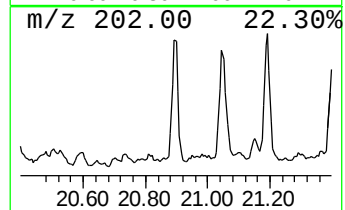
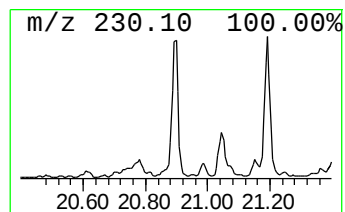
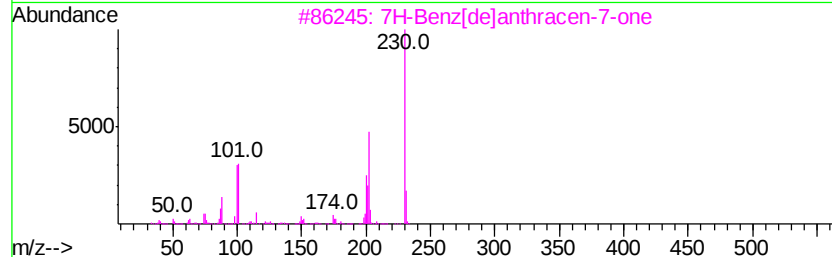
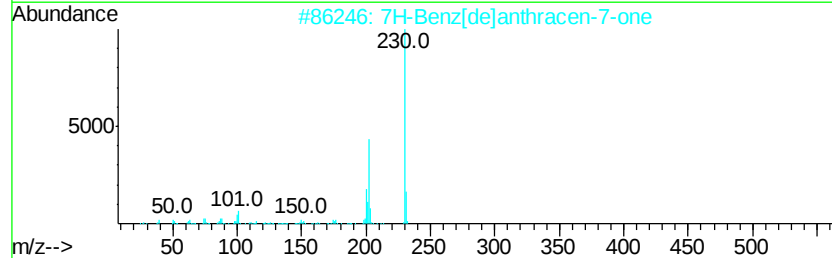
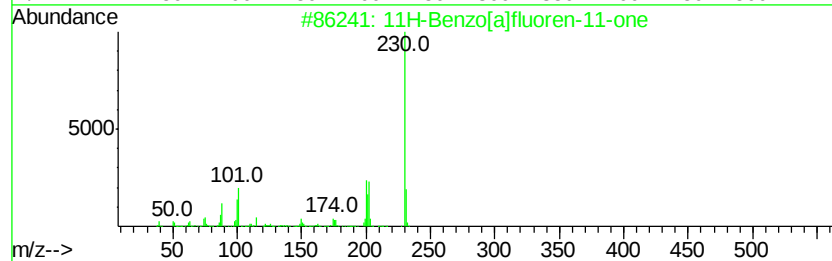
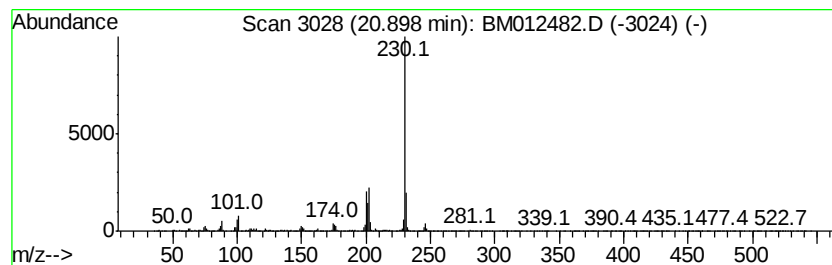
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 11H-Benzo[a]fluoren-11-one Concentration Rank 18

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 20.90 | 2.55 ng/ul | 1725700 | Chrysene-d12     | 21.42 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | 11H-Benzo[a]fluoren-11-one | 230 | C17H10O | 000479-79-8 | 96   |
| 2    |    | 7H-Benz[de]anthracen-7-one | 230 | C17H10O | 000082-05-3 | 90   |
| 3    |    | 7H-Benz[de]anthracen-7-one | 230 | C17H10O | 000082-05-3 | 87   |
| 4    |    | 7H-Benz[de]anthracen-7-one | 230 | C17H10O | 000082-05-3 | 81   |
| 5    |    | 7H-Benz[de]anthracen-7-one | 230 | C17H10O | 000082-05-3 | 81   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\  
Data File : BM012482.D  
Acq On : 27 Oct 2017 12:31  
Operator : SJ/JU  
Sample : I5873-07  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

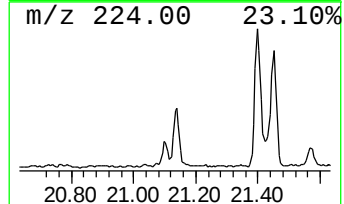
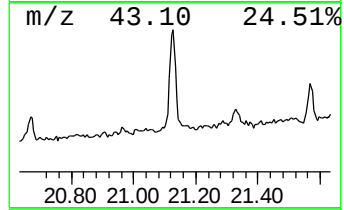
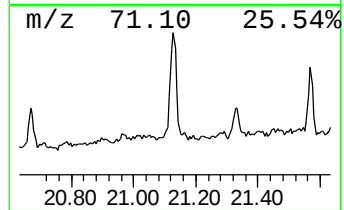
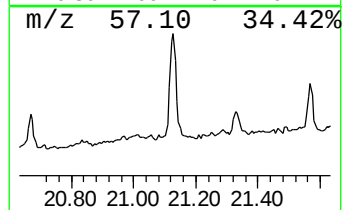
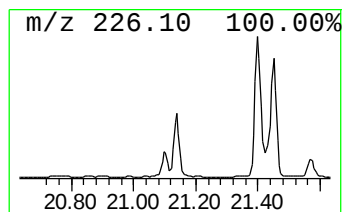
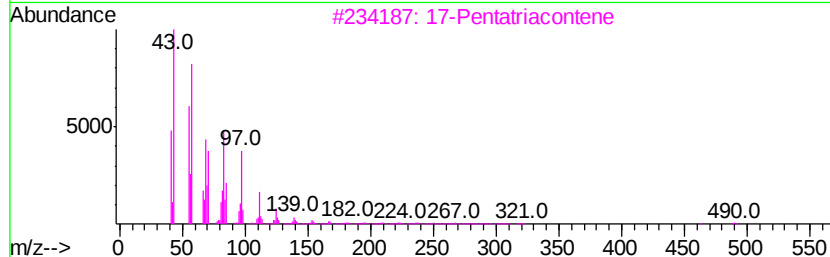
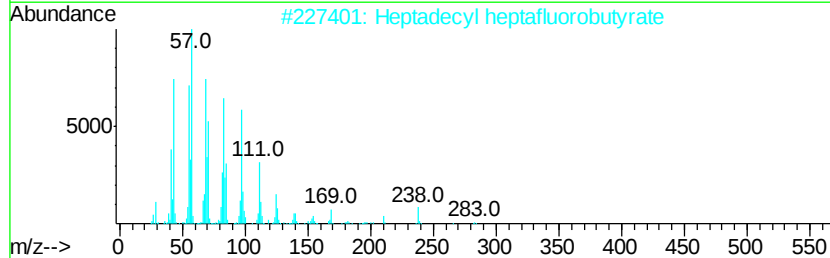
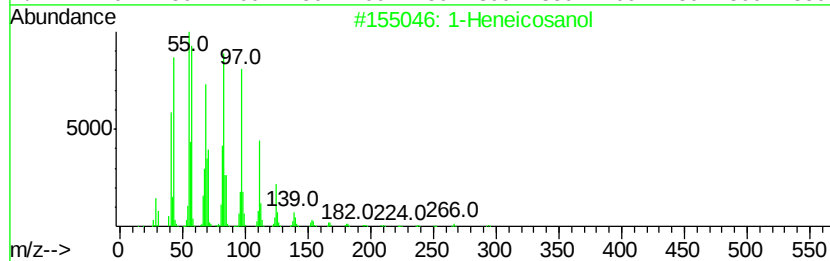
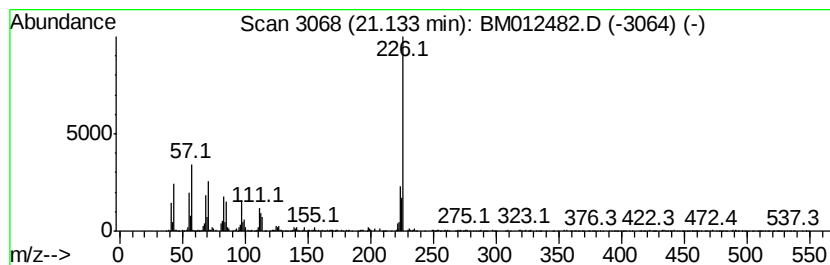
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ClientSampleId :  
B-18(5-6)\_101217

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 18 1-Heneicosanol Concentration Rank 10

| R.T.    | EstConc                       | Area           | Relative to ISTD | R.T.      |
|---------|-------------------------------|----------------|------------------|-----------|
| 21.13   | 3.59 ng/ul                    | 2433890        | Chrysene-d12     | 21.42     |
| Hit# of | 5                             | Tentative ID   | MW MolForm       | CAS# Qual |
| 1       | 1-Heneicosanol                | 312 C21H44O    | 015594-90-8      | 92        |
| 2       | Heptadecyl heptafluorobutyrte | 452 C21H35F7O2 | 959085-66-6      | 90        |
| 3       | 17-Pentatriacontene           | 491 C35H70     | 006971-40-0      | 90        |
| 4       | Behenic alcohol               | 326 C22H46O    | 000661-19-8      | 89        |
| 5       | Nonadecyl trifluoroacetate    | 380 C21H39F3O2 | 1000351-76-3     | 87        |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM102717\  
 Data File : BM012482.D  
 Acq On : 27 Oct 2017 12:31  
 Operator : SJ/JU  
 Sample : I5873-07  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 B-18(5-6)\_101217

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM102417.M  
 Quant Title : SVOA CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |          |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|----------|------|
|                      |       |         |       |          | #                     | RT    | Resp     | Conc |
| unknown-01           | 3.87  | 3.3     | ng/ul | 407649   | 1                     | 7.87  | 2439980  | 20.0 |
| 2-Butanol, 3-(2,2... | 7.80  | 3.4     | ng/ul | 417939   | 1                     | 7.87  | 2439980  | 20.0 |
| Benzene, 1,2-dime... | 12.07 | 21.6    | ng/ul | 3191670  | 2                     | 10.66 | 2958260  | 20.0 |
| Benzene, 1,2-dime... | 13.01 | 3.6     | ng/ul | 667883   | 3                     | 14.50 | 3730300  | 20.0 |
| n-Hexadecanoic acid  | 18.14 | 9.5     | ng/ul | 2045150  | 4                     | 17.24 | 4306570  | 20.0 |
| Anthracene, 1-met... | 18.16 | 5.6     | ng/ul | 1209900  | 4                     | 17.24 | 4306570  | 20.0 |
| 1H-Cyclopropa[1]p... | 18.24 | 2.6     | ng/ul | 567123   | 4                     | 17.24 | 4306570  | 20.0 |
| 4H-Cyclopenta[def... | 18.31 | 6.8     | ng/ul | 1467230  | 4                     | 17.24 | 4306570  | 20.0 |
| Phenanthrene, 4-m... | 18.34 | 2.6     | ng/ul | 566528   | 4                     | 17.24 | 4306570  | 20.0 |
| Naphthalene, 2-ph... | 18.62 | 3.7     | ng/ul | 801997   | 4                     | 17.24 | 4306570  | 20.0 |
| Phenanthrene, 2,5... | 19.03 | 3.5     | ng/ul | 752952   | 4                     | 17.24 | 4306570  | 20.0 |
| Cyclopenta(def)ph... | 19.17 | 4.4     | ng/ul | 955499   | 4                     | 17.24 | 4306570  | 20.0 |
| Pyrene, 1-methyl-    | 19.98 | 2.6     | ng/ul | 1764720  | 5                     | 21.42 | 13554200 | 20.0 |
| 11H-Benzo[a]fluor... | 20.90 | 2.5     | ng/ul | 1725700  | 5                     | 21.42 | 13554200 | 20.0 |
| 1-Heneicosanol       | 21.13 | 3.6     | ng/ul | 2433890  | 5                     | 21.42 | 13554200 | 20.0 |