

Data Path : Z:\HPCHEM1\BNA M\DATA\BM043018\  
 Data File : BM014878.D  
 Acq On : 30 Apr 2018 21:55  
 Operator : SJ/JU  
 Sample : J2551-07  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 F1B58

## 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-BM041918.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         | 3.346       | 5             | 10          | 24           | rVB      | 887632         | 1422012       | 16.80%          | 1.216%        |
| 2         | 3.646       | 56            | 61          | 70           | rVB      | 131283         | 191372        | 2.26%           | 0.164%        |
| 3         | 6.434       | 530           | 535         | 541          | rBV      | 64626          | 104975        | 1.24%           | 0.090%        |
| 4         | 7.622       | 730           | 737         | 745          | rBV      | 452202         | 742681        | 8.77%           | 0.635%        |
| 5         | 7.798       | 759           | 767         | 775          | rBV      | 1469778        | 2364839       | 27.94%          | 2.022%        |
| 6         | 8.016       | 797           | 804         | 813          | rBV      | 1636967        | 2632138       | 31.10%          | 2.250%        |
| 7         | 8.128       | 818           | 823         | 832          | rVB      | 157464         | 259550        | 3.07%           | 0.222%        |
| 8         | 8.498       | 878           | 886         | 900          | rBV      | 1539993        | 2587175       | 30.57%          | 2.212%        |
| 9         | 8.610       | 900           | 905         | 911          | rVB      | 58439          | 91629         | 1.08%           | 0.078%        |
| 10        | 8.881       | 944           | 951         | 959          | rVB      | 209670         | 343961        | 4.06%           | 0.294%        |
| 11        | 9.045       | 972           | 979         | 987          | rBV      | 401476         | 673633        | 7.96%           | 0.576%        |
| 12        | 9.192       | 997           | 1004        | 1015         | rBV      | 1196888        | 1961384       | 23.17%          | 1.677%        |
| 13        | 9.357       | 1025          | 1032        | 1040         | rVB      | 57573          | 97925         | 1.16%           | 0.084%        |
| 14        | 9.681       | 1079          | 1087        | 1097         | rBV      | 1770329        | 3071496       | 36.29%          | 2.626%        |
| 15        | 9.875       | 1114          | 1120        | 1128         | rVB      | 95708          | 159563        | 1.89%           | 0.136%        |
| 16        | 9.998       | 1128          | 1141        | 1147         | rBV      | 226925         | 471636        | 5.57%           | 0.403%        |
| 17        | 10.063      | 1147          | 1152        | 1160         | rVB      | 47668          | 85594         | 1.01%           | 0.073%        |
| 18        | 10.410      | 1200          | 1211        | 1223         | rBV      | 1400941        | 2359065       | 27.87%          | 2.017%        |
| 19        | 10.722      | 1258          | 1264        | 1278         | rBV4     | 68174          | 143613        | 1.70%           | 0.123%        |
| 20        | 10.945      | 1294          | 1302        | 1312         | rBV      | 2037847        | 3491713       | 41.26%          | 2.985%        |
| 21        | 11.339      | 1361          | 1369        | 1374         | rBV      | 1991508        | 3519983       | 41.59%          | 3.009%        |
| 22        | 11.392      | 1374          | 1378        | 1385         | rVV      | 1653768        | 2734216       | 32.31%          | 2.337%        |
| 23        | 11.469      | 1385          | 1391        | 1393         | rVV      | 63857          | 111731        | 1.32%           | 0.096%        |
| 24        | 11.510      | 1393          | 1398        | 1406         | rVB      | 743024         | 1321376       | 15.61%          | 1.130%        |
| 25        | 12.439      | 1549          | 1556        | 1563         | rBV      | 59674          | 111586        | 1.32%           | 0.095%        |
| 26        | 12.863      | 1622          | 1628        | 1640         | rVB2     | 124672         | 253729        | 3.00%           | 0.217%        |
| 27        | 12.963      | 1640          | 1645        | 1651         | rBV      | 132309         | 205641        | 2.43%           | 0.176%        |
| 28        | 13.180      | 1672          | 1682        | 1691         | rVB      | 1520349        | 2521125       | 29.79%          | 2.155%        |
| 29        | 13.945      | 1806          | 1812        | 1821         | rBV      | 1137979        | 1658887       | 19.60%          | 1.418%        |
| 30        | 14.075      | 1828          | 1834        | 1841         | rBV2     | 74099          | 126223        | 1.49%           | 0.108%        |
| 31        | 14.239      | 1857          | 1862        | 1865         | rBV2     | 55806          | 99228         | 1.17%           | 0.085%        |
| 32        | 14.280      | 1865          | 1869        | 1876         | rVB3     | 70497          | 141920        | 1.68%           | 0.121%        |
| 33        | 14.427      | 1888          | 1894        | 1899         | rBV2     | 132225         | 225436        | 2.66%           | 0.193%        |
| 34        | 14.492      | 1899          | 1905        | 1912         | rVB      | 3529868        | 4875455       | 57.60%          | 4.168%        |

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## Integration Parameters: LSCINT.P

Integrator: RTE  
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 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-BM041918.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |         |        |
|----|--------|------|------|------|------|---------|---------|---------|--------|
| 35 | 14.804 | 1950 | 1958 | 1970 | rBV  | 4068264 | 6060470 | 71.61%  | 5.181% |
| 36 | 15.063 | 1997 | 2002 | 2003 | rBV  | 115424  | 142801  | 1.69%   | 0.122% |
| 37 | 15.104 | 2003 | 2009 | 2014 | rVV2 | 2613963 | 4101827 | 48.46%  | 3.506% |
| 38 | 15.163 | 2014 | 2019 | 2025 | rVV  | 1514662 | 2282886 | 26.97%  | 1.951% |
| 39 | 15.227 | 2025 | 2030 | 2033 | rVV  | 557198  | 847327  | 10.01%  | 0.724% |
| 40 | 15.263 | 2033 | 2036 | 2050 | rVB  | 310534  | 510573  | 6.03%   | 0.436% |
| 41 | 15.492 | 2068 | 2075 | 2081 | rBV  | 2599826 | 3723520 | 43.99%  | 3.183% |
| 42 | 16.080 | 2165 | 2175 | 2180 | rBV  | 4897028 | 7121017 | 84.14%  | 6.087% |
| 43 | 16.133 | 2180 | 2184 | 2188 | rVV  | 1239278 | 1754785 | 20.73%  | 1.500% |
| 44 | 16.180 | 2188 | 2192 | 2201 | rVB  | 1484039 | 2019425 | 23.86%  | 1.726% |
| 45 | 16.316 | 2208 | 2215 | 2222 | rVB4 | 68841   | 140203  | 1.66%   | 0.120% |
| 46 | 16.416 | 2228 | 2232 | 2239 | rVB  | 117389  | 181570  | 2.15%   | 0.155% |
| 47 | 16.563 | 2252 | 2257 | 2268 | rVB3 | 139102  | 317286  | 3.75%   | 0.271% |
| 48 | 16.804 | 2285 | 2298 | 2309 | rBV  | 1247666 | 1961329 | 23.17%  | 1.677% |
| 49 | 17.433 | 2400 | 2405 | 2407 | rBV2 | 93865   | 133404  | 1.58%   | 0.114% |
| 50 | 17.468 | 2407 | 2411 | 2418 | rVB  | 576028  | 788990  | 9.32%   | 0.674% |
| 51 | 17.639 | 2436 | 2440 | 2442 | rVV  | 132534  | 191319  | 2.26%   | 0.164% |
| 52 | 17.668 | 2442 | 2445 | 2450 | rVB2 | 83483   | 124238  | 1.47%   | 0.106% |
| 53 | 17.815 | 2461 | 2470 | 2474 | rBV  | 2978040 | 4404036 | 52.03%  | 3.765% |
| 54 | 17.857 | 2474 | 2477 | 2481 | rVV  | 1769201 | 2349032 | 27.75%  | 2.008% |
| 55 | 17.915 | 2481 | 2487 | 2498 | rVV  | 5186699 | 7509670 | 88.73%  | 6.420% |
| 56 | 18.010 | 2498 | 2503 | 2508 | rVB2 | 72121   | 117779  | 1.39%   | 0.101% |
| 57 | 18.204 | 2531 | 2536 | 2542 | rVB  | 898783  | 1292116 | 15.27%  | 1.105% |
| 58 | 18.351 | 2557 | 2561 | 2566 | rVB  | 63870   | 84920   | 1.00%   | 0.073% |
| 59 | 18.433 | 2569 | 2575 | 2578 | rBV6 | 51663   | 95374   | 1.13%   | 0.082% |
| 60 | 18.639 | 2605 | 2610 | 2614 | rBV3 | 90502   | 135747  | 1.60%   | 0.116% |
| 61 | 18.721 | 2619 | 2624 | 2628 | rVB3 | 73379   | 121675  | 1.44%   | 0.104% |
| 62 | 18.857 | 2642 | 2647 | 2655 | rBV2 | 564367  | 869158  | 10.27%  | 0.743% |
| 63 | 19.139 | 2691 | 2695 | 2701 | rVB2 | 65025   | 101304  | 1.20%   | 0.087% |
| 64 | 19.821 | 2808 | 2811 | 2816 | rVB  | 145091  | 181562  | 2.15%   | 0.155% |
| 65 | 19.880 | 2818 | 2821 | 2831 | rVV4 | 80402   | 124857  | 1.48%   | 0.107% |
| 66 | 20.151 | 2861 | 2867 | 2879 | rBV  | 6202540 | 8116347 | 95.90%  | 6.938% |
| 67 | 21.903 | 3160 | 3165 | 3172 | rBV  | 3642089 | 4663806 | 55.10%  | 3.987% |
| 68 | 24.221 | 3552 | 3559 | 3570 | rVB  | 4121271 | 8463716 | 100.00% | 7.235% |
| 69 | 24.380 | 3579 | 3586 | 3599 | rVB  | 2290183 | 4785064 | 56.54%  | 4.090% |

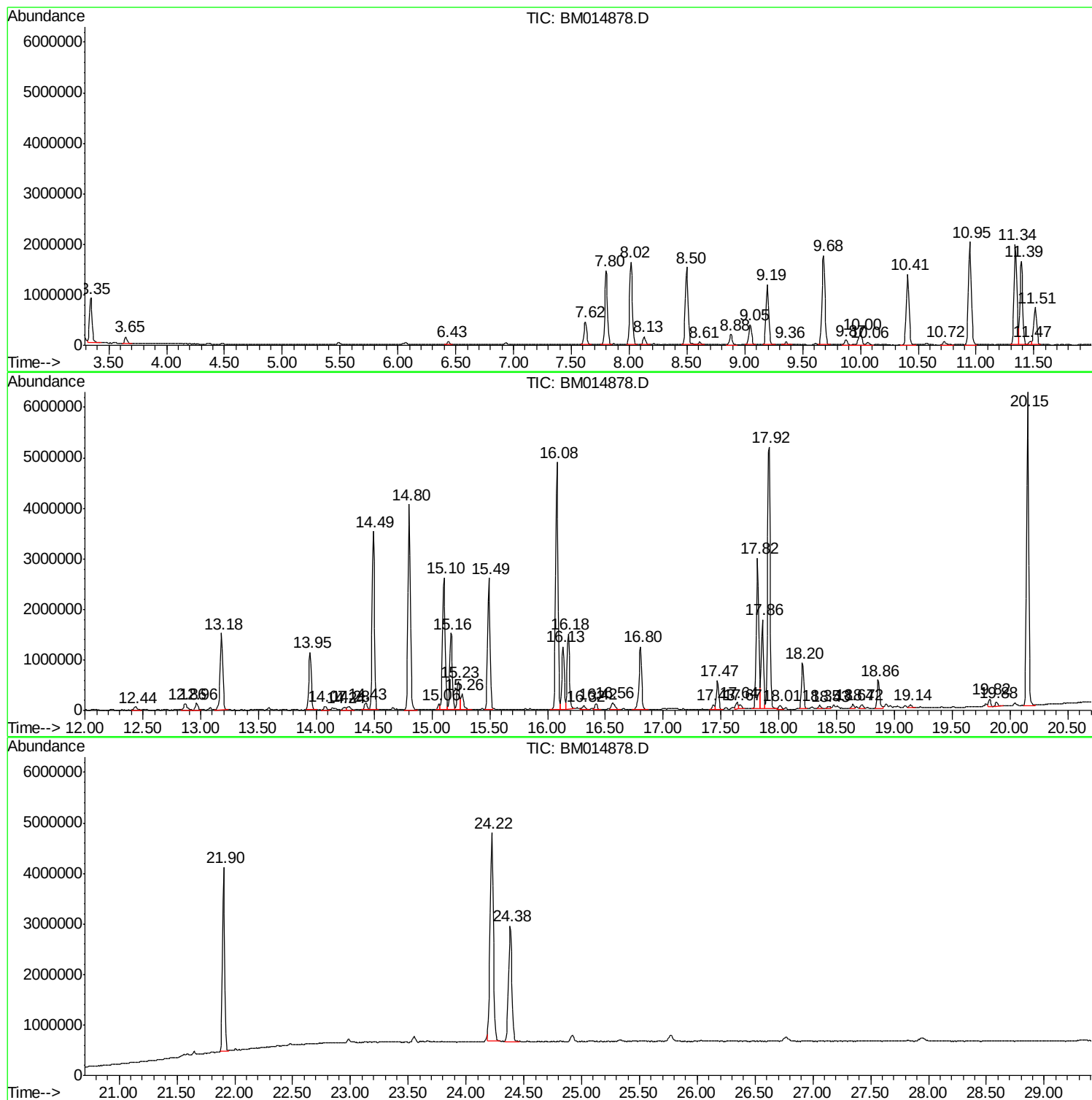
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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M  
Quant Title : SVOA CALIBRATION

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



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Misc :  
ALS Vial : 21 Sample Multiplier: 1

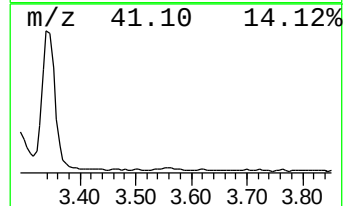
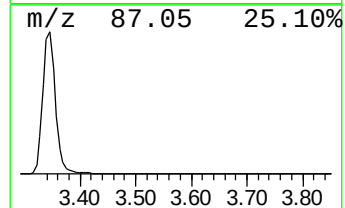
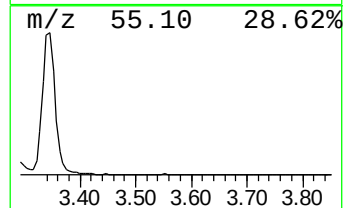
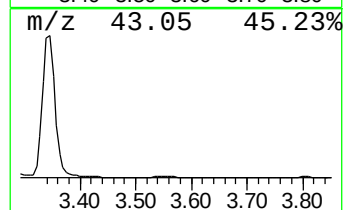
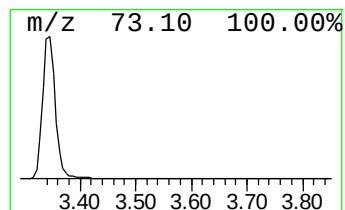
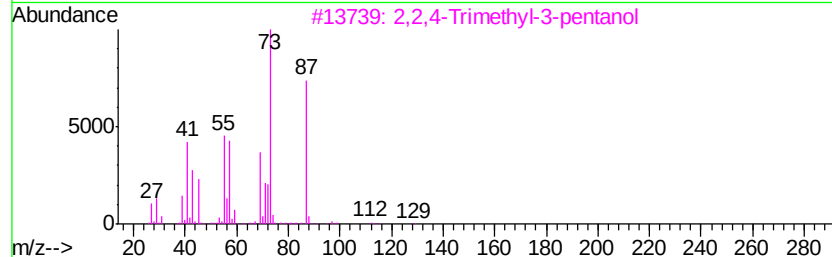
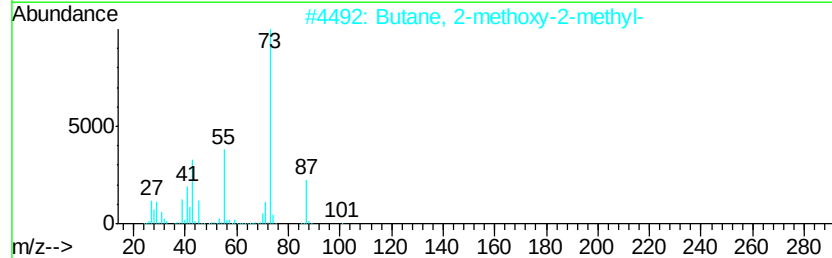
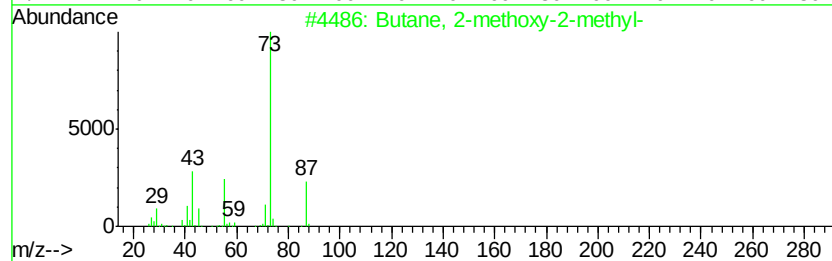
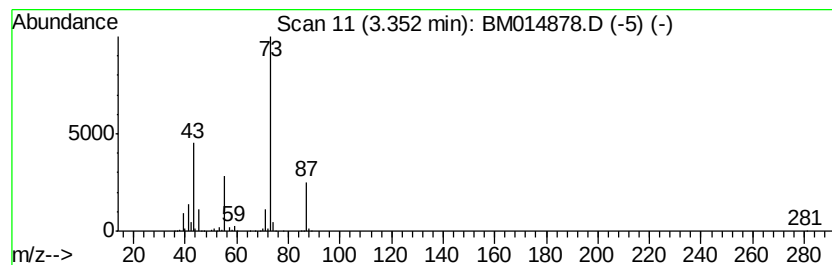
Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

| R.T.    | EstConc                     | Area         | Relative to ISTD       | R.T.        |      |      |
|---------|-----------------------------|--------------|------------------------|-------------|------|------|
| 3.35    | 10.99 ng/ul                 | 1422010      | 1,4-Dichlorobenzene-d4 | 8.50        |      |      |
| Hit# of | 5                           | Tentative ID | MW                     | MolForm     | CAS# | Qual |
| 1       | Butane, 2-methoxy-2-methyl- | 102          | C6H14O                 | 000994-05-8 | 83   |      |
| 2       | Butane, 2-methoxy-2-methyl- | 102          | C6H14O                 | 000994-05-8 | 78   |      |
| 3       | 2,2,4-Trimethyl-3-pentanol  | 130          | C8H18O                 | 005162-48-1 | 40   |      |
| 4       | 2,2,4-Trimethyl-3-pentanol  | 130          | C8H18O                 | 005162-48-1 | 40   |      |
| 5       | Silane, tetramethyl-        | 88           | C4H12Si                | 000075-76-3 | 38   |      |



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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M  
Quant Title : SVOA CALIBRATION

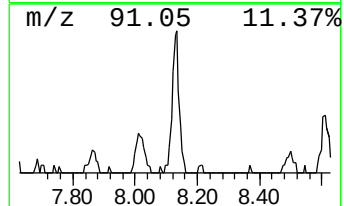
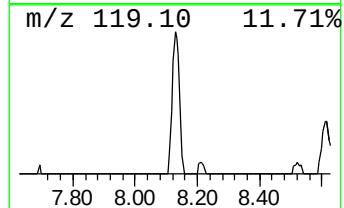
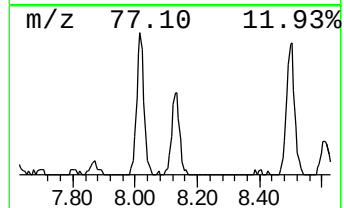
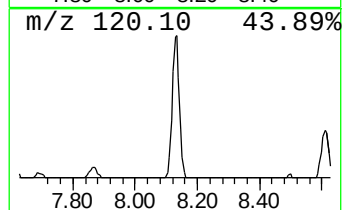
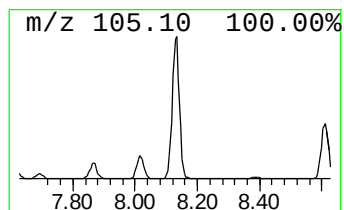
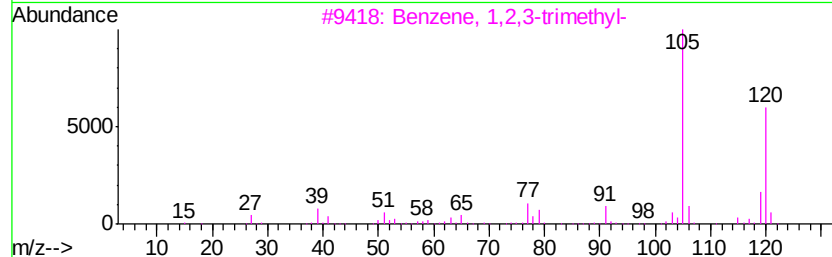
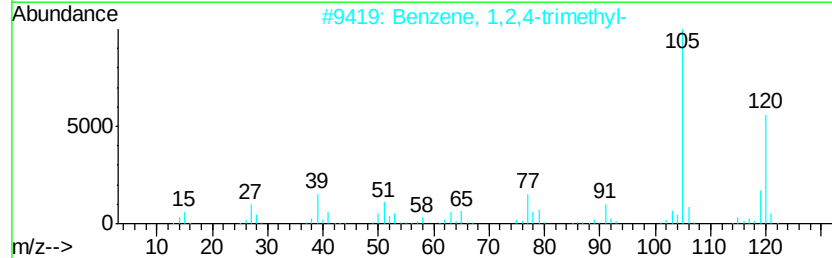
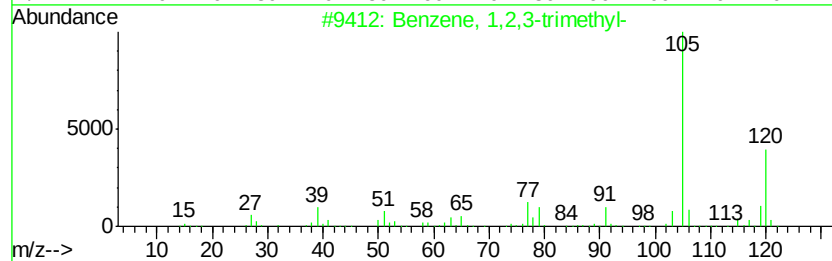
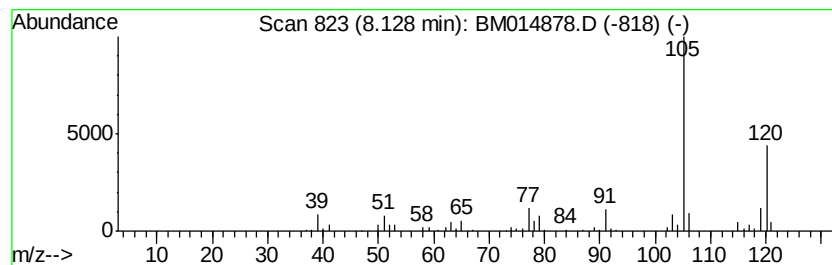
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 Benzene, 1,2,3-trimethyl- Concentration Rank 9

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.13 | 2.01 ng/ul | 259550 | 1,4-Dichlorobenzene-d4 | 8.50 |

| Hit# of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|---------|---|---------------------------|-----|---------|-------------|------|
| 1       |   | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 97   |
| 2       |   | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 95   |
| 3       |   | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 95   |
| 4       |   | Mesitylene                | 120 | C9H12   | 000108-67-8 | 94   |
| 5       |   | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 94   |



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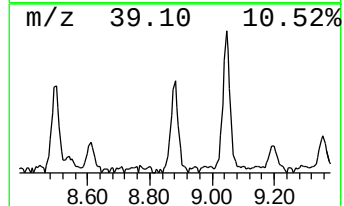
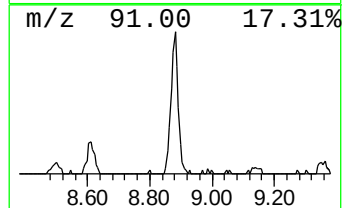
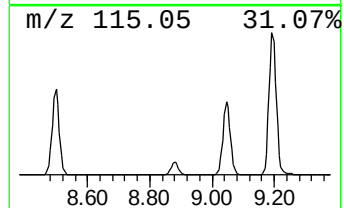
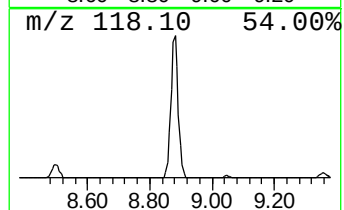
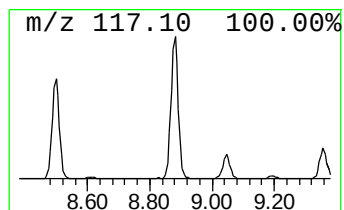
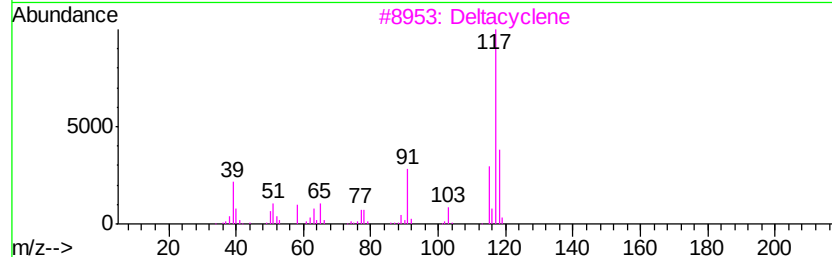
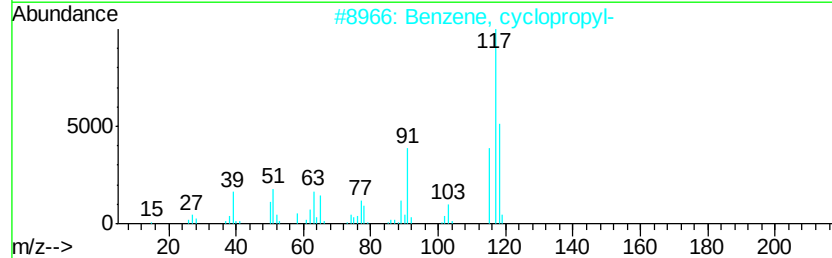
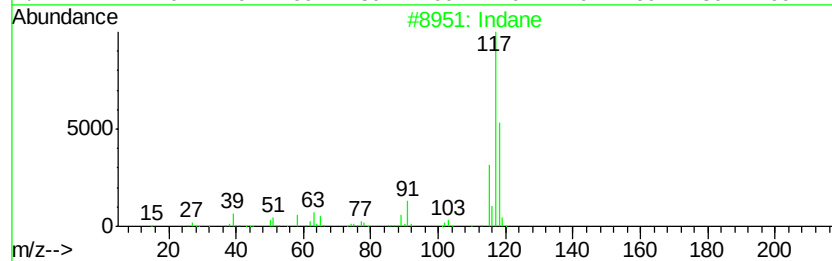
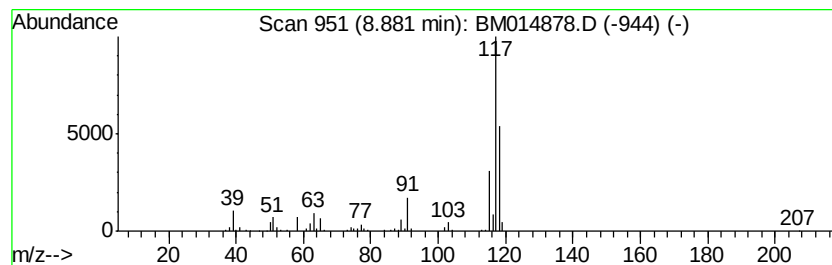
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Indane Concentration Rank 8

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.88 | 2.66 ng/ul | 343961 | 1,4-Dichlorobenzene-d4 | 8.50 |

| Hit# of | 5                     | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|-----------------------|--------------|-----|---------|-------------|------|
| 1       | Indane                |              | 118 | C9H10   | 000496-11-7 | 93   |
| 2       | Benzene, cyclopropyl- |              | 118 | C9H10   | 000873-49-4 | 81   |
| 3       | Deltacyclene          |              | 118 | C9H10   | 007785-10-6 | 76   |
| 4       | Benzene, cyclopropyl- |              | 118 | C9H10   | 000873-49-4 | 74   |
| 5       | Indane                |              | 118 | C9H10   | 000496-11-7 | 74   |



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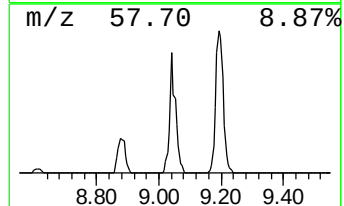
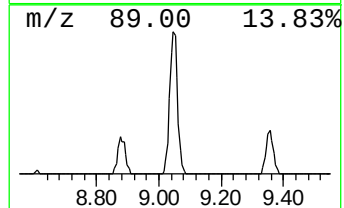
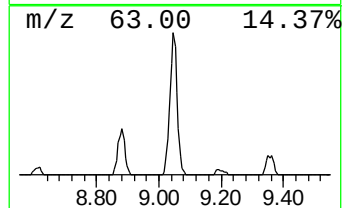
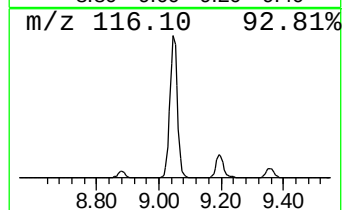
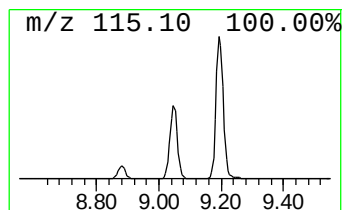
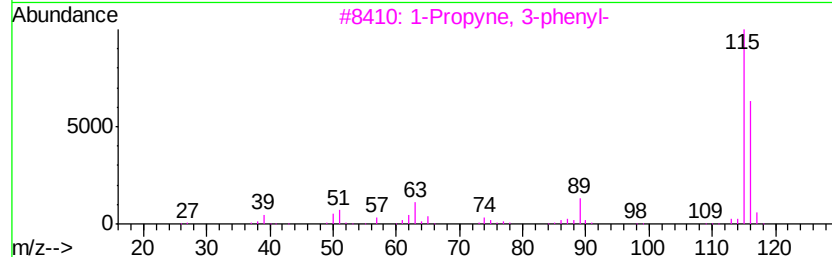
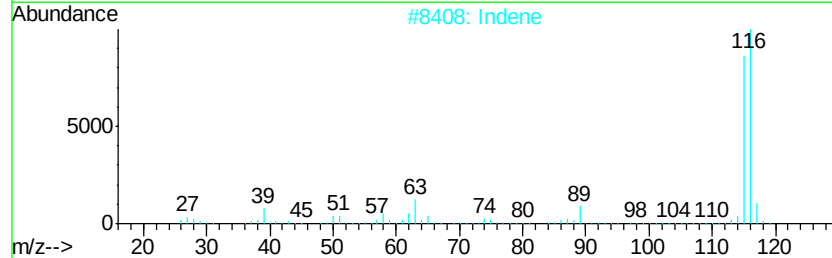
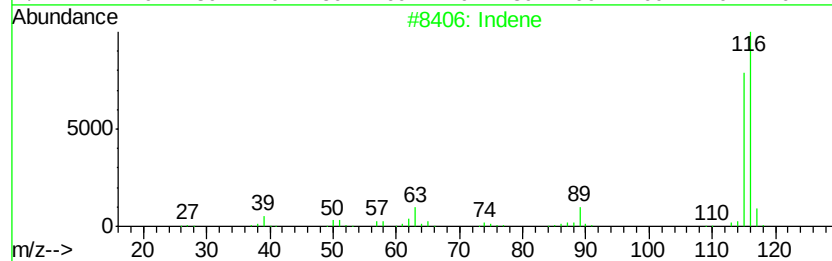
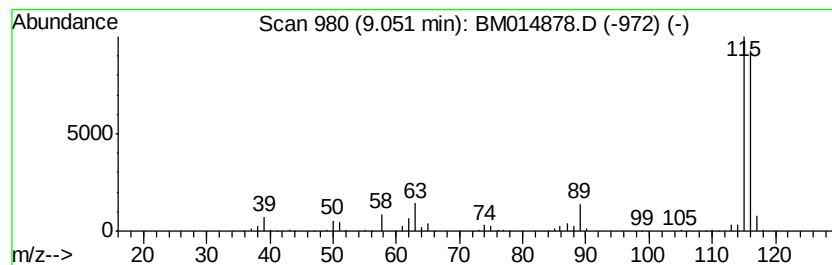
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BNA\_M  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 Indene Concentration Rank 4

| R.T.    | EstConc                 | Area         | Relative to ISTD       | R.T.           |
|---------|-------------------------|--------------|------------------------|----------------|
| 9.05    | 5.21 ng/ul              | 673633       | 1,4-Dichlorobenzene-d4 | 8.50           |
| Hit# of | 5                       | Tentative ID | MW MolForm             | CAS# Qual      |
| 1       | Indene                  |              | 116 C9H8               | 000095-13-6 97 |
| 2       | Indene                  |              | 116 C9H8               | 000095-13-6 97 |
| 3       | 1-Propyne, 3-phenyl-    |              | 116 C9H8               | 010147-11-2 95 |
| 4       | Indene                  |              | 116 C9H8               | 000095-13-6 95 |
| 5       | 3-Methylphenylacetylene |              | 116 C9H8               | 000766-82-5 94 |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM043018\  
Data File : BM014878.D  
Acq On : 30 Apr 2018 21:55  
Operator : SJ/JU  
Sample : J2551-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F1B58

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M  
Quant Title : SVOA CALIBRATION

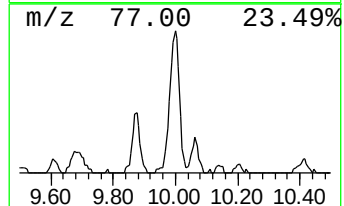
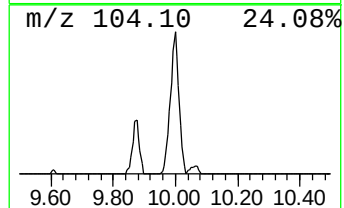
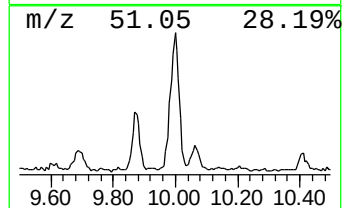
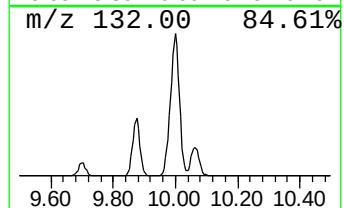
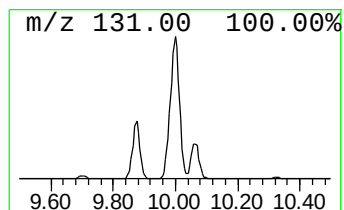
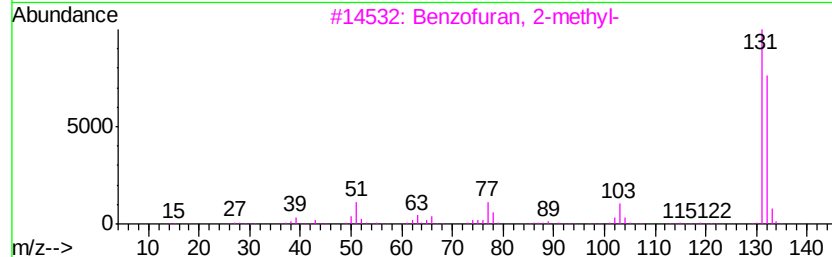
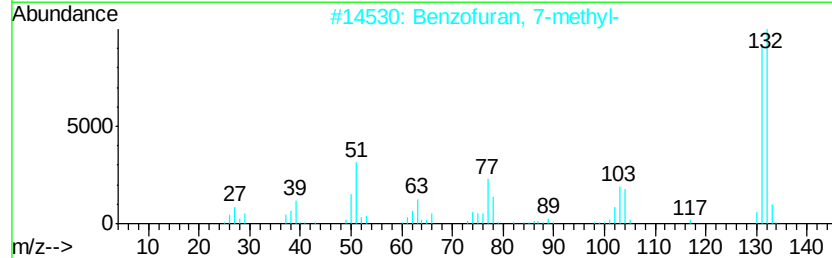
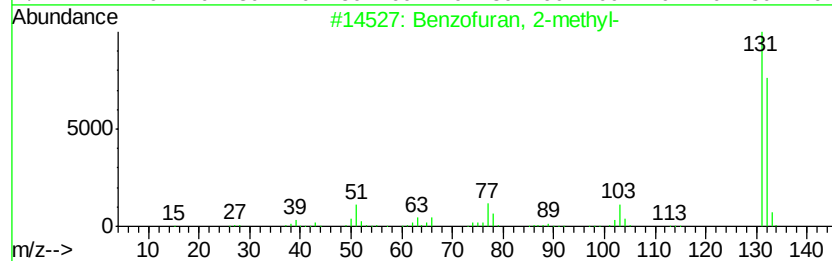
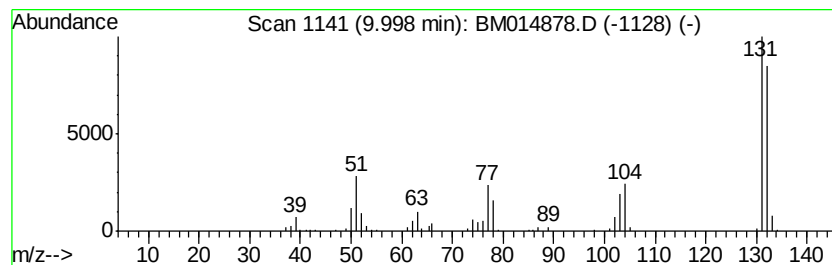
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Benzofuran, 2-methyl- Concentration Rank 7

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 10.00 | 2.68 ng/ul | 471636 | Napthalene-d8    | 11.34 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzofuran, 2-methyl-  | 132 | C9H8O   | 004265-25-2 | 94   |
| 2    |    | Benzofuran, 7-methyl-  | 132 | C9H8O   | 017059-52-8 | 94   |
| 3    |    | Benzofuran, 2-methyl-  | 132 | C9H8O   | 004265-25-2 | 94   |
| 4    |    | 3-Phenyl-2-propyn-1-ol | 132 | C9H8O   | 001504-58-1 | 91   |
| 5    |    | 2-Propenal, 3-phenyl-  | 132 | C9H8O   | 000104-55-2 | 87   |





Data Path : Z:\HPCHEM1\BNA M\DATA\BM043018\  
Data File : BM014878.D  
Acq On : 30 Apr 2018 21:55  
Operator : SJ/JU  
Sample : J2551-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F1B58

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M  
Quant Title : SVOA CALIBRATION

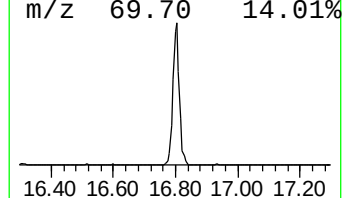
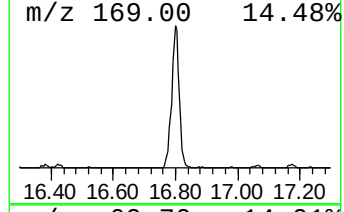
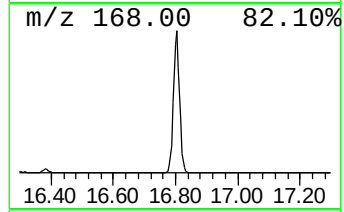
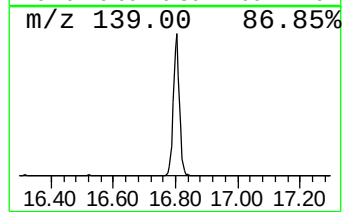
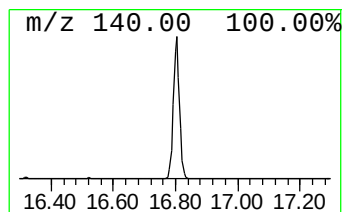
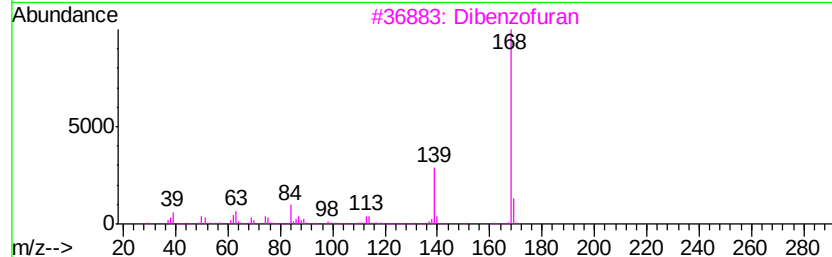
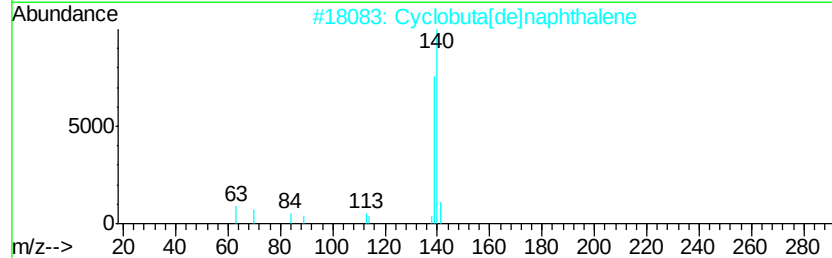
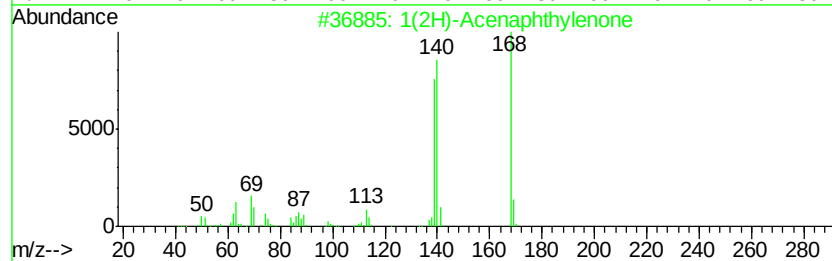
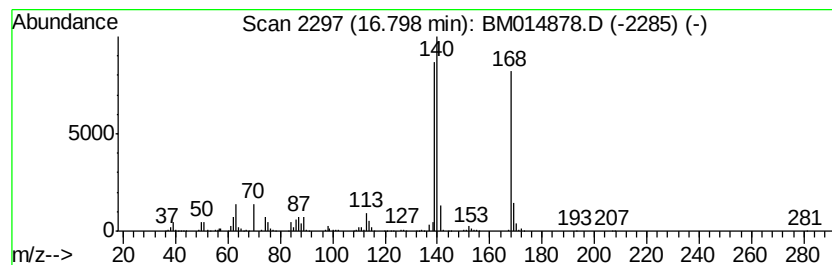
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 1(2H)-Acenaphthylenone Concentration Rank 3

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 16.80 | 8.91 ng/ul | 1961330 | Phenanthrene-d10 | 17.82 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 1(2H)-Acenaphthylenone              | 168 | C12H8O   | 002235-15-6 | 96   |
| 2    |    | Cyclobuta[de]naphthalene            | 140 | C11H8    | 024973-91-9 | 64   |
| 3    |    | Dibenzofuran                        | 168 | C12H8O   | 000132-64-9 | 64   |
| 4    |    | 7(4H)-Benzothiazolone, 2-amino-5... | 168 | C7H8N2OS | 017583-10-7 | 50   |
| 5    |    | 2-Amino-5,6-dihydro-4H-cyclopent... | 140 | C6H8N2S  | 053051-97-1 | 47   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM043018\  
Data File : BM014878.D  
Acq On : 30 Apr 2018 21:55  
Operator : SJ/JU  
Sample : J2551-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F1B58

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M  
Quant Title : SVOA CALIBRATION

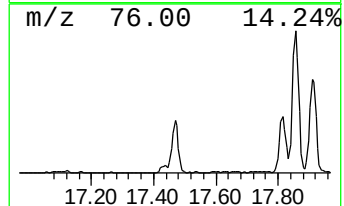
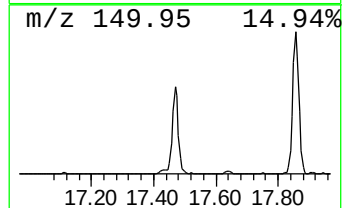
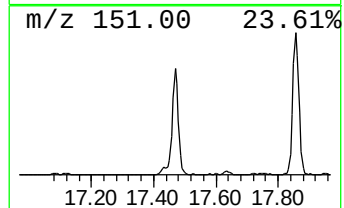
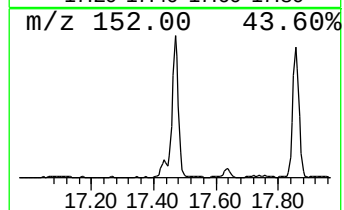
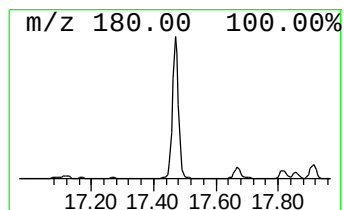
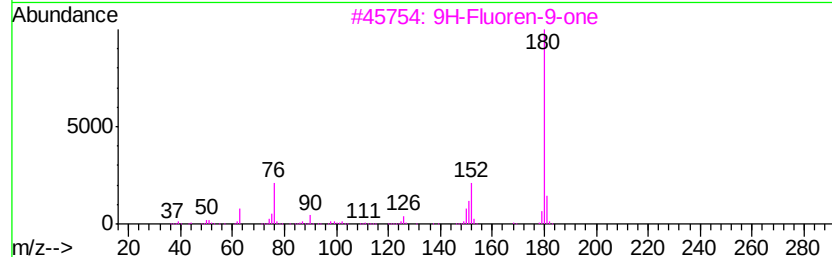
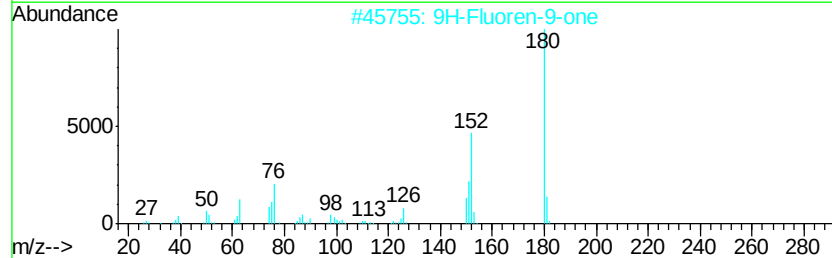
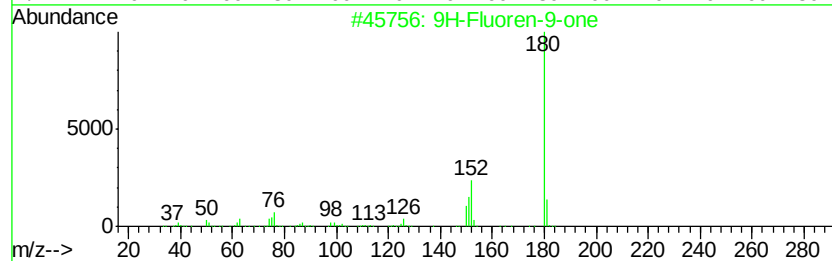
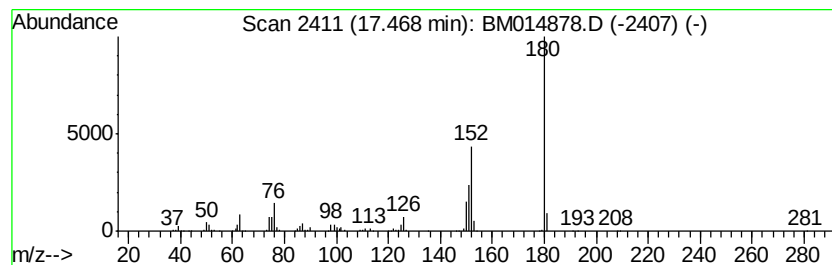
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 9H-Fluoren-9-one Concentration Rank 6

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.47 | 3.58 ng/ul | 788990 | Phenanthrene-d10 | 17.82 |

| Hit# | of | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------|-----|---------|-------------|------|
| 1    | 5  | 9H-Fluoren-9-one  | 180 | C13H8O  | 000486-25-9 | 95   |
| 2    |    | 9H-Fluoren-9-one  | 180 | C13H8O  | 000486-25-9 | 91   |
| 3    |    | 9H-Fluoren-9-one  | 180 | C13H8O  | 000486-25-9 | 91   |
| 4    |    | 9H-Fluoren-9-one  | 180 | C13H8O  | 000486-25-9 | 76   |
| 5    |    | Benzo[h]cinnoline | 180 | C12H8N2 | 000230-31-9 | 58   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM043018\  
Data File : BM014878.D  
Acq On : 30 Apr 2018 21:55  
Operator : SJ/JU  
Sample : J2551-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

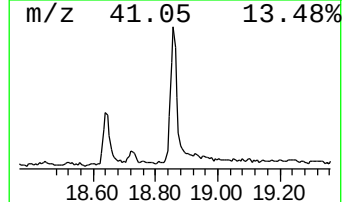
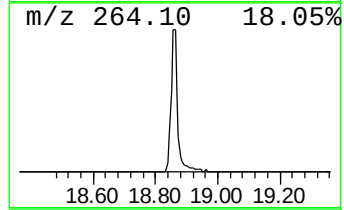
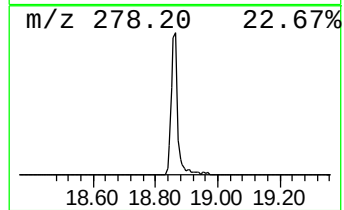
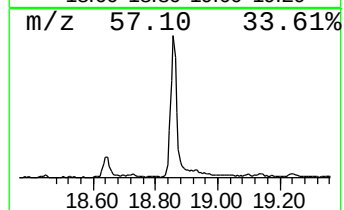
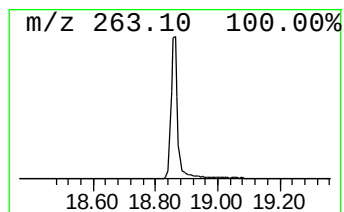
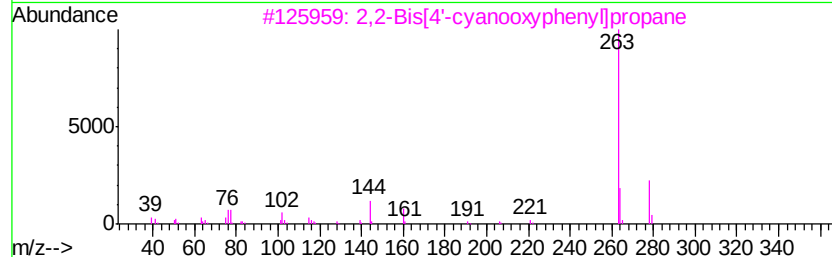
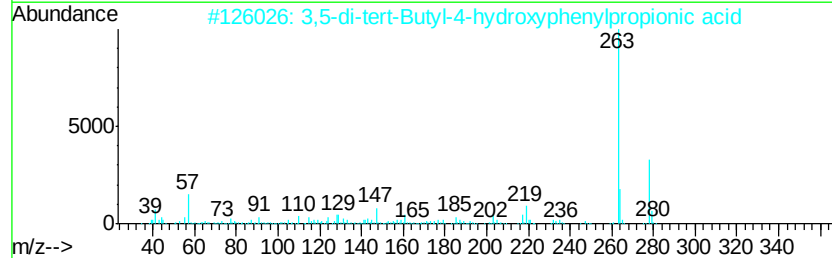
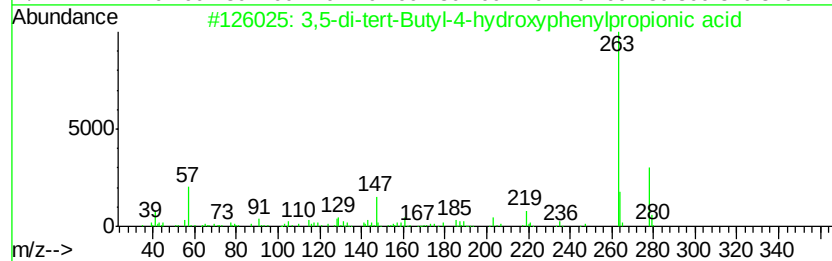
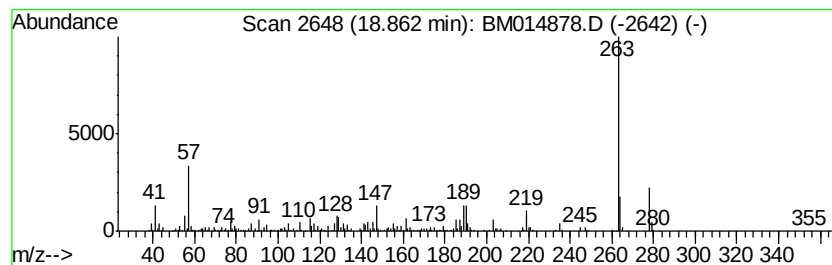
Instrument :  
BNA\_M  
ClientSampleId :  
F1B58

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 9 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 5

| R.T.    | EstConc                              | Area         | Relative to ISTD | R.T.         |      |      |
|---------|--------------------------------------|--------------|------------------|--------------|------|------|
| 18.86   | 3.95 ng/ul                           | 869158       | Phenanthrene-d10 | 17.82        |      |      |
| Hit# of | 5                                    | Tentative ID | MW               | MolForm      | CAS# | Qual |
| 1       | 3,5-di-tert-Butyl-4-hydroxyphenyl... | 278          | C17H26O3         | 020170-32-5  | 94   |      |
| 2       | 3,5-di-tert-Butyl-4-hydroxyphenyl... | 278          | C17H26O3         | 020170-32-5  | 86   |      |
| 3       | 2,2-Bis[4'-cyanoxyphenyl]propane     | 278          | C17H14N2O2       | 1000322-13-3 | 64   |      |
| 4       | Fenoxon sulfoxide                    | 278          | C10H15O5PS       | 006552-13-2  | 62   |      |
| 5       | 4,6-Dinitro-1,1,3,3,5-pentamethy...  | 278          | C14H18N2O4       | 000116-66-5  | 59   |      |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM043018\  
Data File : BM014878.D  
Acq On : 30 Apr 2018 21:55  
Operator : SJ/JU  
Sample : J2551-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F1B58

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM041918.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 |
| Butane, 2-methoxy... | 3.35  | 11.0    | ng/ul | 1422010  | 1                     | 8.50  | 2587180 | 20.0 |
| Benzene, 1,2,3-tr... | 8.13  | 2.0     | ng/ul | 259550   | 1                     | 8.50  | 2587180 | 20.0 |
| Indane               | 8.88  | 2.7     | ng/ul | 343961   | 1                     | 8.50  | 2587180 | 20.0 |
| Indene               | 9.05  | 5.2     | ng/ul | 673633   | 1                     | 8.50  | 2587180 | 20.0 |
| Benzofuran, 2-met... | 10.00 | 2.7     | ng/ul | 471636   | 2                     | 11.34 | 3519980 | 20.0 |
| 1(2H)-Acenaphthyl... | 16.80 | 8.9     | ng/ul | 1961330  | 4                     | 17.82 | 4404040 | 20.0 |
| 9H-Fluoren-9-one     | 17.47 | 3.6     | ng/ul | 788990   | 4                     | 17.82 | 4404040 | 20.0 |
| 3,5-di-tert-Butyl... | 18.86 | 4.0     | ng/ul | 869158   | 4                     | 17.82 | 4404040 | 20.0 |