

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101119\  
 Data File : BM023142.D  
 Acq On : 11 Oct 2019 22:32  
 Operator : JU  
 Sample : K5262-15  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BFBB6

## Integration Parameters: LSCINT.P

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

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM092719MA.M  
 Title : SVOA CALIBRATION

Signal : TIC

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|--------|----------|------------|----------|-----------|-------|-------------|------------|--------------|------------|
| 1      | 3.252    | 42         | 45       | 57        | rVB   | 39466       | 58739      | 0.69%        | 0.126%     |
| 2      | 3.981    | 165        | 169      | 173       | rVB   | 8715        | 11087      | 0.13%        | 0.024%     |
| 3      | 5.846    | 481        | 486      | 490       | rBV5  | 5878        | 9948       | 0.12%        | 0.021%     |
| 4      | 5.887    | 490        | 493      | 502       | rVV2  | 14811       | 27433      | 0.32%        | 0.059%     |
| 5      | 6.940    | 665        | 672      | 680       | rBV   | 152817      | 244504     | 2.89%        | 0.524%     |
| 6      | 7.104    | 693        | 700      | 722       | rBV   | 495257      | 822604     | 9.73%        | 1.763%     |
| 7      | 7.304    | 727        | 734      | 746       | rVB   | 614665      | 963858     | 11.40%       | 2.066%     |
| 8      | 7.769    | 804        | 813      | 819       | rBV   | 623955      | 995706     | 11.78%       | 2.134%     |
| 9      | 7.839    | 819        | 825      | 835       | rVB   | 37858       | 70099      | 0.83%        | 0.150%     |
| 10     | 8.475    | 925        | 933      | 944       | rBV   | 427430      | 708503     | 8.38%        | 1.519%     |
| 11     | 8.592    | 948        | 953      | 960       | rVB   | 26440       | 42854      | 0.51%        | 0.092%     |
| 12     | 8.863    | 991        | 999      | 1003      | rBV   | 43006       | 68655      | 0.81%        | 0.147%     |
| 13     | 8.922    | 1003       | 1009     | 1021      | rVB   | 633291      | 1039707    | 12.30%       | 2.229%     |
| 14     | 9.051    | 1027       | 1031     | 1034      | rBV2  | 6596        | 9413       | 0.11%        | 0.020%     |
| 15     | 9.092    | 1034       | 1038     | 1048      | rVB2  | 11632       | 19896      | 0.24%        | 0.043%     |
| 16     | 9.292    | 1068       | 1072     | 1082      | rBV5  | 11326       | 27399      | 0.32%        | 0.059%     |
| 17     | 9.639    | 1123       | 1131     | 1147      | rBV   | 536148      | 888313     | 10.51%       | 1.904%     |
| 18     | 9.816    | 1155       | 1161     | 1162      | rBV4  | 5883        | 8853       | 0.10%        | 0.019%     |
| 19     | 10.181   | 1215       | 1223     | 1242      | rBV   | 792601      | 1347769    | 15.95%       | 2.889%     |
| 20     | 10.339   | 1245       | 1250     | 1255      | rVB3  | 6095        | 9977       | 0.12%        | 0.021%     |
| 21     | 10.469   | 1268       | 1272     | 1279      | rVB5  | 10625       | 20610      | 0.24%        | 0.044%     |
| 22     | 10.557   | 1279       | 1287     | 1294      | rBV   | 805369      | 1371160    | 16.22%       | 2.939%     |
| 23     | 10.616   | 1294       | 1297     | 1301      | rVB2  | 9069        | 12816      | 0.15%        | 0.027%     |
| 24     | 10.869   | 1333       | 1340     | 1344      | rBV4  | 5161        | 8773       | 0.10%        | 0.019%     |
| 25     | 11.216   | 1394       | 1399     | 1410      | rVB   | 7814        | 13883      | 0.16%        | 0.030%     |
| 26     | 11.357   | 1417       | 1423     | 1429      | rBV2  | 14976       | 26485      | 0.31%        | 0.057%     |
| 27     | 11.904   | 1507       | 1516     | 1523      | rBV3  | 8540        | 18185      | 0.22%        | 0.039%     |
| 28     | 12.145   | 1552       | 1557     | 1562      | rBV   | 11650       | 18420      | 0.22%        | 0.039%     |
| 29     | 12.763   | 1654       | 1662     | 1670      | rBV   | 30077       | 48699      | 0.58%        | 0.104%     |
| 30     | 13.016   | 1697       | 1705     | 1712      | rBV   | 67056       | 99107      | 1.17%        | 0.212%     |
| 31     | 13.310   | 1748       | 1755     | 1764      | rBV   | 414593      | 688285     | 8.14%        | 1.475%     |
| 32     | 13.721   | 1820       | 1825     | 1832      | rVB   | 163344      | 203109     | 2.40%        | 0.435%     |
| 33     | 13.816   | 1834       | 1841     | 1851      | rBV   | 1321014     | 1814053    | 21.46%       | 3.889%     |
| 34     | 14.098   | 1879       | 1889     | 1899      | rBV   | 1624317     | 2315202    | 27.39%       | 4.963%     |

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Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM092719MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |         |         |
|----|--------|------|------|------|------|---------|---------|---------|---------|
| 35 | 14.404 | 1935 | 1941 | 1949 | rVB  | 1115258 | 1656016 | 19.59%  | 3.550%  |
| 36 | 14.616 | 1972 | 1977 | 1988 | rBV  | 79506   | 147514  | 1.75%   | 0.316%  |
| 37 | 14.868 | 2016 | 2020 | 2024 | rBV2 | 7438    | 10063   | 0.12%   | 0.022%  |
| 38 | 15.151 | 2062 | 2068 | 2076 | rVV  | 508871  | 682741  | 8.08%   | 1.463%  |
| 39 | 15.227 | 2076 | 2081 | 2086 | rVB2 | 41928   | 65001   | 0.77%   | 0.139%  |
| 40 | 15.398 | 2103 | 2110 | 2118 | rBV  | 2007033 | 2766365 | 32.73%  | 5.930%  |
| 41 | 15.521 | 2125 | 2131 | 2139 | rBV  | 491609  | 704178  | 8.33%   | 1.509%  |
| 42 | 15.639 | 2147 | 2151 | 2156 | rBV3 | 21131   | 27496   | 0.33%   | 0.059%  |
| 43 | 15.762 | 2166 | 2172 | 2179 | rBV  | 30183   | 43260   | 0.51%   | 0.093%  |
| 44 | 15.951 | 2195 | 2204 | 2214 | rBV  | 19511   | 41172   | 0.49%   | 0.088%  |
| 45 | 16.180 | 2240 | 2243 | 2246 | rBV2 | 7846    | 9343    | 0.11%   | 0.020%  |
| 46 | 16.657 | 2316 | 2324 | 2336 | rBV  | 105622  | 170606  | 2.02%   | 0.366%  |
| 47 | 16.927 | 2367 | 2370 | 2377 | rVB4 | 5198    | 8794    | 0.10%   | 0.019%  |
| 48 | 17.145 | 2401 | 2407 | 2415 | rVB2 | 1258610 | 1754623 | 20.76%  | 3.761%  |
| 49 | 17.245 | 2417 | 2424 | 2434 | rBV  | 2147630 | 2968486 | 35.12%  | 6.363%  |
| 50 | 17.433 | 2452 | 2456 | 2460 | rVB  | 16145   | 19038   | 0.23%   | 0.041%  |
| 51 | 17.815 | 2516 | 2521 | 2527 | rBV  | 340686  | 417798  | 4.94%   | 0.896%  |
| 52 | 18.115 | 2568 | 2572 | 2576 | rBV  | 30491   | 38261   | 0.45%   | 0.082%  |
| 53 | 18.186 | 2579 | 2584 | 2588 | rBV  | 16815   | 26838   | 0.32%   | 0.058%  |
| 54 | 18.298 | 2600 | 2603 | 2609 | rVB8 | 5984    | 8503    | 0.10%   | 0.018%  |
| 55 | 19.174 | 2748 | 2752 | 2758 | rBV  | 24019   | 33544   | 0.40%   | 0.072%  |
| 56 | 19.427 | 2792 | 2795 | 2799 | rVB  | 18747   | 19613   | 0.23%   | 0.042%  |
| 57 | 19.480 | 2800 | 2804 | 2808 | rBV  | 134930  | 156148  | 1.85%   | 0.335%  |
| 58 | 19.539 | 2808 | 2814 | 2824 | rVV  | 2646066 | 3539322 | 41.87%  | 7.587%  |
| 59 | 20.468 | 2964 | 2972 | 2986 | rBV  | 5782095 | 8452604 | 100.00% | 18.119% |
| 60 | 20.574 | 2987 | 2990 | 2996 | rVB  | 70176   | 105124  | 1.24%   | 0.225%  |
| 61 | 21.033 | 3064 | 3068 | 3076 | rBV  | 260076  | 349530  | 4.14%   | 0.749%  |
| 62 | 21.327 | 3113 | 3118 | 3124 | rBV  | 1681492 | 2117284 | 25.05%  | 4.539%  |
| 63 | 23.438 | 3471 | 3477 | 3490 | rVB  | 2078107 | 4000553 | 47.33%  | 8.575%  |
| 64 | 23.580 | 3495 | 3501 | 3510 | rVB  | 1192559 | 2277516 | 26.94%  | 4.882%  |

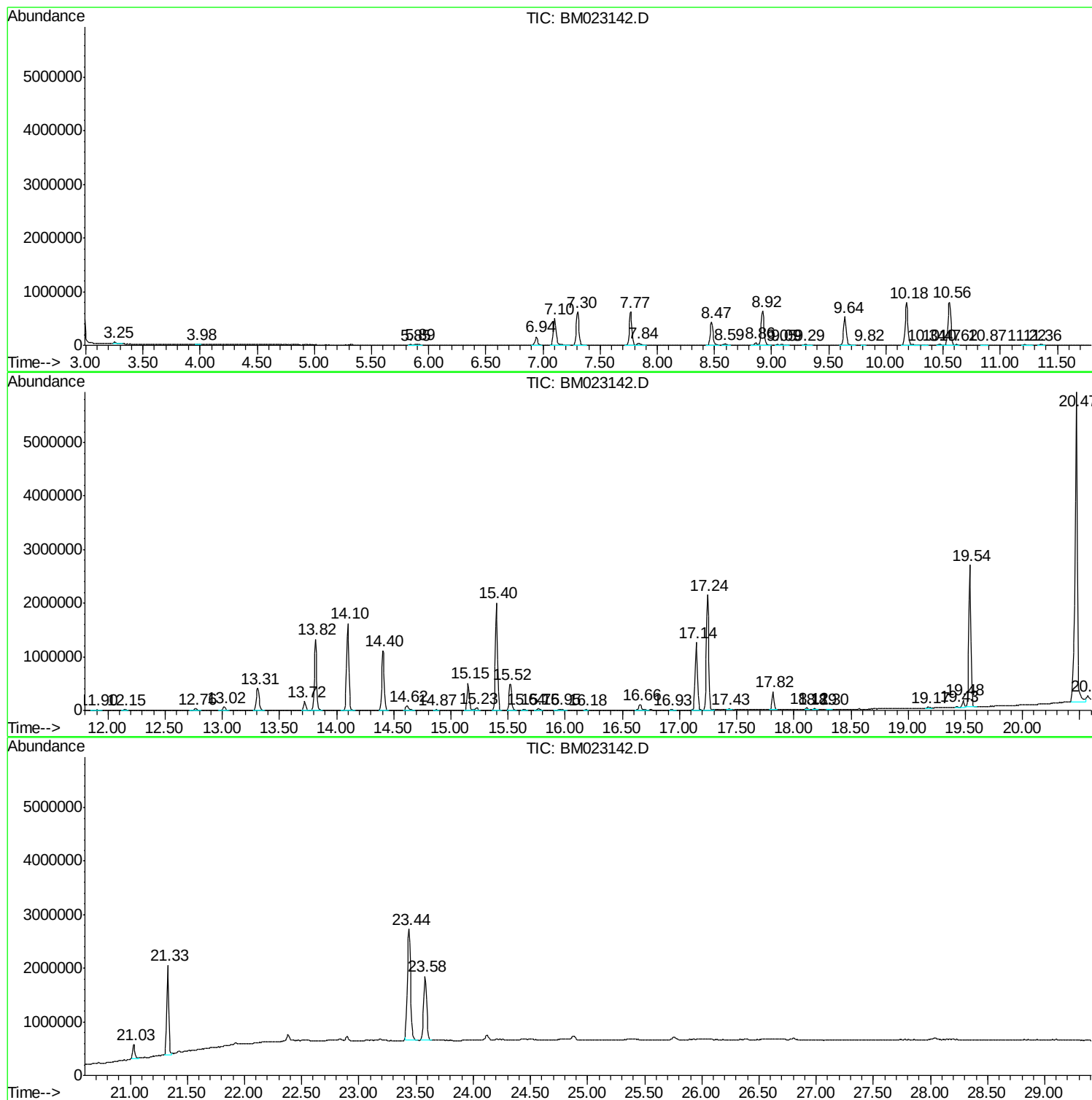
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BNA\_M  
ClientSampleId :  
BFBB6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM092719MA.M  
Quant Title : SVOA CALIBRATION

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



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Data File : BM023142.D  
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ALS Vial : 23 Sample Multiplier: 1

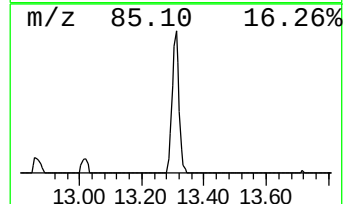
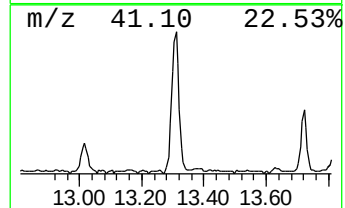
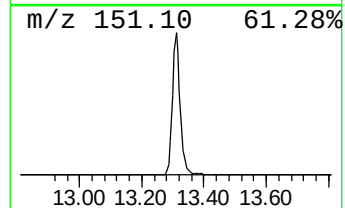
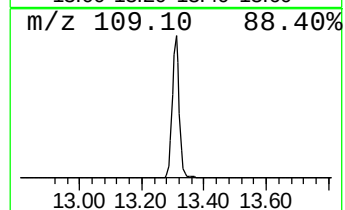
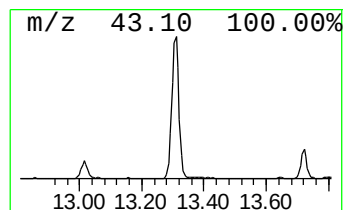
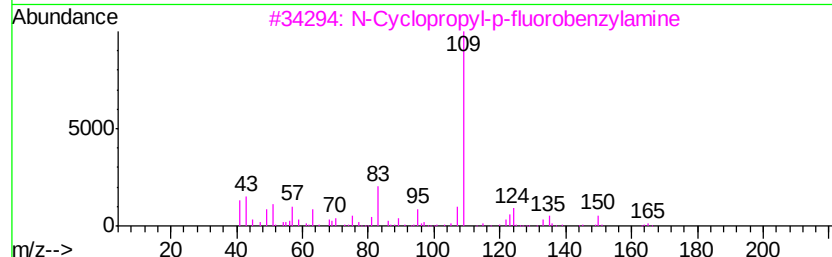
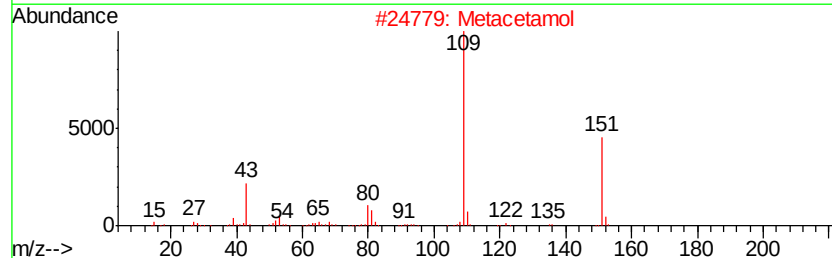
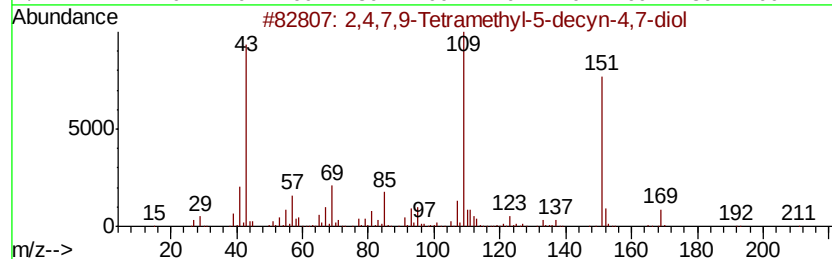
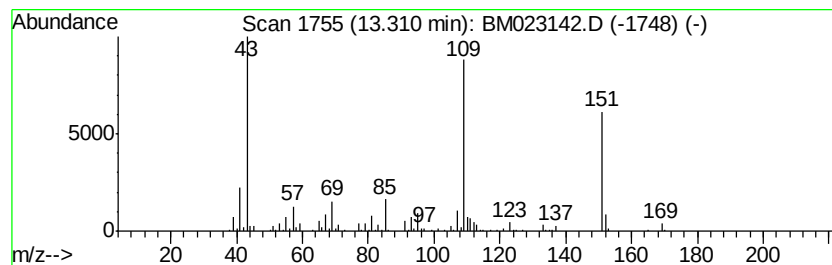
Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM092719MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 2

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 13.31     | 8.31 ng/ul                          | 688285 | Acenaphthene-d10 | 14.40        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 2,4,7,9-Tetramethyl-5-decyn-4,7-... | 226    | C14H26O2         | 000126-86-3  | 91   |
| 2         | Metacetamol                         | 151    | C8H9NO2          | 000621-42-1  | 53   |
| 3         | N-Cvclopropyl-p-fluorobenzylamine   | 165    | C10H12FN         | 1000250-88-9 | 47   |
| 4         | Acetamide, N-(2-hydroxyphenyl)-     | 151    | C8H9NO2          | 000614-80-2  | 42   |
| 5         | 3-Pyridinol, 4-methyl-, acetate ... | 151    | C8H9NO2          | 001006-96-8  | 42   |



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM092719MA.M  
Quant Title : SVOA CALIBRATION

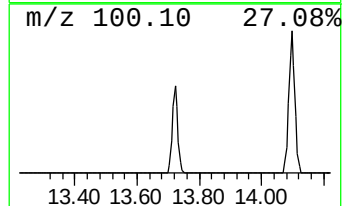
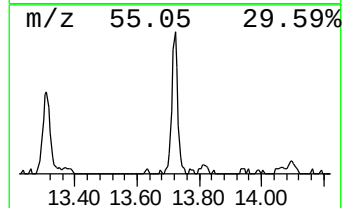
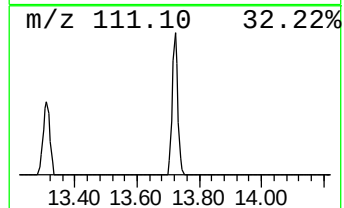
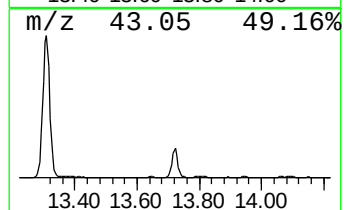
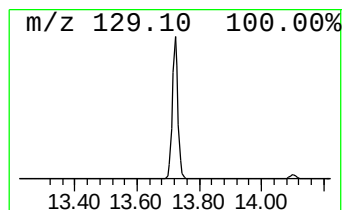
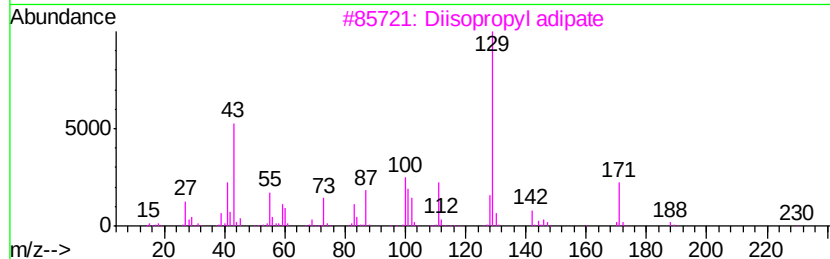
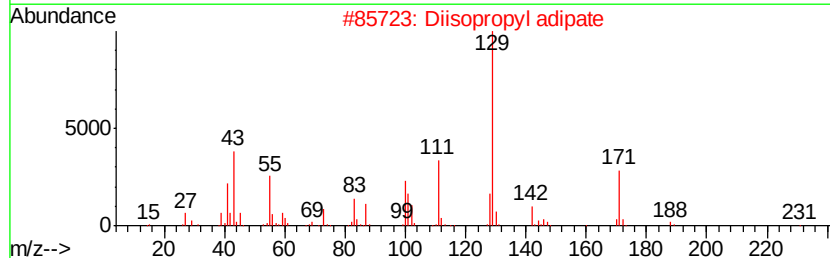
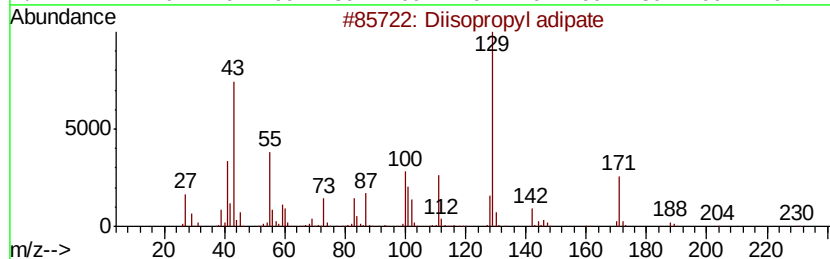
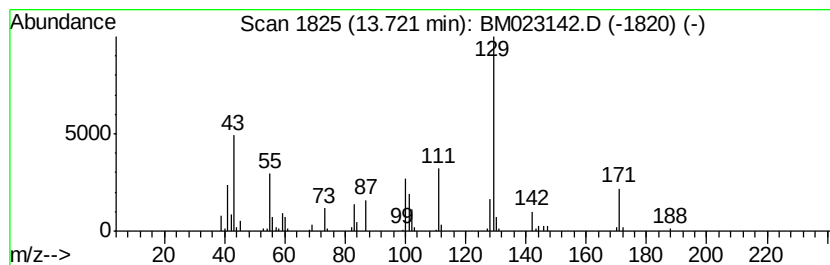
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Diisopropyl adipate Concentration Rank 6

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.72 | 2.45 ng/ul | 203109 | Acenaphthene-d10 | 14.40 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Diisopropyl adipate                 | 230 | C12H22O4 | 006938-94-9  | 91   |
| 2    |    | Diisopropyl adipate                 | 230 | C12H22O4 | 006938-94-9  | 91   |
| 3    |    | Diisopropyl adipate                 | 230 | C12H22O4 | 006938-94-9  | 91   |
| 4    |    | Adipic acid, isobutyl 2-propyl e... | 244 | C13H24O4 | 1000324-42-9 | 90   |
| 5    |    | Adipic acid, butyl 2-propyl ester   | 244 | C13H24O4 | 1000324-43-0 | 90   |



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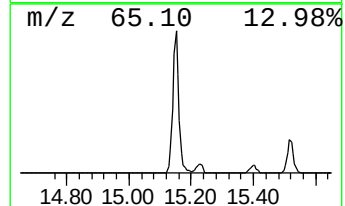
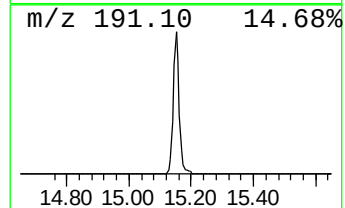
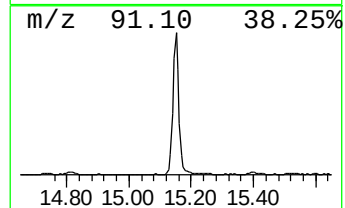
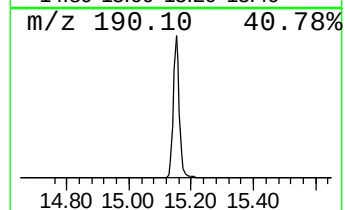
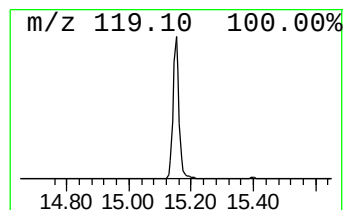
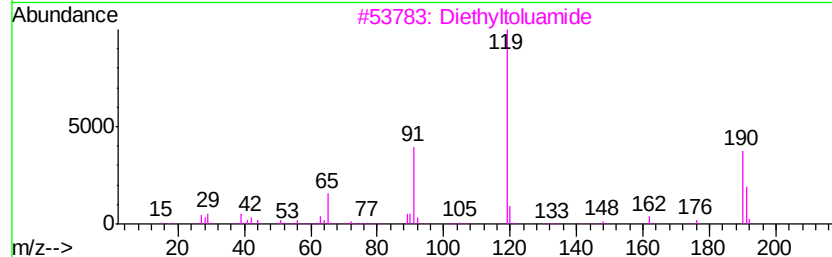
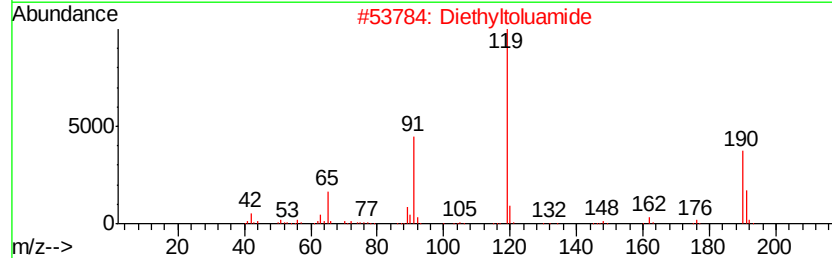
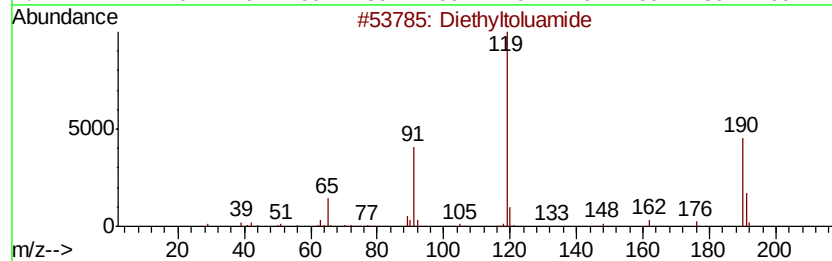
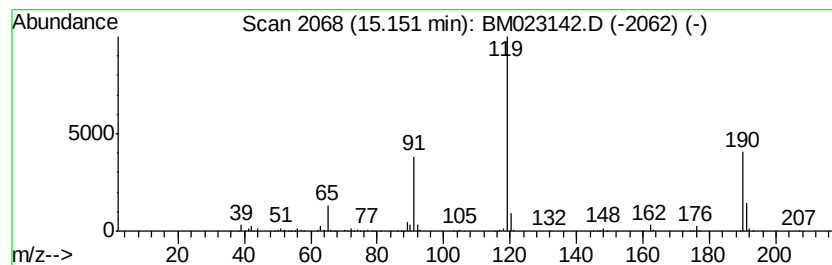
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TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Diethyltoluamide Concentration Rank 3

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.           |
|---------|-------------------------------------|--------------|------------------|----------------|
| 15.15   | 8.25 ng/ul                          | 682741       | Acenaphthene-d10 | 14.40          |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual      |
| 1       | Diethyltoluamide                    | 191          | C12H17NO         | 000134-62-3 97 |
| 2       | Diethyltoluamide                    | 191          | C12H17NO         | 000134-62-3 94 |
| 3       | Diethyltoluamide                    | 191          | C12H17NO         | 000134-62-3 60 |
| 4       | Benzamide, N-(1,1-dimethylethyl)... | 191          | C12H17NO         | 042498-32-8 58 |
| 5       | N-[1-(Dimethylamino)-2,2,2-trich... | 308          | C12H15Cl3N2O     | 300814-41-9 53 |



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Quant Title : SVOA CALIBRATION

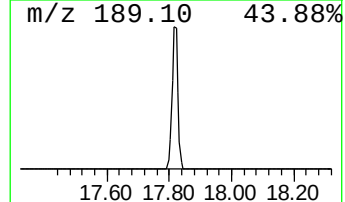
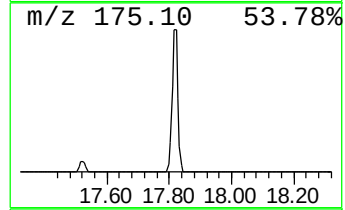
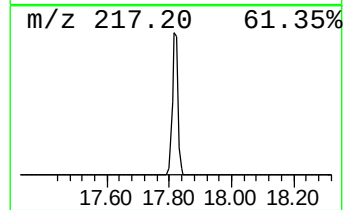
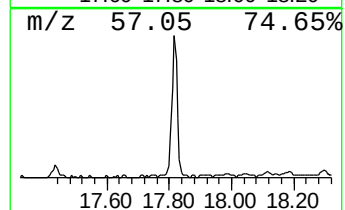
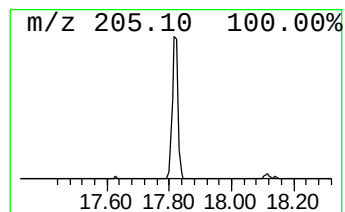
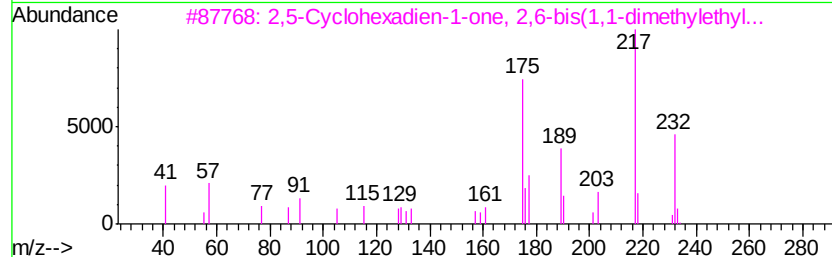
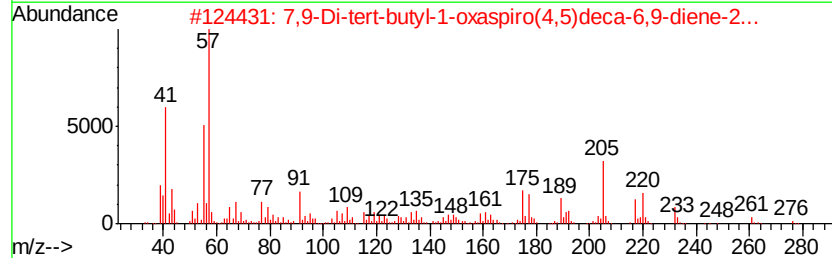
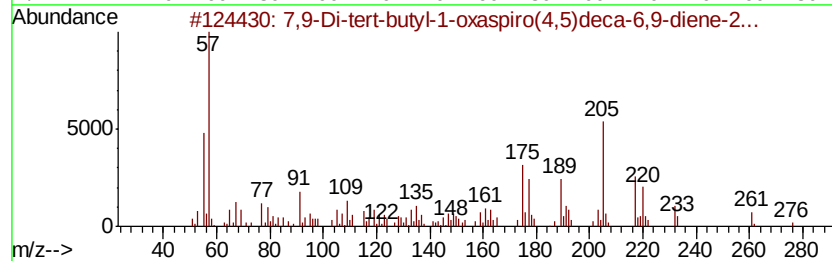
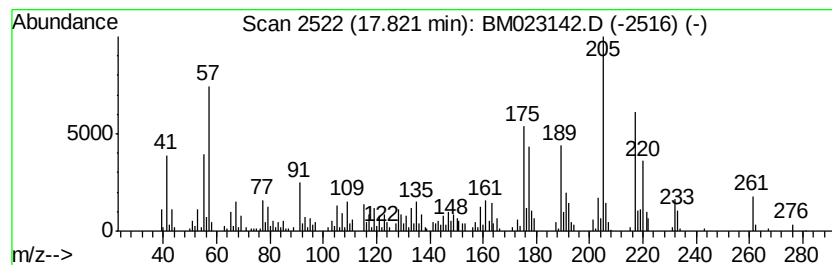
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 4

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.82 | 4.76 ng/ul | 417798 | Phenanthrene-d10 | 17.14 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 7,9-Di-tert-butyl-1-oxaspiro(4,5... | 276 | C17H24O3     | 082304-66-3 | 99      |      |      |
| 2    | 7,9-Di-tert-butyl-1-oxaspiro(4,5... | 276 | C17H24O3     | 082304-66-3 | 98      |      |      |
| 3    | 2,5-Cyclohexadien-1-one, 2,6-bis... | 232 | C16H24O      | 006738-27-8 | 49      |      |      |
| 4    | Ethanone, 1-(5,6,7,8-tetrahydro...  | 220 | C14H20O2     | 071596-88-8 | 45      |      |      |
| 5    | 2,5-di-tert-Butyl-1,4-benzoquinone  | 220 | C14H20O2     | 002460-77-7 | 44      |      |      |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101119\  
Data File : BM023142.D  
Acq On : 11 Oct 2019 22:32  
Operator : JU  
Sample : K5262-15  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BFBB6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM092719MA.M  
Quant Title : SVOA CALIBRATION

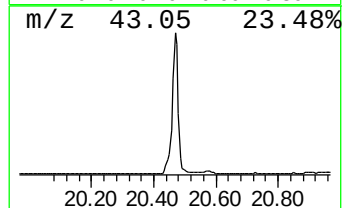
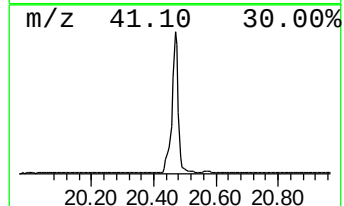
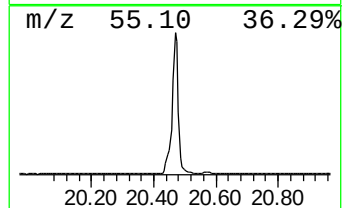
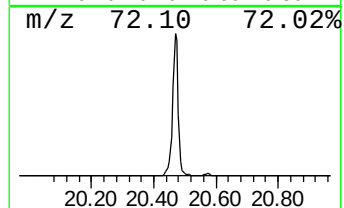
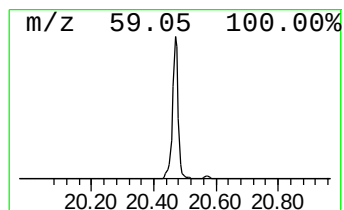
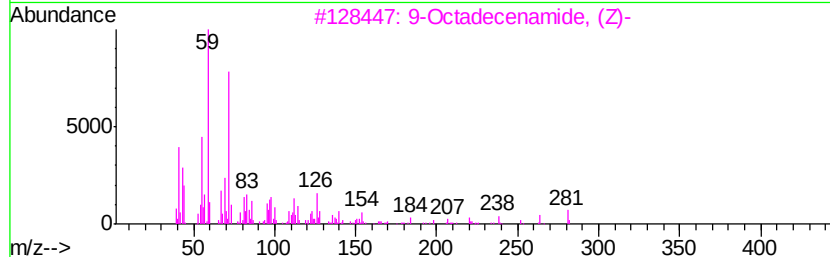
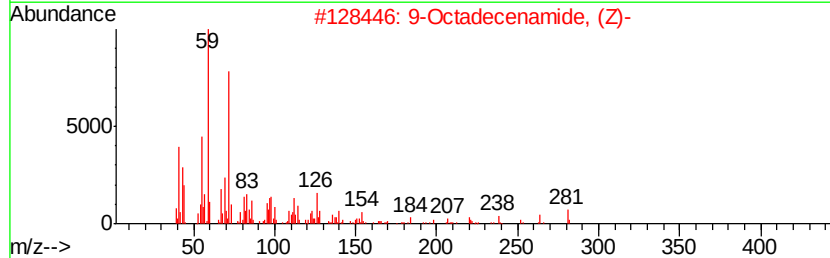
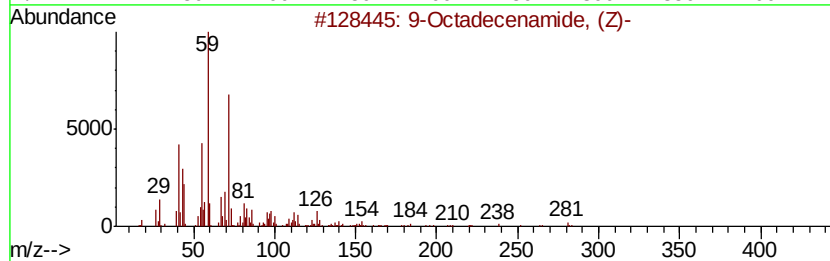
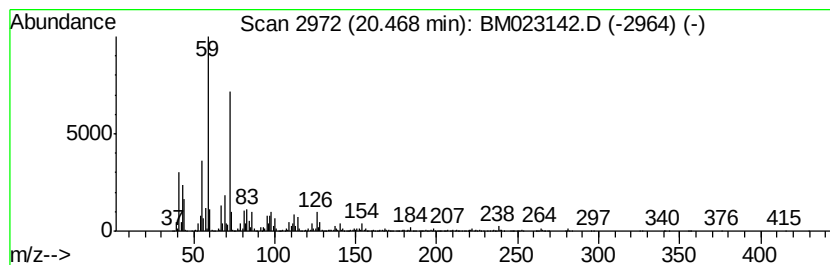
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 9-Octadecenamide, (Z)- Concentration Rank 1

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 20.47 | 79.84 ng/ul | 8452600 | Chrysene-d12     | 21.33 |

| Hit# | of                     | 5 | Tentative ID | MW  | MolForm  | CAS#        | Qual |
|------|------------------------|---|--------------|-----|----------|-------------|------|
| 1    | 9-Octadecenamide, (Z)- |   |              | 281 | C18H35NO | 000301-02-0 | 99   |
| 2    | 9-Octadecenamide, (Z)- |   |              | 281 | C18H35NO | 000301-02-0 | 95   |
| 3    | 9-Octadecenamide, (Z)- |   |              | 281 | C18H35NO | 000301-02-0 | 93   |
| 4    | 9-Octadecenamide, (Z)- |   |              | 281 | C18H35NO | 000301-02-0 | 87   |
| 5    | 9-Octadecenamide, (Z)- |   |              | 281 | C18H35NO | 000301-02-0 | 83   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101119\  
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Acq On : 11 Oct 2019 22:32  
Operator : JU  
Sample : K5262-15  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BFBB6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM092719MA.M  
Quant Title : SVOA CALIBRATION

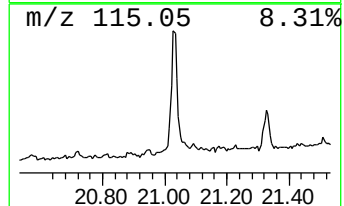
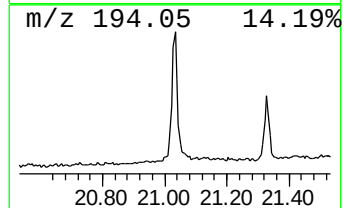
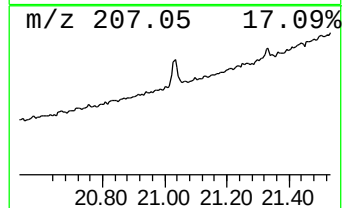
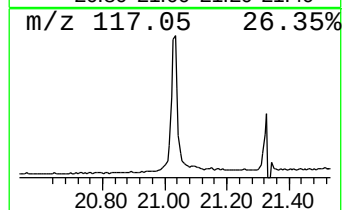
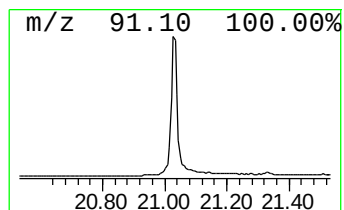
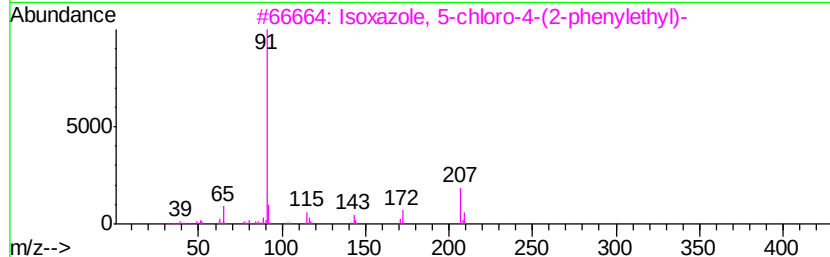
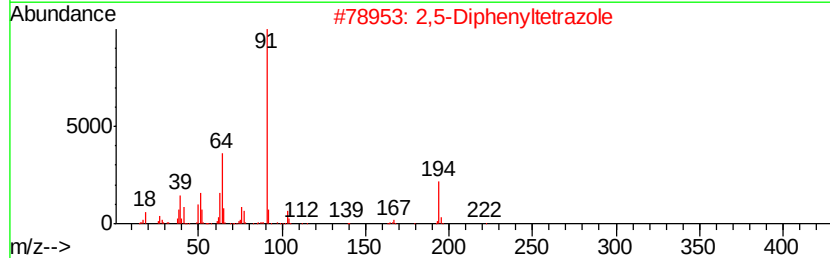
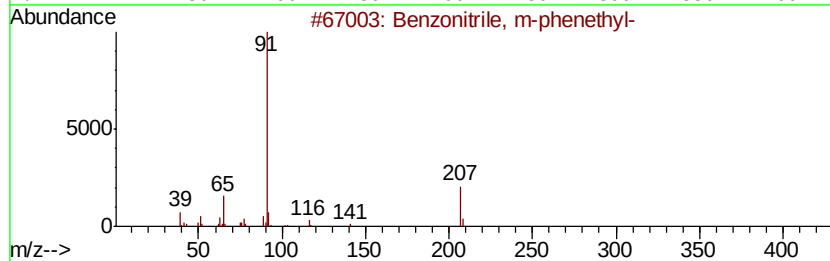
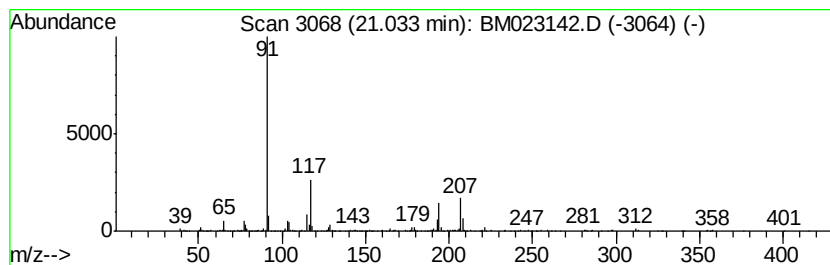
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 unknown-01 Concentration Rank 5

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 21.03 | 3.30 ng/ul | 349530 | Chrysene-d12     | 21.33 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | Benzonitrile, m-phenethyl-          | 207 | C15H13N     | 034176-91-5  | 27   |
| 2    |    | 2,5-Diphenyltetrazole               | 222 | C13H10N4    | 018039-45-7  | 25   |
| 3    |    | Isoxazole, 5-chloro-4-(2-phenyle... | 207 | C11H10ClNO  | 1000362-09-0 | 16   |
| 4    |    | 7-Benzyloxy-3-phenyl-4H-chromen-... | 328 | C22H16O3    | 004253-05-8  | 11   |
| 5    |    | Alpha-phenyl-alpha-tropylacetal...  | 378 | C22H22N2O2S | 022532-16-7  | 10   |



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Instrument :  
BNA\_M  
ClientSampleId :  
BFBB6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM092719MA.M  
Quant Title : SVOA CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| 2,4,7,9-Tetrameth... | 13.31 | 8.3     | ng/ul | 688285   | 3                     | 14.40 | 1656020 | 20.0 |
| Diisopropyl adipate  | 13.72 | 2.5     | ng/ul | 203109   | 3                     | 14.40 | 1656020 | 20.0 |
| Diethyltoluamide     | 15.15 | 8.3     | ng/ul | 682741   | 3                     | 14.40 | 1656020 | 20.0 |
| 7,9-Di-tert-butyl... | 17.82 | 4.8     | ng/ul | 417798   | 4                     | 17.14 | 1754620 | 20.0 |
| 9-Octadecenamide,... | 20.47 | 79.8    | ng/ul | 8452600  | 5                     | 21.33 | 2117280 | 20.0 |
| unknown-01           | 21.03 | 3.3     | ng/ul | 349530   | 5                     | 21.33 | 2117280 | 20.0 |