

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012221\  
 Data File : BN013455.D  
 Acq On : 21 Jan 2021 17:55  
 Operator : CG/JU  
 Sample : M1050-02  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 X2X34

## 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\_N\METHODS\SFAM-EPA-BN012221.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         | 3.187       | 13            | 17          | 20           | rBV      | 141405         | 181009        | 0.14%           | 0.043%        |
| 2         | 3.217       | 20            | 22          | 31           | rVB      | 242404         | 339286        | 0.27%           | 0.080%        |
| 3         | 3.569       | 79            | 82          | 106          | rBV      | 1690955        | 2424290       | 1.92%           | 0.570%        |
| 4         | 3.899       | 130           | 138         | 145          | rBV3     | 45527391       | 126457827     | 100.00%         | 29.724%       |
| 5         | 4.764       | 281           | 285         | 295          | rBV      | 872640         | 1196200       | 0.95%           | 0.281%        |
| 6         | 6.740       | 614           | 621         | 640          | rBV2     | 4421072        | 9791465       | 7.74%           | 2.301%        |
| 7         | 6.928       | 647           | 653         | 664          | rBV      | 1623397        | 2562797       | 2.03%           | 0.602%        |
| 8         | 7.122       | 679           | 686         | 689          | rBV      | 2336271        | 3799836       | 3.00%           | 0.893%        |
| 9         | 7.152       | 689           | 691         | 700          | rVB      | 1523740        | 1988913       | 1.57%           | 0.467%        |
| 10        | 7.581       | 757           | 764         | 774          | rVB      | 1789076        | 2808131       | 2.22%           | 0.660%        |
| 11        | 8.116       | 845           | 855         | 861          | rBV      | 1297748        | 3180058       | 2.51%           | 0.747%        |
| 12        | 8.163       | 861           | 863         | 870          | rVB      | 145624         | 183089        | 0.14%           | 0.043%        |
| 13        | 8.287       | 877           | 884         | 894          | rBV      | 2376729        | 3806863       | 3.01%           | 0.895%        |
| 14        | 8.387       | 894           | 901         | 910          | rBV      | 3563738        | 5541395       | 4.38%           | 1.302%        |
| 15        | 8.652       | 939           | 946         | 952          | rBV      | 1026023        | 1679975       | 1.33%           | 0.395%        |
| 16        | 8.728       | 952           | 959         | 962          | rVV      | 1882087        | 3070084       | 2.43%           | 0.722%        |
| 17        | 8.769       | 962           | 966         | 975          | rVB      | 2247384        | 3449178       | 2.73%           | 0.811%        |
| 18        | 9.040       | 1005          | 1012        | 1022         | rBV      | 1195894        | 1727852       | 1.37%           | 0.406%        |
| 19        | 9.440       | 1066          | 1080        | 1088         | rBV      | 1560570        | 2525035       | 2.00%           | 0.594%        |
| 20        | 9.552       | 1092          | 1099        | 1105         | rVB      | 110272         | 170278        | 0.13%           | 0.040%        |
| 21        | 9.775       | 1129          | 1137        | 1146         | rBV      | 1688269        | 2474370       | 1.96%           | 0.582%        |
| 22        | 9.969       | 1163          | 1170        | 1181         | rBV      | 2518388        | 4072747       | 3.22%           | 0.957%        |
| 23        | 10.346      | 1227          | 1234        | 1239         | rBV      | 2309649        | 4022059       | 3.18%           | 0.945%        |
| 24        | 10.399      | 1239          | 1243        | 1250         | rVV      | 2763738        | 4330795       | 3.42%           | 1.018%        |
| 25        | 10.481      | 1250          | 1257        | 1272         | rVB      | 1663828        | 2780034       | 2.20%           | 0.653%        |
| 26        | 11.063      | 1351          | 1356        | 1362         | rVB      | 156827         | 235031        | 0.19%           | 0.055%        |
| 27        | 11.263      | 1381          | 1390        | 1394         | rBV      | 629455         | 1407593       | 1.11%           | 0.331%        |
| 28        | 11.646      | 1447          | 1455        | 1466         | rBV      | 2833721        | 4643193       | 3.67%           | 1.091%        |
| 29        | 11.887      | 1488          | 1496        | 1502         | rBV      | 1176792        | 1852891       | 1.47%           | 0.436%        |
| 30        | 12.240      | 1547          | 1556        | 1565         | rVB      | 5300756        | 8106643       | 6.41%           | 1.905%        |
| 31        | 12.634      | 1614          | 1623        | 1629         | rBV      | 1362427        | 2055589       | 1.63%           | 0.483%        |
| 32        | 12.704      | 1629          | 1635        | 1645         | rVB      | 1357921        | 2064641       | 1.63%           | 0.485%        |
| 33        | 13.087      | 1687          | 1700        | 1708         | rBV      | 5961475        | 9336627       | 7.38%           | 2.195%        |
| 34        | 13.422      | 1748          | 1757        | 1762         | rBV2     | 106904         | 175483        | 0.14%           | 0.041%        |

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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: 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\_N\METHODS\SFAM-EPA-BN012221.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |          |       |        |
|----|--------|------|------|------|------|---------|----------|-------|--------|
| 35 | 13.634 | 1787 | 1793 | 1797 | rBV  | 3639523 | 5065983  | 4.01% | 1.191% |
| 36 | 13.681 | 1797 | 1801 | 1813 | rVB  | 3191302 | 4343772  | 3.43% | 1.021% |
| 37 | 13.910 | 1830 | 1840 | 1842 | rBV  | 4482609 | 6919384  | 5.47% | 1.626% |
| 38 | 13.940 | 1842 | 1845 | 1854 | rVB  | 3783749 | 5120978  | 4.05% | 1.204% |
| 39 | 14.216 | 1884 | 1892 | 1898 | rBV  | 3035058 | 4528275  | 3.58% | 1.064% |
| 40 | 14.281 | 1898 | 1903 | 1914 | rVB  | 2734735 | 3859357  | 3.05% | 0.907% |
| 41 | 14.439 | 1921 | 1930 | 1950 | rBV2 | 3710362 | 6691523  | 5.29% | 1.573% |
| 42 | 14.622 | 1950 | 1961 | 1968 | rVB  | 2881212 | 4154384  | 3.29% | 0.976% |
| 43 | 14.734 | 1973 | 1980 | 1987 | rBV  | 3797714 | 4956561  | 3.92% | 1.165% |
| 44 | 14.845 | 1992 | 1999 | 2011 | rBV  | 1305181 | 1844429  | 1.46% | 0.434% |
| 45 | 15.063 | 2030 | 2036 | 2048 | rVB  | 6018848 | 7702226  | 6.09% | 1.810% |
| 46 | 15.216 | 2055 | 2062 | 2067 | rBV  | 5249602 | 7503189  | 5.93% | 1.764% |
| 47 | 15.275 | 2067 | 2072 | 2077 | rVV  | 8087099 | 11195475 | 8.85% | 2.631% |
| 48 | 15.334 | 2077 | 2082 | 2095 | rVV  | 1412514 | 1879960  | 1.49% | 0.442% |
| 49 | 16.169 | 2217 | 2224 | 2236 | rVB  | 6206907 | 7843094  | 6.20% | 1.844% |
| 50 | 16.275 | 2236 | 2242 | 2249 | rBV  | 1837663 | 2534173  | 2.00% | 0.596% |
| 51 | 16.622 | 2294 | 2301 | 2303 | rBV  | 2735986 | 3933968  | 3.11% | 0.925% |
| 52 | 16.645 | 2303 | 2305 | 2315 | rVB  | 3405309 | 4245226  | 3.36% | 0.998% |
| 53 | 16.969 | 2354 | 2360 | 2363 | rBV  | 3197341 | 4505150  | 3.56% | 1.059% |
| 54 | 17.010 | 2363 | 2367 | 2372 | rVV  | 6341343 | 8779709  | 6.94% | 2.064% |
| 55 | 17.069 | 2372 | 2377 | 2389 | rVB  | 5213670 | 7148023  | 5.65% | 1.680% |
| 56 | 17.451 | 2435 | 2442 | 2449 | rBV  | 4498473 | 5882068  | 4.65% | 1.383% |
| 57 | 17.592 | 2461 | 2466 | 2469 | rBV  | 203705  | 265589   | 0.21% | 0.062% |
| 58 | 17.957 | 2521 | 2528 | 2533 | rBV  | 3639827 | 4193284  | 3.32% | 0.986% |
| 59 | 19.033 | 2700 | 2711 | 2719 | rBV  | 4638997 | 6369328  | 5.04% | 1.497% |
| 60 | 19.369 | 2762 | 2768 | 2778 | rBV  | 5474654 | 7520649  | 5.95% | 1.768% |
| 61 | 20.327 | 2926 | 2931 | 2937 | rBV  | 3043624 | 3304099  | 2.61% | 0.777% |
| 62 | 21.116 | 3060 | 3065 | 3068 | rBV  | 2255027 | 2475986  | 1.96% | 0.582% |
| 63 | 21.163 | 3068 | 3073 | 3078 | rVV2 | 5266069 | 10227791 | 8.09% | 2.404% |
| 64 | 21.210 | 3078 | 3081 | 3087 | rVB  | 3966806 | 4588961  | 3.63% | 1.079% |
| 65 | 21.980 | 3207 | 3212 | 3219 | rBV  | 6417503 | 7059131  | 5.58% | 1.659% |
| 66 | 22.233 | 3252 | 3255 | 3259 | rVB2 | 166234  | 203624   | 0.16% | 0.048% |
| 67 | 22.704 | 3329 | 3335 | 3338 | rBV  | 2462183 | 3833171  | 3.03% | 0.901% |
| 68 | 22.745 | 3338 | 3342 | 3353 | rVB  | 4641204 | 7670544  | 6.07% | 1.803% |
| 69 | 23.215 | 3415 | 3422 | 3429 | rBV  | 3871419 | 7049193  | 5.57% | 1.657% |
| 70 | 23.286 | 3430 | 3434 | 3438 | rVV  | 429011  | 666998   | 0.53% | 0.157% |
| 71 | 23.357 | 3439 | 3446 | 3455 | rVB2 | 2363105 | 4255606  | 3.37% | 1.000% |

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## Integration Parameters: LSCINT.P

Integrator: RTE  
Smoothing : OFF Filtering: 5  
Sampling : 1 Min Area: 0.1 % of largest Peak  
Start Thrs: 0.2 Max Peaks: 100  
Stop Thrs : 0 Peak Location: TOP

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN012221.M  
Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |       |        |
|----|--------|------|------|------|------|---------|---------|-------|--------|
| 72 | 23.915 | 3536 | 3541 | 3548 | rVB  | 427057  | 746122  | 0.59% | 0.175% |
| 73 | 24.645 | 3659 | 3665 | 3676 | rVB  | 380091  | 804989  | 0.64% | 0.189% |
| 74 | 25.539 | 3804 | 3817 | 3830 | rVB3 | 2877079 | 9132244 | 7.22% | 2.147% |

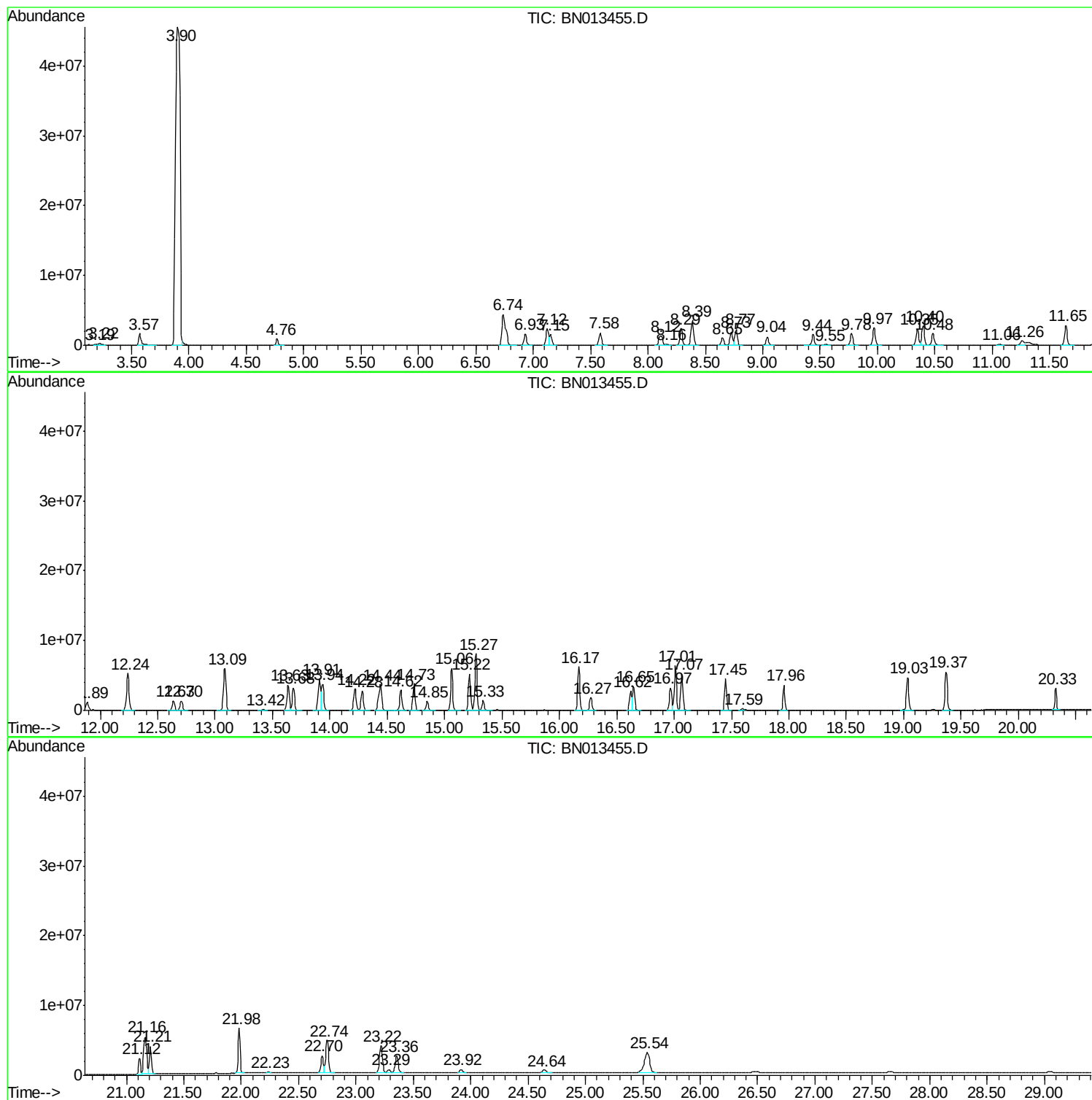
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN012221.M  
Quant Title : SVOA CALIBRATION

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



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ALS Vial : 12 Sample Multiplier: 1

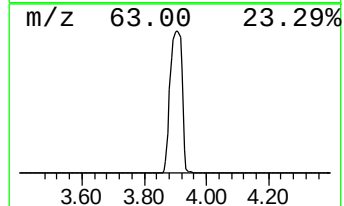
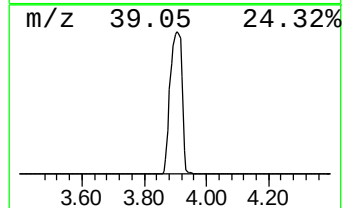
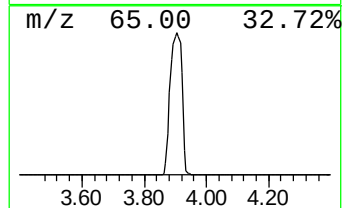
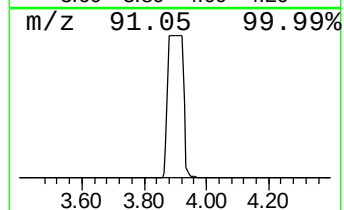
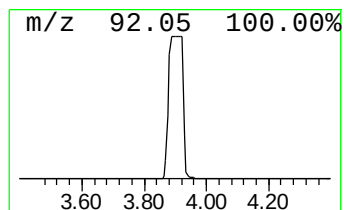
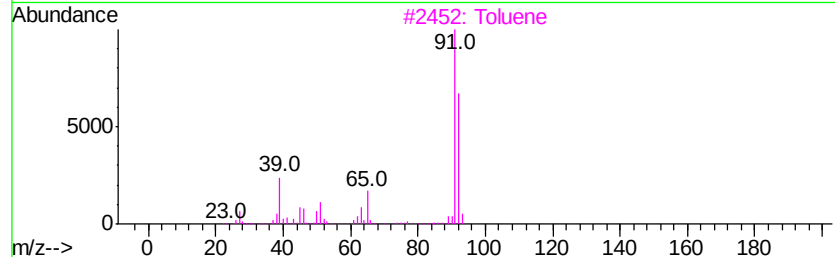
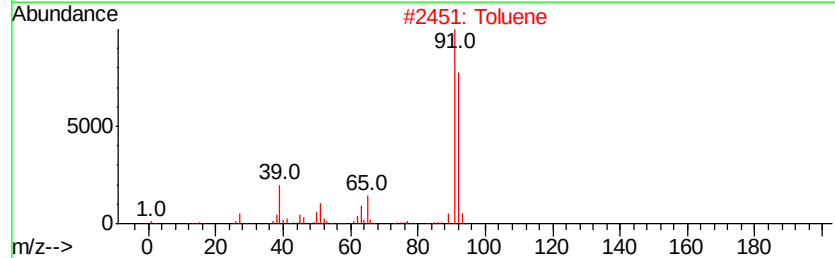
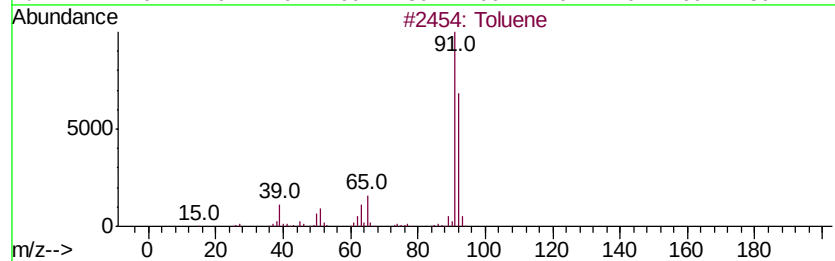
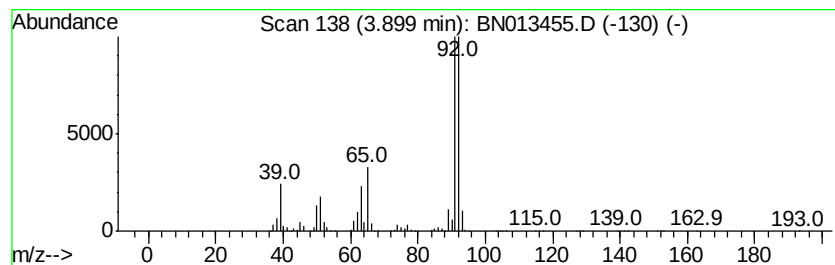
Instrument :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN012221.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 Toluene Concentration Rank 1

| R.T.    | EstConc                       | Area         | Relative to ISTD       | R.T.           |
|---------|-------------------------------|--------------|------------------------|----------------|
| 3.90    | 900.66 ng/ul                  | 126458000    | 1,4-Dichlorobenzene-d4 | 7.59           |
| Hit# of | 5                             | Tentative ID | MW MolForm             | CAS# Qual      |
| 1       | Toluene                       |              | 92 C7H8                | 000108-88-3 93 |
| 2       | Toluene                       |              | 92 C7H8                | 000108-88-3 91 |
| 3       | Toluene                       |              | 92 C7H8                | 000108-88-3 90 |
| 4       | Toluene                       |              | 92 C7H8                | 000108-88-3 87 |
| 5       | 1,5-Hexadien-3-yne, 2-methyl- |              | 92 C7H8                | 000820-54-2 80 |



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Quant Title : SVOA CALIBRATION

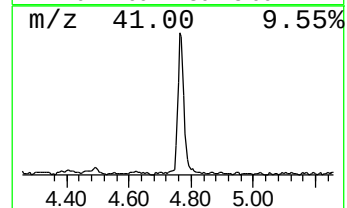
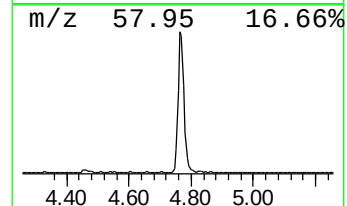
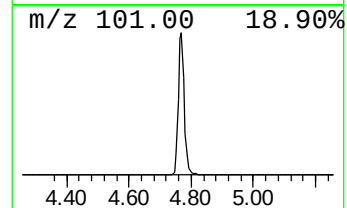
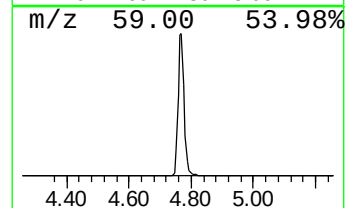
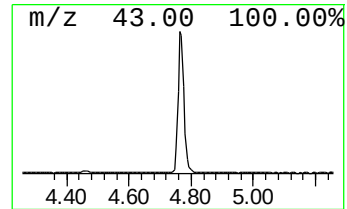
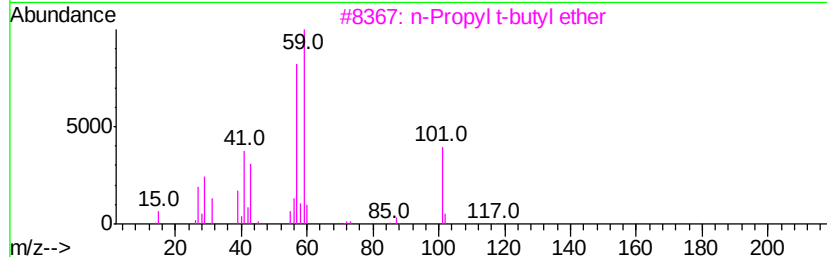
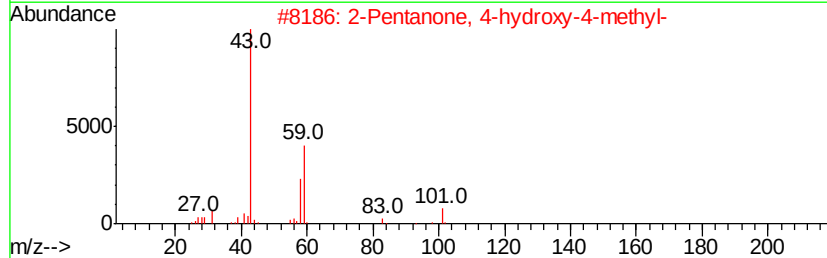
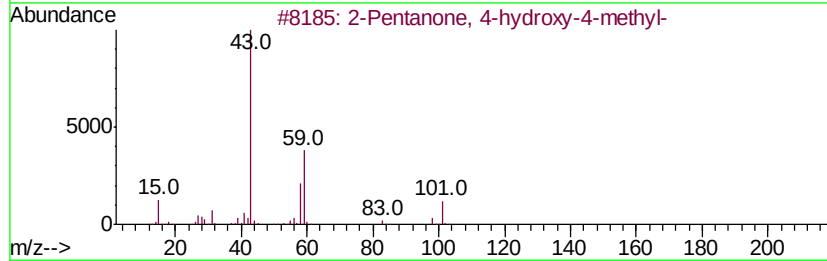
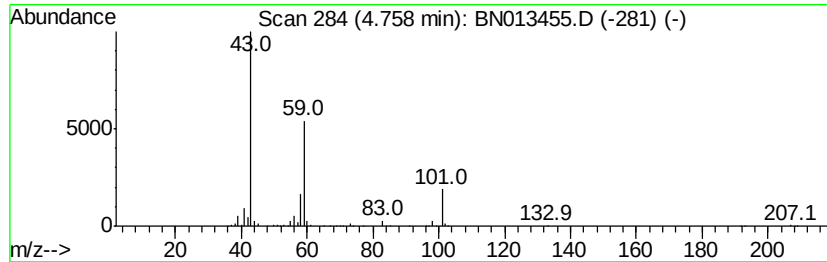
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

| R.T. | EstConc    | Area    | Relative to ISTD       | R.T. |
|------|------------|---------|------------------------|------|
| 4.76 | 8.52 ng/ul | 1196200 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2  | 50   |
| 2    |    | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2  | 40   |
| 3    |    | n-Propyl t-butyl ether              | 116 | C7H16O   | 1000334-81-9 | 33   |
| 4    |    | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2  | 32   |
| 5    |    | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9  | 23   |



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Quant Title : SVOA CALIBRATION

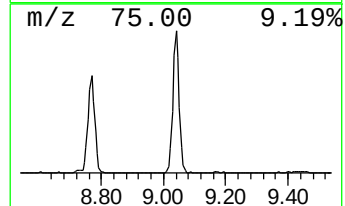
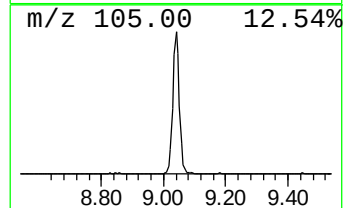
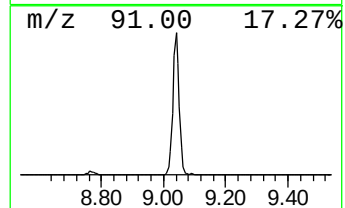
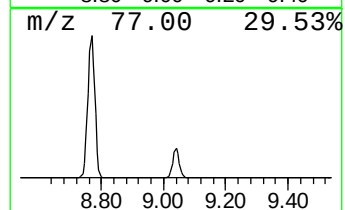
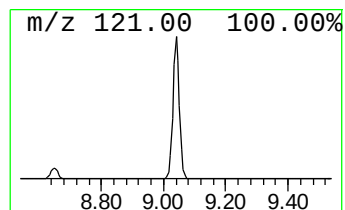
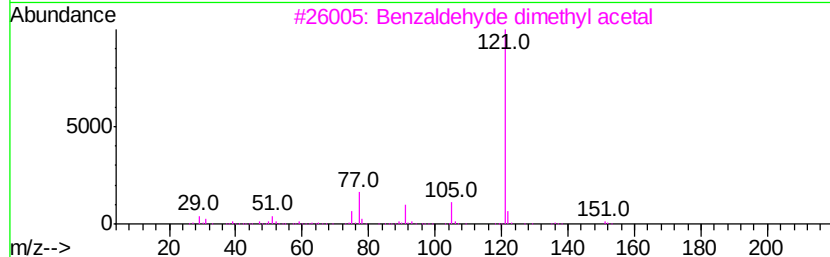
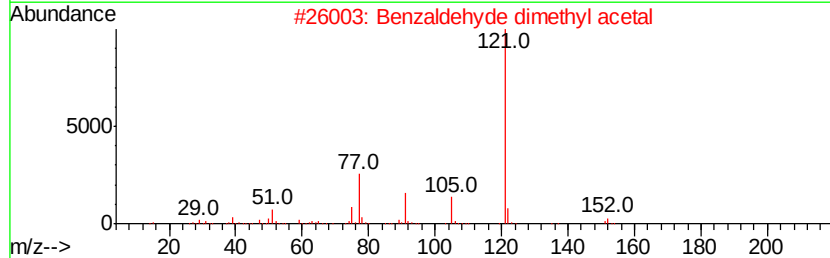
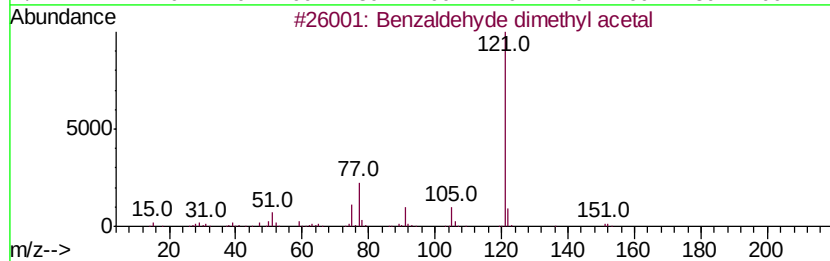
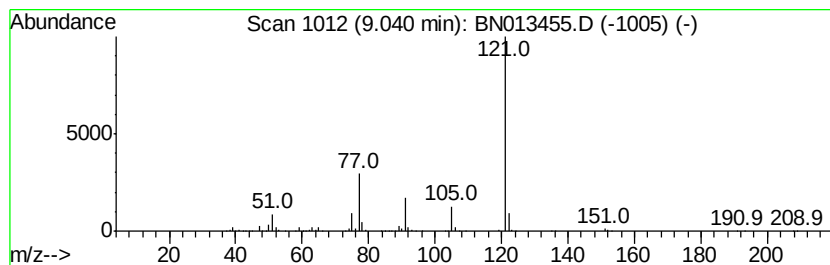
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 Benzaldehyde dimethyl acetal Concentration Rank 5

| R.T. | EstConc    | Area    | Relative to ISTD | R.T.  |
|------|------------|---------|------------------|-------|
| 9.04 | 8.59 ng/ul | 1727850 | Napthalene-d8    | 10.35 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Benzaldehyde dimethyl acetal        | 152 | C9H12O2  | 001125-88-8  | 90   |
| 2    |    | Benzaldehyde dimethyl acetal        | 152 | C9H12O2  | 001125-88-8  | 90   |
| 3    |    | Benzaldehyde dimethyl acetal        | 152 | C9H12O2  | 001125-88-8  | 86   |
| 4    |    | Benzeneacetic acid, .alpha.-meth... | 180 | C10H12O3 | 056143-21-6  | 78   |
| 5    |    | Benzene, (3-iodo-1-methoxypropyl)-  | 276 | C10H13IO | 1000327-31-9 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012221\  
Data File : BN013455.D  
Acq On : 21 Jan 2021 17:55  
Operator : CG/JU  
Sample : M1050-02  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

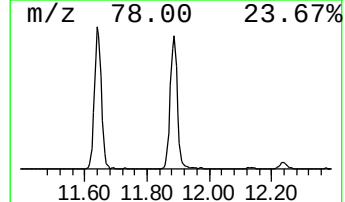
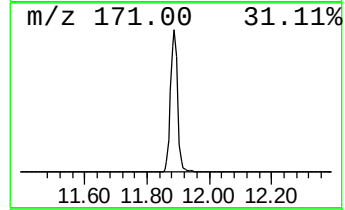
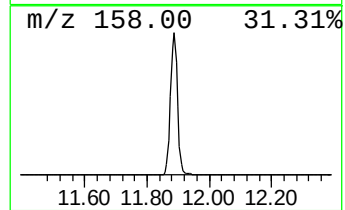
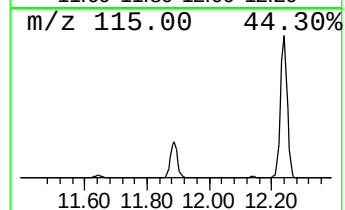
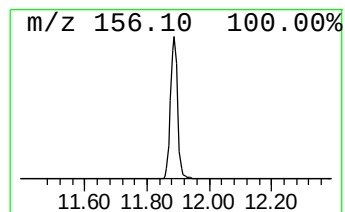
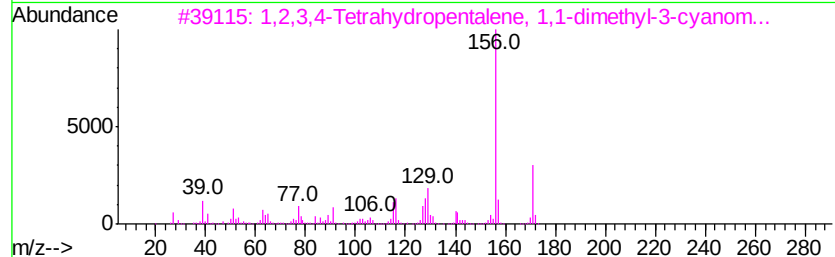
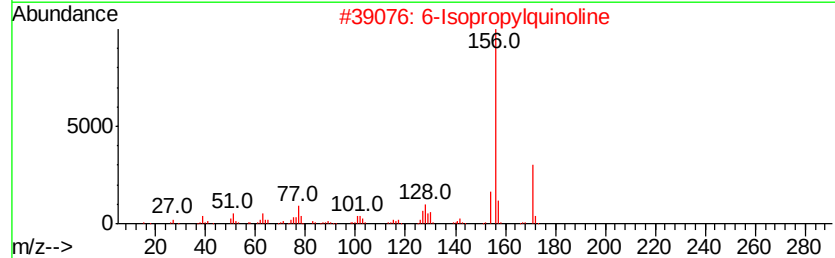
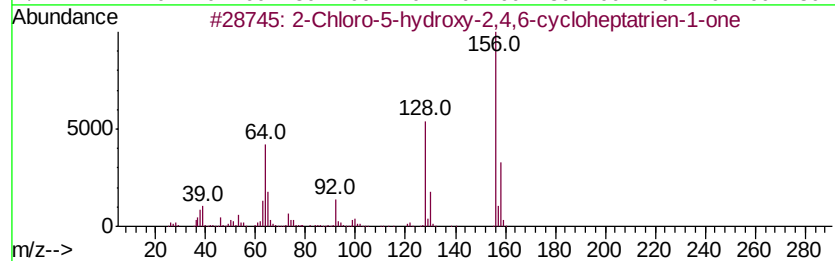
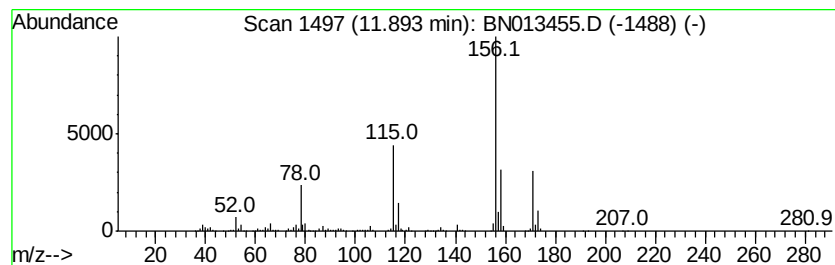
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ClientSampleId :  
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 unknown-01 Concentration Rank 4

| R.T.    | EstConc                             | Area           | Relative to ISTD | R.T.      |
|---------|-------------------------------------|----------------|------------------|-----------|
| 11.89   | 9.21 ng/ul                          | 1852890        | Napthalene-d8    | 10.35     |
| Hit# of | 5                                   | Tentative ID   | MW MolForm       | CAS# Qual |
| 1       | 2-Chloro-5-hydroxy-2,4,6-cyclohe... | 156 C7H5ClO2   | 007009-16-7      | 38        |
| 2       | 6-Isopropylquinoline                | 171 C12H13N    | 000135-79-5      | 37        |
| 3       | 1,2,3,4-Tetrahydropentalene, 1,1... | 171 C12H13N    | 1000217-18-6     | 37        |
| 4       | 3,5-Xylenol, 4-chloro-, acetate     | 198 C10H11ClO2 | 022012-58-4      | 32        |
| 5       | Bicyclo[3.2.0]hept-2-ene, 4-isop... | 171 C12H13N    | 1000217-15-6     | 32        |





Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012221\  
Data File : BN013455.D  
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Operator : CG/JU  
Sample : M1050-02  
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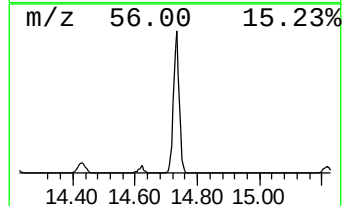
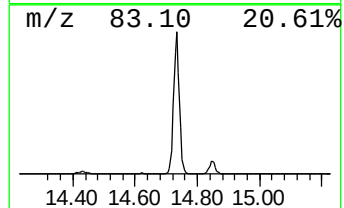
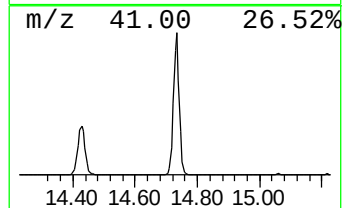
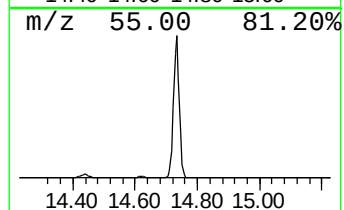
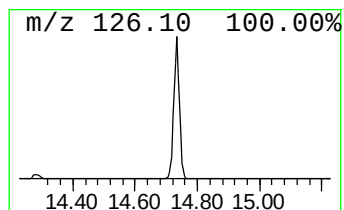
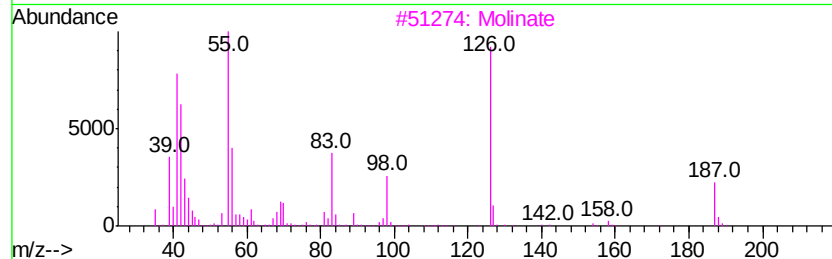
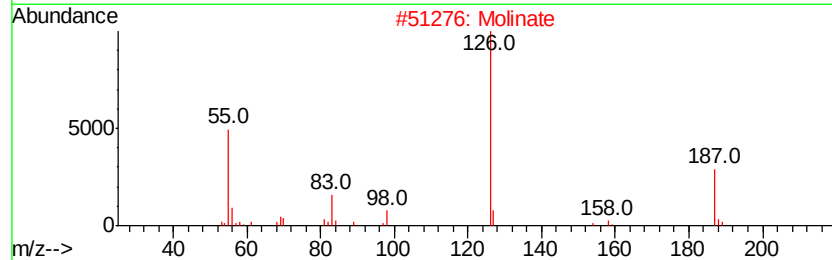
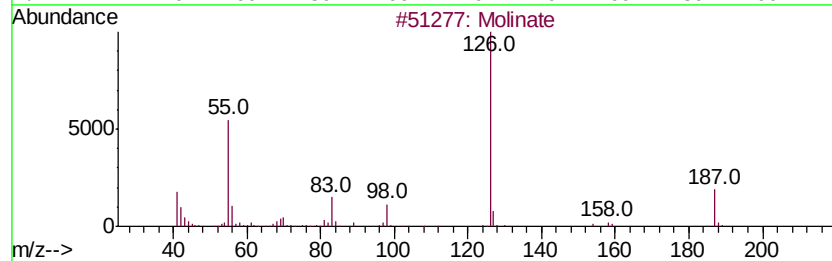
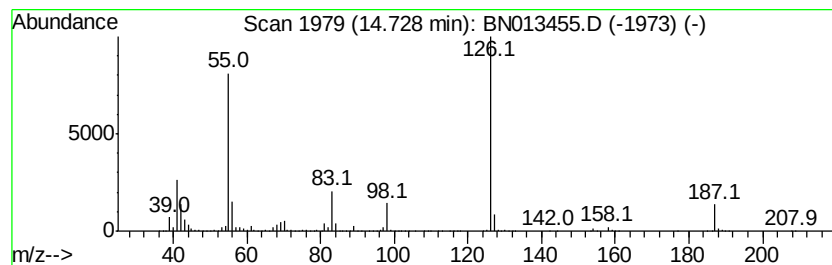
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 Molinate Concentration Rank 3

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.           |
|---------|-------------------------------------|--------------|------------------|----------------|
| 14.73   | 21.89 ng/ul                         | 4956560      | Acenaphthene-d10 | 14.22          |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual      |
| 1       | Molinate                            |              | 187 C9H17NOS     | 002212-67-1 97 |
| 2       | Molinate                            |              | 187 C9H17NOS     | 002212-67-1 96 |
| 3       | Molinate                            |              | 187 C9H17NOS     | 002212-67-1 64 |
| 4       | Thymine                             |              | 126 C5H6N2O2     | 000065-71-4 59 |
| 5       | 1,2,4-Cyclopentanetrione, 3-methyl- |              | 126 C6H6O3       | 004505-54-8 52 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012221\  
Data File : BN013455.D  
Acq On : 21 Jan 2021 17:55  
Operator : CG/JU  
Sample : M1050-02  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

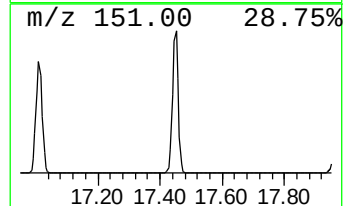
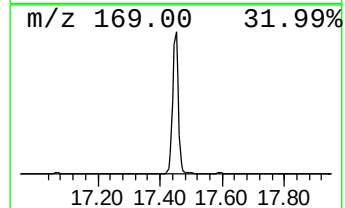
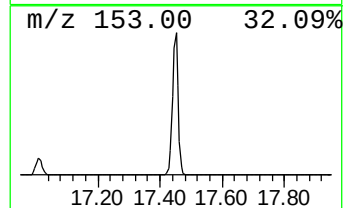
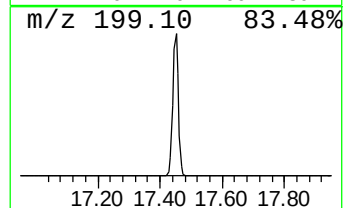
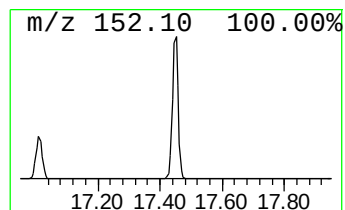
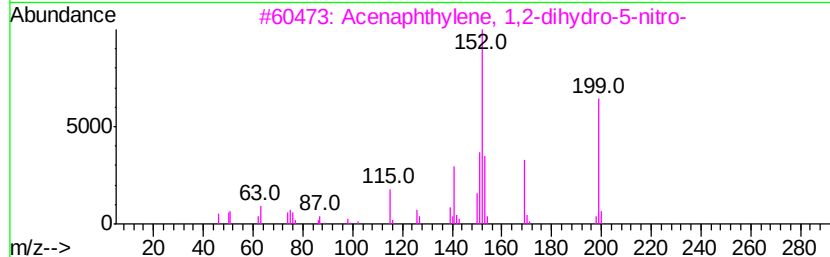
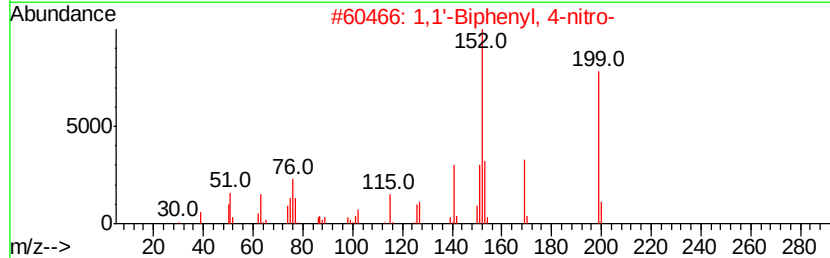
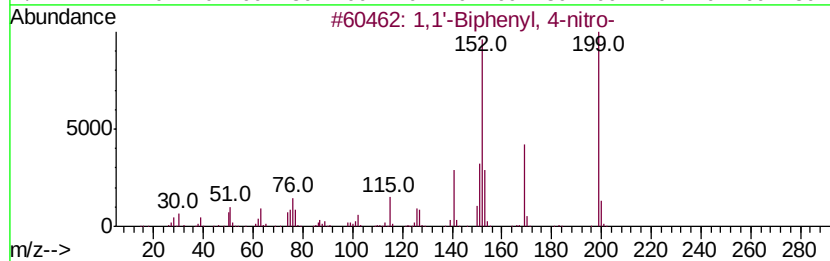
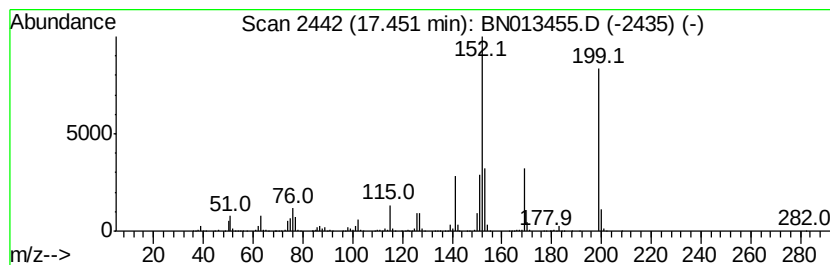
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN012221.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 6 1,1'-Biphenyl, 4-nitro- Concentration Rank 2

| R.T.    | EstConc     | Area                                | Relative to ISTD | R.T.     |             |      |
|---------|-------------|-------------------------------------|------------------|----------|-------------|------|
| 17.45   | 26.11 ng/ul | 5882070                             | Phenanthrene-d10 | 16.97    |             |      |
| Hit# of | 5           | Tentative ID                        | MW               | MolForm  | CAS#        | Qual |
| 1       |             | 1,1'-Biphenyl, 4-nitro-             | 199              | C12H9NO2 | 000092-93-3 | 98   |
| 2       |             | 1,1'-Biphenyl, 4-nitro-             | 199              | C12H9NO2 | 000092-93-3 | 98   |
| 3       |             | Acenaphthylene, 1,2-dihydro-5-ni... | 199              | C12H9NO2 | 000602-87-9 | 94   |
| 4       |             | 1,1'-Biphenyl, 3-nitro-             | 199              | C12H9NO2 | 002113-58-8 | 81   |
| 5       |             | 1,1'-Biphenyl, 3-nitro-             | 199              | C12H9NO2 | 002113-58-8 | 68   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012221\  
Data File : BN013455.D  
Acq On : 21 Jan 2021 17:55  
Operator : CG/JU  
Sample : M1050-02  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
X2X34

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Quant Title : SVOA CALIBRATION

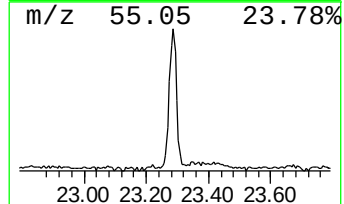
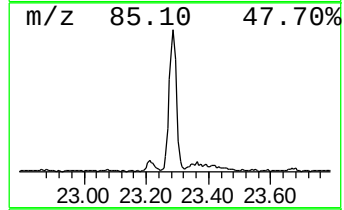
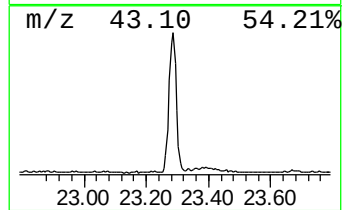
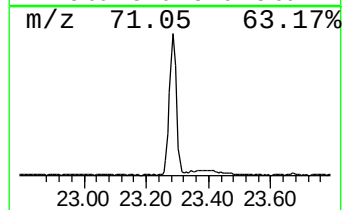
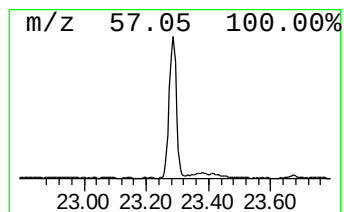
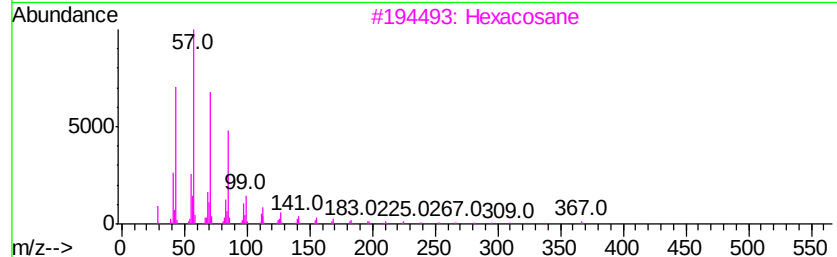
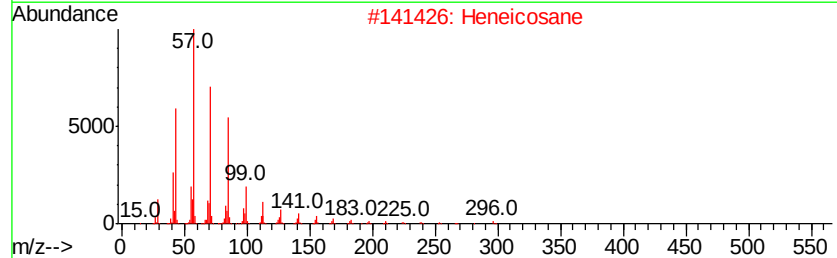
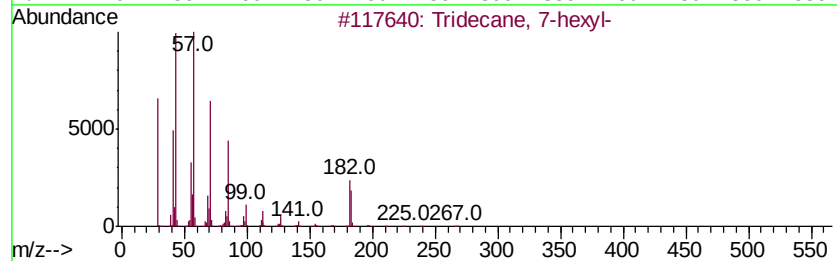
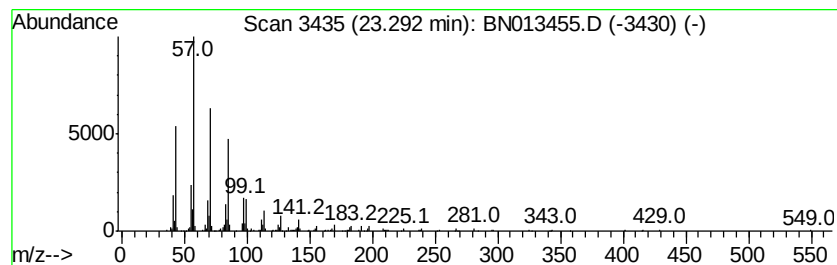
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 23.29 | 3.13 ng/ul | 666998 | Perylene-d12     | 23.36 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Tridecane, 7-hexyl-   | 268 | C19H40  | 007225-66-3 | 90   |
| 2    |    | Heneicosane           | 296 | C21H44  | 000629-94-7 | 90   |
| 3    |    | Hexacosane            | 366 | C26H54  | 000630-01-3 | 90   |
| 4    |    | Decane, 3,8-dimethyl- | 170 | C12H26  | 017312-55-9 | 87   |
| 5    |    | Octadecane            | 254 | C18H38  | 000593-45-3 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012221\  
Data File : BN013455.D  
Acq On : 21 Jan 2021 17:55  
Operator : CG/JU  
Sample : M1050-02  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
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Quant Title : SVOA CALIBRATION

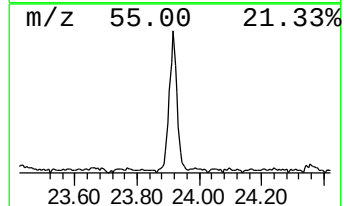
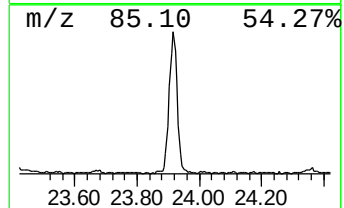
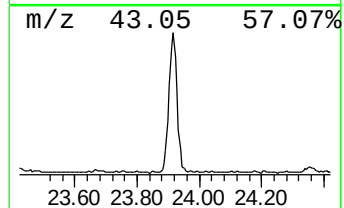
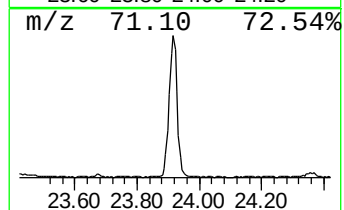
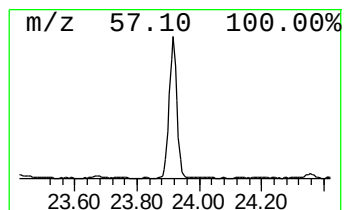
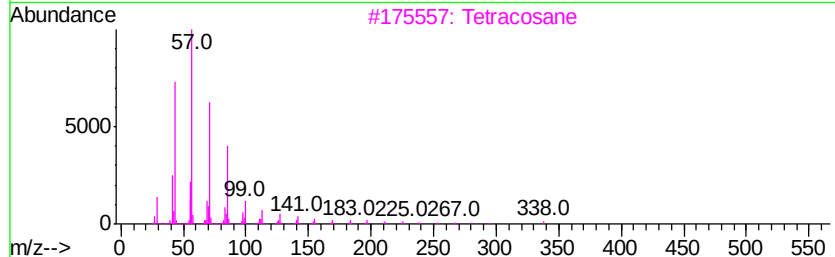
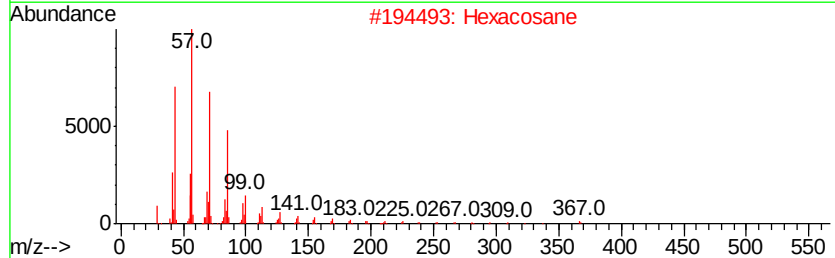
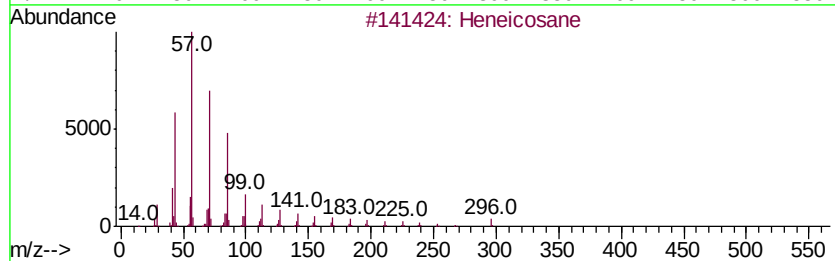
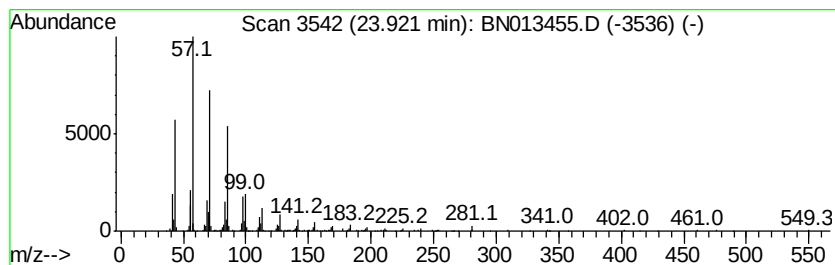
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 23.92 | 3.51 ng/ul | 746122 | Perylene-d12     | 23.36 |

| Hit# | of | 5 | Tentative ID       | MW  | MolForm | CAS#         | Qual |
|------|----|---|--------------------|-----|---------|--------------|------|
| 1    |    |   | Heneicosane        | 296 | C21H44  | 000629-94-7  | 98   |
| 2    |    |   | Hexacosane         | 366 | C26H54  | 000630-01-3  | 93   |
| 3    |    |   | Tetracosane        | 338 | C24H50  | 000646-31-1  | 90   |
| 4    |    |   | 2-methyloctacosane | 408 | C29H60  | 1000376-72-8 | 90   |
| 5    |    |   | Heneicosane        | 296 | C21H44  | 000629-94-7  | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012221\  
 Data File : BN013455.D  
 Acq On : 21 Jan 2021 17:55  
 Operator : CG/JU  
 Sample : M1050-02  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

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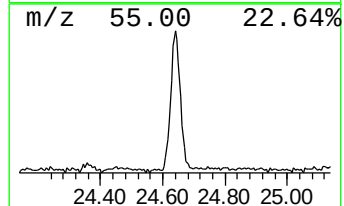
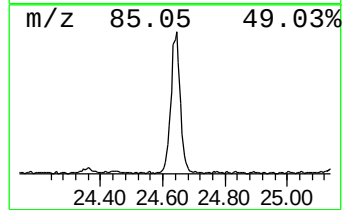
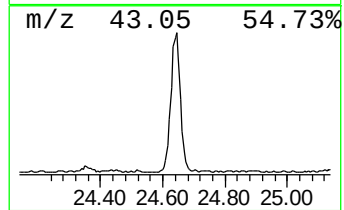
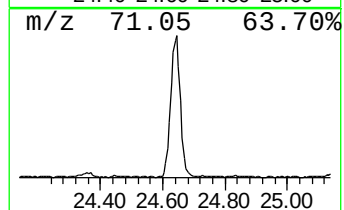
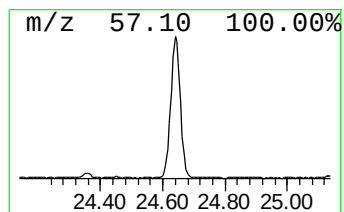
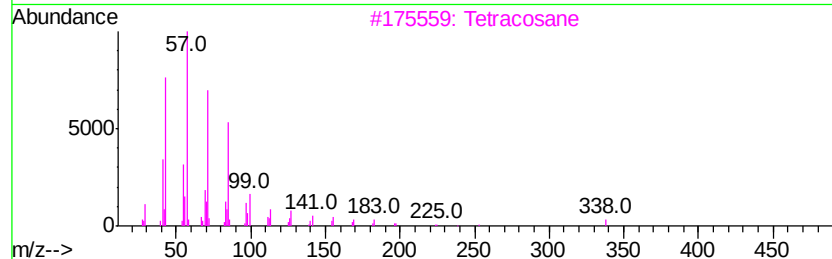
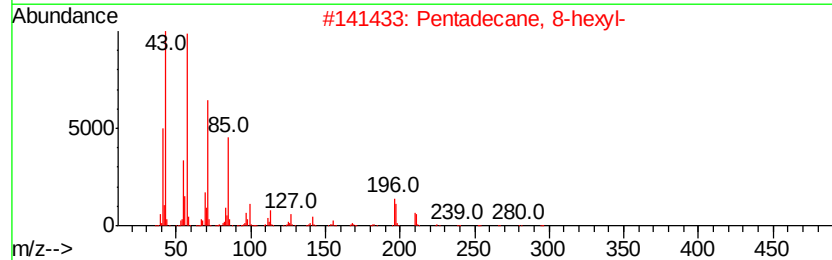
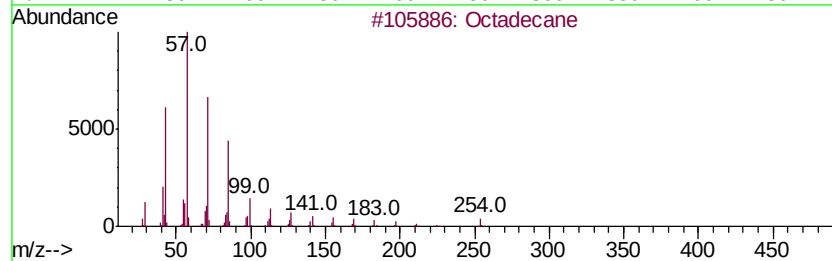
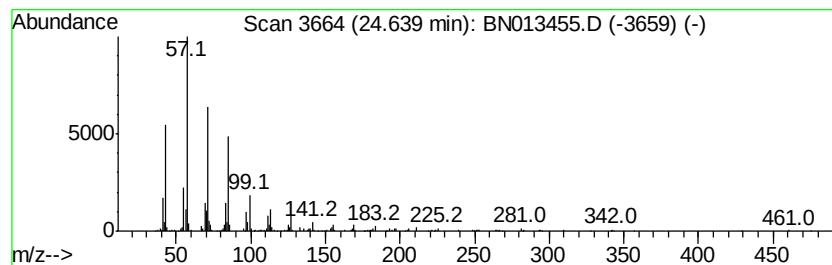
TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

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 Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 24.64 | 3.78 ng/ul | 804989 | Perylene-d12     | 23.36 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Octadecane            | 254 | C18H38  | 000593-45-3 | 96   |
| 2    |    | Pentadecane, 8-hexyl- | 296 | C21H44  | 013475-75-7 | 90   |
| 3    |    | Tetracosane           | 338 | C24H50  | 000646-31-1 | 90   |
| 4    |    | Eicosane              | 282 | C20H42  | 000112-95-8 | 90   |
| 5    |    | Hexacosane            | 366 | C26H54  | 000630-01-3 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN012221\  
Data File : BN013455.D  
Acq On : 21 Jan 2021 17:55  
Operator : CG/JU  
Sample : M1050-02  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
X2X34

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN012221.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 |
| Toluene              | 3.90  | 900.7   | ng/ul | 126458000 | 1                     | 7.59  | 2808130 | 20.0 |
| 2-Pentanone, 4-hy... | 4.76  | 8.5     | ng/ul | 1196200   | 1                     | 7.59  | 2808130 | 20.0 |
| Benzaldehyde dime... | 9.04  | 8.6     | ng/ul | 1727850   | 2                     | 10.35 | 4022060 | 20.0 |
| unknown-01           | 11.89 | 9.2     | ng/ul | 1852890   | 2                     | 10.35 | 4022060 | 20.0 |
| Molinate             | 14.73 | 21.9    | ng/ul | 4956560   | 3                     | 14.22 | 4528280 | 20.0 |
| 1,1'-Biphenyl, 4-... | 17.45 | 26.1    | ng/ul | 5882070   | 4                     | 16.97 | 4505150 | 20.0 |
| (DEL) Alkane: Str... | 23.29 | 3.1     | ng/ul | 666998    | 6                     | 23.36 | 4255610 | 20.0 |
| (DEL) Alkane: Str... | 23.92 | 3.5     | ng/ul | 746122    | 6                     | 23.36 | 4255610 | 20.0 |
| (DEL) Alkane: Str... | 24.64 | 3.8     | ng/ul | 804989    | 6                     | 23.36 | 4255610 | 20.0 |