

Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
 Data File : BG019343.D  
 Acq On : 24 Oct 2015 19:45  
 Operator : UM/NP  
 Sample : G4058-11  
 Misc :  
 ALS Vial : 33 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 E5LR9

## 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:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG102315.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         | 3.207       | 56            | 63          | 72           | rBV      | 70937          | 133381        | 7.09%           | 0.419%        |
| 2         | 3.290       | 72            | 77          | 91           | rVB      | 457316         | 667527        | 35.49%          | 2.094%        |
| 3         | 4.347       | 251           | 257         | 264          | rVB      | 105126         | 157096        | 8.35%           | 0.493%        |
| 4         | 7.373       | 765           | 772         | 782          | rVB      | 172040         | 294538        | 15.66%          | 0.924%        |
| 5         | 7.549       | 796           | 802         | 810          | rVV      | 117638         | 199481        | 10.61%          | 0.626%        |
| 6         | 7.767       | 831           | 839         | 847          | rVB      | 151730         | 264000        | 14.04%          | 0.828%        |
| 7         | 7.843       | 847           | 852         | 857          | rBV3     | 7877           | 15046         | 0.80%           | 0.047%        |
| 8         | 8.237       | 911           | 919         | 926          | rVV      | 160474         | 278042        | 14.78%          | 0.872%        |
| 9         | 8.930       | 1024          | 1037        | 1046         | rBV2     | 206421         | 360133        | 19.15%          | 1.130%        |
| 10        | 9.406       | 1111          | 1118        | 1125         | rVV      | 165006         | 306889        | 16.32%          | 0.963%        |
| 11        | 9.465       | 1125          | 1128        | 1136         | rVV5     | 12064          | 24218         | 1.29%           | 0.076%        |
| 12        | 10.135      | 1234          | 1242        | 1249         | rBV2     | 161965         | 293963        | 15.63%          | 0.922%        |
| 13        | 10.334      | 1270          | 1276        | 1281         | rBV3     | 11542          | 18424         | 0.98%           | 0.058%        |
| 14        | 10.675      | 1325          | 1334        | 1341         | rVB      | 230824         | 407808        | 21.68%          | 1.280%        |
| 15        | 11.016      | 1388          | 1392        | 1395         | rBV2     | 15090          | 24499         | 1.30%           | 0.077%        |
| 16        | 11.069      | 1395          | 1401        | 1406         | rVV      | 221941         | 390011        | 20.74%          | 1.224%        |
| 17        | 11.122      | 1406          | 1410        | 1413         | rVB2     | 10555          | 16168         | 0.86%           | 0.051%        |
| 18        | 11.192      | 1416          | 1422        | 1430         | rBV      | 167719         | 321289        | 17.08%          | 1.008%        |
| 19        | 11.815      | 1521          | 1528        | 1533         | rBV3     | 16096          | 27177         | 1.45%           | 0.085%        |
| 20        | 12.215      | 1591          | 1596        | 1600         | rBV5     | 11913          | 22211         | 1.18%           | 0.070%        |
| 21        | 12.450      | 1631          | 1636        | 1642         | rVB3     | 21600          | 34688         | 1.84%           | 0.109%        |
| 22        | 12.567      | 1650          | 1656        | 1662         | rBV6     | 7888           | 16058         | 0.85%           | 0.050%        |
| 23        | 12.708      | 1673          | 1680        | 1686         | rBV4     | 28312          | 50855         | 2.70%           | 0.160%        |
| 24        | 12.920      | 1711          | 1716        | 1723         | rBV2     | 17515          | 28249         | 1.50%           | 0.089%        |
| 25        | 13.119      | 1744          | 1750        | 1757         | rVB8     | 41955          | 71914         | 3.82%           | 0.226%        |
| 26        | 13.348      | 1781          | 1789        | 1794         | rBV2     | 23078          | 36894         | 1.96%           | 0.116%        |
| 27        | 13.648      | 1836          | 1840        | 1841         | rBV3     | 23966          | 25744         | 1.37%           | 0.081%        |
| 28        | 13.666      | 1841          | 1843        | 1849         | rVV4     | 28748          | 46016         | 2.45%           | 0.144%        |
| 29        | 13.766      | 1855          | 1860        | 1866         | rVB4     | 24885          | 41546         | 2.21%           | 0.130%        |
| 30        | 13.983      | 1892          | 1897        | 1900         | rBV7     | 11174          | 18893         | 1.00%           | 0.059%        |
| 31        | 14.024      | 1900          | 1904        | 1909         | rVV7     | 16576          | 28433         | 1.51%           | 0.089%        |
| 32        | 14.195      | 1927          | 1933        | 1936         | rBV6     | 31262          | 61674         | 3.28%           | 0.194%        |
| 33        | 14.253      | 1937          | 1943        | 1948         | rVV      | 463147         | 736467        | 39.16%          | 2.311%        |
| 34        | 14.300      | 1948          | 1951        | 1959         | rVB      | 142149         | 200824        | 10.68%          | 0.630%        |

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 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:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG102315.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |        |         |        |        |
|----|--------|------|------|------|------|--------|---------|--------|--------|
| 35 | 14.418 | 1966 | 1971 | 1973 | rBV6 | 12279  | 19602   | 1.04%  | 0.062% |
| 36 | 14.453 | 1973 | 1977 | 1985 | rVV6 | 31113  | 56968   | 3.03%  | 0.179% |
| 37 | 14.553 | 1987 | 1994 | 2002 | rVV  | 483775 | 749903  | 39.87% | 2.353% |
| 38 | 14.694 | 2015 | 2018 | 2021 | rVV4 | 16930  | 26158   | 1.39%  | 0.082% |
| 39 | 14.735 | 2021 | 2025 | 2030 | rVV  | 40978  | 53817   | 2.86%  | 0.169% |
| 40 | 14.858 | 2036 | 2046 | 2059 | rBV2 | 427914 | 768944  | 40.89% | 2.413% |
| 41 | 14.994 | 2062 | 2069 | 2071 | rBV  | 78249  | 129640  | 6.89%  | 0.407% |
| 42 | 15.023 | 2071 | 2074 | 2081 | rVB  | 141790 | 226245  | 12.03% | 0.710% |
| 43 | 15.176 | 2096 | 2100 | 2107 | rVB8 | 30081  | 56839   | 3.02%  | 0.178% |
| 44 | 15.240 | 2107 | 2111 | 2121 | rVB5 | 79592  | 143662  | 7.64%  | 0.451% |
| 45 | 15.329 | 2122 | 2126 | 2128 | rBV5 | 13538  | 19269   | 1.02%  | 0.060% |
| 46 | 15.393 | 2128 | 2137 | 2143 | rBV2 | 66269  | 123887  | 6.59%  | 0.389% |
| 47 | 15.470 | 2147 | 2150 | 2155 | rVB7 | 16425  | 19052   | 1.01%  | 0.060% |
| 48 | 15.517 | 2155 | 2158 | 2161 | rVB4 | 9804   | 14293   | 0.76%  | 0.045% |
| 49 | 15.564 | 2161 | 2166 | 2169 | rBV7 | 11604  | 22774   | 1.21%  | 0.071% |
| 50 | 15.616 | 2170 | 2175 | 2181 | rBV7 | 22662  | 51261   | 2.73%  | 0.161% |
| 51 | 15.681 | 2181 | 2186 | 2190 | rVB2 | 70102  | 97543   | 5.19%  | 0.306% |
| 52 | 15.728 | 2190 | 2194 | 2198 | rBV7 | 17156  | 28843   | 1.53%  | 0.090% |
| 53 | 15.781 | 2198 | 2203 | 2208 | rBV  | 187331 | 259777  | 13.81% | 0.815% |
| 54 | 15.846 | 2208 | 2214 | 2219 | rBV  | 611610 | 951377  | 50.59% | 2.985% |
| 55 | 15.945 | 2225 | 2231 | 2237 | rVB  | 258646 | 390782  | 20.78% | 1.226% |
| 56 | 16.057 | 2245 | 2250 | 2251 | rBV4 | 20236  | 27852   | 1.48%  | 0.087% |
| 57 | 16.139 | 2260 | 2264 | 2267 | rBV5 | 34836  | 56073   | 2.98%  | 0.176% |
| 58 | 16.269 | 2281 | 2286 | 2292 | rBV  | 147690 | 214691  | 11.42% | 0.674% |
| 59 | 16.539 | 2328 | 2332 | 2335 | rBV  | 79410  | 101554  | 5.40%  | 0.319% |
| 60 | 16.668 | 2349 | 2354 | 2358 | rBV5 | 44547  | 75283   | 4.00%  | 0.236% |
| 61 | 16.868 | 2383 | 2388 | 2398 | rBV3 | 187917 | 332987  | 17.71% | 1.045% |
| 62 | 16.968 | 2401 | 2405 | 2409 | rVB2 | 60080  | 74533   | 3.96%  | 0.234% |
| 63 | 17.244 | 2449 | 2452 | 2457 | rVB4 | 40266  | 55648   | 2.96%  | 0.175% |
| 64 | 17.332 | 2460 | 2467 | 2471 | rBV3 | 122391 | 203904  | 10.84% | 0.640% |
| 65 | 17.526 | 2498 | 2500 | 2503 | rBV4 | 29932  | 30410   | 1.62%  | 0.095% |
| 66 | 17.596 | 2505 | 2512 | 2517 | rVV2 | 610988 | 1082710 | 57.57% | 3.397% |
| 67 | 17.638 | 2517 | 2519 | 2523 | rVB  | 74345  | 82915   | 4.41%  | 0.260% |
| 68 | 17.696 | 2523 | 2529 | 2534 | rBV  | 718657 | 1075085 | 57.16% | 3.373% |
| 69 | 17.737 | 2534 | 2536 | 2540 | rVB2 | 71036  | 71143   | 3.78%  | 0.223% |
| 70 | 17.914 | 2561 | 2566 | 2570 | rVV7 | 31094  | 54431   | 2.89%  | 0.171% |
| 71 | 18.072 | 2589 | 2593 | 2597 | rBV  | 158857 | 188993  | 10.05% | 0.593% |

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Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG102315.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |        |
|-----|--------|------|------|------|------|---------|---------|---------|--------|
| 72  | 18.243 | 2619 | 2622 | 2625 | rBV3 | 41330   | 45505   | 2.42%   | 0.143% |
| 73  | 18.466 | 2655 | 2660 | 2664 | rBV  | 535216  | 707612  | 37.62%  | 2.220% |
| 74  | 18.754 | 2705 | 2709 | 2714 | rBV  | 235833  | 319801  | 17.00%  | 1.003% |
| 75  | 18.807 | 2715 | 2718 | 2722 | rVB  | 83781   | 115445  | 6.14%   | 0.362% |
| 76  | 18.995 | 2748 | 2750 | 2756 | rBV7 | 37423   | 59684   | 3.17%   | 0.187% |
| 77  | 19.377 | 2811 | 2815 | 2822 | rVB2 | 336861  | 437751  | 23.28%  | 1.374% |
| 78  | 19.629 | 2855 | 2858 | 2863 | rVB  | 183184  | 282298  | 15.01%  | 0.886% |
| 79  | 19.723 | 2870 | 2874 | 2881 | rVB2 | 230580  | 297335  | 15.81%  | 0.933% |
| 80  | 19.923 | 2905 | 2908 | 2910 | rBV2 | 99476   | 127895  | 6.80%   | 0.401% |
| 81  | 19.964 | 2910 | 2915 | 2925 | rVB3 | 908631  | 1880700 | 100.00% | 5.901% |
| 82  | 20.446 | 2994 | 2997 | 2999 | rBV4 | 93261   | 131782  | 7.01%   | 0.413% |
| 83  | 20.475 | 2999 | 3002 | 3007 | rVB  | 442419  | 519913  | 27.64%  | 1.631% |
| 84  | 20.793 | 3053 | 3056 | 3059 | rVB3 | 124980  | 139685  | 7.43%   | 0.438% |
| 85  | 20.981 | 3081 | 3088 | 3095 | rVB2 | 480663  | 949761  | 50.50%  | 2.980% |
| 86  | 21.498 | 3171 | 3176 | 3181 | rVB2 | 408650  | 640722  | 34.07%  | 2.010% |
| 87  | 21.662 | 3200 | 3204 | 3210 | rVB  | 285085  | 422330  | 22.46%  | 1.325% |
| 88  | 21.886 | 3236 | 3242 | 3249 | rBV  | 1022860 | 1691462 | 89.94%  | 5.307% |
| 89  | 22.038 | 3263 | 3268 | 3272 | rBV3 | 366226  | 772006  | 41.05%  | 2.422% |
| 90  | 22.079 | 3273 | 3275 | 3280 | rVB2 | 383405  | 483225  | 25.69%  | 1.516% |
| 91  | 22.737 | 3381 | 3387 | 3392 | rBV  | 697898  | 1241795 | 66.03%  | 3.896% |
| 92  | 22.937 | 3416 | 3421 | 3429 | rVB2 | 274618  | 535421  | 28.47%  | 1.680% |
| 93  | 23.243 | 3469 | 3473 | 3478 | rVB4 | 111041  | 216454  | 11.51%  | 0.679% |
| 94  | 23.396 | 3494 | 3499 | 3501 | rBV5 | 127986  | 240621  | 12.79%  | 0.755% |
| 95  | 23.437 | 3502 | 3506 | 3516 | rVB4 | 423784  | 847700  | 45.07%  | 2.660% |
| 96  | 24.377 | 3660 | 3666 | 3675 | rVB2 | 455152  | 1050320 | 55.85%  | 3.296% |
| 97  | 25.047 | 3773 | 3780 | 3796 | rVB  | 461045  | 1368316 | 72.76%  | 4.293% |
| 98  | 25.287 | 3813 | 3821 | 3829 | rVB2 | 377750  | 1015923 | 54.02%  | 3.188% |
| 99  | 25.987 | 3932 | 3940 | 3950 | rVB  | 324154  | 924390  | 49.15%  | 2.900% |
| 100 | 26.627 | 4043 | 4049 | 4067 | rVB  | 227285  | 846149  | 44.99%  | 2.655% |

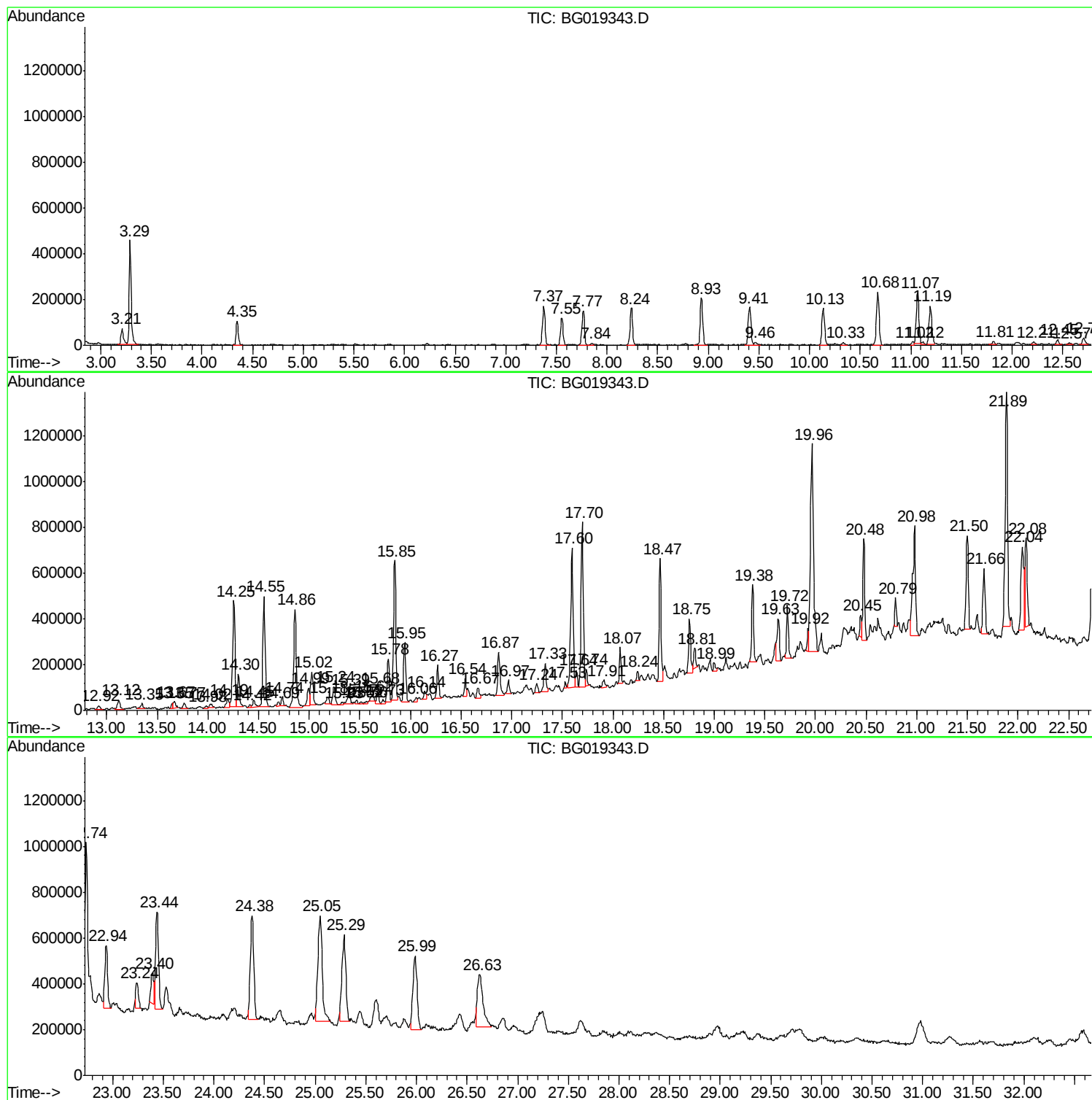
Sum of corrected areas: 31871005

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P



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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

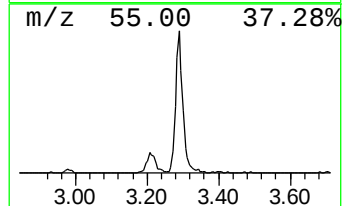
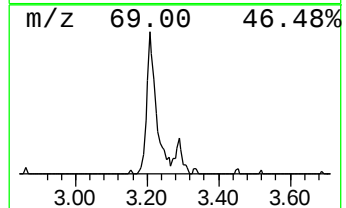
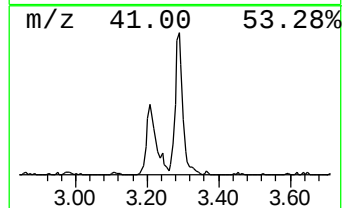
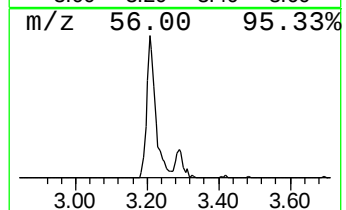
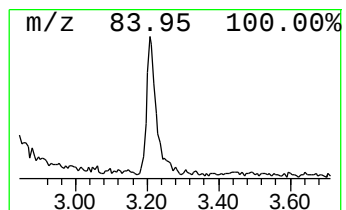
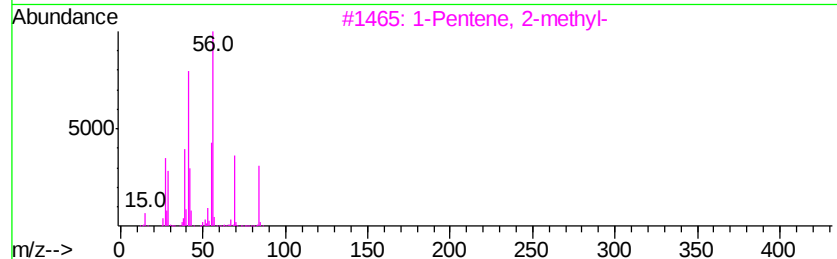
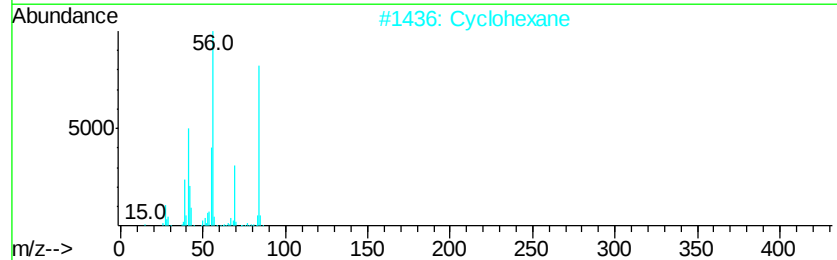
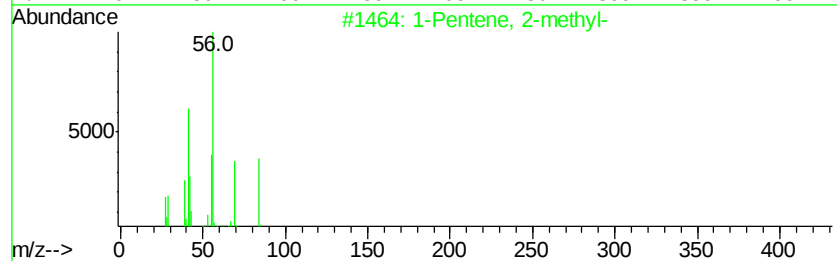
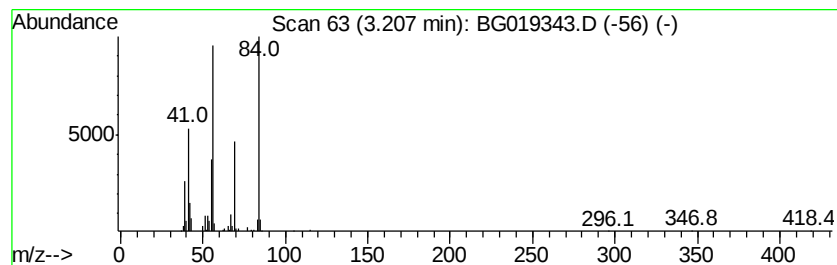
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 (DEL) Alkane: Cyclic3.21 Concentration Rank 10

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 3.21 | 9.59 ng/ul | 133381 | 1,4-Dichlorobenzene-d4 | 8.24 |

| Hit# | of | 5 | Tentative ID         | MW | MolForm | CAS#        | Qual |
|------|----|---|----------------------|----|---------|-------------|------|
| 1    |    |   | 1-Pentene, 2-methyl- | 84 | C6H12   | 000763-29-1 | 90   |
| 2    |    |   | Cyclohexane          | 84 | C6H12   | 000110-82-7 | 87   |
| 3    |    |   | 1-Pentene, 2-methyl- | 84 | C6H12   | 000763-29-1 | 80   |
| 4    |    |   | Cyclohexane          | 84 | C6H12   | 000110-82-7 | 72   |
| 5    |    |   | Cyclohexane          | 84 | C6H12   | 000110-82-7 | 72   |



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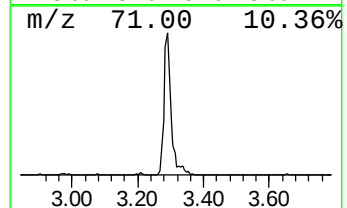
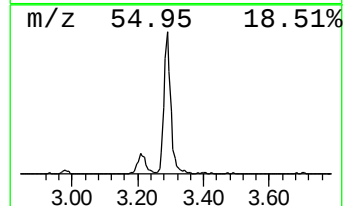
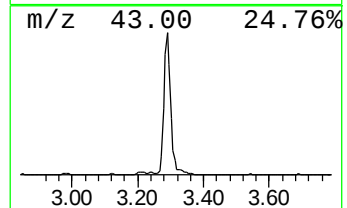
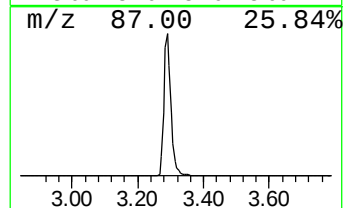
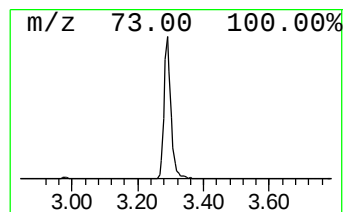
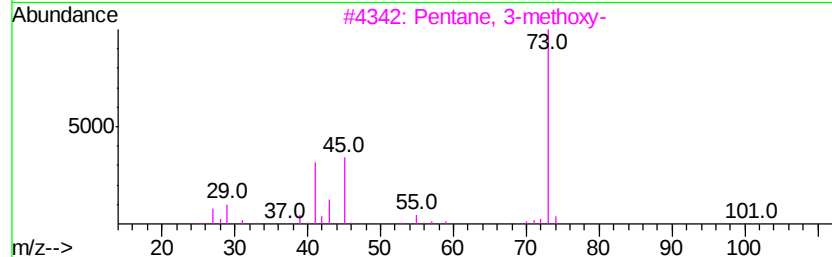
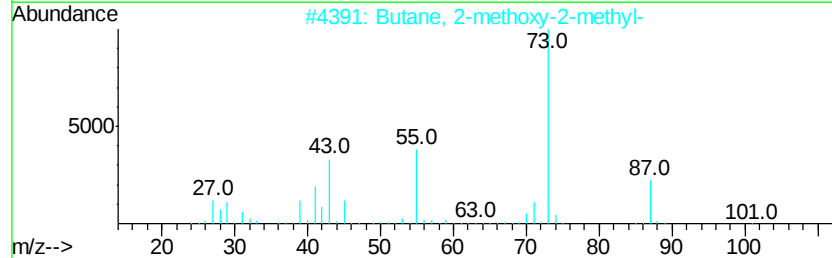
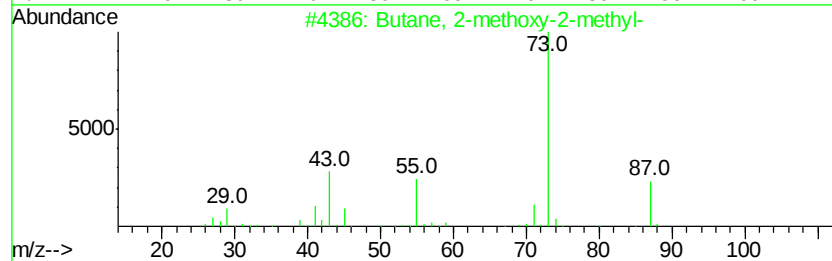
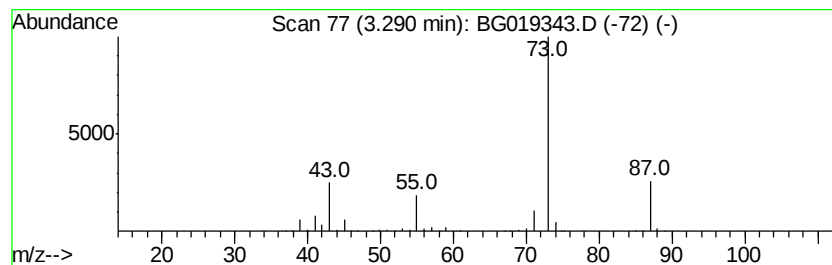
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 3.29 | 48.02 ng/ul | 667527 | 1,4-Dichlorobenzene-d4 | 8.24 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 72   |
| 2    |    | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 64   |
| 3    |    | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 17   |
| 4    |    | Acetamide, N-ethyl-         | 87  | C4H9NO  | 000625-50-3 | 12   |
| 5    |    | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 9    |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

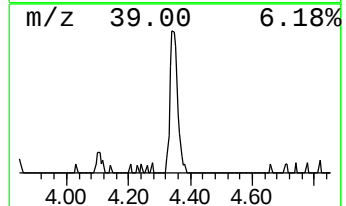
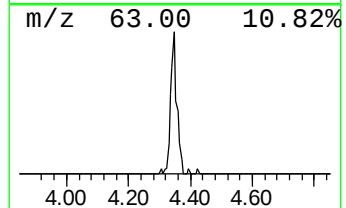
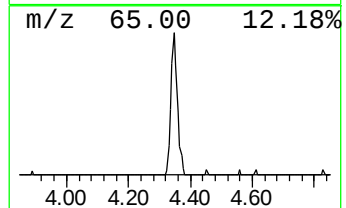
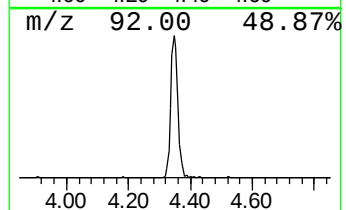
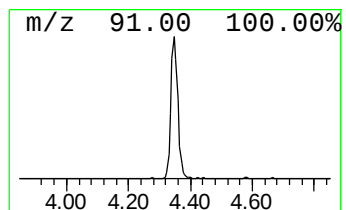
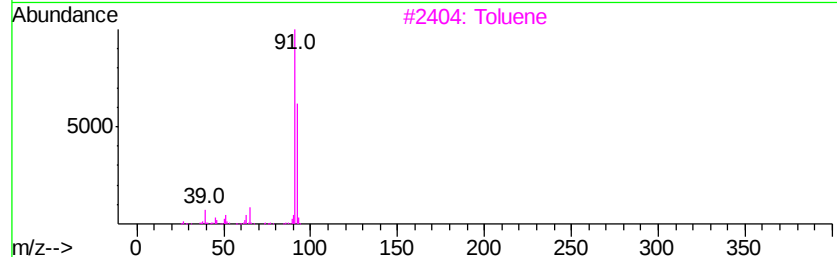
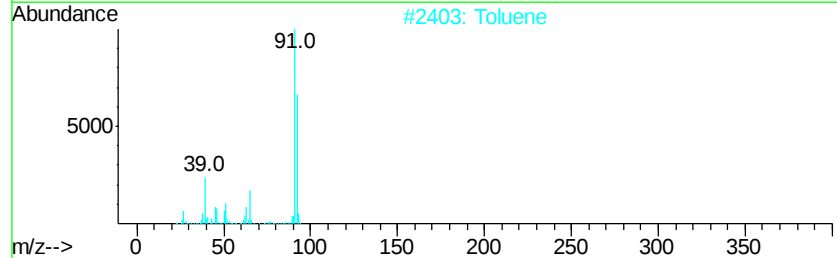
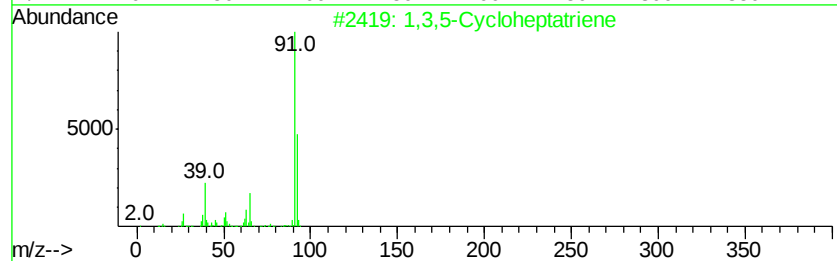
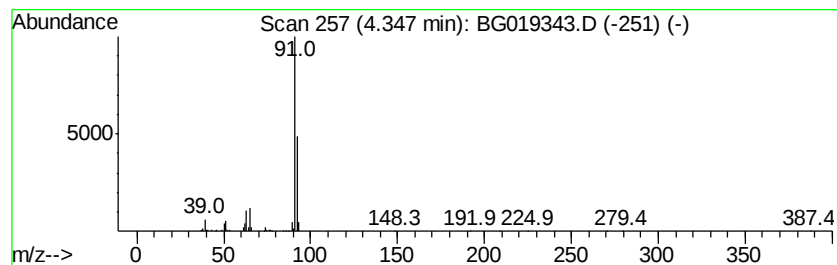
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 1,3,5-Cycloheptatriene Concentration Rank 7

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 4.35 | 11.30 ng/ul | 157096 | 1,4-Dichlorobenzene-d4 | 8.24 |

| Hit# | of | Tentative ID           | MW | MolForm | CAS#        | Qual |
|------|----|------------------------|----|---------|-------------|------|
| 1    | 5  | 1,3,5-Cycloheptatriene | 92 | C7H8    | 000544-25-2 | 90   |
| 2    |    | Toluene                | 92 | C7H8    | 000108-88-3 | 87   |
| 3    |    | Toluene                | 92 | C7H8    | 000108-88-3 | 87   |
| 4    |    | Toluene                | 92 | C7H8    | 000108-88-3 | 81   |
| 5    |    | Toluene                | 92 | C7H8    | 000108-88-3 | 81   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

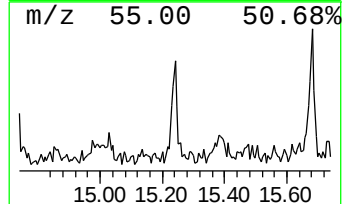
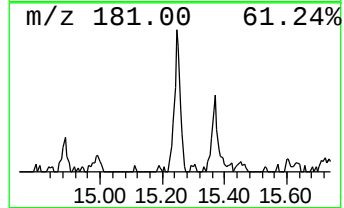
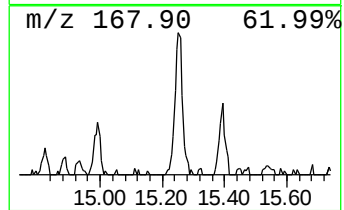
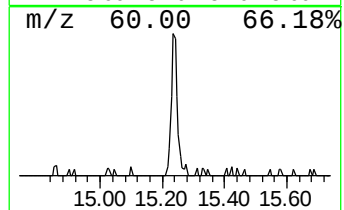
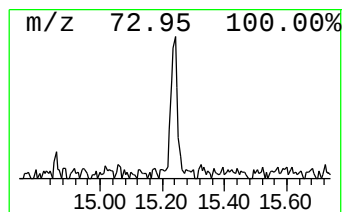
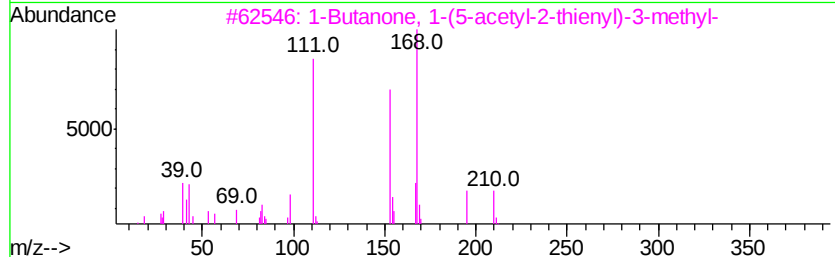
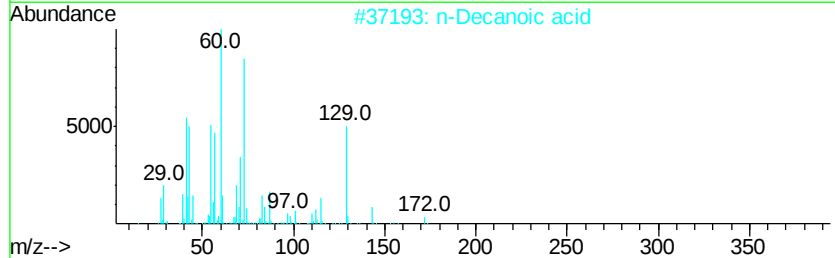
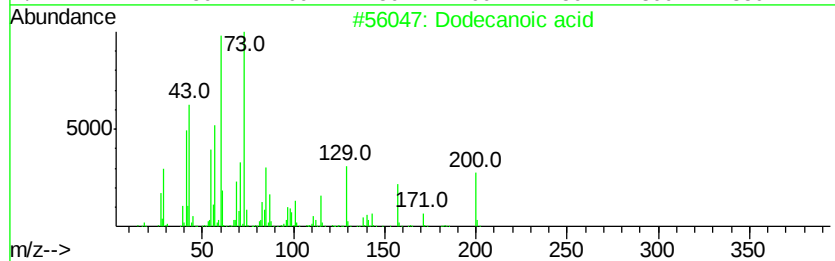
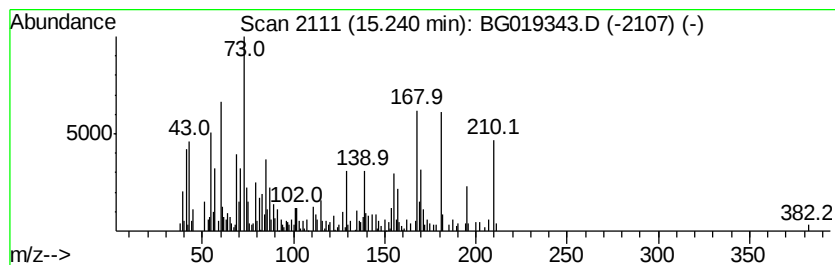
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 unknown-01 Concentration Rank 24

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.24 | 3.74 ng/ul | 143662 | Acenaphthene-d10 | 14.86 |

| Hit# | of | Tentative ID                                 | MW  | MolForm   | CAS#        | Qual |
|------|----|----------------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Dodecanoic acid                              | 200 | C12H24O2  | 000143-07-7 | 40   |
| 2    |    | n-Decanoic acid                              | 172 | C10H20O2  | 000334-48-5 | 25   |
| 3    |    | 1-Butanone, 1-(5-acetyl-2-thienyl)-3-methyl- | 210 | C11H14O2S | 018282-21-8 | 22   |
| 4    |    | Allo-Inositol, hexaacetate                   | 432 | C18H24O12 | 029267-04-7 | 16   |
| 5    |    | Inositol, hexaacetate, D-chiro-              | 432 | C18H24O12 | 029267-03-6 | 12   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

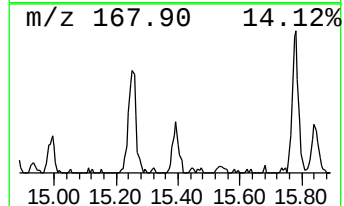
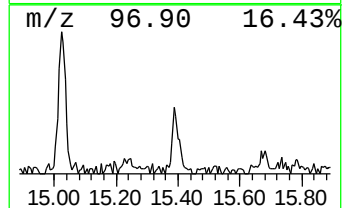
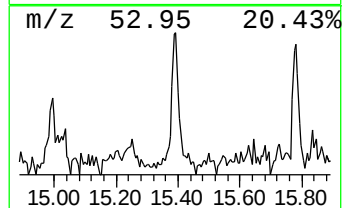
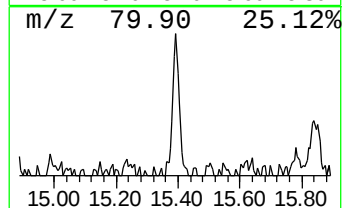
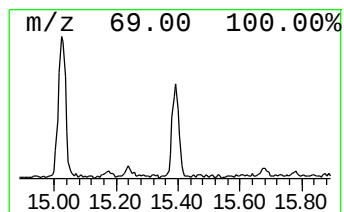
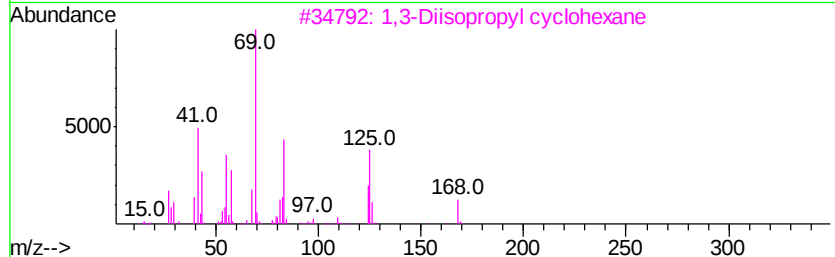
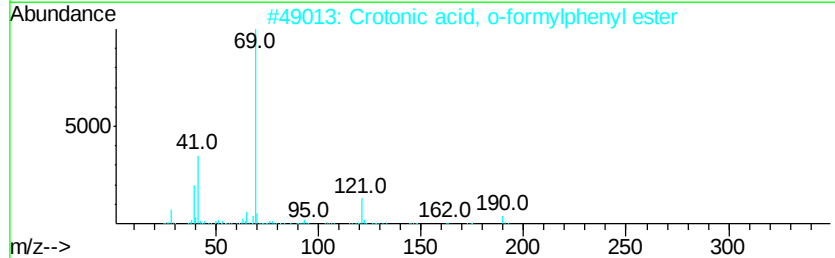
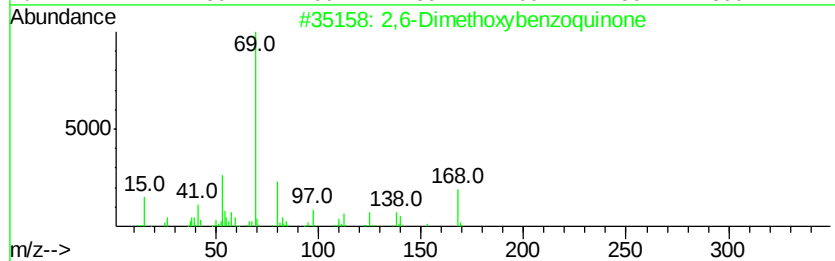
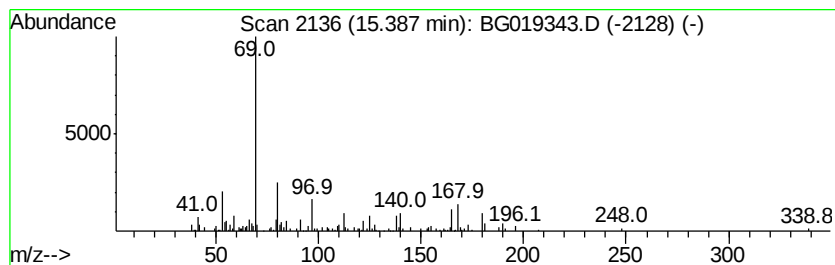
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 2,6-Dimethoxybenzoquinone Concentration Rank 26

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.39 | 3.22 ng/ul | 123887 | Acenaphthene-d10 | 14.86 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2,6-Dimethoxybenzoquinone           | 168 | C8H8O4   | 000530-55-2 | 68   |
| 2    |    | Crotonic acid, o-formylphenyl ester | 190 | C11H10O3 | 114636-69-0 | 38   |
| 3    |    | 1,3-Diisopropyl cyclohexane         | 168 | C12H24   | 007045-70-7 | 37   |
| 4    |    | 1-Hexyn-3-ol, 3-methyl-             | 112 | C7H12O   | 004339-05-3 | 37   |
| 5    |    | Methanone, dicyclopropyl-           | 110 | C7H10O   | 001121-37-5 | 32   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

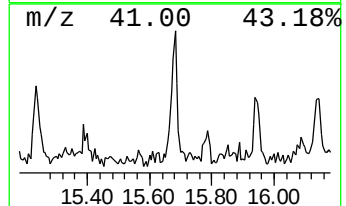
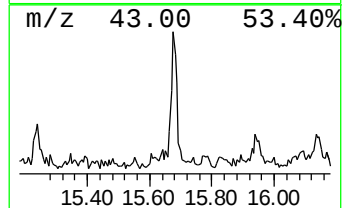
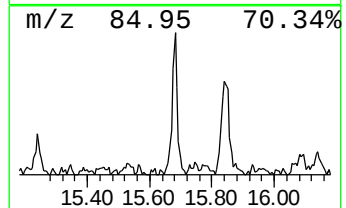
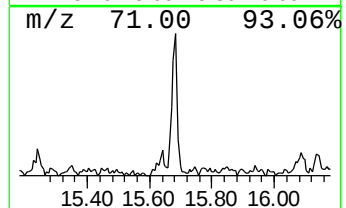
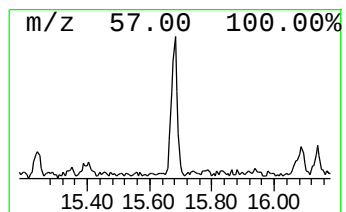
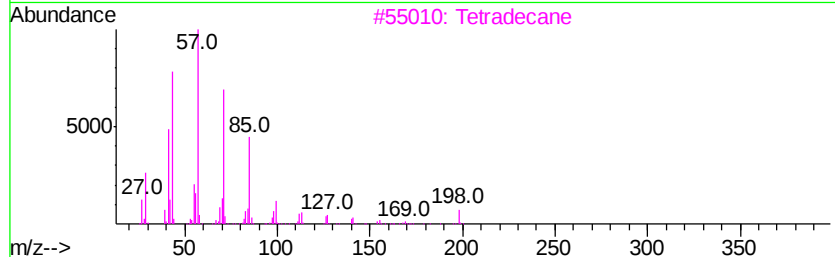
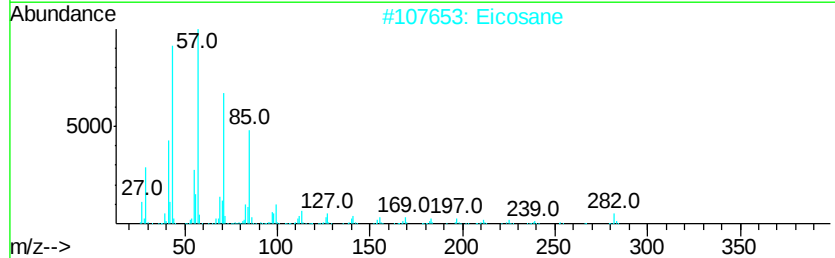
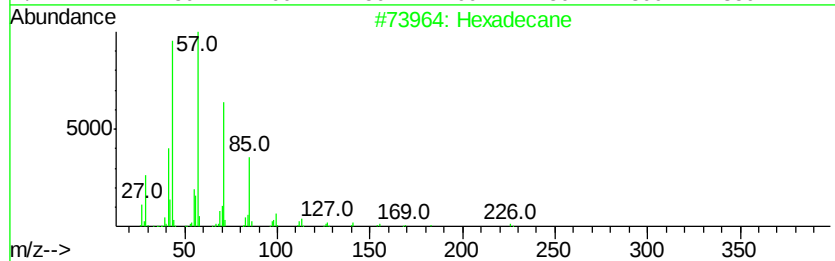
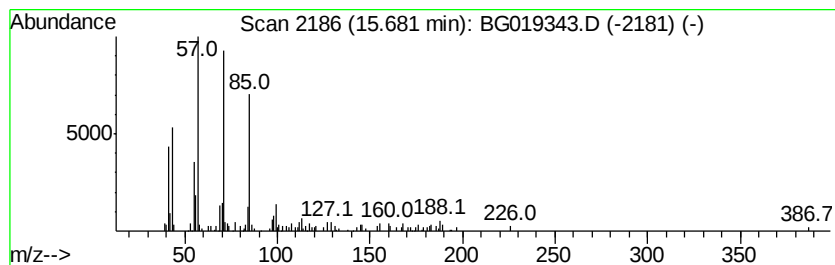
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 29

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 15.68 | 2.54 ng/ul | 97543 | Acenaphthene-d10 | 14.86 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 93   |
| 2    |    |   | Eicosane     | 282 | C20H42  | 000112-95-8 | 86   |
| 3    |    |   | Tetradecane  | 198 | C14H30  | 000629-59-4 | 86   |
| 4    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 76   |
| 5    |    |   | Pentadecane  | 212 | C15H32  | 000629-62-9 | 52   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

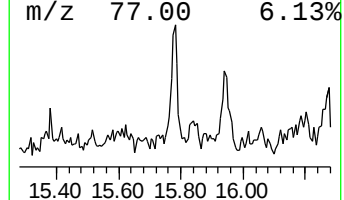
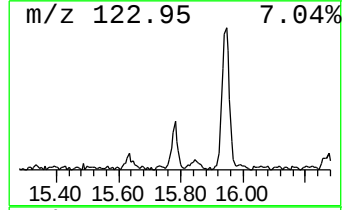
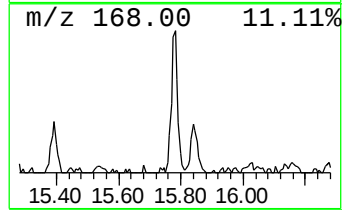
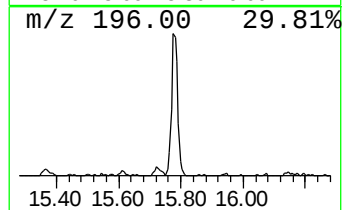
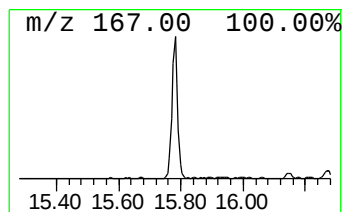
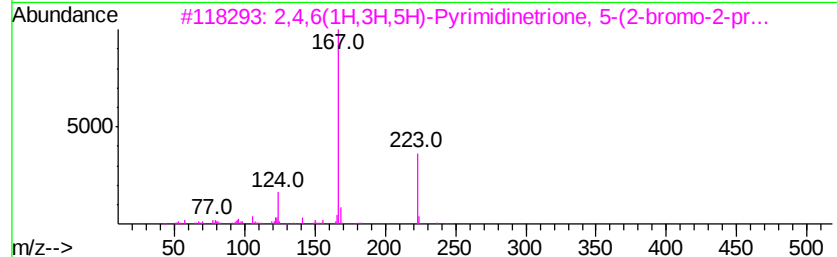
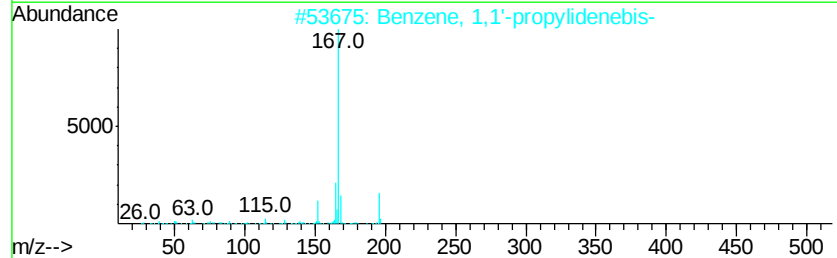
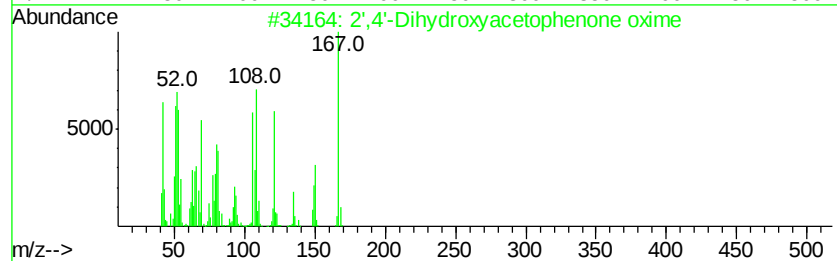
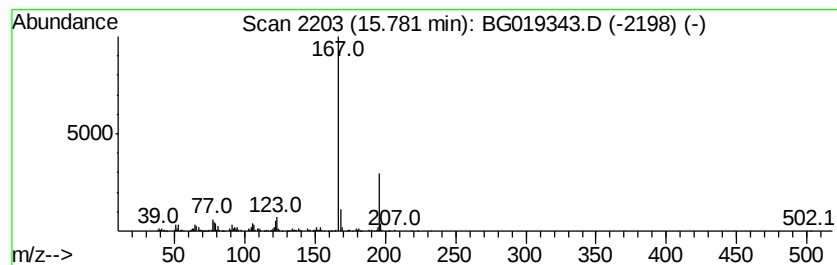
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 2',4'-Dihydroxyacetophenone... Concentration Rank 14

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.78 | 6.76 ng/ul | 259777 | Acenaphthene-d10 | 14.86 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|--------------|--------------|------|
| 1    |    |   | 2',4'-Dihydroxyacetophenone oxime   | 167 | C8H9NO3      | 1000212-89-5 | 52   |
| 2    |    |   | Benzene, 1,1'-propylidenebis-       | 196 | C15H16       | 001530-03-6  | 50   |
| 3    |    |   | 2,4,6(1H,3H,5H)-Pyrimidinetrione... | 302 | C11H15BrN2O3 | 001142-70-7  | 45   |
| 4    |    |   | Benzene, 1,1'-[(2-phenylethoxy)m... | 288 | C21H20O      | 067810-88-2  | 42   |
| 5    |    |   | Cyclopentylmethylphenylphosphine... | 208 | C12H17OP     | 254100-71-5  | 40   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

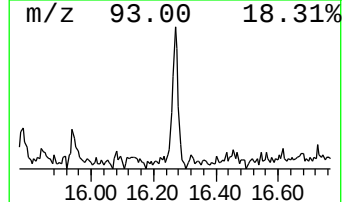
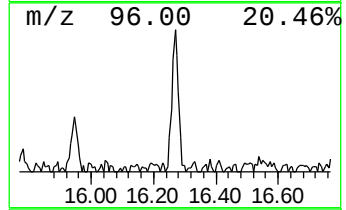
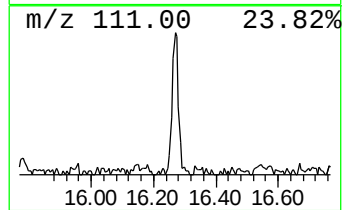
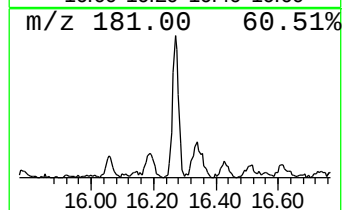
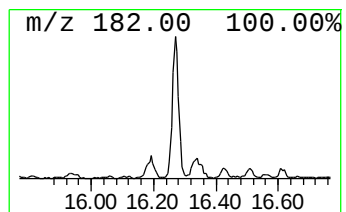
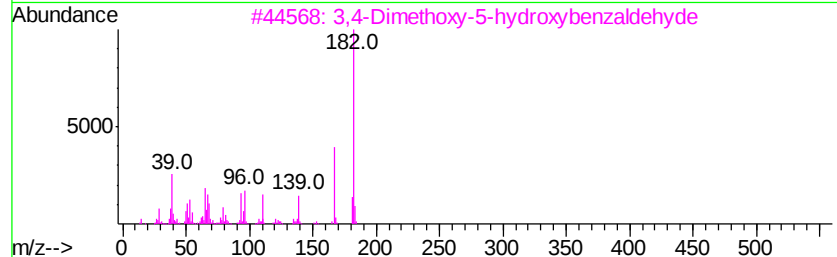
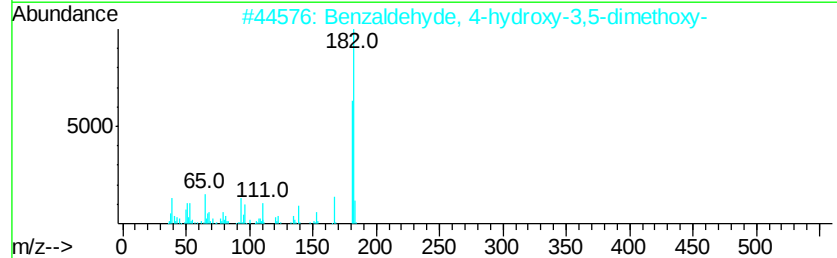
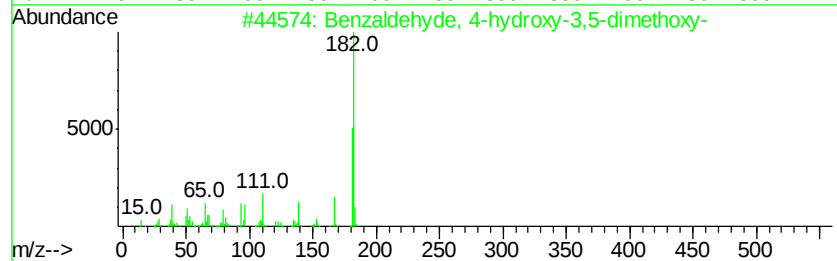
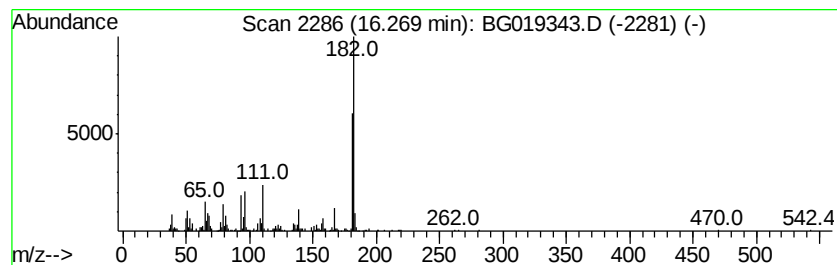
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Benzaldehyde, 4-hydroxy-3,5... Concentration Rank 22

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.27 | 3.97 ng/ul | 214691 | Phenanthrene-d10 | 17.60 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Benzaldehyde, 4-hydroxy-3,5-dime... | 182 | C9H10O4  | 000134-96-3  | 81   |
| 2    |    | Benzaldehyde, 4-hydroxy-3,5-dime... | 182 | C9H10O4  | 000134-96-3  | 68   |
| 3    |    | 3,4-Dimethoxy-5-hydroxybenzaldehyd  | 182 | C9H10O4  | 029865-90-5  | 52   |
| 4    |    | 7-Phosphabicyclo[2.2.1]hept-2-en... | 182 | C11H19P  | 1000151-83-5 | 47   |
| 5    |    | 9H-Pyrido[3,4-b]indole, 1-methyl-   | 182 | C12H10N2 | 000486-84-0  | 43   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

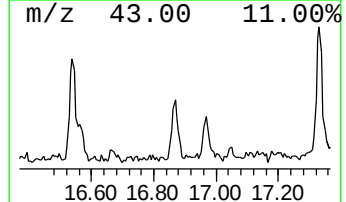
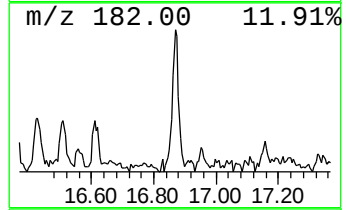
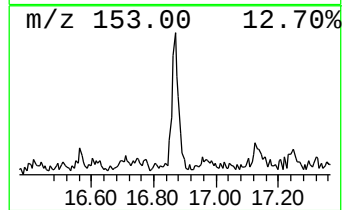
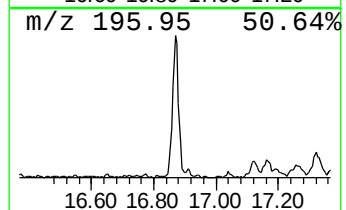
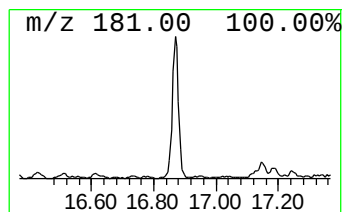
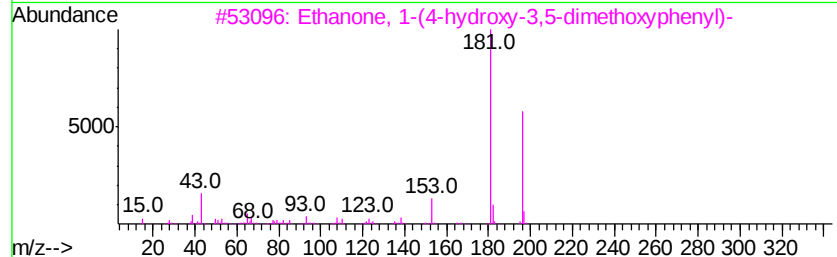
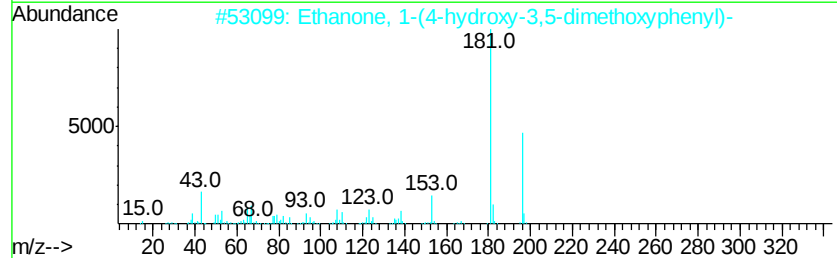
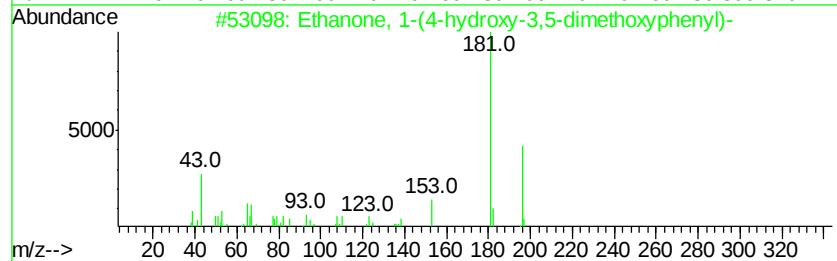
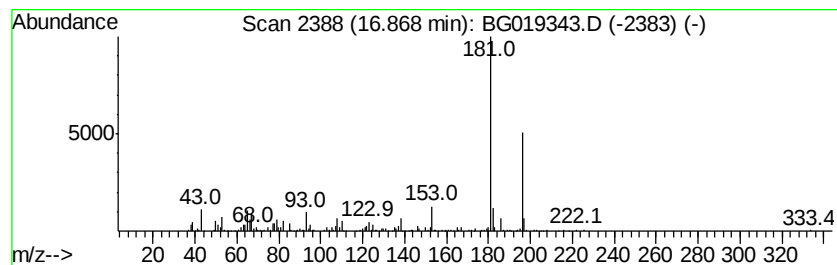
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Ethanone, 1-(4-hydroxy-3,5-... Concentration Rank 16

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.87 | 6.15 ng/ul | 332987 | Phenanthrene-d10 | 17.60 |

| Hit# | of | Tentative ID                                 | MW  | MolForm   | CAS#        | Qual |
|------|----|----------------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Ethanone, 1-(4-hydroxy-3,5-dimethoxyphenyl)- | 196 | C10H12O4  | 002478-38-8 | 96   |
| 2    |    | Ethanone, 1-(4-hydroxy-3,5-dimethoxyphenyl)- | 196 | C10H12O4  | 002478-38-8 | 94   |
| 3    |    | Ethanone, 1-(4-hydroxy-3,5-dimethoxyphenyl)- | 196 | C10H12O4  | 002478-38-8 | 93   |
| 4    |    | 2-Chloro-2-methyl-1-oxa-2-sila-1-phenylene   | 196 | C9H9ClOSi | 073060-34-1 | 64   |
| 5    |    | 1,1'-Biphenyl, 4-(1-methylethyl)-            | 196 | C15H16    | 007116-95-2 | 50   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

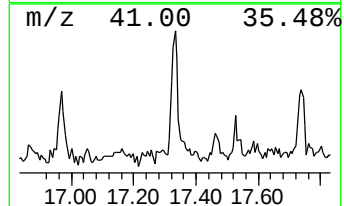
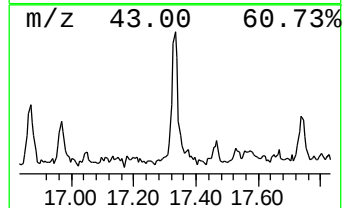
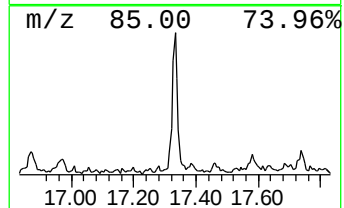
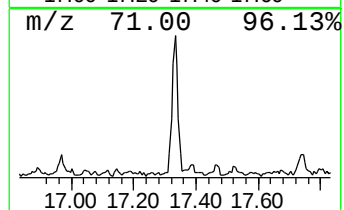
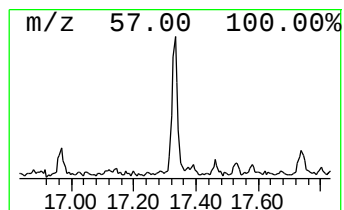
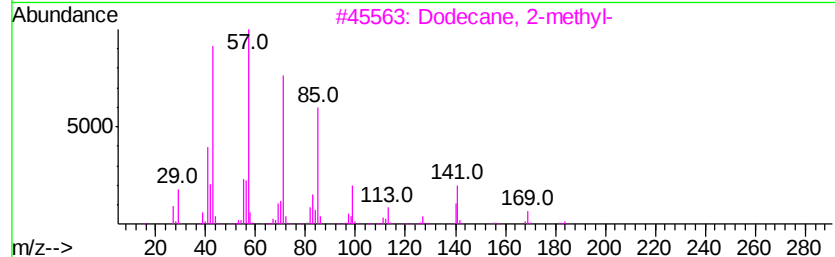
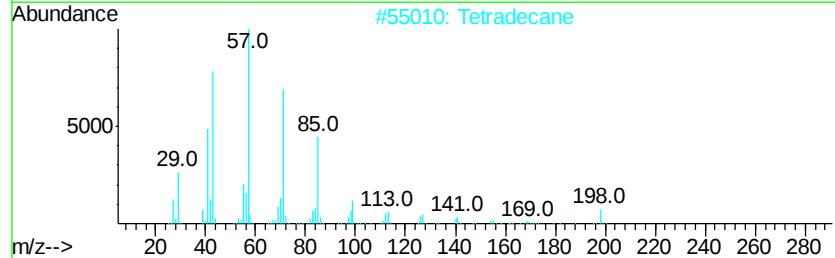
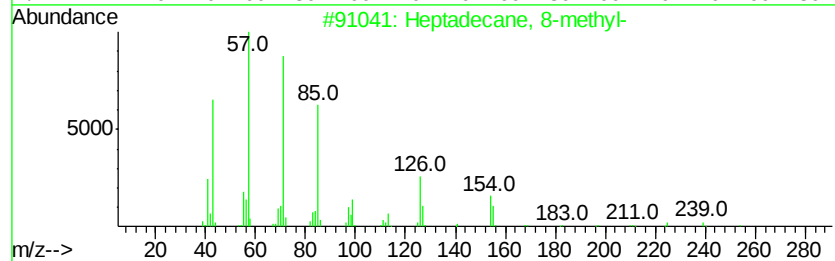
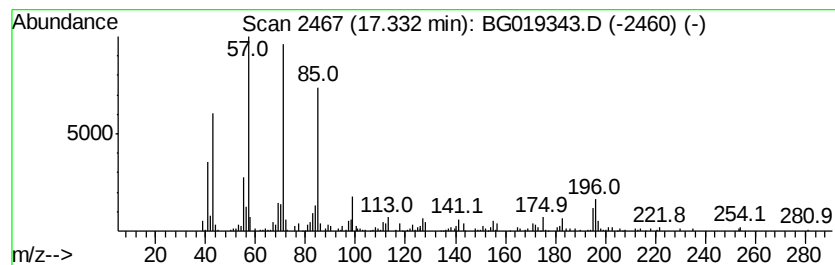
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 23

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.33 | 3.77 ng/ul | 203904 | Phenanthrene-d10 | 17.60 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Heptadecane, 8-methyl-    | 254 | C18H38  | 013287-23-5 | 90   |
| 2    |    | Tetradecane               | 198 | C14H30  | 000629-59-4 | 87   |
| 3    |    | Dodecane, 2-methyl-       | 184 | C13H28  | 001560-97-0 | 83   |
| 4    |    | Nonane, 5-butyl-          | 184 | C13H28  | 017312-63-9 | 72   |
| 5    |    | Hexadecane, 7,9-dimethyl- | 254 | C18H38  | 021164-95-4 | 52   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

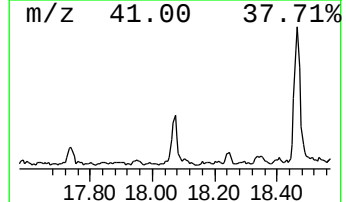
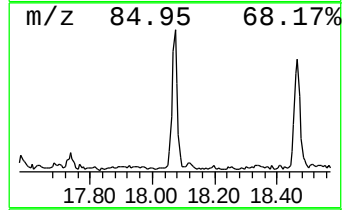
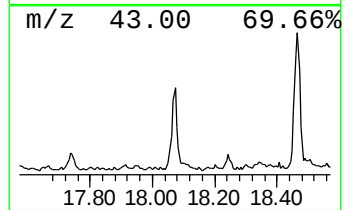
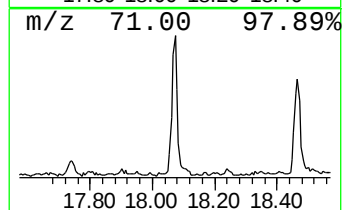
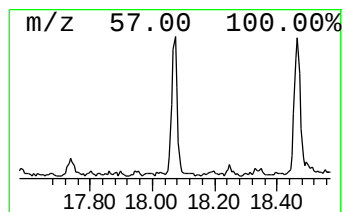
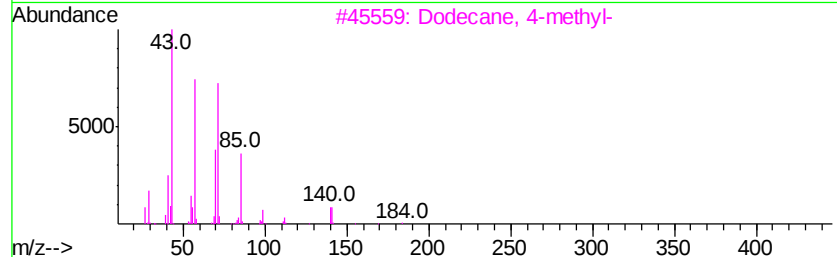
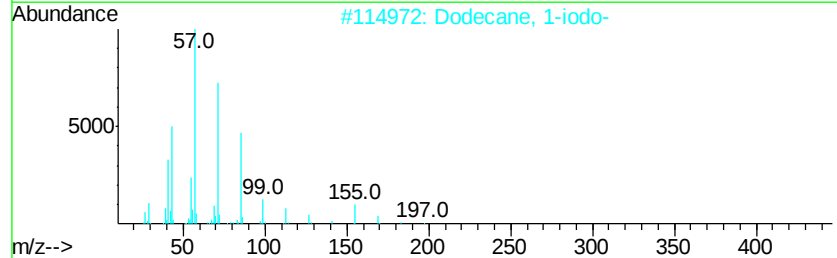
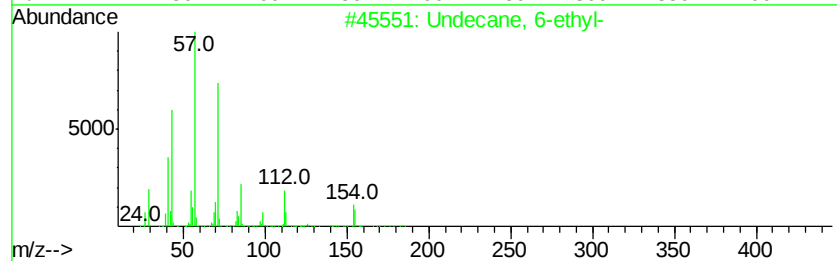
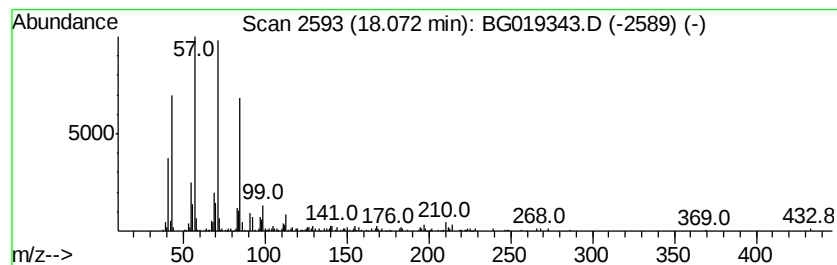
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 25

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.07 | 3.49 ng/ul | 188993 | Phenanthrene-d10 | 17.60 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Undecane, 6-ethyl-       | 184 | C13H28  | 017312-60-6 | 81   |
| 2    |    |   | Dodecane, 1-iodo-        | 296 | C12H25I | 004292-19-7 | 80   |
| 3    |    |   | Dodecane, 4-methyl-      | 184 | C13H28  | 006117-97-1 | 80   |
| 4    |    |   | 10-Methylnonadecane      | 282 | C20H42  | 056862-62-5 | 80   |
| 5    |    |   | Decane, 2,3,5-trimethyl- | 184 | C13H28  | 062238-11-3 | 78   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

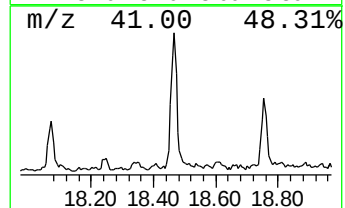
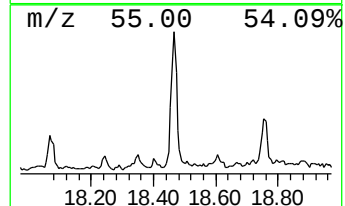
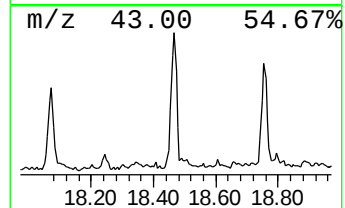
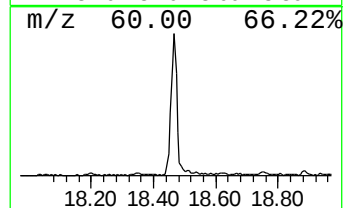
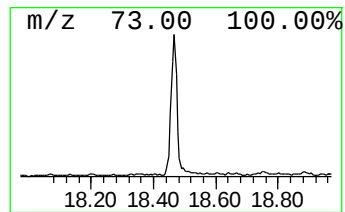
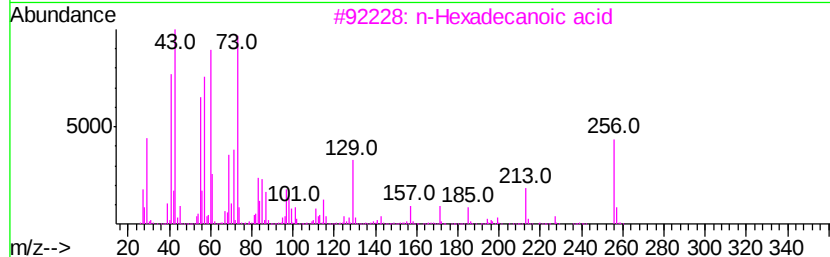
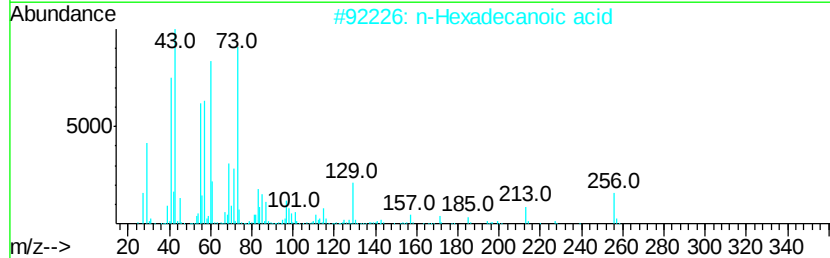
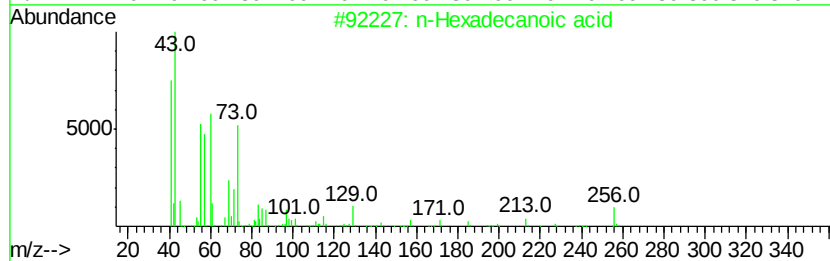
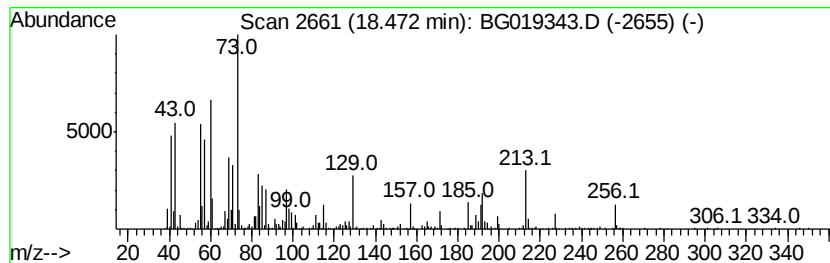
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 n-Hexadecanoic acid Concentration Rank 6

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 18.47 | 13.07 ng/ul | 707612 | Phenanthrene-d10 | 17.60 |

| Hit# | of | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------|-----|----------|-------------|------|
| 1    | 5  | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 95   |
| 2    |    | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 94   |
| 3    |    | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 83   |
| 4    |    | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 64   |
| 5    |    | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 52   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

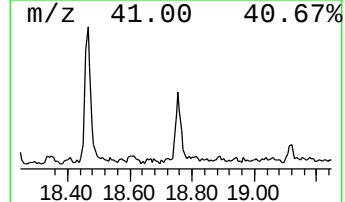
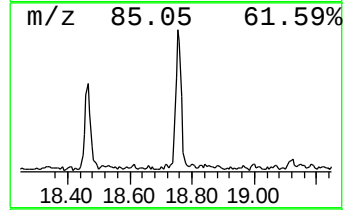
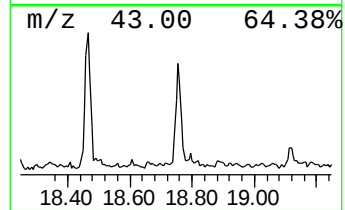
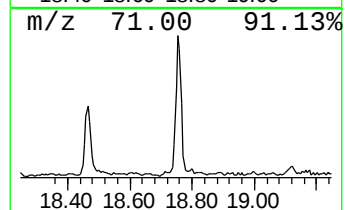
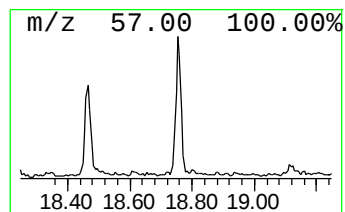
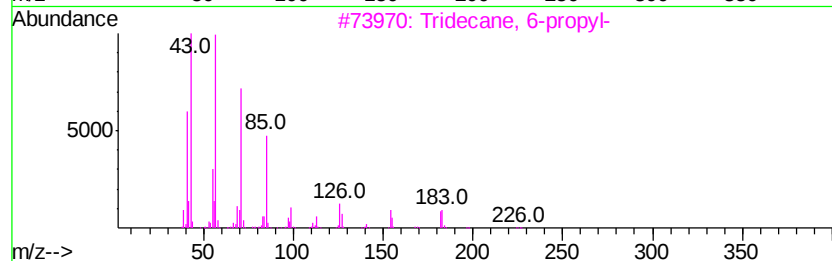
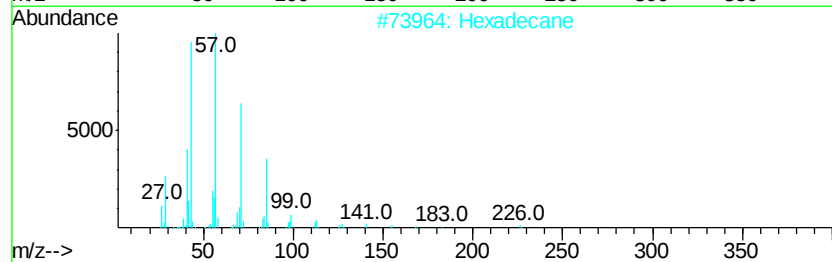
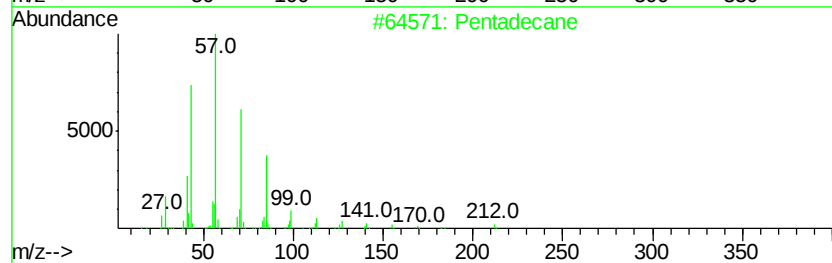
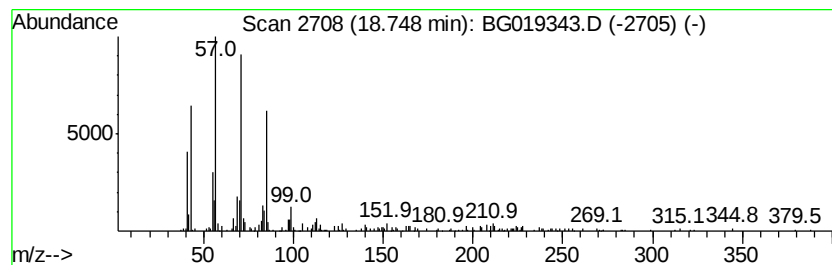
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 18

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.75 | 5.91 ng/ul | 319801 | Phenanthrene-d10 | 17.60 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Pentadecane          | 212 | C15H32  | 000629-62-9 | 90   |
| 2    |    |   | Hexadecane           | 226 | C16H34  | 000544-76-3 | 90   |
| 3    |    |   | Tridecane, 6-propyl- | 226 | C16H34  | 055045-10-8 | 87   |
| 4    |    |   | Hexadecane           | 226 | C16H34  | 000544-76-3 | 81   |
| 5    |    |   | Docosane             | 310 | C22H46  | 000629-97-0 | 78   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

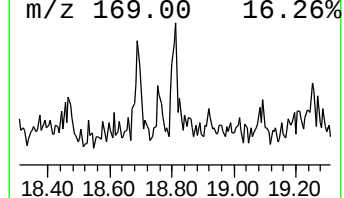
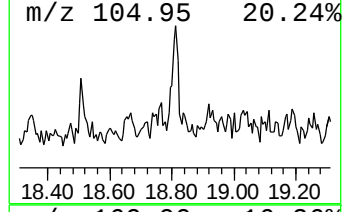
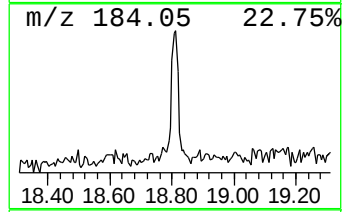
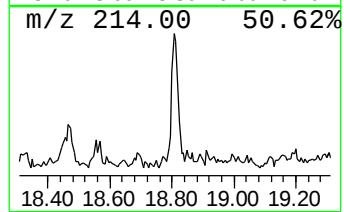
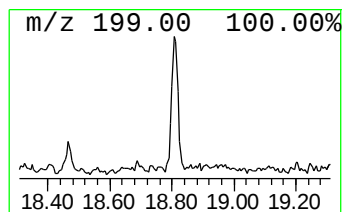
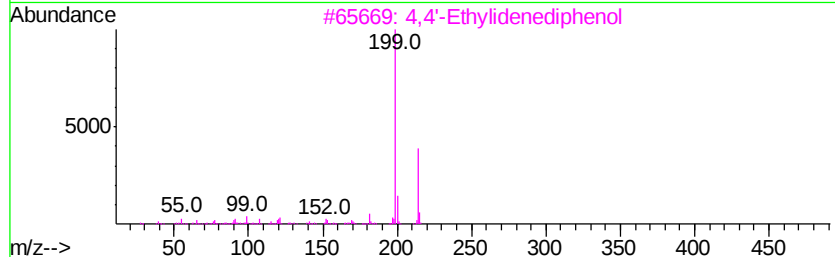
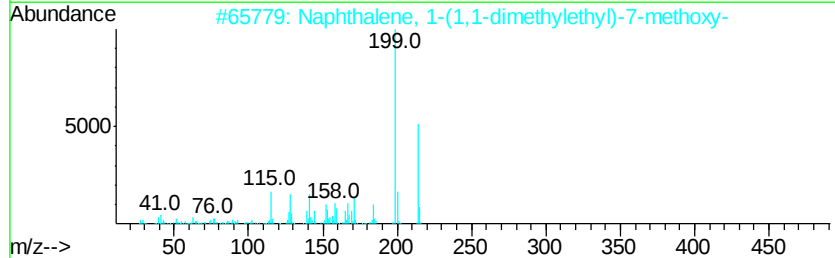
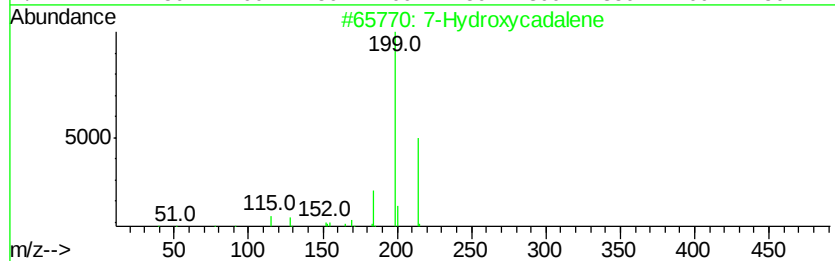
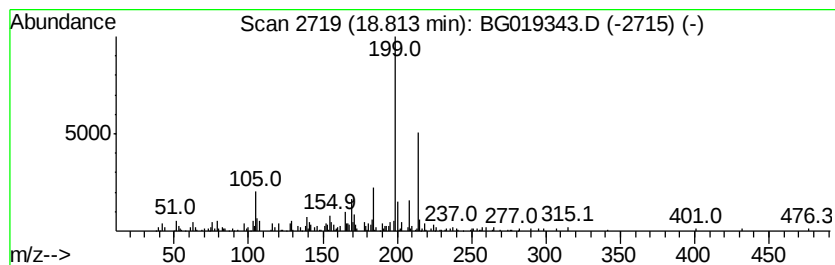
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 7-Hydroxycadalene Concentration Rank 30

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.81 | 2.13 ng/ul | 115445 | Phenanthrene-d10 | 17.60 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 7-Hydroxycadalene                   | 214 | C15H18O  | 002102-75-2  | 87   |
| 2    |    |   | Naphthalene, 1-(1,1-dimethylethy... | 214 | C15H18O  | 060683-42-3  | 68   |
| 3    |    |   | 4,4'-Ethylidenediphenol             | 214 | C14H14O2 | 002081-08-5  | 53   |
| 4    |    |   | 1,2,4-Benzenetriamine, N1-phenyl-   | 199 | C12H13N3 | 000136-17-4  | 49   |
| 5    |    |   | 5,7,7-Trimethyl-4,5,7,8-tetrahyd... | 214 | C15H18O  | 1000147-53-3 | 47   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

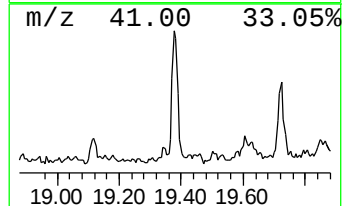
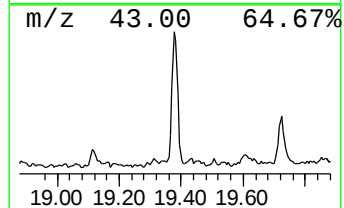
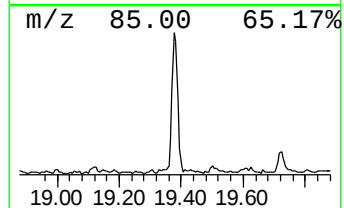
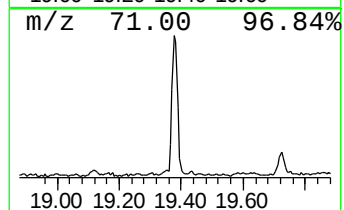
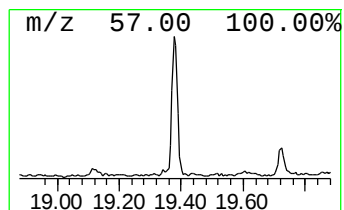
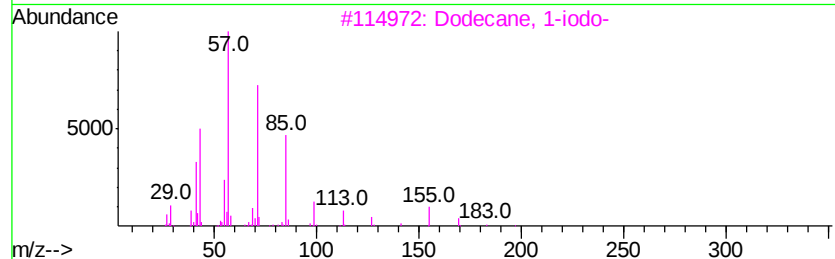
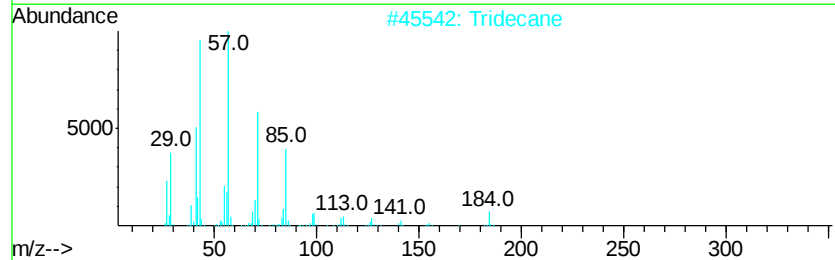
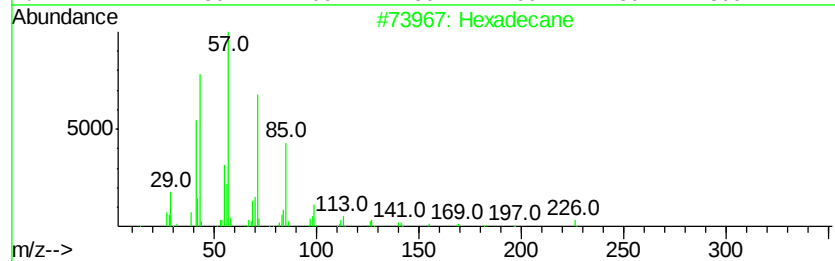
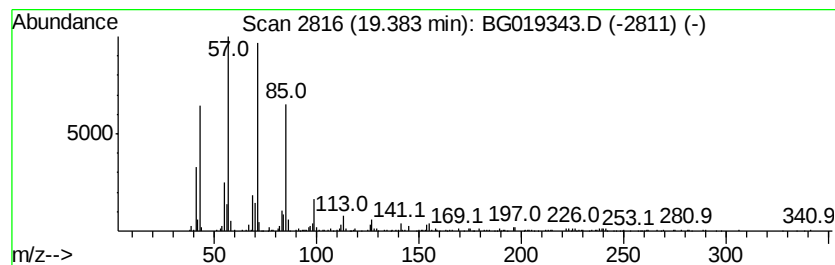
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 12

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.38 | 8.09 ng/ul | 437751 | Phenanthrene-d10 | 17.60 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane          | 226 | C16H34  | 000544-76-3 | 83   |
| 2    |    | Tridecane           | 184 | C13H28  | 000629-50-5 | 80   |
| 3    |    | Dodecane, 1-iodo-   | 296 | C12H25I | 004292-19-7 | 80   |
| 4    |    | Dodecane, 4-methyl- | 184 | C13H28  | 006117-97-1 | 80   |
| 5    |    | Heptacosane         | 380 | C27H56  | 000593-49-7 | 64   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

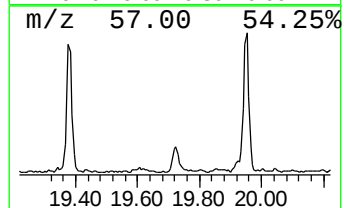
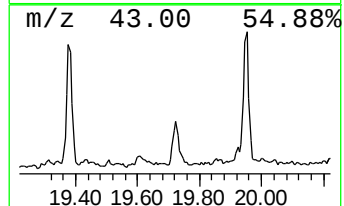
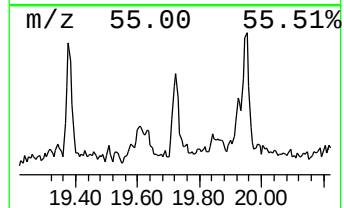
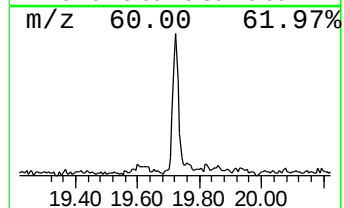
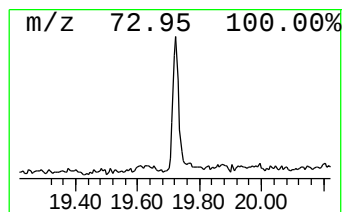
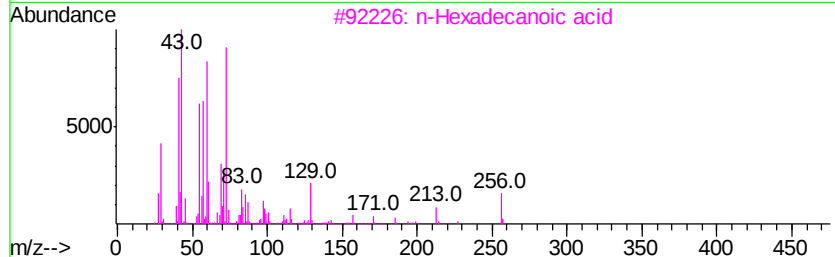
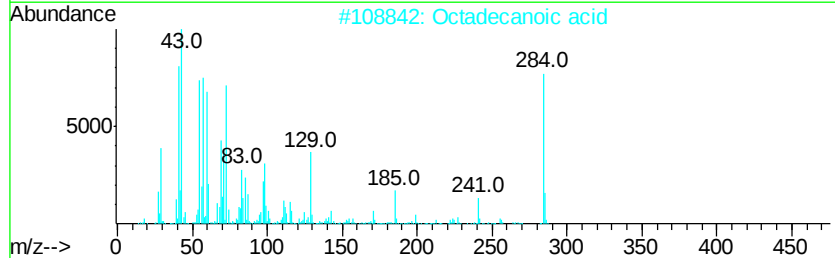
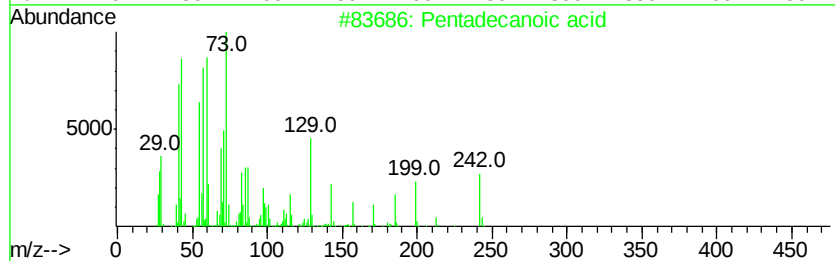
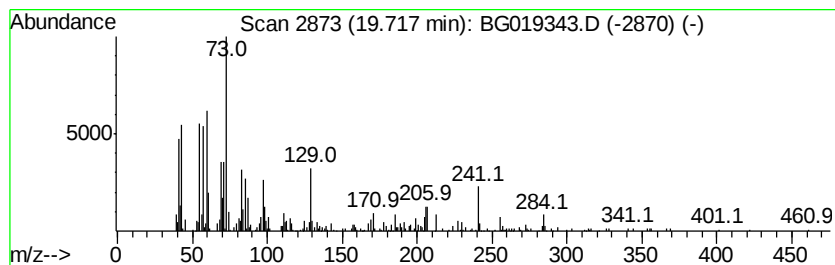
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Pentadecanoic acid Concentration Rank 20

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.72 | 5.49 ng/ul | 297335 | Phenanthrene-d10 | 17.60 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|---------------------|-----|----------|-------------|------|
| 1    |    |   | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 78   |
| 2    |    |   | Octadecanoic acid   | 284 | C18H36O2 | 000057-11-4 | 60   |
| 3    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 55   |
| 4    |    |   | Octadecanoic acid   | 284 | C18H36O2 | 000057-11-4 | 53   |
| 5    |    |   | Octadecanoic acid   | 284 | C18H36O2 | 000057-11-4 | 38   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

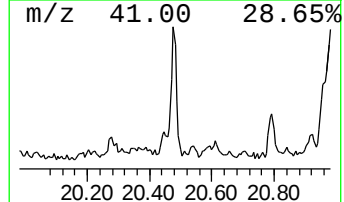
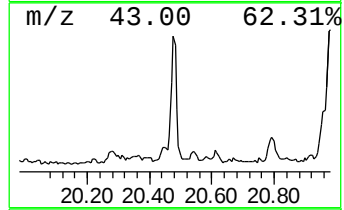
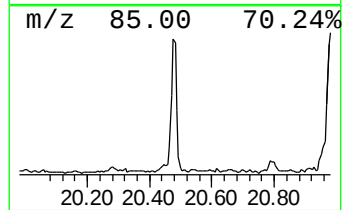
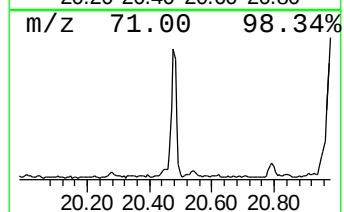
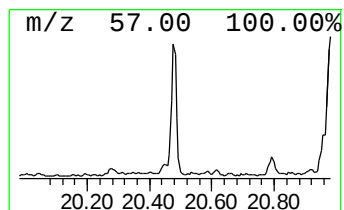
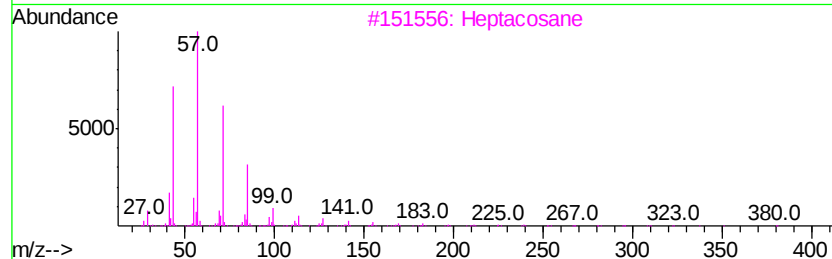
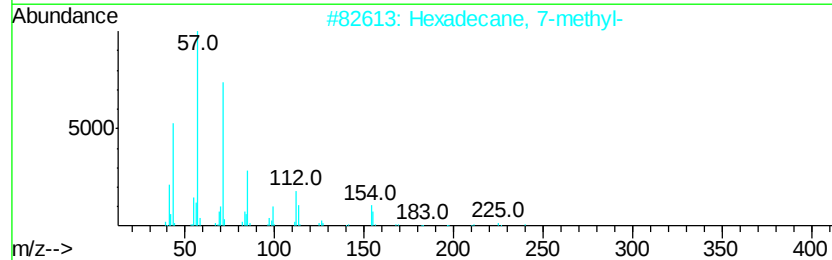
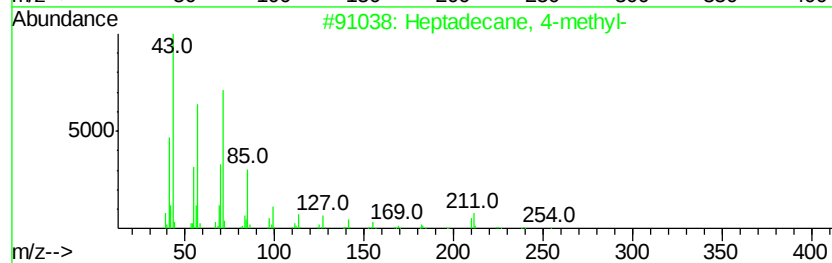
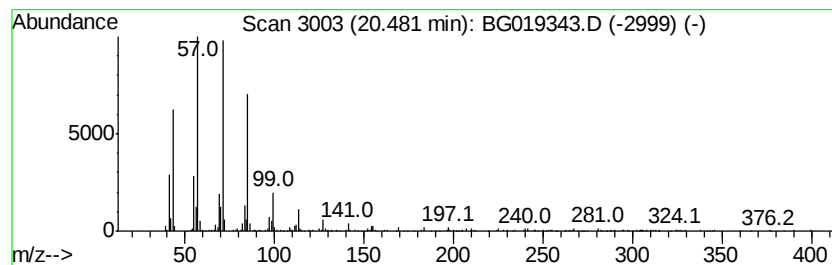
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 17

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.48 | 6.15 ng/ul | 519913 | Chrysene-d12     | 21.89 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane, 4-methyl- | 254 | C18H38  | 026429-11-8 | 80   |
| 2    |    |   | Hexadecane, 7-methyl-  | 240 | C17H36  | 026730-20-1 | 76   |
| 3    |    |   | Heptacosane            | 380 | C27H56  | 000593-49-7 | 74   |
| 4    |    |   | Nonane, 3,7-dimethyl-  | 156 | C11H24  | 017302-32-8 | 68   |
| 5    |    |   | Tridecane, 4-methyl-   | 198 | C14H30  | 026730-12-1 | 68   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

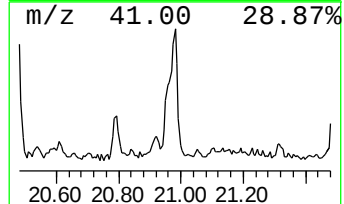
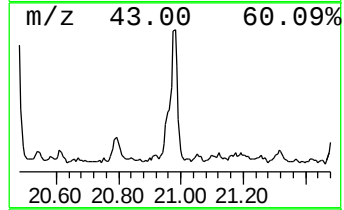
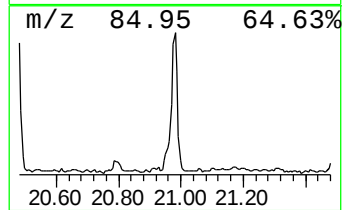
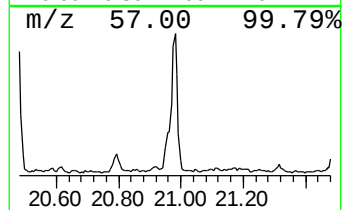
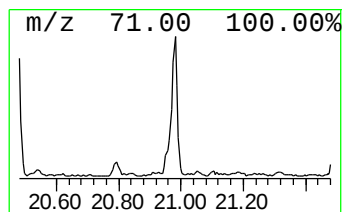
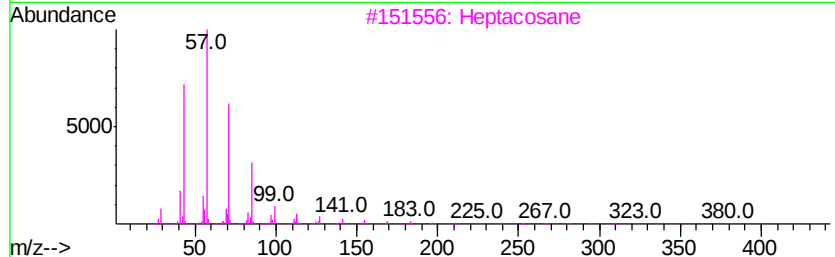
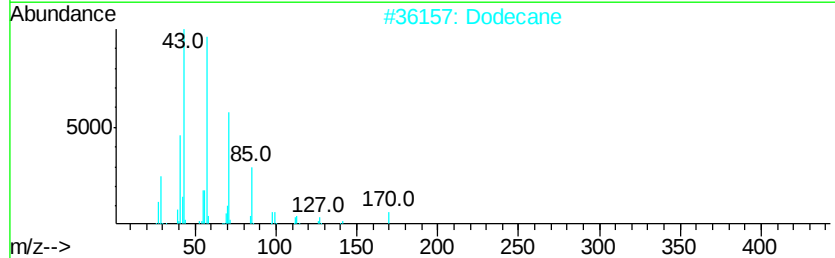
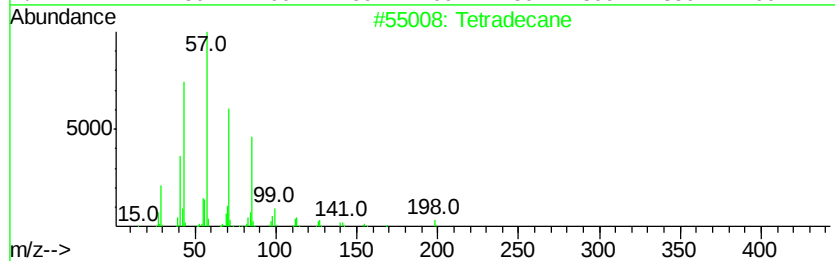
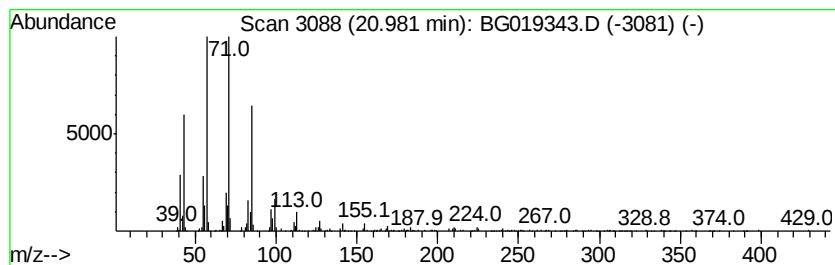
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 20.98 | 11.23 ng/ul | 949761 | Chrysene-d12     | 21.89 |

| Hit# | of | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------|-----|---------|-------------|------|
| 1    | 5  | Tetradecane       | 198 | C14H30  | 000629-59-4 | 93   |
| 2    |    | Dodecane          | 170 | C12H26  | 000112-40-3 | 87   |
| 3    |    | Heptacosane       | 380 | C27H56  | 000593-49-7 | 83   |
| 4    |    | Dodecane, 1-iodo- | 296 | C12H25I | 004292-19-7 | 80   |
| 5    |    | Tetracosane       | 338 | C24H50  | 000646-31-1 | 64   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

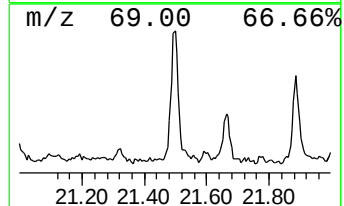
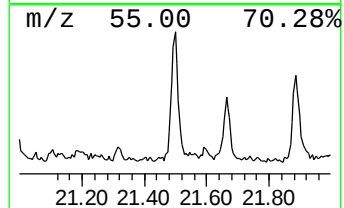
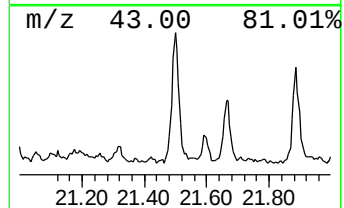
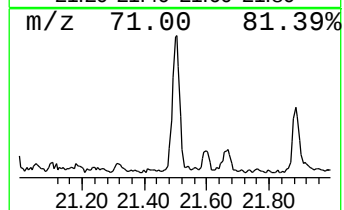
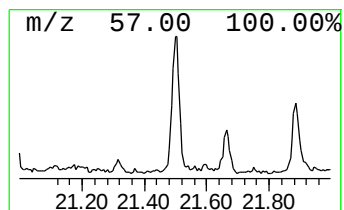
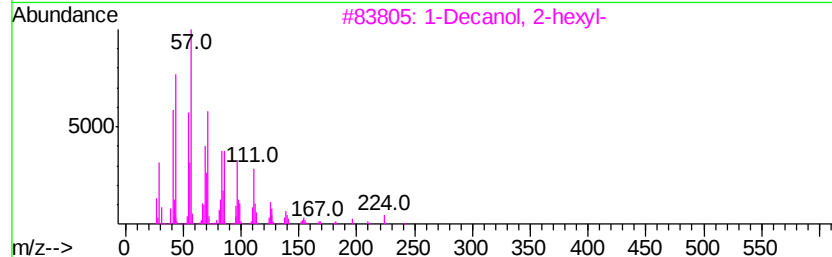
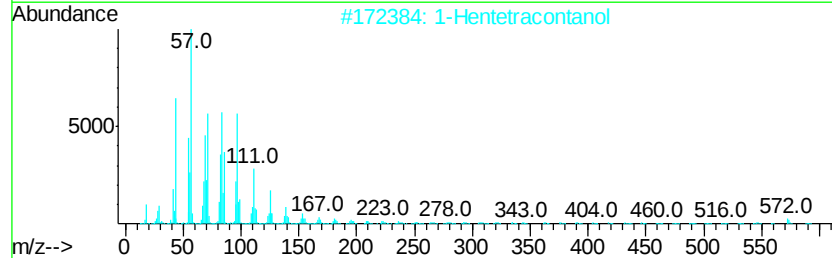
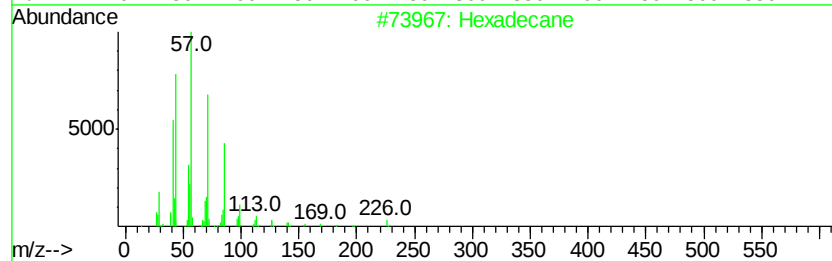
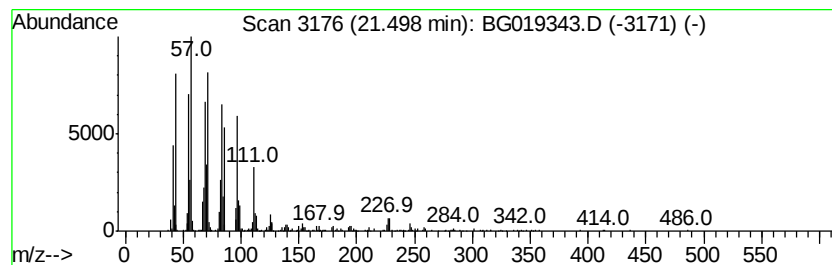
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 13

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 21.50 | 7.58 ng/ul | 640722 | Chrysene-d12     | 21.89 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Hexadecane                          | 226 | C16H34   | 000544-76-3 | 86   |
| 2    |    |   | 1-Hentetracontanol                  | 593 | C41H84O  | 040710-42-7 | 86   |
| 3    |    |   | 1-Decanol, 2-hexyl-                 | 242 | C16H34O  | 002425-77-6 | 74   |
| 4    |    |   | Decanoic acid, 5-ethyl-3,5,9-tri... | 256 | C16H32O2 | 056247-63-3 | 52   |
| 5    |    |   | Cyclotetracosane                    | 336 | C24H48   | 000297-03-0 | 49   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

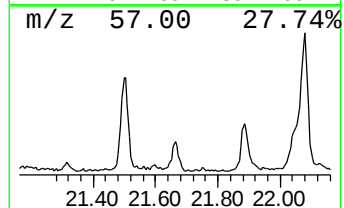
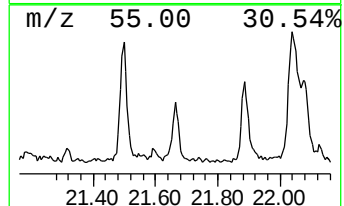
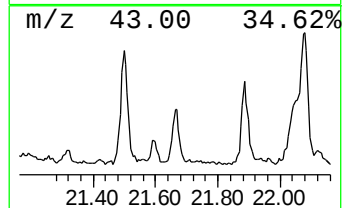
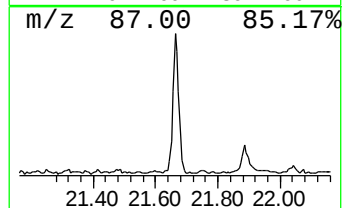
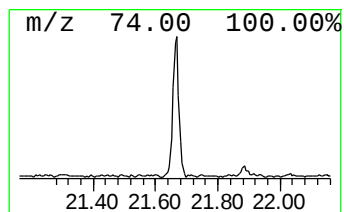
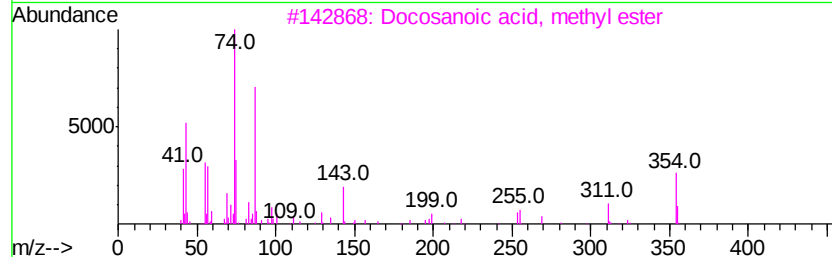
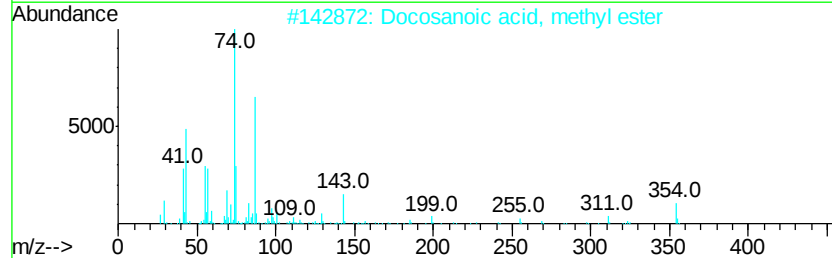
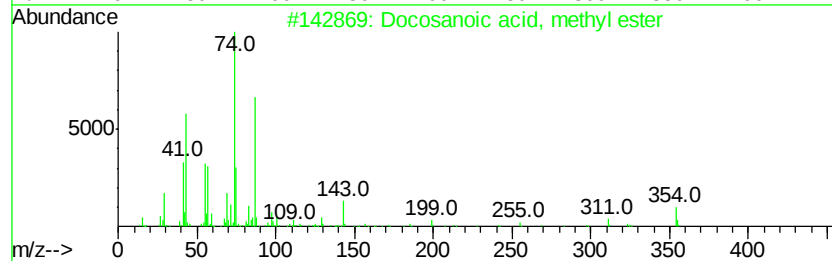
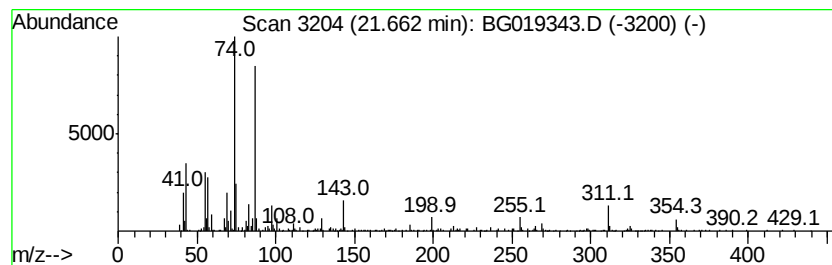
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 Docosanoic acid, methyl ester Concentration Rank 21

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 21.66 | 4.99 ng/ul | 422330 | Chrysene-d12     | 21.89 |

| Hit# | of | Tentative ID                    | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------------------|-----|----------|-------------|------|
| 1    | 5  | Docosanoic acid, methyl ester   | 354 | C23H46O2 | 000929-77-1 | 95   |
| 2    |    | Docosanoic acid, methyl ester   | 354 | C23H46O2 | 000929-77-1 | 94   |
| 3    |    | Docosanoic acid, methyl ester   | 354 | C23H46O2 | 000929-77-1 | 93   |
| 4    |    | Octadecanoic acid, methyl ester | 298 | C19H38O2 | 000112-61-8 | 91   |
| 5    |    | Docosanoic acid, methyl ester   | 354 | C23H46O2 | 000929-77-1 | 89   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

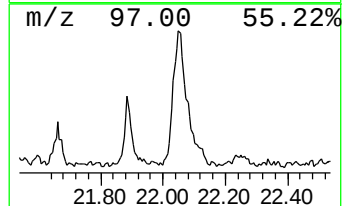
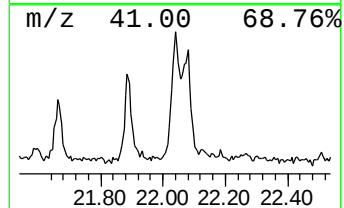
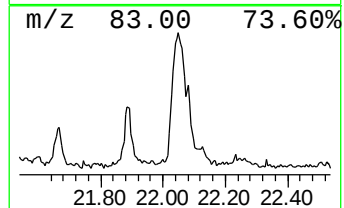
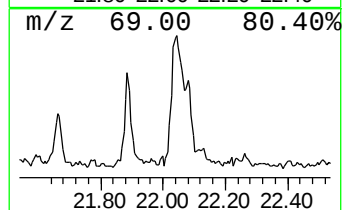
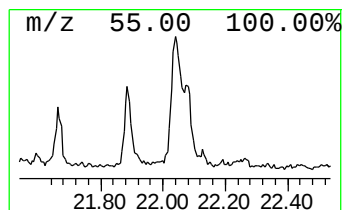
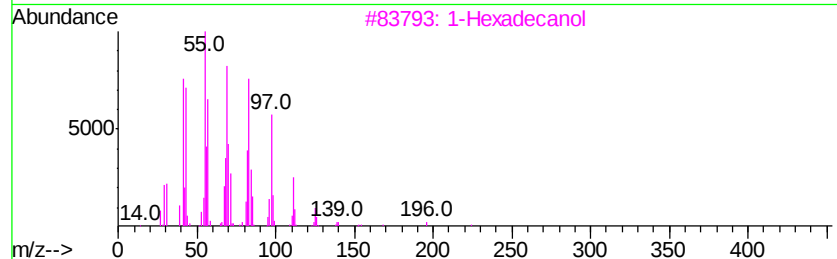
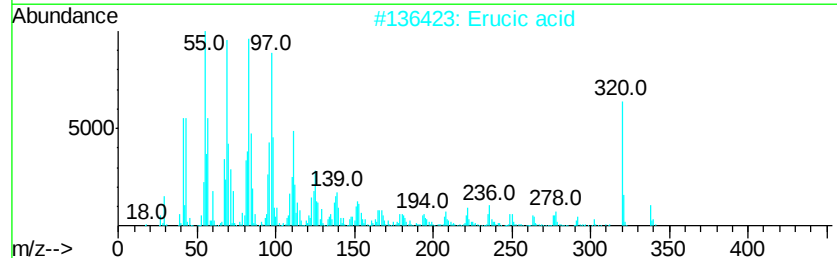
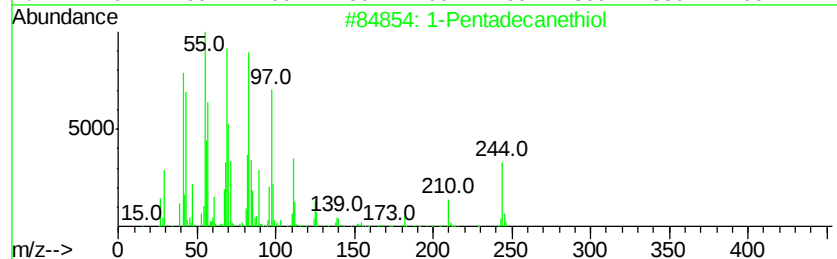
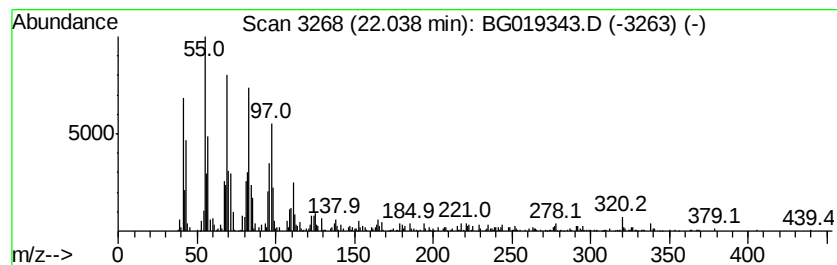
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 1-Pentadecanethiol Concentration Rank 11

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 22.04 | 9.13 ng/ul | 772006 | Chrysene-d12     | 21.89 |

| Hit# | of | 5 | Tentative ID       | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------|-----|----------|-------------|------|
| 1    |    |   | 1-Pentadecanethiol | 244 | C15H32S  | 025276-70-4 | 91   |
| 2    |    |   | Erucic acid        | 338 | C22H42O2 | 000112-86-7 | 87   |
| 3    |    |   | 1-Hexadecanol      | 242 | C16H34O  | 036653-82-4 | 83   |
| 4    |    |   | 9-Octadecene, (E)- | 252 | C18H36   | 007206-25-9 | 81   |
| 5    |    |   | 5-Octadecene, (E)- | 252 | C18H36   | 007206-21-5 | 72   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

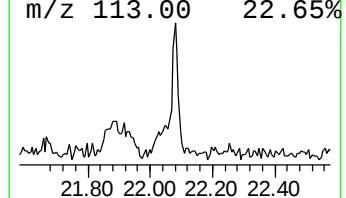
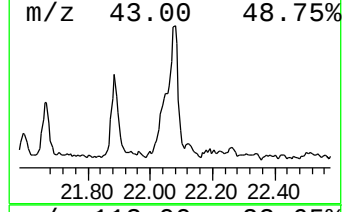
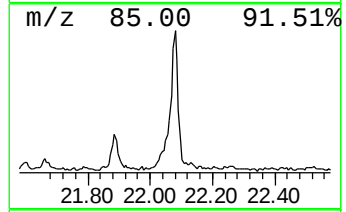
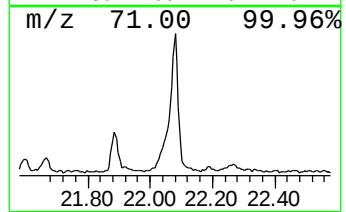
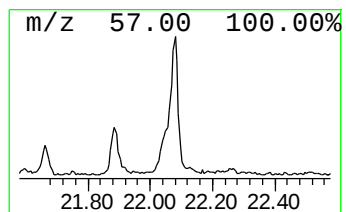
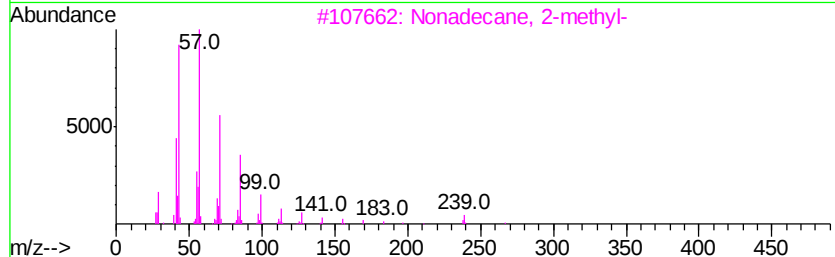
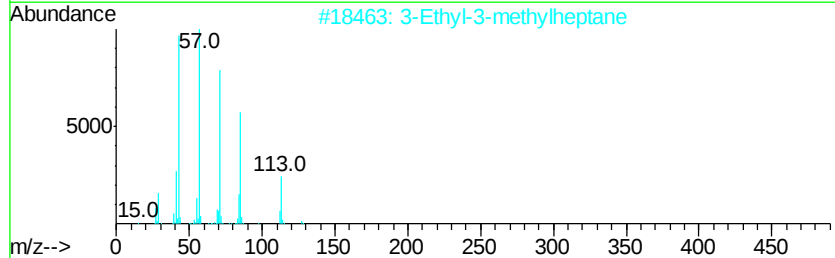
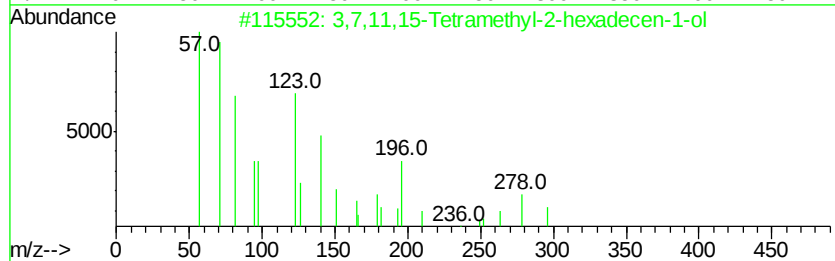
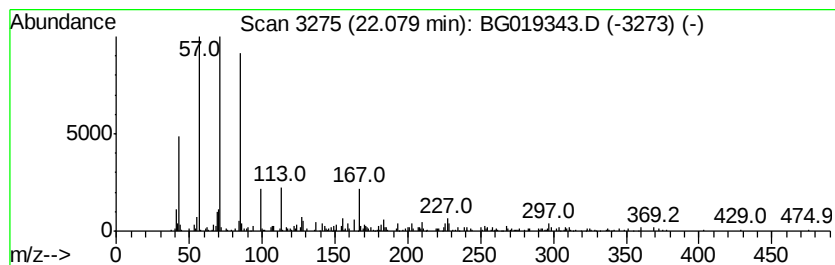
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 3,7,11,15-Tetramethyl-2-hex... Concentration Rank 19

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 22.08 | 5.71 ng/ul | 483225 | Chrysene-d12     | 21.89 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 3,7,11,15-Tetramethyl-2-hexadec... | 296 | C20H400 | 102608-53-7 | 94   |
| 2    |    |   | 3-Ethyl-3-methylheptane            | 142 | C10H22  | 017302-01-1 | 59   |
| 3    |    |   | Nonadecane, 2-methyl-              | 282 | C20H42  | 001560-86-7 | 45   |
| 4    |    |   | Octadecane, 1-iodo-                | 380 | C18H37I | 000629-93-6 | 45   |
| 5    |    |   | Heptacosane                        | 380 | C27H56  | 000593-49-7 | 43   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

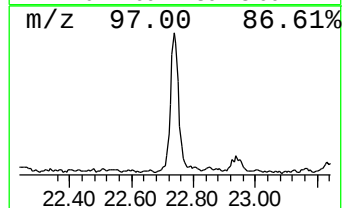
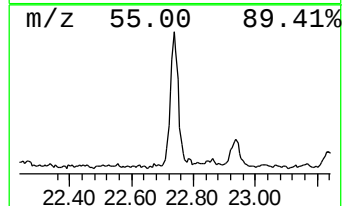
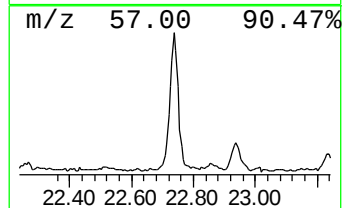
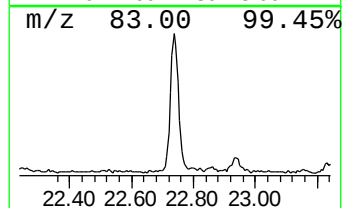
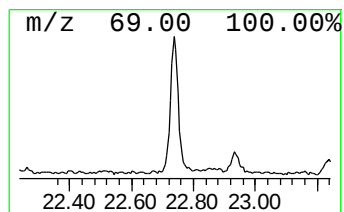
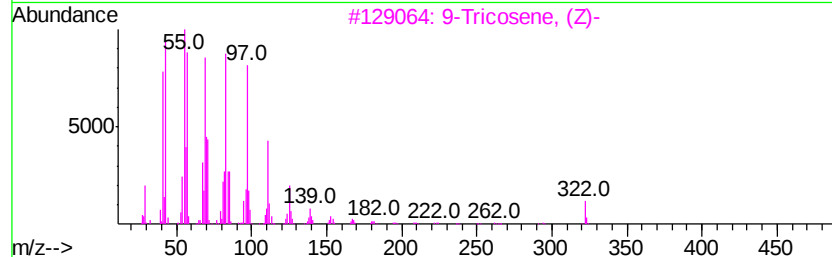
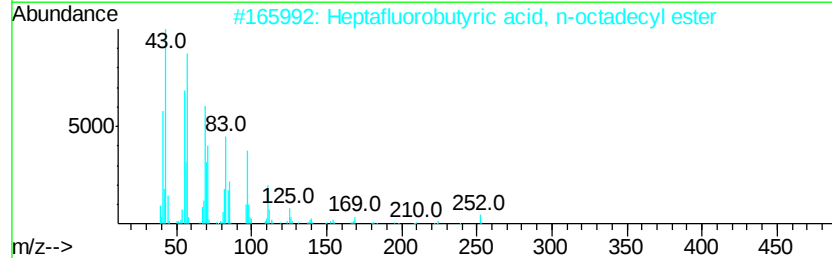
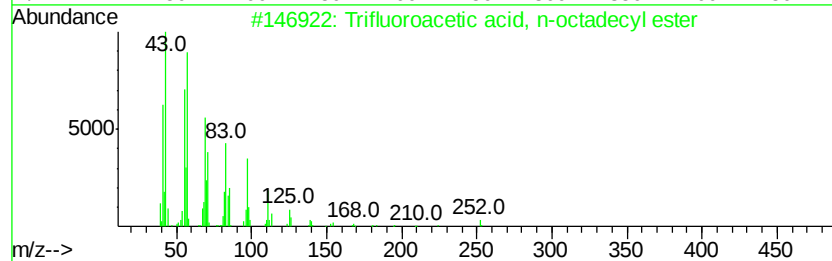
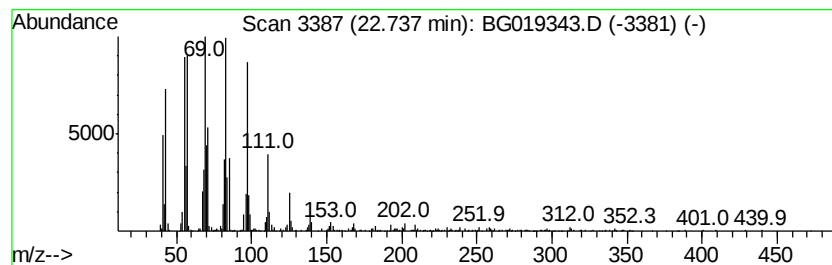
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 23 Trifluoroacetic acid, n-oct... Concentration Rank 5

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 22.74 | 14.68 ng/ul | 1241800 | Chrysene-d12     | 21.89 |

| Hit# | of | Tentative ID                         | MW  | MolForm    | CAS#        | Qual |
|------|----|--------------------------------------|-----|------------|-------------|------|
| 1    | 5  | Trifluoroacetic acid, n-octadecyl... | 366 | C20H37F3O2 | 079392-43-1 | 91   |
| 2    |    | Heptafluorobutyric acid, n-octad...  | 466 | C22H37F7O2 | 000400-57-7 | 91   |
| 3    |    | 9-Tricosene, (Z)-                    | 322 | C23H46     | 027519-02-4 | 91   |
| 4    |    | 1-Heneicosyl formate                 | 340 | C22H44O2   | 077899-03-7 | 90   |
| 5    |    | 1-Octadecene                         | 252 | C18H36     | 000112-88-9 | 83   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

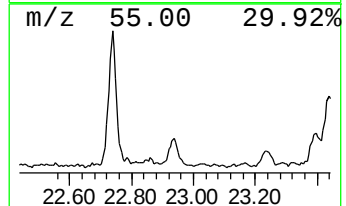
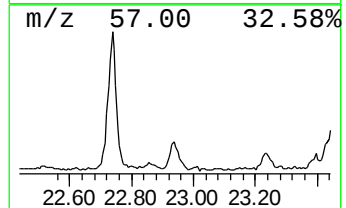
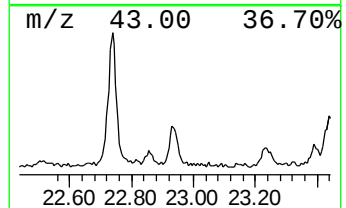
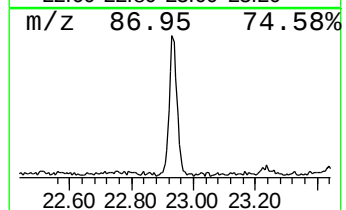
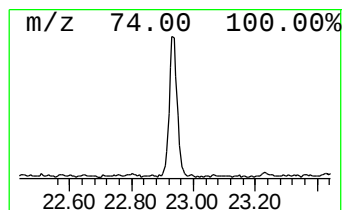
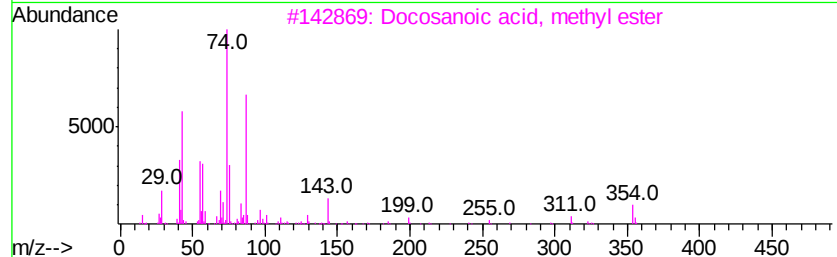
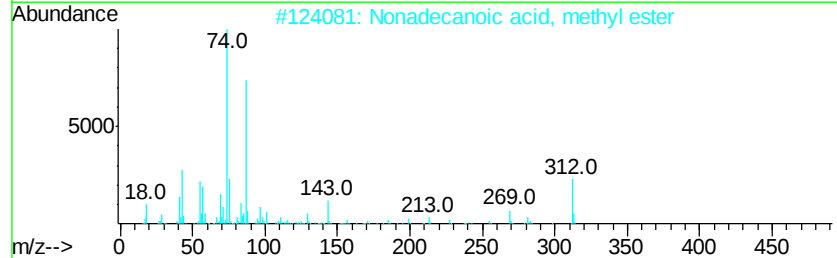
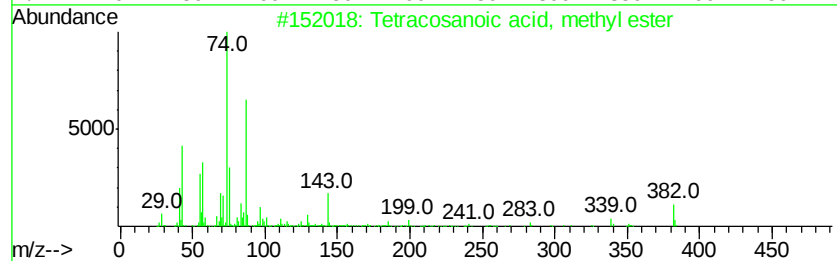
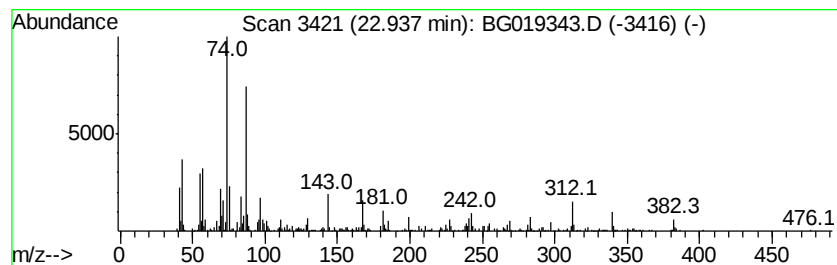
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 24 Tetracosanoic acid, methyl ... Concentration Rank 15

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 22.94 | 6.33 ng/ul | 535421 | Chrysene-d12     | 21.89 |

| Hit# | of | Tentative ID                     | MW  | MolForm  | CAS#        | Qual |
|------|----|----------------------------------|-----|----------|-------------|------|
| 1    | 5  | Tetracosanoic acid, methyl ester | 382 | C25H50O2 | 002442-49-1 | 97   |
| 2    |    | Nonadecanoic acid, methyl ester  | 312 | C20H40O2 | 001731-94-8 | 95   |
| 3    |    | Docosanoic acid, methyl ester    | 354 | C23H46O2 | 000929-77-1 | 95   |
| 4    |    | Nonadecanoic acid, methyl ester  | 312 | C20H40O2 | 001731-94-8 | 93   |
| 5    |    | Docosanoic acid, methyl ester    | 354 | C23H46O2 | 000929-77-1 | 93   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

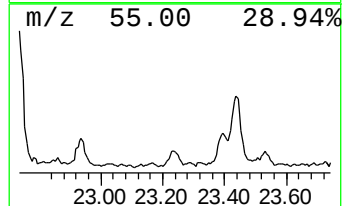
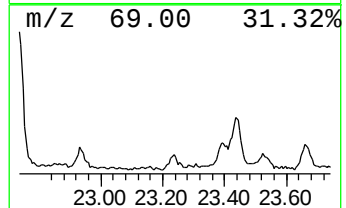
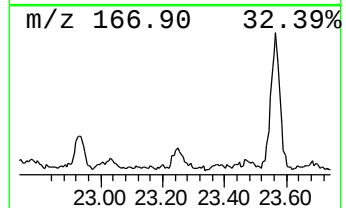
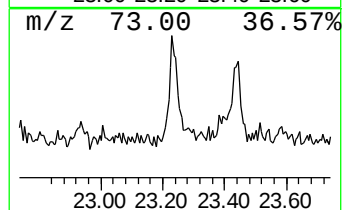
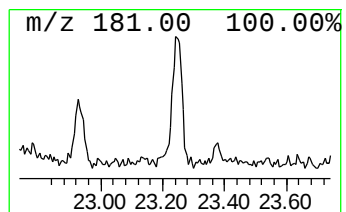
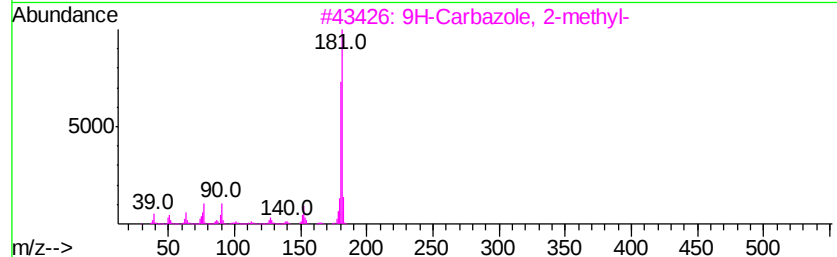
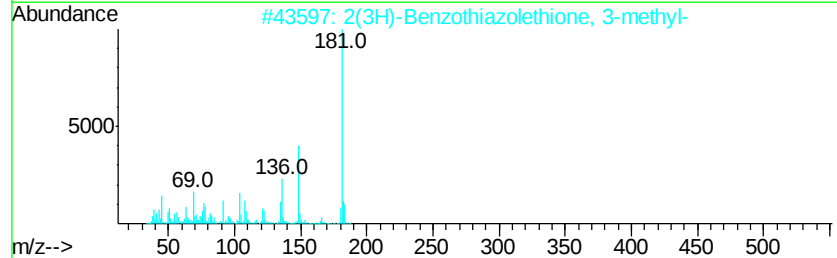
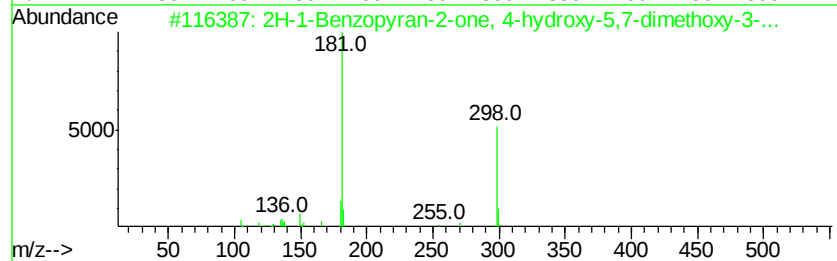
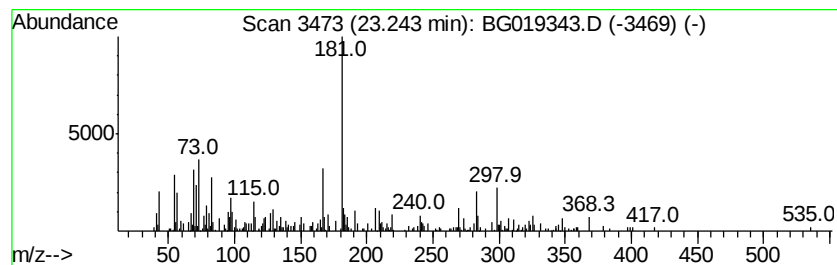
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 25 unknown-02 Concentration Rank 28

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 23.24 | 2.56 ng/ul | 216454 | Chrysene-d12     | 21.89 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 2H-1-Benzopyran-2-one, 4-hydroxy... | 298 | C17H14O5 | 004222-00-8  | 40   |
| 2    |    | 2(3H)-Benzothiazolethione, 3-met... | 181 | C8H7NS2  | 002254-94-6  | 38   |
| 3    |    | 9H-Carbazole, 2-methyl-             | 181 | C13H11N  | 003652-91-3  | 35   |
| 4    |    | Benzaldehyde, 4-[1-[4-(acetyloxy... | 374 | C20H22O7 | 055525-48-9  | 35   |
| 5    |    | 8-Methoxy-[1,2,4]triazolo[4,3-a]... | 181 | C7H7N3OS | 1000186-80-8 | 35   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

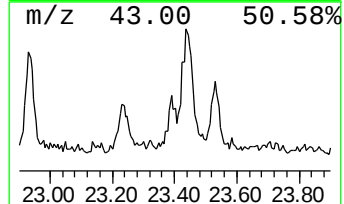
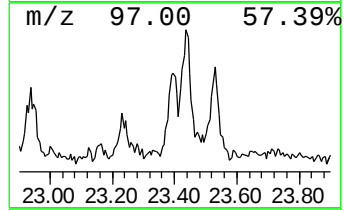
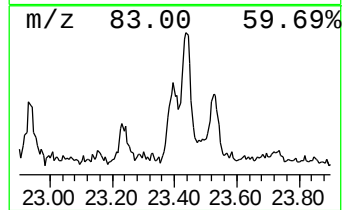
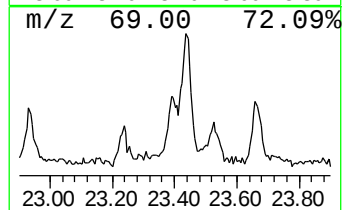
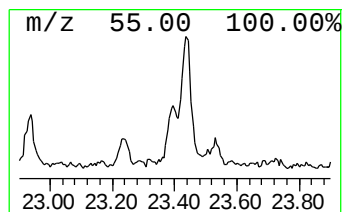
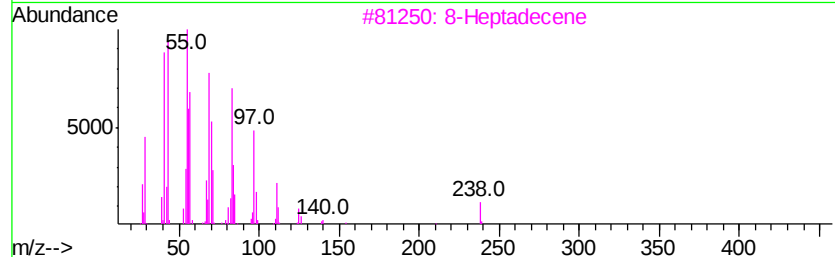
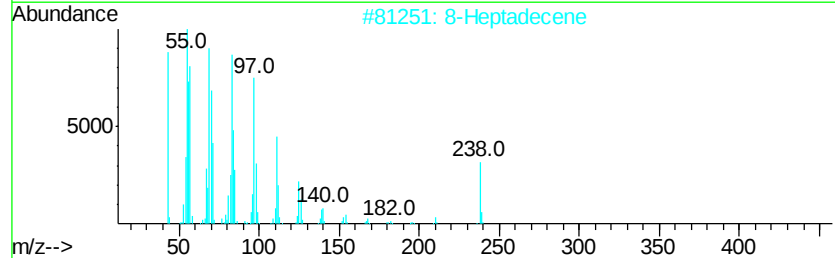
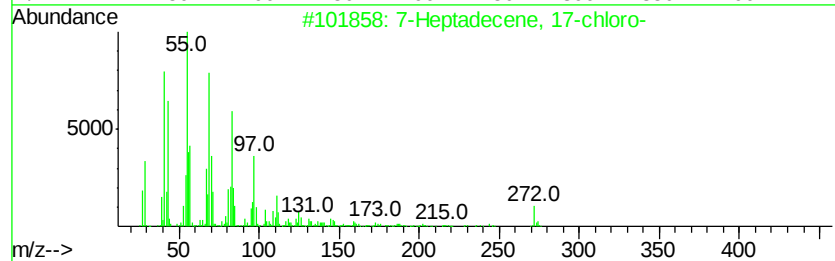
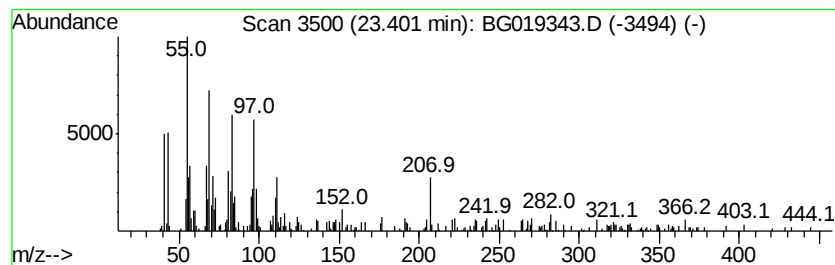
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 26 7-Heptadecene, 17-chloro- Concentration Rank 27

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 23.40 | 2.85 ng/ul | 240621 | Chrysene-d12     | 21.89 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm  | CAS#        | Qual |
|------|----|---|---------------------------|-----|----------|-------------|------|
| 1    |    |   | 7-Heptadecene, 17-chloro- | 272 | C17H33Cl | 056554-79-1 | 90   |
| 2    |    |   | 8-Heptadecene             | 238 | C17H34   | 002579-04-6 | 87   |
| 3    |    |   | 8-Heptadecene             | 238 | C17H34   | 054290-12-9 | 87   |
| 4    |    |   | 8-Heptadecene             | 238 | C17H34   | 054290-12-9 | 80   |
| 5    |    |   | Cycloeicosane             | 280 | C20H40   | 000296-56-0 | 64   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

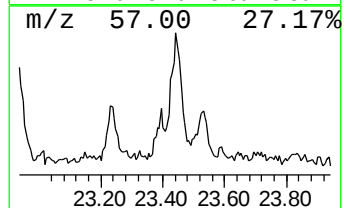
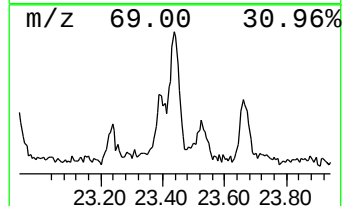
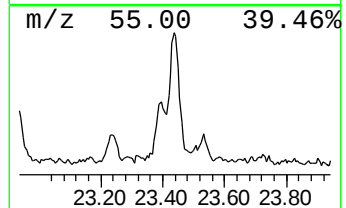
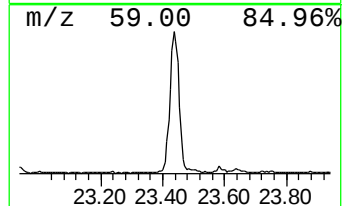
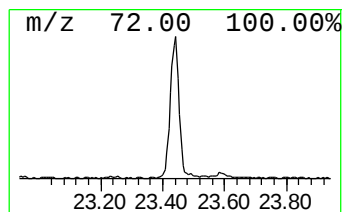
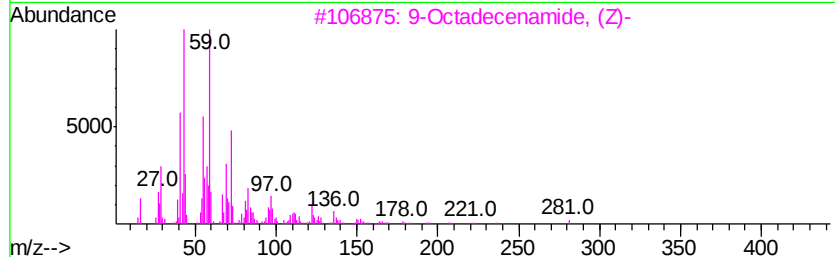
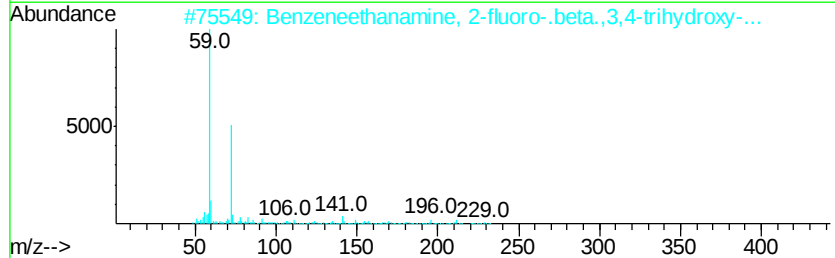
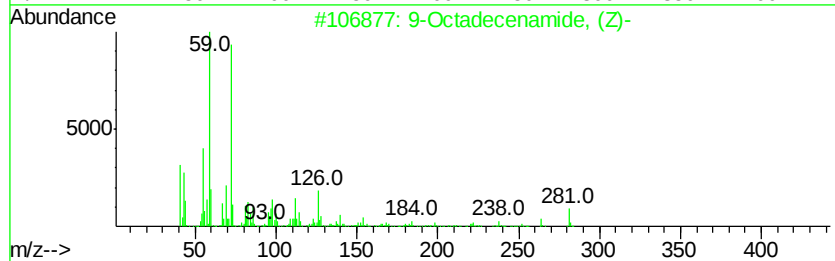
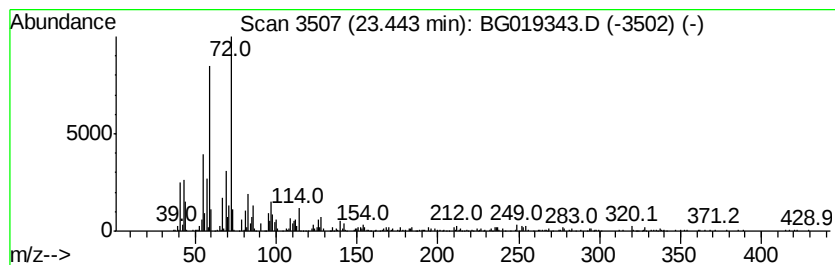
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 27 9-Octadecenamide, (Z)- Concentration Rank 9

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 23.44 | 10.02 ng/ul | 847700 | Chrysene-d12     | 21.89 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | 9-Octadecenamide, (Z)-              | 281 | C18H35NO   | 000301-02-0 | 59   |
| 2    |    | Benzeneethanamine, 2-fluoro-.bet... | 229 | C11H16FN03 | 061338-98-5 | 50   |
| 3    |    | 9-Octadecenamide, (Z)-              | 281 | C18H35NO   | 000301-02-0 | 41   |
| 4    |    | Tetradecanamide                     | 227 | C14H29NO   | 000638-58-4 | 38   |
| 5    |    | Decanamide-                         | 171 | C10H21NO   | 002319-29-1 | 37   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

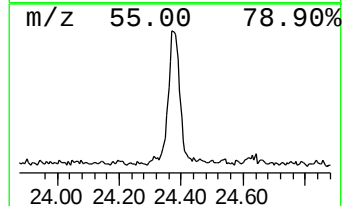
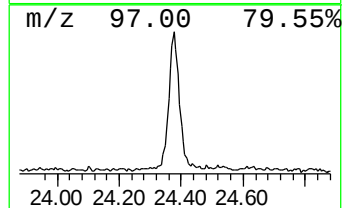
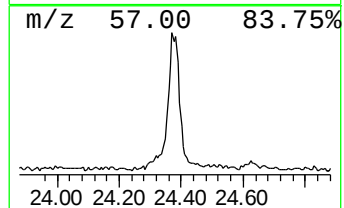
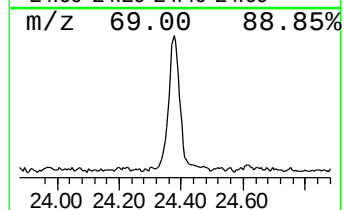
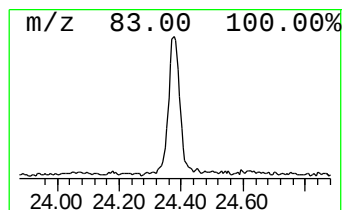
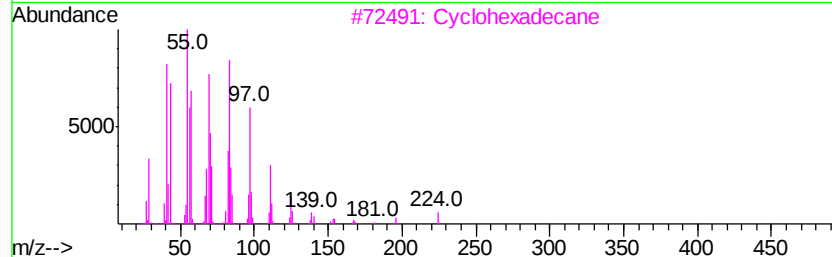
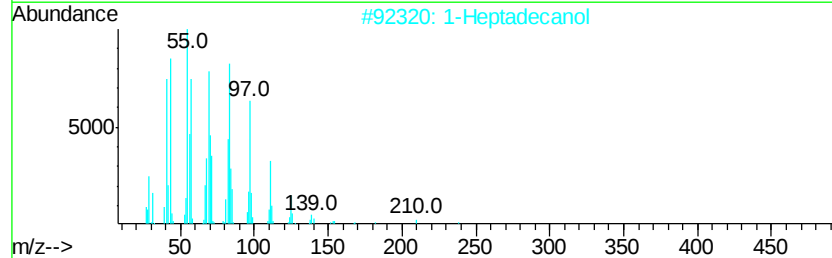
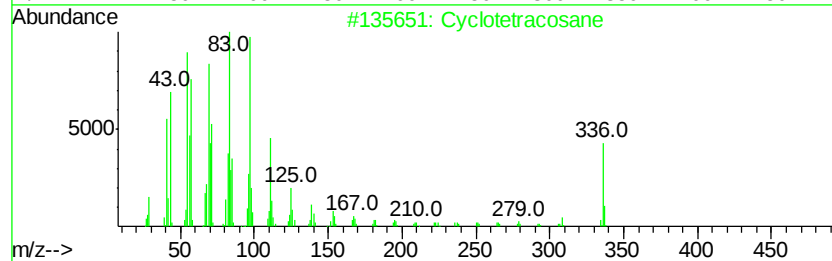
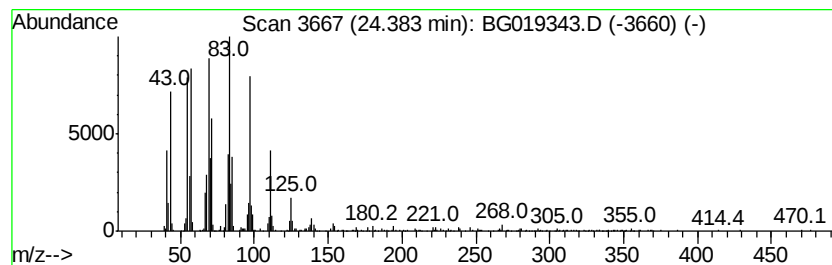
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 28 (DEL) Alkane: Cyclic24.38 Concentration Rank 2

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 24.38 | 20.68 ng/ul | 1050320 | Perylene-d12     | 25.29 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|-------------------------------------|-----|-------------|-------------|------|
| 1    | 5  | Cyclotetracosane                    | 336 | C24H48      | 000297-03-0 | 96   |
| 2    |    | 1-Heptadecanol                      | 256 | C17H36O     | 001454-85-9 | 90   |
| 3    |    | Cyclohexadecane                     | 224 | C16H32      | 000295-65-8 | 87   |
| 4    |    | Trichloroacetic acid, hexadecyl ... | 386 | C18H33Cl3O2 | 074339-54-1 | 87   |
| 5    |    | 1-Docosene                          | 308 | C22H44      | 001599-67-3 | 76   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

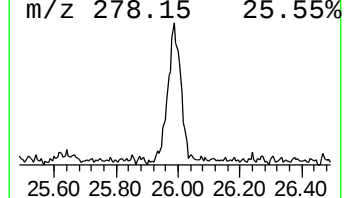
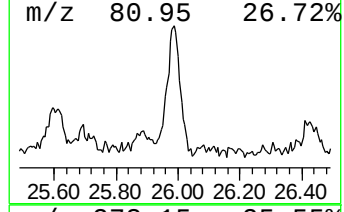
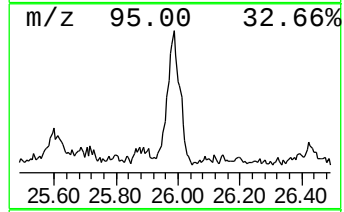
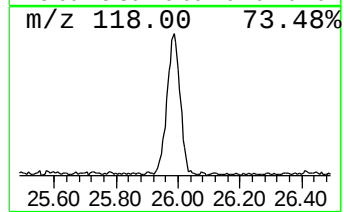
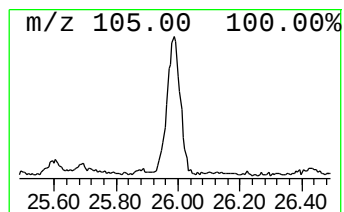
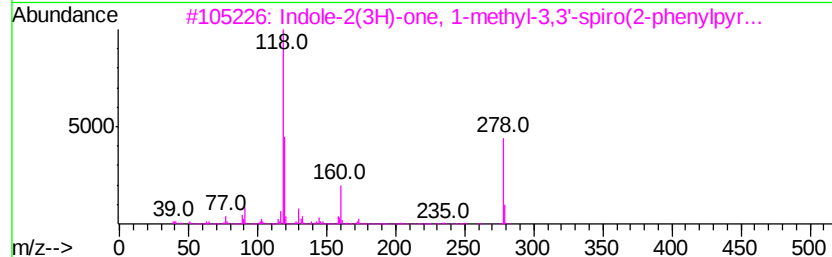
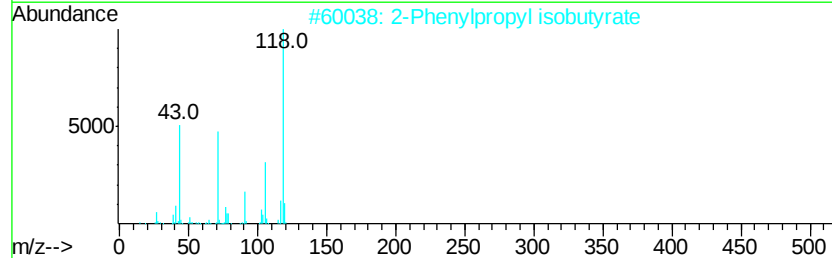
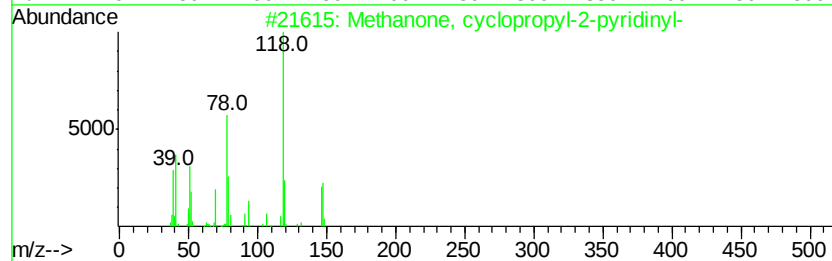
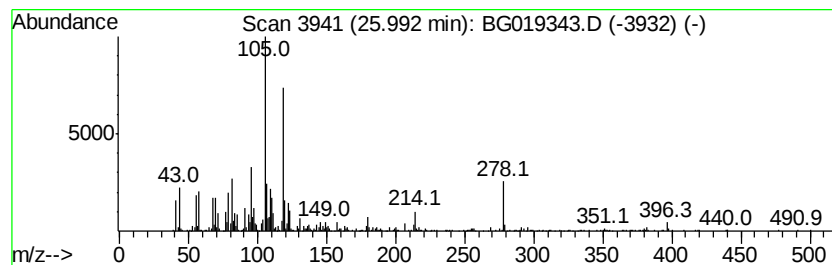
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 29 unknown-03 Concentration Rank 3

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 25.99 | 18.20 ng/ul | 924390 | Perylene-d12     | 25.29 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Methanone, cyclopropyl-2-pyridinyl- | 147 | C9H9NO    | 057276-28-5 | 35   |
| 2    |    | 2-Phenylpropyl isobutyrate          | 206 | C13H18O2  | 065813-53-8 | 22   |
| 3    |    | Indole-2(3H)-one, 1-methyl-3,3'-... | 278 | C18H18N2O | 132962-61-9 | 17   |
| 4    |    | Methanamine, N-(phenylmethyl)-      | 119 | C8H9N     | 000622-29-7 | 16   |
| 5    |    | Benzoic acid, 2,5-dimethyl-, (2,... | 268 | C18H20O2  | 055000-49-2 | 12   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG102315\  
Data File : BG019343.D  
Acq On : 24 Oct 2015 19:45  
Operator : UM/NP  
Sample : G4058-11  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
E5LR9

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102315.M  
Quant Title : SVOA CALIBRATION

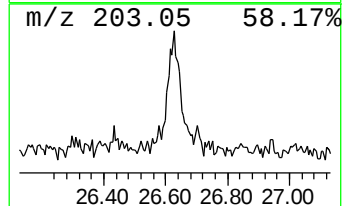
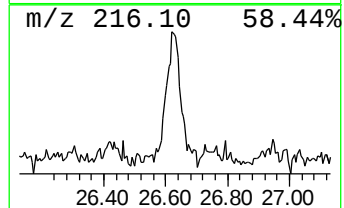
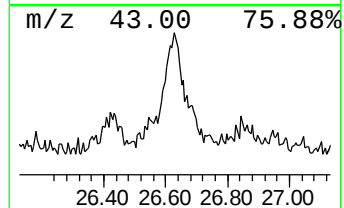
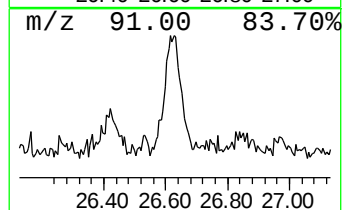
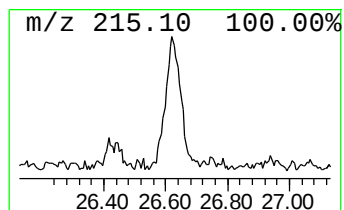
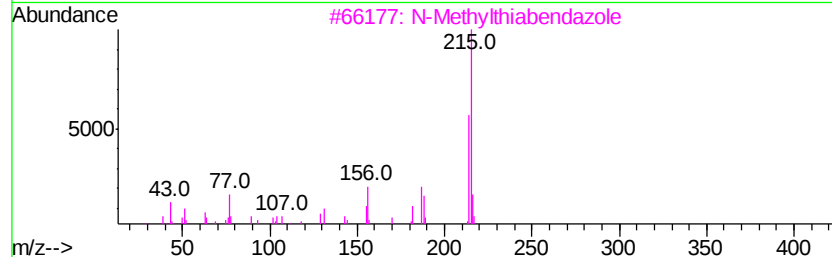
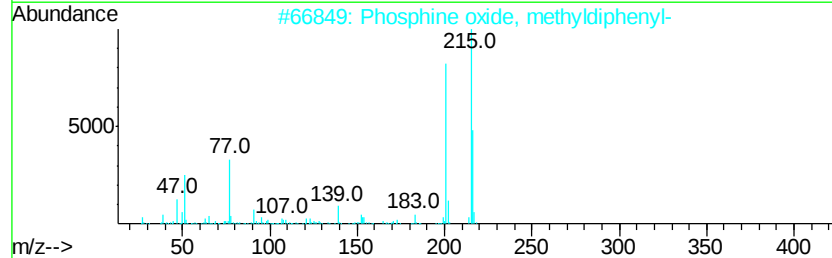
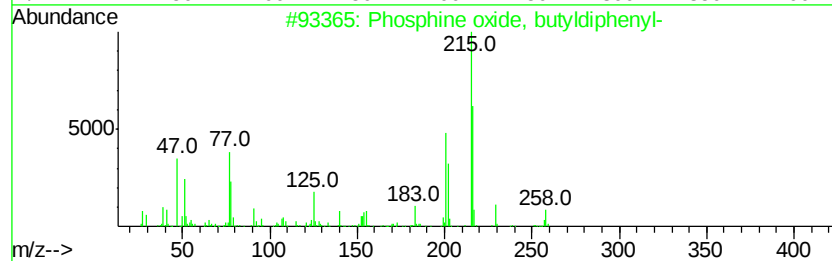
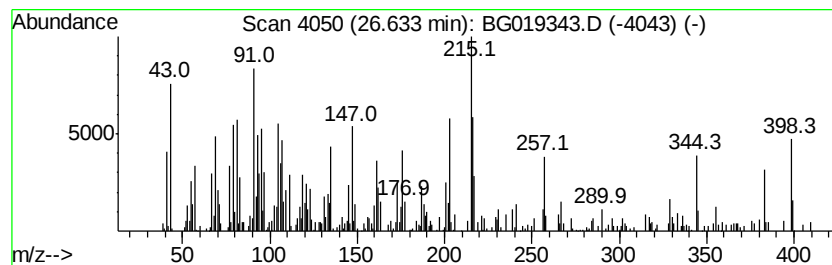
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 30 unknown-04 Concentration Rank 4

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 26.63 | 16.66 ng/ul | 846149 | Perylene-d12     | 25.29 |

| Hit# | of | Tentative ID                     | MW  | MolForm  | CAS#        | Qual |
|------|----|----------------------------------|-----|----------|-------------|------|
| 1    | 5  | Phosphine oxide, butyldiphenyl-  | 258 | C16H19OP | 004233-13-0 | 30   |
| 2    |    | Phosphine oxide, methyldiphenyl- | 216 | C13H13OP | 002129-89-7 | 18   |
| 3    |    | N-Methylthiabendazole            | 215 | C11H9N3S | 032594-70-0 | 18   |
| 4    |    | Phosphine oxide, methyldiphenyl- | 216 | C13H13OP | 002129-89-7 | 14   |
| 5    |    | 2-Phenoxathinamine               | 215 | C12H9NOS | 010350-07-9 | 11   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG102315\  
 Data File : BG019343.D  
 Acq On : 24 Oct 2015 19:45  
 Operator : UM/NP  
 Sample : G4058-11  
 Misc :  
 ALS Vial : 33 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 E5LR9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG102315.M  
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name      | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|-----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                       |       |         |       |          | #                     | RT    | Resp    | Conc |
| (DEL) Alkane: Cyc...  | 3.21  | 9.6     | ng/ul | 133381   | 1                     | 8.24  | 278042  | 20.0 |
| Butane, 2-methoxy...  | 3.29  | 48.0    | ng/ul | 667527   | 1                     | 8.24  | 278042  | 20.0 |
| 1,3,5-Cycloheptat...  | 4.35  | 11.3    | ng/ul | 157096   | 1                     | 8.24  | 278042  | 20.0 |
| unknown-01            | 15.24 | 3.7     | ng/ul | 143662   | 3                     | 14.86 | 768944  | 20.0 |
| 2,6-Dimethoxybenz...  | 15.39 | 3.2     | ng/ul | 123887   | 3                     | 14.86 | 768944  | 20.0 |
| (DEL) Alkane: Str...  | 15.68 | 2.5     | ng/ul | 97543    | 3                     | 14.86 | 768944  | 20.0 |
| 2',4'-Dihydroxyac...  | 15.78 | 6.8     | ng/ul | 259777   | 3                     | 14.86 | 768944  | 20.0 |
| Benzaldehyde, 4-h...  | 16.27 | 4.0     | ng/ul | 214691   | 4                     | 17.60 | 1082710 | 20.0 |
| Ethanone, 1-(4-hy...  | 16.87 | 6.2     | ng/ul | 332987   | 4                     | 17.60 | 1082710 | 20.0 |
| (DEL) Alkane: Str...  | 17.33 | 3.8     | ng/ul | 203904   | 4                     | 17.60 | 1082710 | 20.0 |
| (DEL) Alkane: Str...  | 18.07 | 3.5     | ng/ul | 188993   | 4                     | 17.60 | 1082710 | 20.0 |
| n-Hexadecanoic acid   | 18.47 | 13.1    | ng/ul | 707612   | 4                     | 17.60 | 1082710 | 20.0 |
| (DEL) Alkane: Str...  | 18.75 | 5.9     | ng/ul | 319801   | 4                     | 17.60 | 1082710 | 20.0 |
| 7-Hydroxycadalene     | 18.81 | 2.1     | ng/ul | 115445   | 4                     | 17.60 | 1082710 | 20.0 |
| (DEL) Alkane: Str...  | 19.38 | 8.1     | ng/ul | 437751   | 4                     | 17.60 | 1082710 | 20.0 |
| Pentadecanoic acid    | 19.72 | 5.5     | ng/ul | 297335   | 4                     | 17.60 | 1082710 | 20.0 |
| (DEL) Alkane: Str...  | 20.48 | 6.2     | ng/ul | 519913   | 5                     | 21.89 | 1691460 | 20.0 |
| (DEL) Alkane: Str...  | 20.98 | 11.2    | ng/ul | 949761   | 5                     | 21.89 | 1691460 | 20.0 |
| (DEL) Alkane: Str...  | 21.50 | 7.6     | ng/ul | 640722   | 5                     | 21.89 | 1691460 | 20.0 |
| Docosanoic acid, ...  | 21.66 | 5.0     | ng/ul | 422330   | 5                     | 21.89 | 1691460 | 20.0 |
| 1-Pentadecanethiol    | 22.04 | 9.1     | ng/ul | 772006   | 5                     | 21.89 | 1691460 | 20.0 |
| 3,7,11,15-Tetrame...  | 22.08 | 5.7     | ng/ul | 483225   | 5                     | 21.89 | 1691460 | 20.0 |
| Trifluoroacetic a...  | 22.74 | 14.7    | ng/ul | 1241800  | 5                     | 21.89 | 1691460 | 20.0 |
| Tetracosanoic aci...  | 22.94 | 6.3     | ng/ul | 535421   | 5                     | 21.89 | 1691460 | 20.0 |
| unknown-02            | 23.24 | 2.6     | ng/ul | 216454   | 5                     | 21.89 | 1691460 | 20.0 |
| 7-Heptadecene, 17...  | 23.40 | 2.9     | ng/ul | 240621   | 5                     | 21.89 | 1691460 | 20.0 |
| 9-Octadecenamide, ... | 23.44 | 10.0    | ng/ul | 847700   | 5                     | 21.89 | 1691460 | 20.0 |
| (DEL) Alkane: Cyc...  | 24.38 | 20.7    | ng/ul | 1050320  | 6                     | 25.29 | 1015920 | 20.0 |
| unknown-03            | 25.99 | 18.2    | ng/ul | 924390   | 6                     | 25.29 | 1015920 | 20.0 |
| unknown-04            | 26.63 | 16.7    | ng/ul | 846149   | 6                     | 25.29 | 1015920 | 20.0 |