

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
 Data File : BG061841.D  
 Acq On : 2 Jul 2024 11:02  
 Operator : MA/JU  
 Sample : P2963-07  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 A4CQ8

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062024.MA.M  
 Title : SVOA CALIBRATION

Signal : TIC: BG061841.D\data.ms

| 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         | 3.458       | 26            | 29          | 34           | rVB3     | 17176          | 25839         | 0.79%           | 0.058%        |
| 2         | 3.734       | 71            | 76          | 81           | rBV2     | 31508          | 50073         | 1.53%           | 0.112%        |
| 3         | 3.987       | 112           | 119         | 126          | rVB      | 62478          | 98955         | 3.03%           | 0.221%        |
| 4         | 4.216       | 152           | 158         | 165          | rVV2     | 37226          | 69983         | 2.14%           | 0.157%        |
| 5         | 4.774       | 249           | 253         | 258          | rVV      | 21837          | 37951         | 1.16%           | 0.085%        |
| 6         | 5.133       | 307           | 314         | 324          | rBV      | 75719          | 130630        | 4.00%           | 0.292%        |
| 7         | 5.227       | 324           | 330         | 338          | rBV4     | 19292          | 36988         | 1.13%           | 0.083%        |
| 8         | 7.219       | 661           | 669         | 677          | rVB      | 432520         | 772675        | 23.65%          | 1.729%        |
| 9         | 7.383       | 688           | 697         | 704          | rBV      | 291806         | 493744        | 15.12%          | 1.105%        |
| 10        | 7.589       | 725           | 732         | 738          | rVV      | 404700         | 690462        | 21.14%          | 1.545%        |
| 11        | 7.671       | 738           | 746         | 748          | rVV4     | 17987          | 35395         | 1.08%           | 0.079%        |
| 12        | 8.059       | 801           | 812         | 818          | rVV      | 285480         | 502537        | 15.38%          | 1.125%        |
| 13        | 8.611       | 901           | 906         | 915          | rVV9     | 9364           | 28284         | 0.87%           | 0.063%        |
| 14        | 8.764       | 925           | 932         | 938          | rVV2     | 475222         | 824626        | 25.25%          | 1.845%        |
| 15        | 8.881       | 947           | 952         | 960          | rVB3     | 30420          | 55826         | 1.71%           | 0.125%        |
| 16        | 9.222       | 1003          | 1010        | 1017         | rVV      | 416593         | 736209        | 22.54%          | 1.647%        |
| 17        | 9.275       | 1017          | 1019        | 1024         | rVV2     | 23782          | 27148         | 0.83%           | 0.061%        |
| 18        | 9.945       | 1126          | 1133        | 1143         | rBV      | 380349         | 660987        | 20.24%          | 1.479%        |
| 19        | 10.097      | 1152          | 1159        | 1165         | rVV8     | 12125          | 26952         | 0.83%           | 0.060%        |
| 20        | 10.491      | 1217          | 1226        | 1233         | rVV      | 503096         | 925276        | 28.33%          | 2.071%        |
| 21        | 10.826      | 1278          | 1283        | 1284         | rBV2     | 17564          | 24146         | 0.74%           | 0.054%        |
| 22        | 10.867      | 1284          | 1290        | 1296         | rVV      | 422737         | 757567        | 23.19%          | 1.695%        |
| 23        | 10.920      | 1296          | 1299        | 1304         | rVB      | 75611          | 112401        | 3.44%           | 0.252%        |
| 24        | 11.008      | 1307          | 1314        | 1320         | rBV2     | 45055          | 93507         | 2.86%           | 0.209%        |
| 25        | 11.390      | 1373          | 1379        | 1385         | rBV5     | 17485          | 30885         | 0.95%           | 0.069%        |
| 26        | 11.607      | 1408          | 1416        | 1420         | rBV2     | 30886          | 63192         | 1.93%           | 0.141%        |
| 27        | 11.872      | 1457          | 1461        | 1466         | rVV4     | 15342          | 25449         | 0.78%           | 0.057%        |
| 28        | 12.518      | 1566          | 1571        | 1579         | rVB2     | 57255          | 91552         | 2.80%           | 0.205%        |
| 29        | 12.736      | 1603          | 1608        | 1614         | rVB2     | 42095          | 73084         | 2.24%           | 0.164%        |
| 30        | 13.505      | 1734          | 1739        | 1740         | rBV4     | 21813          | 32334         | 0.99%           | 0.072%        |
| 31        | 13.717      | 1768          | 1775        | 1779         | rBV4     | 19642          | 39538         | 1.21%           | 0.088%        |
| 32        | 13.858      | 1794          | 1799        | 1806         | rVB7     | 31048          | 74047         | 2.27%           | 0.166%        |
| 33        | 14.011      | 1818          | 1825        | 1829         | rBV2     | 46379          | 88577         | 2.71%           | 0.198%        |
| 34        | 14.093      | 1829          | 1839        | 1844         | rBV      | 845442         | 1271496       | 38.93%          | 2.845%        |
| 35        | 14.152      | 1846          | 1849        | 1853         | rVB3     | 25148          | 34503         | 1.06%           | 0.077%        |
| 36        | 14.246      | 1861          | 1865        | 1868         | rBV2     | 22410          | 30451         | 0.93%           | 0.068%        |

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 ClientSampleId :  
 A4CQ8

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062024.MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 14.381 | 1882 | 1888 | 1900 | rVV  | 905246  | 1509143 | 46.20% | 3.377% |
| 38 | 14.569 | 1916 | 1920 | 1925 | rBV4 | 22050   | 37633   | 1.15%  | 0.084% |
| 39 | 14.627 | 1925 | 1930 | 1933 | rVV  | 29156   | 49385   | 1.51%  | 0.111% |
| 40 | 14.686 | 1933 | 1940 | 1946 | rVV  | 636946  | 998533  | 30.57% | 2.235% |
| 41 | 14.751 | 1946 | 1951 | 1958 | rVB  | 166231  | 260713  | 7.98%  | 0.583% |
| 42 | 14.886 | 1967 | 1974 | 1985 | rBV  | 464505  | 803441  | 24.60% | 1.798% |
| 43 | 14.998 | 1985 | 1993 | 1995 | rVV6 | 21727   | 48915   | 1.50%  | 0.109% |
| 44 | 15.033 | 1997 | 1999 | 2003 | rVV5 | 20093   | 27233   | 0.83%  | 0.061% |
| 45 | 15.080 | 2003 | 2007 | 2013 | rVB2 | 133849  | 215312  | 6.59%  | 0.482% |
| 46 | 15.156 | 2015 | 2020 | 2024 | rBV3 | 24686   | 35046   | 1.07%  | 0.078% |
| 47 | 15.297 | 2039 | 2044 | 2049 | rVV6 | 25826   | 49710   | 1.52%  | 0.111% |
| 48 | 15.350 | 2049 | 2053 | 2057 | rVB5 | 24432   | 36356   | 1.11%  | 0.081% |
| 49 | 15.468 | 2067 | 2073 | 2077 | rBV6 | 32796   | 63096   | 1.93%  | 0.141% |
| 50 | 15.515 | 2079 | 2081 | 2085 | rVV3 | 24452   | 29589   | 0.91%  | 0.066% |
| 51 | 15.679 | 2102 | 2109 | 2114 | rVV  | 1129542 | 1771594 | 54.24% | 3.964% |
| 52 | 15.732 | 2114 | 2118 | 2123 | rVV  | 169292  | 261382  | 8.00%  | 0.585% |
| 53 | 15.791 | 2123 | 2128 | 2132 | rVV  | 228768  | 342327  | 10.48% | 0.766% |
| 54 | 15.850 | 2136 | 2138 | 2142 | rVV3 | 44483   | 58862   | 1.80%  | 0.132% |
| 55 | 15.897 | 2142 | 2146 | 2149 | rVV5 | 40267   | 70115   | 2.15%  | 0.157% |
| 56 | 15.938 | 2149 | 2153 | 2157 | rVV7 | 30585   | 65745   | 2.01%  | 0.147% |
| 57 | 15.979 | 2157 | 2160 | 2163 | rVV5 | 21806   | 37879   | 1.16%  | 0.085% |
| 58 | 16.026 | 2165 | 2168 | 2172 | rVV2 | 61159   | 86879   | 2.66%  | 0.194% |
| 59 | 16.173 | 2188 | 2193 | 2201 | rVB2 | 52208   | 112984  | 3.46%  | 0.253% |
| 60 | 16.243 | 2201 | 2205 | 2207 | rBV5 | 16695   | 24178   | 0.74%  | 0.054% |
| 61 | 16.267 | 2207 | 2209 | 2214 | rVB5 | 27236   | 29513   | 0.90%  | 0.066% |
| 62 | 16.402 | 2224 | 2232 | 2237 | rBV6 | 78979   | 180757  | 5.53%  | 0.405% |
| 63 | 16.713 | 2279 | 2285 | 2286 | rBV6 | 38539   | 73019   | 2.24%  | 0.163% |
| 64 | 16.778 | 2293 | 2296 | 2300 | rBV3 | 44562   | 72370   | 2.22%  | 0.162% |
| 65 | 16.966 | 2320 | 2328 | 2330 | rBV8 | 42523   | 87711   | 2.69%  | 0.196% |
| 66 | 17.001 | 2330 | 2334 | 2337 | rVB6 | 39339   | 62100   | 1.90%  | 0.139% |
| 67 | 17.083 | 2344 | 2348 | 2354 | rVB  | 100102  | 156581  | 4.79%  | 0.350% |
| 68 | 17.177 | 2357 | 2364 | 2368 | rVV7 | 90327   | 195658  | 5.99%  | 0.438% |
| 69 | 17.248 | 2372 | 2376 | 2384 | rVB  | 115416  | 193300  | 5.92%  | 0.433% |
| 70 | 17.336 | 2389 | 2391 | 2395 | rVB5 | 23531   | 25968   | 0.79%  | 0.058% |
| 71 | 17.430 | 2401 | 2407 | 2411 | rBV  | 726262  | 1187501 | 36.35% | 2.657% |
| 72 | 17.477 | 2411 | 2415 | 2420 | rVV  | 1858757 | 2734458 | 83.71% | 6.119% |
| 73 | 17.530 | 2420 | 2424 | 2428 | rVV  | 1219726 | 1886035 | 57.74% | 4.221% |
| 74 | 17.565 | 2428 | 2430 | 2435 | rVB  | 368484  | 423155  | 12.95% | 0.947% |
| 75 | 17.718 | 2450 | 2456 | 2457 | rBV6 | 27227   | 54770   | 1.68%  | 0.123% |
| 76 | 17.835 | 2468 | 2476 | 2481 | rBV2 | 162343  | 334078  | 10.23% | 0.748% |

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 ClientSampleId :  
 A4CQ8

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

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Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

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Stop Thrs : 0

Peak Location: TOP

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062024.MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |        |
|-----|--------|------|------|------|------|---------|---------|---------|--------|
| 77  | 17.912 | 2486 | 2489 | 2492 | rVV4 | 46531   | 54052   | 1.65%   | 0.121% |
| 78  | 17.947 | 2492 | 2495 | 2500 | rVV2 | 39147   | 65668   | 2.01%   | 0.147% |
| 79  | 18.018 | 2503 | 2507 | 2514 | rVB4 | 90875   | 213844  | 6.55%   | 0.479% |
| 80  | 18.311 | 2552 | 2557 | 2560 | rBV2 | 388749  | 601386  | 18.41%  | 1.346% |
| 81  | 18.353 | 2560 | 2564 | 2571 | rVB  | 434513  | 633925  | 19.41%  | 1.419% |
| 82  | 18.435 | 2574 | 2578 | 2583 | rVB2 | 123054  | 176886  | 5.42%   | 0.396% |
| 83  | 18.499 | 2583 | 2589 | 2593 | rBV3 | 407803  | 776113  | 23.76%  | 1.737% |
| 84  | 18.535 | 2593 | 2595 | 2602 | rVB  | 237769  | 314646  | 9.63%   | 0.704% |
| 85  | 18.605 | 2603 | 2607 | 2611 | rBV4 | 91965   | 139841  | 4.28%   | 0.313% |
| 86  | 18.799 | 2637 | 2640 | 2644 | rVV  | 184209  | 223073  | 6.83%   | 0.499% |
| 87  | 18.840 | 2644 | 2647 | 2653 | rVB2 | 188253  | 277261  | 8.49%   | 0.620% |
| 88  | 19.046 | 2679 | 2682 | 2686 | rVB5 | 90511   | 122112  | 3.74%   | 0.273% |
| 89  | 19.216 | 2707 | 2711 | 2716 | rBV3 | 207598  | 395844  | 12.12%  | 0.886% |
| 90  | 19.481 | 2751 | 2756 | 2761 | rBV  | 2300004 | 3266441 | 100.00% | 7.310% |
| 91  | 19.533 | 2762 | 2765 | 2769 | rVV6 | 107723  | 151961  | 4.65%   | 0.340% |
| 92  | 19.816 | 2807 | 2813 | 2815 | rBV2 | 1478069 | 2397008 | 73.38%  | 5.364% |
| 93  | 19.845 | 2815 | 2818 | 2825 | rVB  | 1807748 | 2467196 | 75.53%  | 5.521% |
| 94  | 20.333 | 2898 | 2901 | 2905 | rVB3 | 437251  | 592864  | 18.15%  | 1.327% |
| 95  | 20.427 | 2915 | 2917 | 2921 | rVB  | 170246  | 201265  | 6.16%   | 0.450% |
| 96  | 21.337 | 3068 | 3072 | 3076 | rVB2 | 467777  | 595191  | 18.22%  | 1.332% |
| 97  | 21.684 | 3125 | 3131 | 3137 | rBV3 | 1185293 | 2866017 | 87.74%  | 6.414% |
| 98  | 21.743 | 3138 | 3141 | 3149 | rVB  | 996684  | 1613516 | 49.40%  | 3.611% |
| 99  | 22.506 | 3266 | 3271 | 3281 | rVB5 | 456016  | 1157965 | 35.45%  | 2.591% |
| 100 | 23.899 | 3503 | 3508 | 3516 | rBV2 | 553862  | 1616115 | 49.48%  | 3.617% |

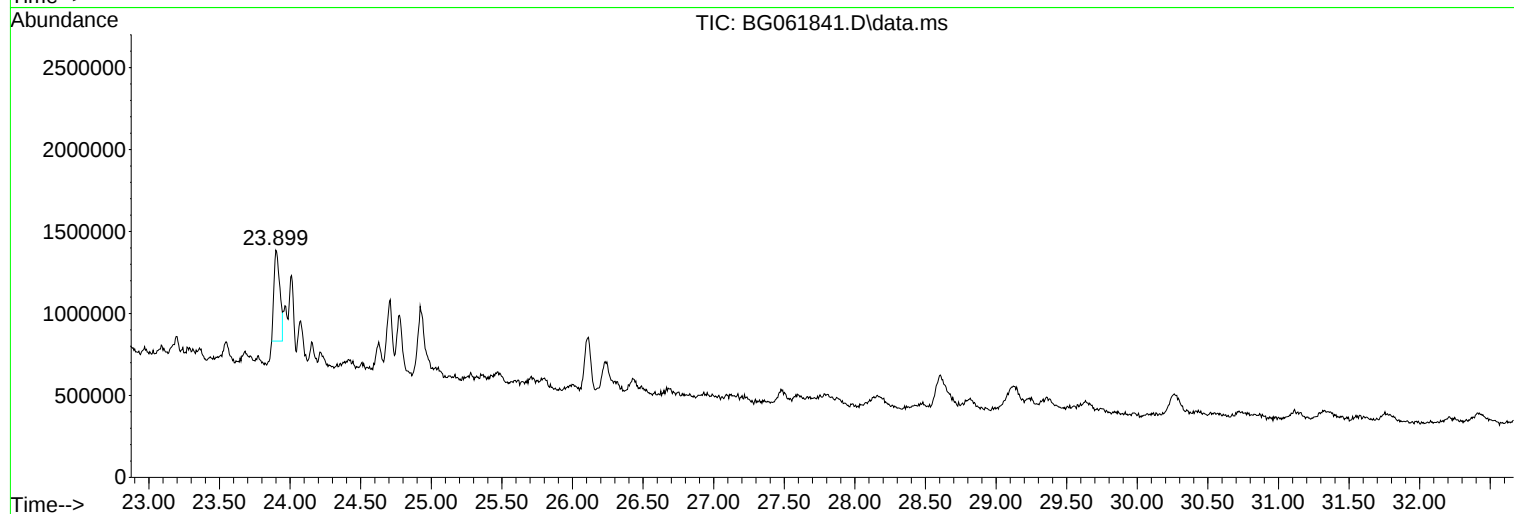
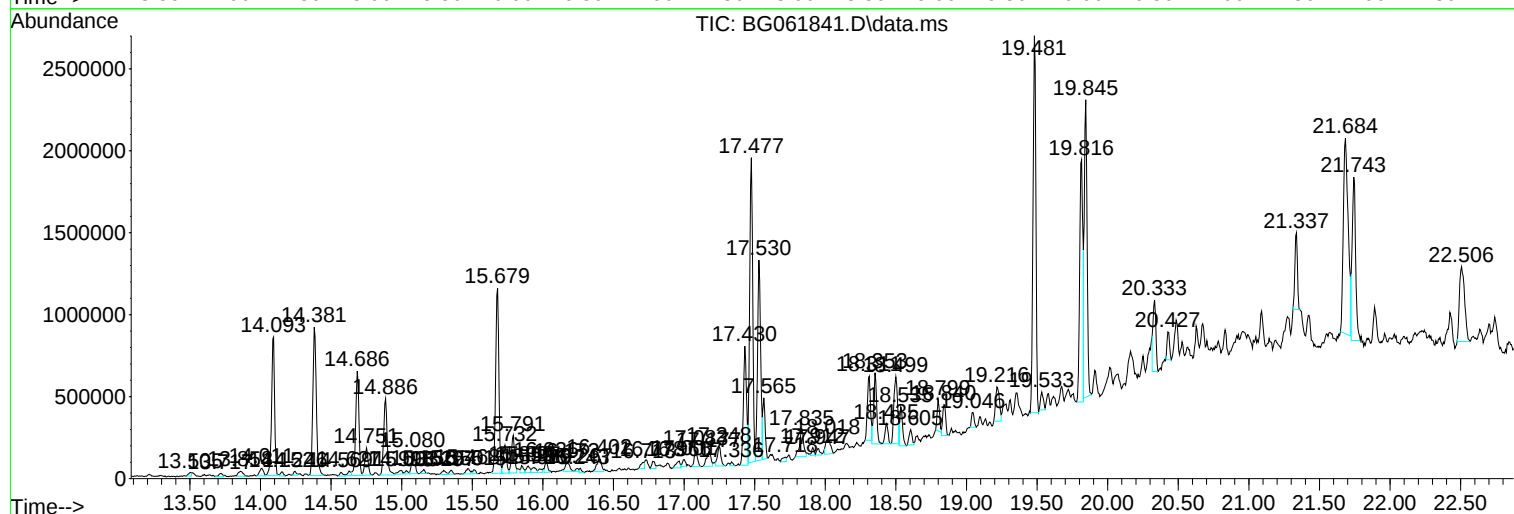
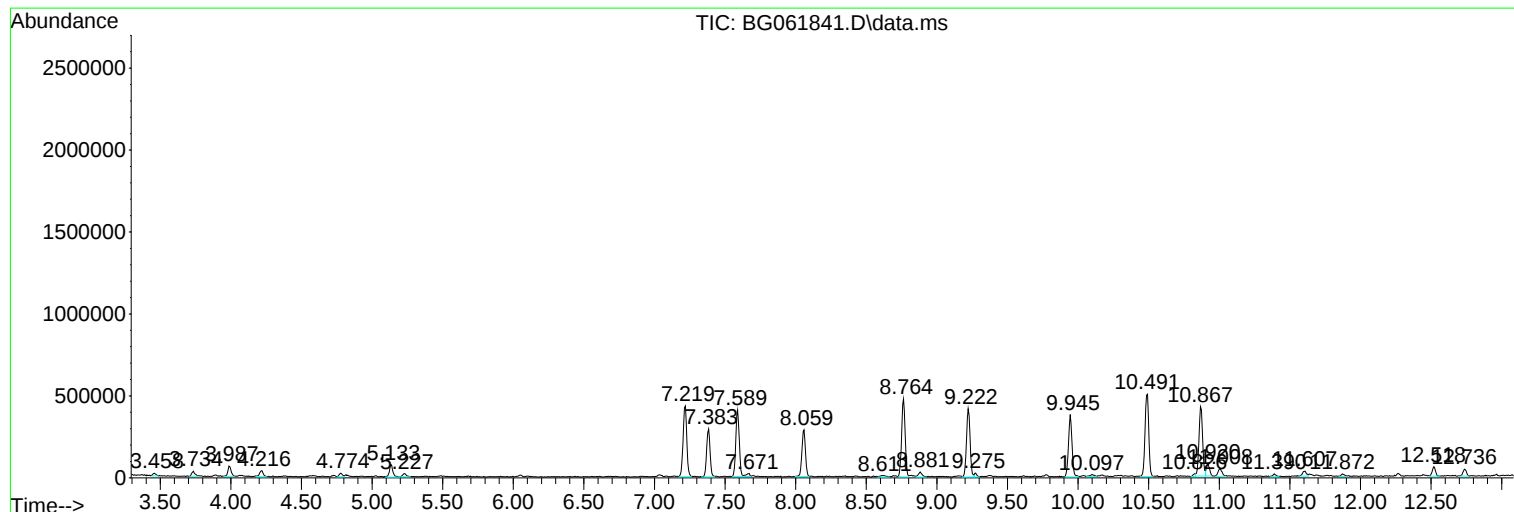
Sum of corrected areas: 44686483

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Instrument :  
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ClientSampleId :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062024.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P



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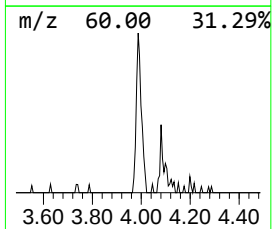
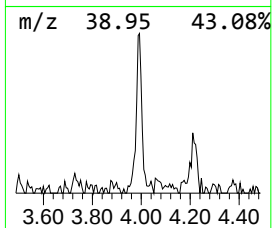
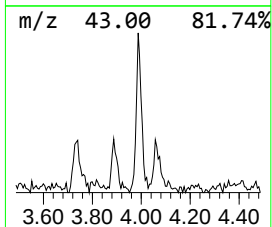
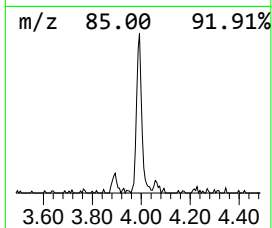
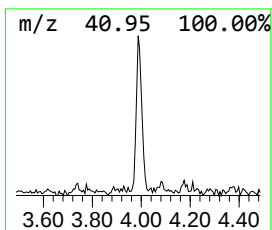
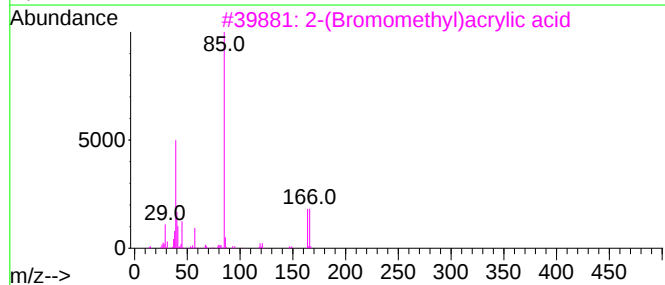
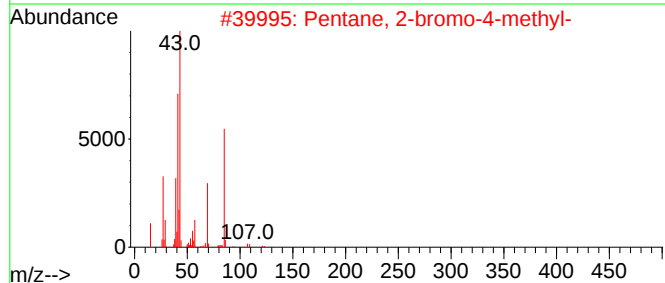
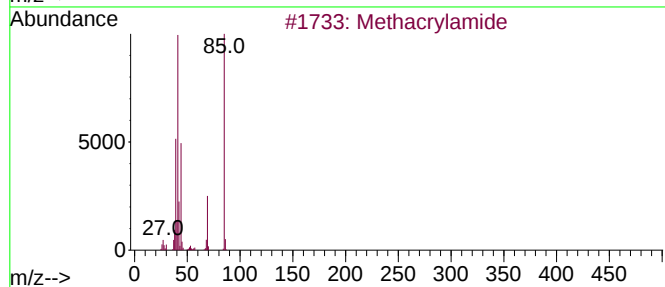
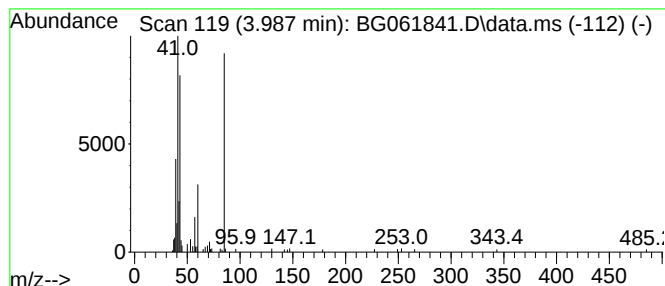
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062024.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Methacrylamide Concentration Rank 11

| R.T.  | EstConc    | Area  | Relative to ISTD       | R.T.  |
|-------|------------|-------|------------------------|-------|
| 3.987 | 3.94 ng/ul | 98955 | 1,4-Dichlorobenzene-d4 | 8.059 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------|-----|----------|-------------|------|
| 1    |    |   | Methacrylamide                | 85  | C4H7NO   | 000079-39-0 | 58   |
| 2    |    |   | Pentane, 2-bromo-4-methyl-    | 164 | C6H13Br  | 030310-22-6 | 37   |
| 3    |    |   | 2-(Bromomethyl)acrylic acid   | 164 | C4H5BrO2 | 072707-66-5 | 33   |
| 4    |    |   | Pentane, 2,4-dimethyl-        | 100 | C7H16    | 000108-08-7 | 28   |
| 5    |    |   | 2-Aminoethanol, N,O-diacetyl- | 145 | C6H11NO3 | 016180-96-4 | 28   |



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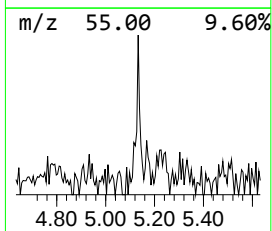
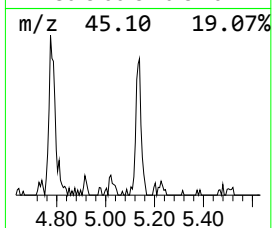
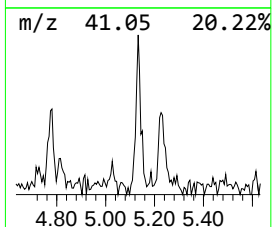
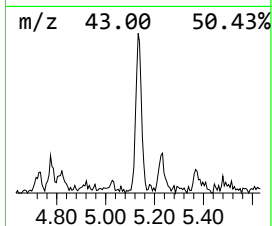
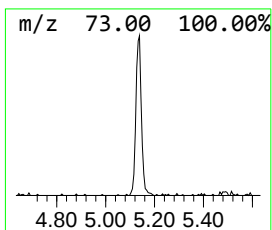
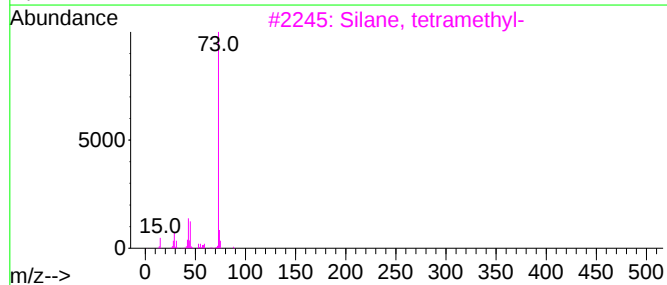
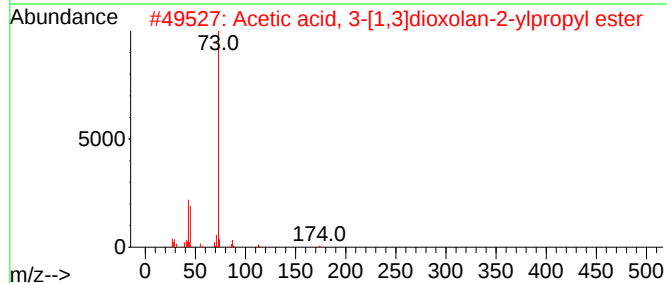
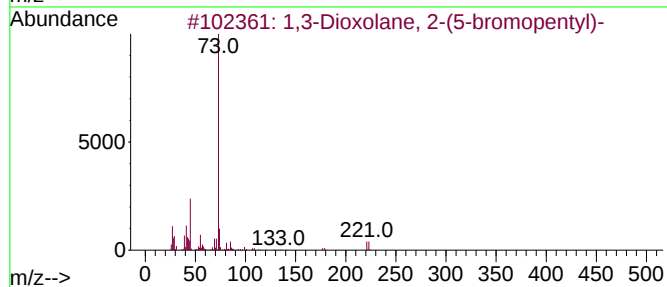
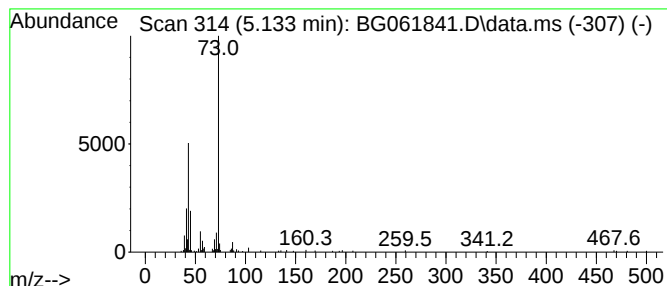
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062024.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 unknown-01 Concentration Rank 7

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 5.133 | 5.20 ng/ul | 130630 | 1,4-Dichlorobenzene-d4 | 8.059 |

| Hit# | of | 5 | Tentative ID                                  | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-----------------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 1,3-Dioxolane, 2-(5-bromopentyl)-             | 222 | C8H15BrO2 | 056741-68-5  | 38   |
| 2    |    |   | Acetic acid, 3-[1,3]dioxolan-2-ylpropyl ester | 174 | C8H14O4   | 1000186-02-3 | 36   |
| 3    |    |   | Silane, tetramethyl-                          | 88  | C4H12Si   | 000075-76-3  | 35   |
| 4    |    |   | 2,2'-Bi-1,3-dioxolane                         | 146 | C6H10O4   | 006705-89-1  | 33   |
| 5    |    |   | Methane, isothiocyanato-                      | 73  | C2H3NS    | 000556-61-6  | 28   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
Acq On : 2 Jul 2024 11:02  
Operator : MA/JU  
Sample : P2963-07  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
A4CQ8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062024.MA.M  
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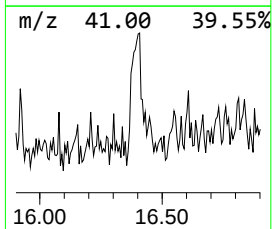
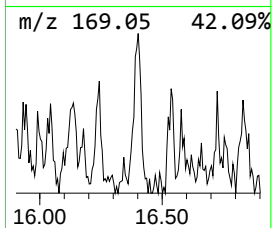
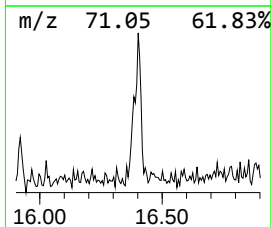
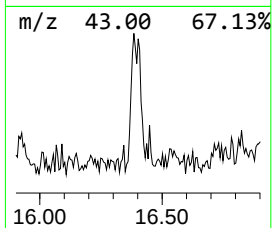
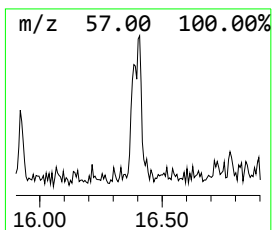
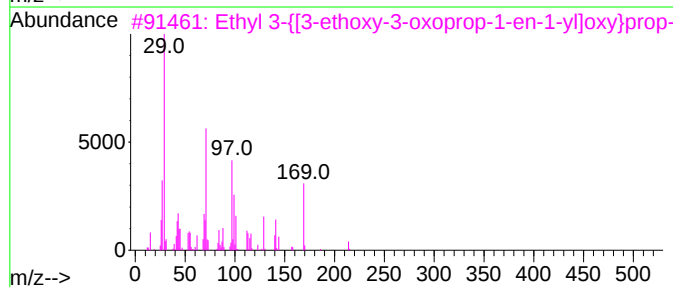
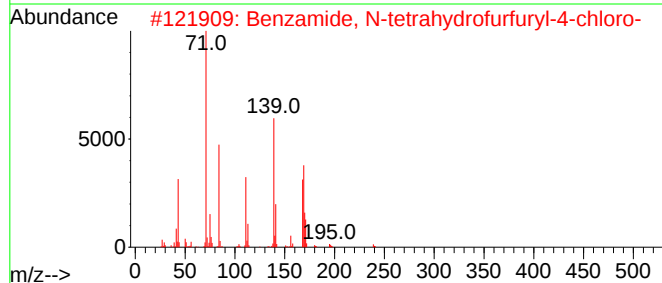
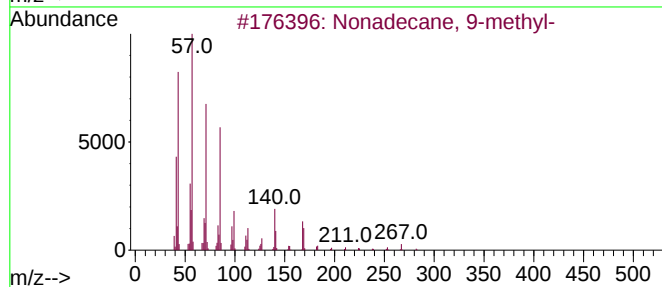
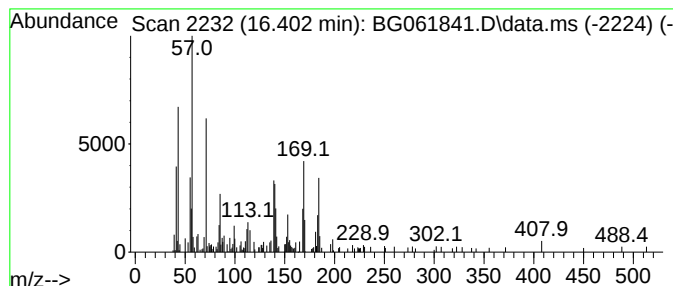
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 16

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.402 | 3.04 ng/ul | 180757 | Phenanthrene-d10 | 17.430 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Nonadecane, 9-methyl-               | 282 | C20H42      | 013287-24-6  | 22   |
| 2    |    |   | Benzamide, N-tetrahydrofurfuryl-... | 239 | C12H14ClNO2 | 1000307-39-4 | 12   |
| 3    |    |   | Ethyl 3-[[3-ethoxy-3-oxoprop-1-e... | 214 | C10H14O5    | 090927-35-8  | 12   |
| 4    |    |   | 5-Fluoro-2-methoxypyridin-4-amin... | 226 | C10H11FN2O3 | 1000503-13-3 | 10   |
| 5    |    |   | Heptadecane, 7-methyl-              | 254 | C18H38      | 020959-33-5  | 10   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
Acq On : 2 Jul 2024 11:02  
Operator : MA/JU  
Sample : P2963-07  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
A4CQ8

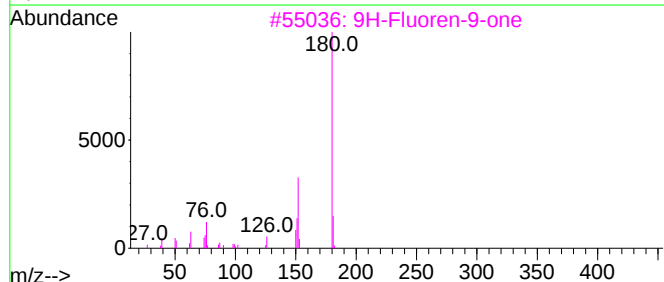
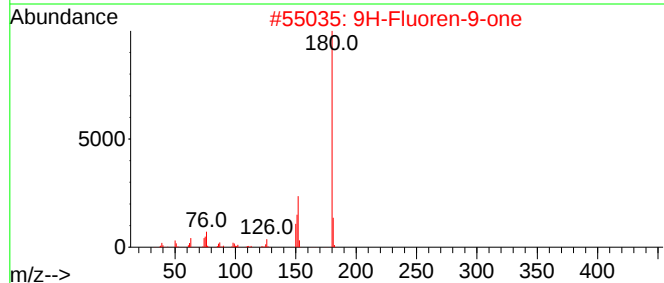
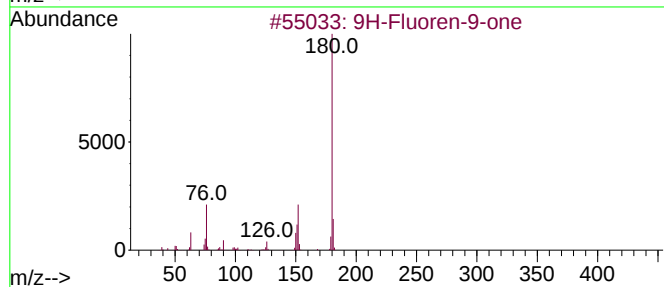
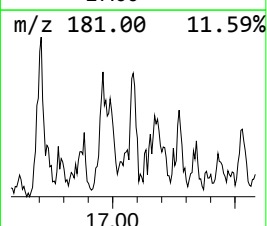
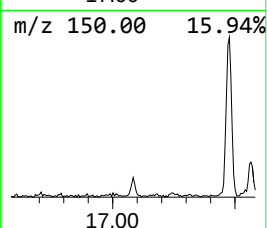
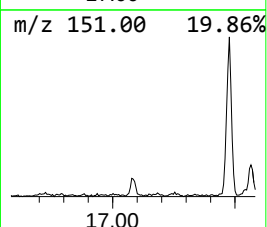
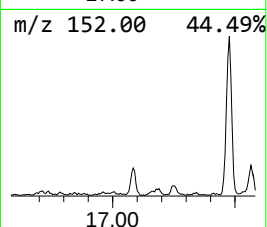
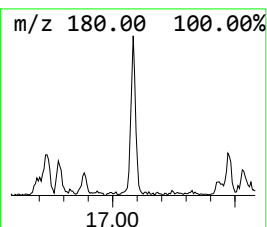
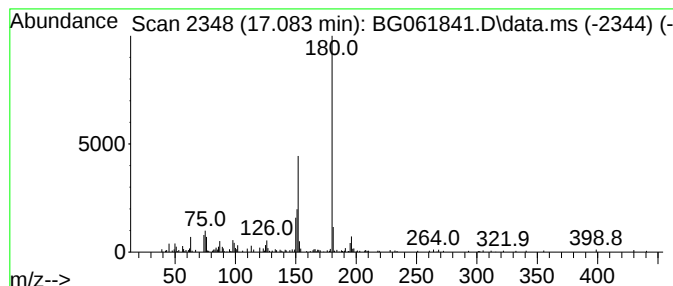
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 9H-Fluoren-9-one Concentration Rank 19

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.083 | 2.64 ng/ul | 156581 | Phenanthrene-d10 | 17.430 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm      | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|--------------|-------------|------|
| 1    |    |   | 9H-Fluoren-9-one                    | 180 | C13H8O       | 000486-25-9 | 87   |
| 2    |    |   | 9H-Fluoren-9-one                    | 180 | C13H8O       | 000486-25-9 | 81   |
| 3    |    |   | 9H-Fluoren-9-one                    | 180 | C13H8O       | 000486-25-9 | 80   |
| 4    |    |   | 9,9-Bis[4-aminophenylsulfonyl]fl... | 476 | C25H20N2O4S2 | 081269-16-1 | 78   |
| 5    |    |   | 9H-Fluoren-9-one                    | 180 | C13H8O       | 000486-25-9 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
Acq On : 2 Jul 2024 11:02  
Operator : MA/JU  
Sample : P2963-07  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
A4CQ8

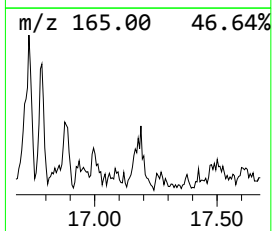
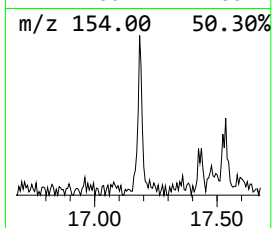
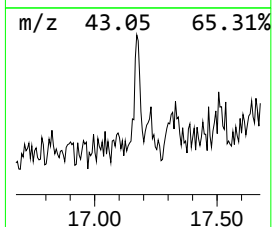
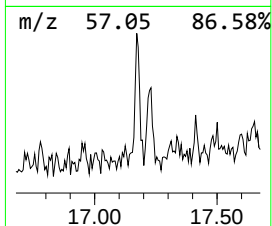
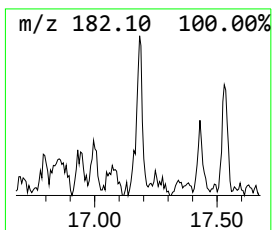
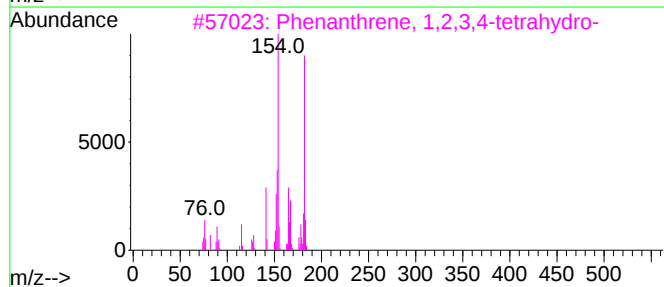
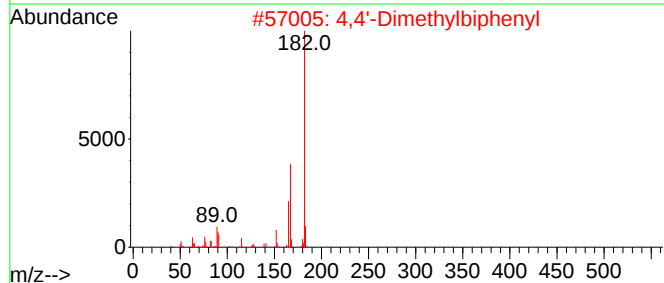
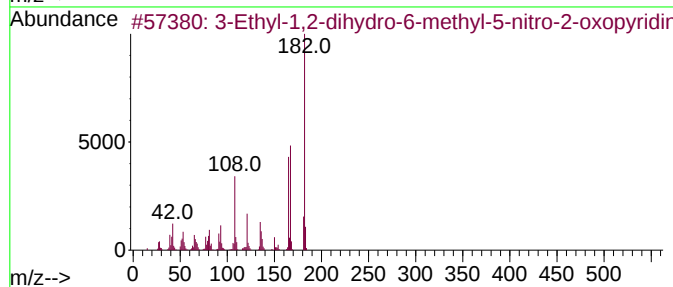
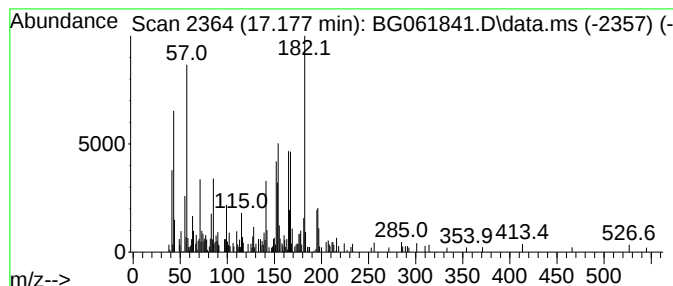
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 unknown-02 Concentration Rank 14

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.177 | 3.30 ng/ul | 195658 | Phenanthrene-d10 | 17.430 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 3-Ethyl-1,2-dihydro-6-methyl-5-n... | 182 | C8H10N2O3  | 1000241-09-2 | 49   |
| 2    |    |   | 4,4'-Dimethylbiphenyl               | 182 | C14H14     | 000613-33-2  | 38   |
| 3    |    |   | Phenanthrene, 1,2,3,4-tetrahydro-   | 182 | C14H14     | 001013-08-7  | 38   |
| 4    |    |   | But-2-enedinitrile, 2-cyano-3-(1... | 182 | C10H6N4    | 1000304-34-1 | 35   |
| 5    |    |   | N-Methyl-N-trifluoroacetyl-L-valine | 227 | C8H12F3NO3 | 1000452-81-7 | 30   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
Acq On : 2 Jul 2024 11:02  
Operator : MA/JU  
Sample : P2963-07  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
A4CQ8

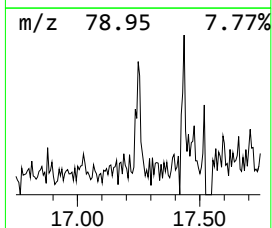
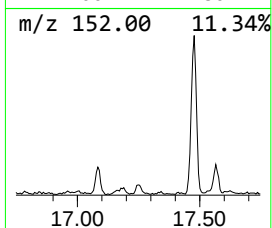
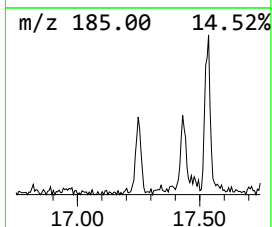
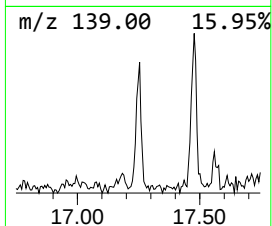
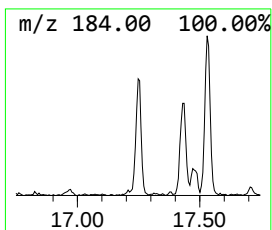
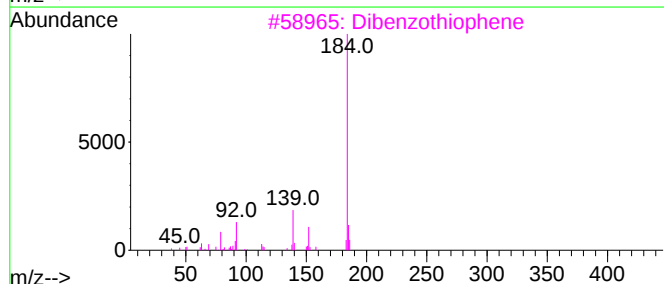
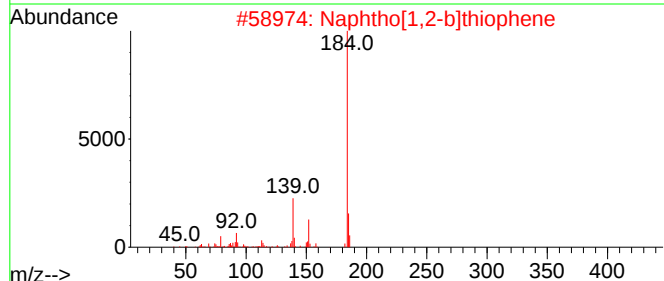
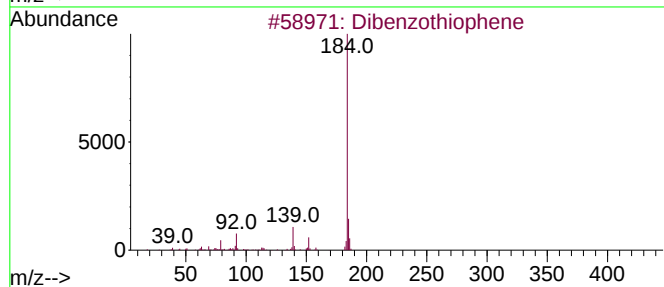
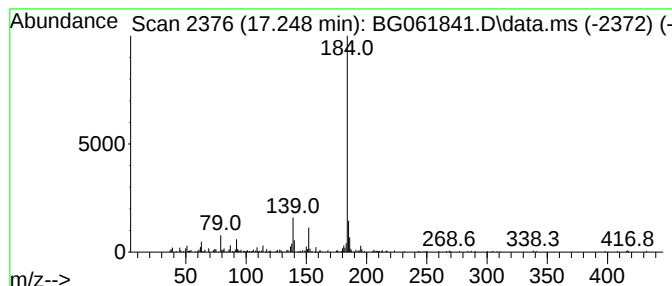
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Dibenzothiophene Concentration Rank 15

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.248 | 3.26 ng/ul | 193300 | Phenanthrene-d10 | 17.430 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Dibenzothiophene        | 184 | C12H8S  | 000132-65-0 | 90   |
| 2    |    |   | Naphtho[1,2-b]thiophene | 184 | C12H8S  | 000234-41-3 | 90   |
| 3    |    |   | Dibenzothiophene        | 184 | C12H8S  | 000132-65-0 | 90   |
| 4    |    |   | Dibenzothiophene        | 184 | C12H8S  | 000132-65-0 | 90   |
| 5    |    |   | Naphtho[2,1-b]thiophene | 184 | C12H8S  | 000233-02-3 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
Acq On : 2 Jul 2024 11:02  
Operator : MA/JU  
Sample : P2963-07  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

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

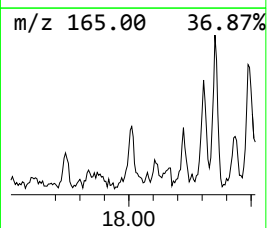
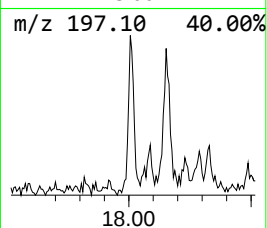
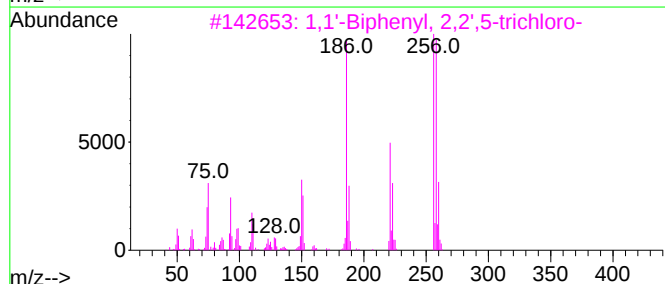
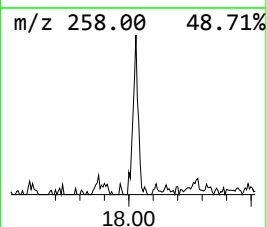
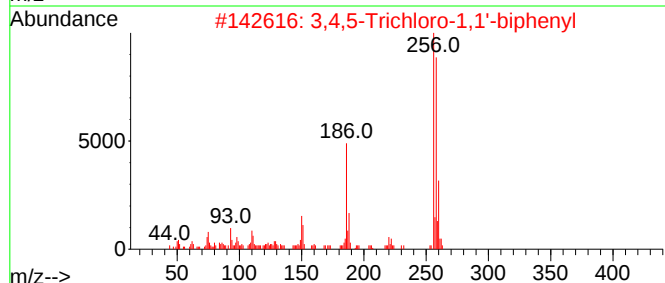
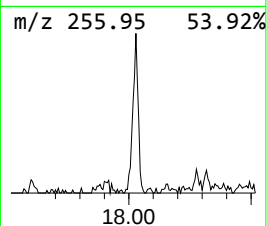
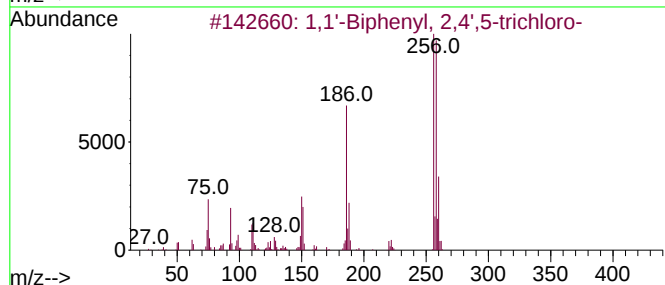
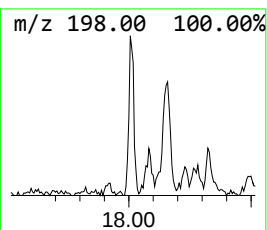
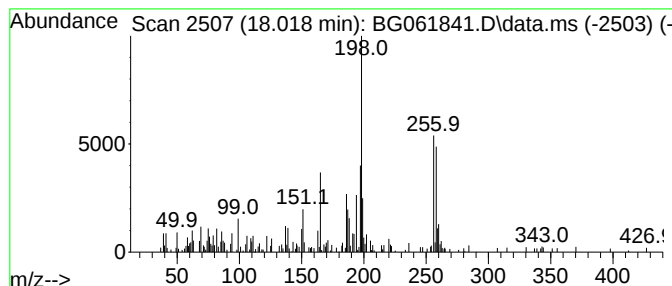
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 1,1'-Biphenyl, 2,4',5-trich... Concentration Rank 13

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.018 | 3.60 ng/ul | 213844 | Phenanthrene-d10 | 17.430 |

| Hit# | of                               | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|----------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1,1'-Biphenyl, 2,4',5-trichloro- | 256 | C12H7Cl3     | 016606-02-3 | 70      |      |      |
| 2    | 3,4,5-Trichloro-1,1'-biphenyl    | 256 | C12H7Cl3     | 053555-66-1 | 50      |      |      |
| 3    | 1,1'-Biphenyl, 2,2',5-trichloro- | 256 | C12H7Cl3     | 037680-65-2 | 38      |      |      |
| 4    | 3,3',4-Trichloro-1,1'-biphenyl   | 256 | C12H7Cl3     | 037680-69-6 | 30      |      |      |
| 5    | 3,4,5-Trichloro-1,1'-biphenyl    | 256 | C12H7Cl3     | 053555-66-1 | 30      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
Acq On : 2 Jul 2024 11:02  
Operator : MA/JU  
Sample : P2963-07  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
A4CQ8

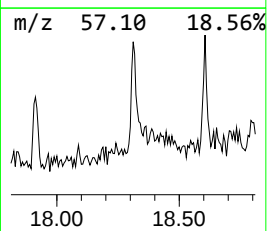
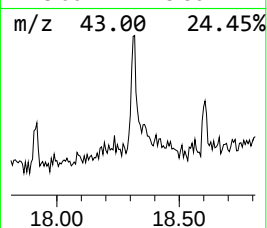
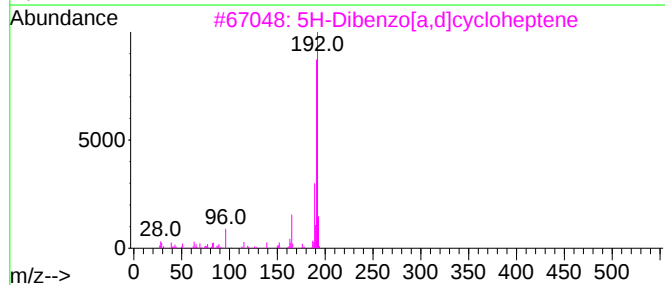
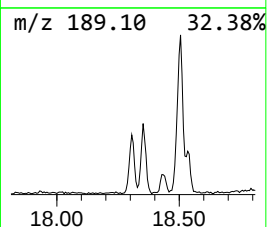
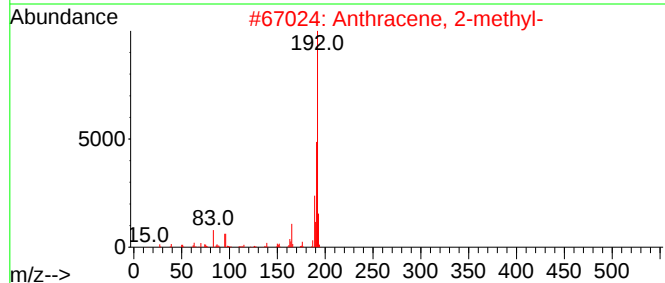
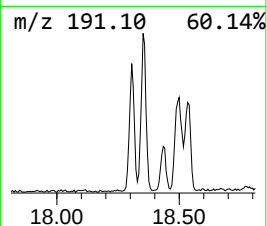
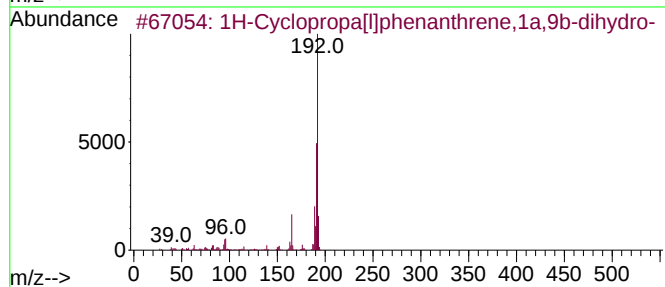
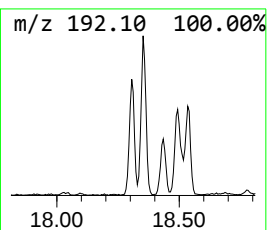
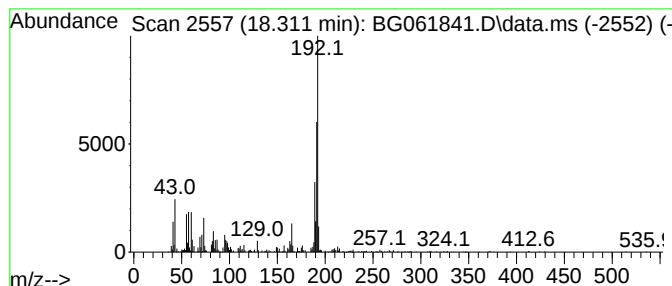
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 1H-Cyclopropa[l]phenanthren... Concentration Rank 3

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 18.311 | 10.13 ng/ul | 601386 | Phenanthrene-d10 | 17.430 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Cyclopropa[l]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 86   |
| 2    |    |   | Anthracene, 2-methyl-               | 192 | C15H12  | 000613-12-7 | 86   |
| 3    |    |   | 5H-Dibenzo[a,d]cycloheptene         | 192 | C15H12  | 000256-81-5 | 76   |
| 4    |    |   | Anthracene, 9-methyl-               | 192 | C15H12  | 000779-02-2 | 76   |
| 5    |    |   | Anthracene, 9-methyl-               | 192 | C15H12  | 000779-02-2 | 76   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
Acq On : 2 Jul 2024 11:02  
Operator : MA/JU  
Sample : P2963-07  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
A4CQ8

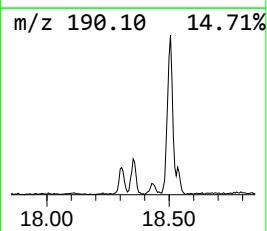
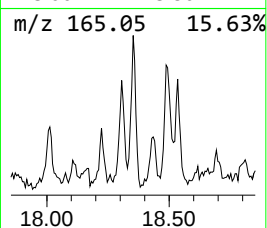
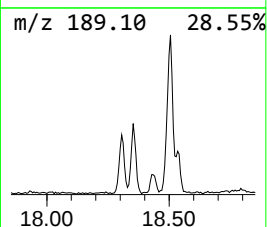
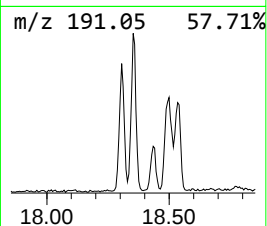
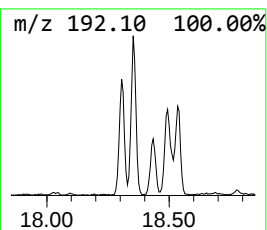
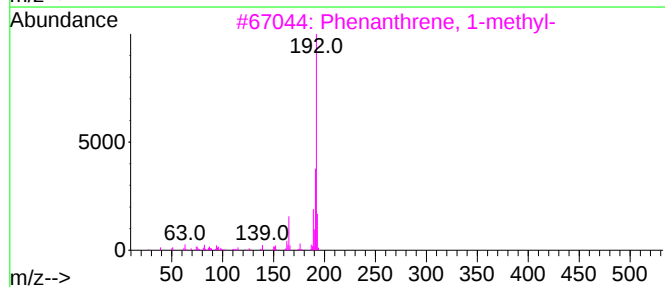
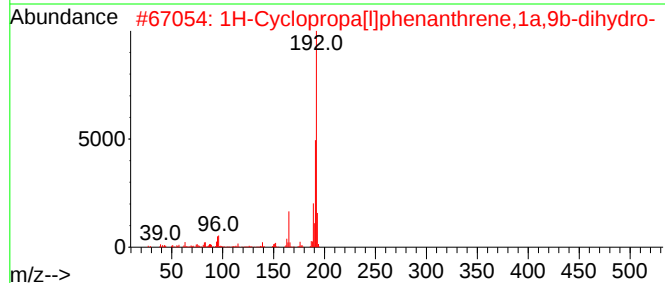
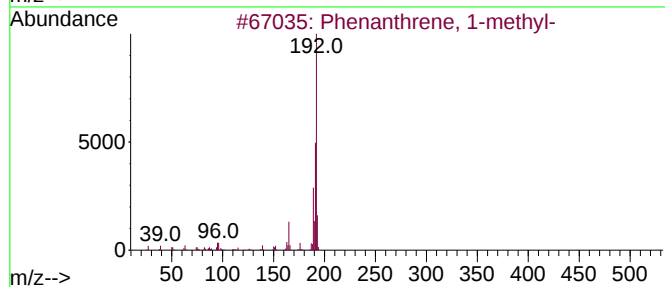
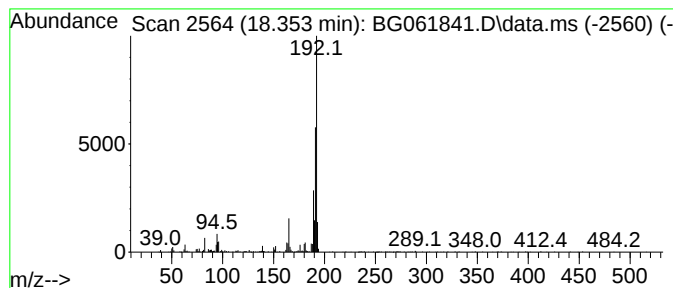
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Phenanthrene, 1-methyl- Concentration Rank 2

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 18.352 | 10.68 ng/ul | 633925 | Phenanthrene-d10 | 17.430 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 97   |
| 2    |      | 1H-Cyclopropa[1]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 96   |
| 3    |      | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 93   |
| 4    |      | Anthracene, 1-methyl-               | 192 | C15H12  | 000610-48-0 | 93   |
| 5    |      | Anthracene, 2-methyl-               | 192 | C15H12  | 000613-12-7 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
Acq On : 2 Jul 2024 11:02  
Operator : MA/JU  
Sample : P2963-07  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

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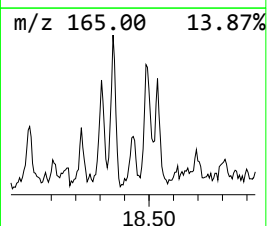
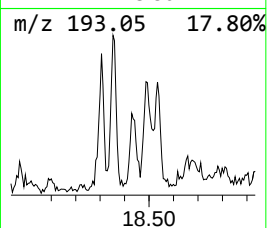
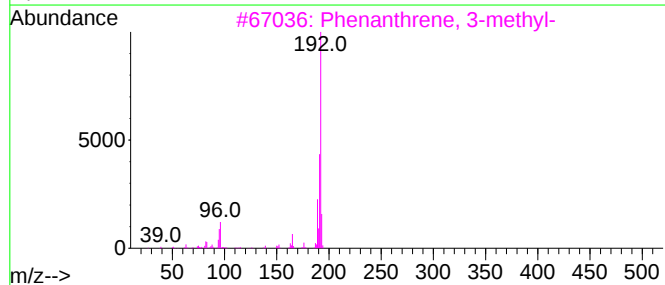
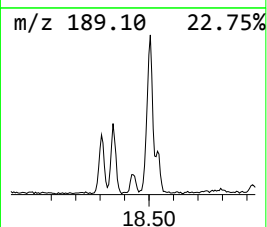
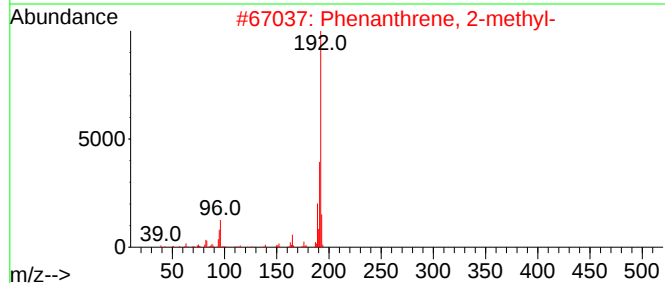
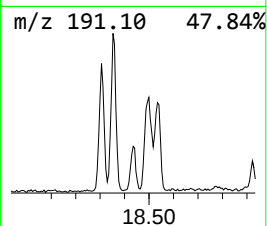
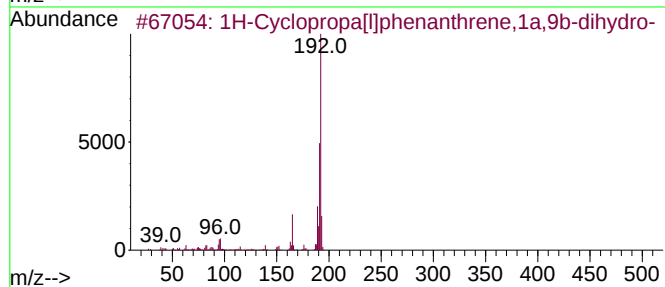
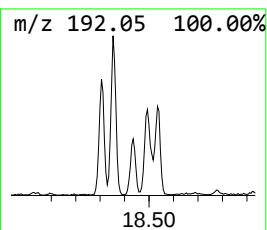
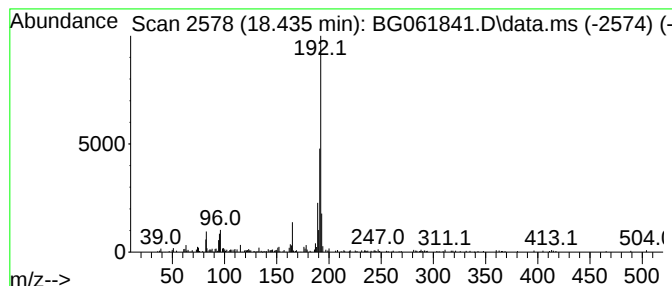
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Phenanthrene, 2-methyl- Concentration Rank 17

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.435 | 2.98 ng/ul | 176886 | Phenanthrene-d10 | 17.430 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | 1H-Cyclopropa[l]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 95   |
| 2    |      | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 94   |
| 3    |      | Phenanthrene, 3-methyl-             | 192 | C15H12  | 000832-71-3 | 94   |
| 4    |      | Phenanthrene, 9-methyl-             | 192 | C15H12  | 000883-20-5 | 94   |
| 5    |      | Phenanthrene, 4-methyl-             | 192 | C15H12  | 000832-64-4 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
Acq On : 2 Jul 2024 11:02  
Operator : MA/JU  
Sample : P2963-07  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

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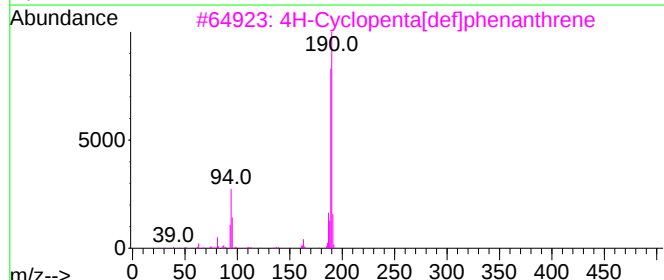
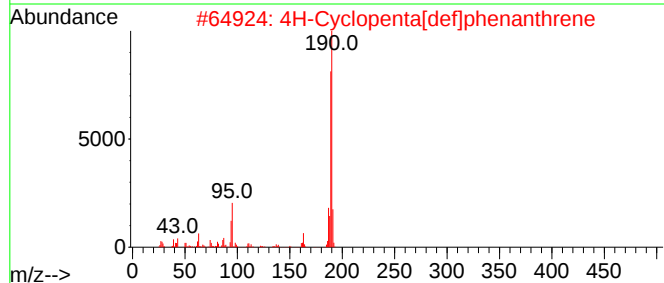
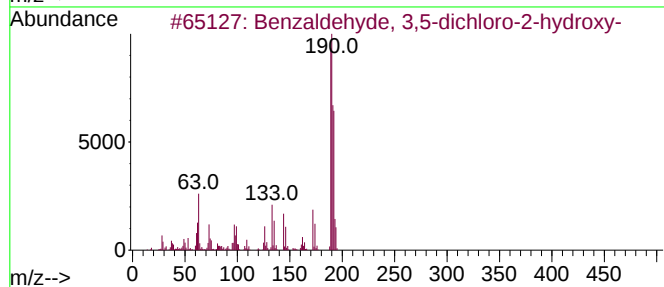
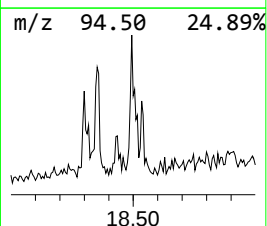
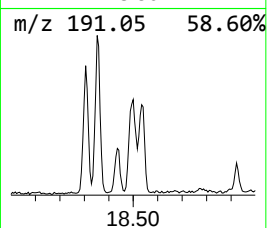
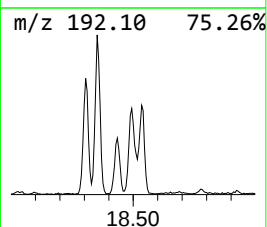
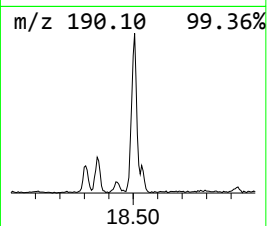
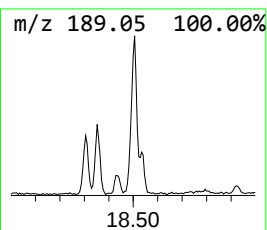
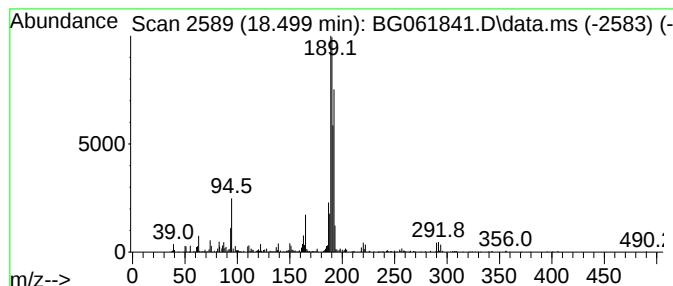
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Benzaldehyde, 3,5-dichloro-... Concentration Rank 1

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 18.499 | 13.07 ng/ul | 776113 | Phenanthrene-d10 | 17.430 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | Benzaldehyde, 3,5-dichloro-2-hyd... | 190 | C7H4Cl2O2 | 000090-60-8 | 72   |
| 2    |    |   | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10    | 000203-64-5 | 55   |
| 3    |    |   | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10    | 000203-64-5 | 49   |
| 4    |    |   | Benzaldehyde, 3,5-dichloro-2-hyd... | 190 | C7H4Cl2O2 | 000090-60-8 | 43   |
| 5    |    |   | Anthracene, 9-methyl-               | 192 | C15H12    | 000779-02-2 | 30   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
Acq On : 2 Jul 2024 11:02  
Operator : MA/JU  
Sample : P2963-07  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

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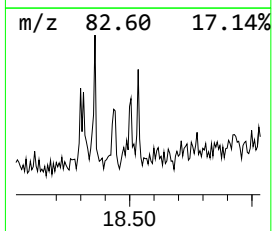
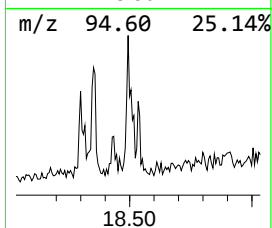
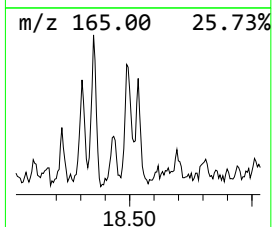
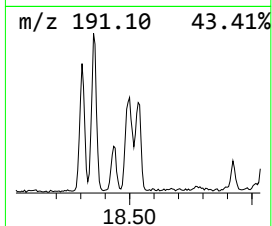
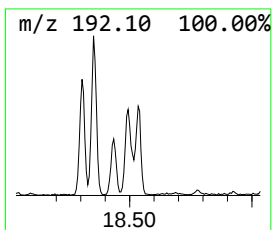
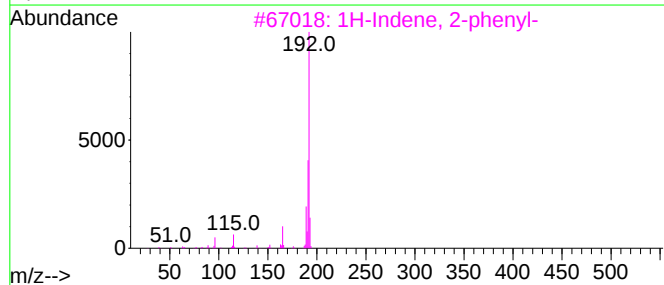
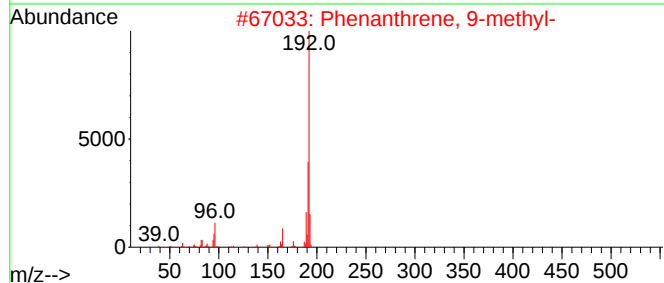
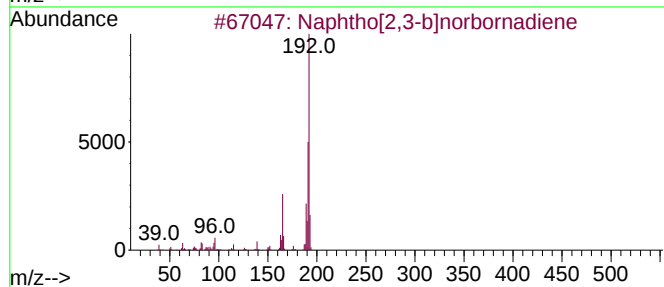
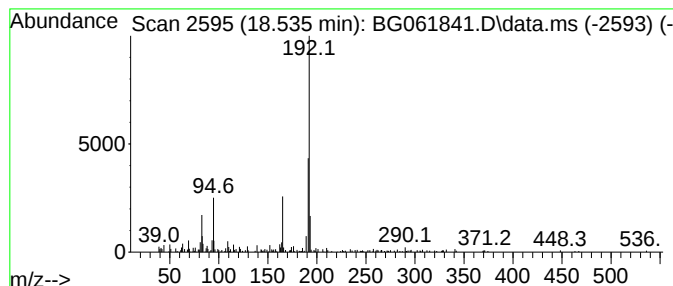
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 Naphtho[2,3-b]norbornadiene Concentration Rank 6

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.535 | 5.30 ng/ul | 314646 | Phenanthrene-d10 | 17.430 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphtho[2,3-b]norbornadiene | 192 | C15H12  | 107426-38-0 | 81   |
| 2    |    |   | Phenanthrene, 9-methyl-     | 192 | C15H12  | 000883-20-5 | 74   |
| 3    |    |   | 1H-Indene, 2-phenyl-        | 192 | C15H12  | 004505-48-0 | 72   |
| 4    |    |   | Phenanthrene, 2-methyl-     | 192 | C15H12  | 002531-84-2 | 72   |
| 5    |    |   | Anthracene, 1-methyl-       | 192 | C15H12  | 000610-48-0 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
Acq On : 2 Jul 2024 11:02  
Operator : MA/JU  
Sample : P2963-07  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

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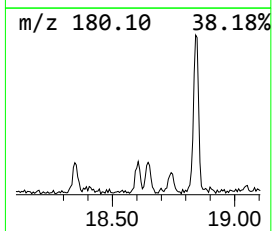
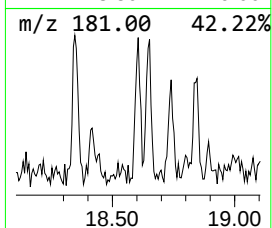
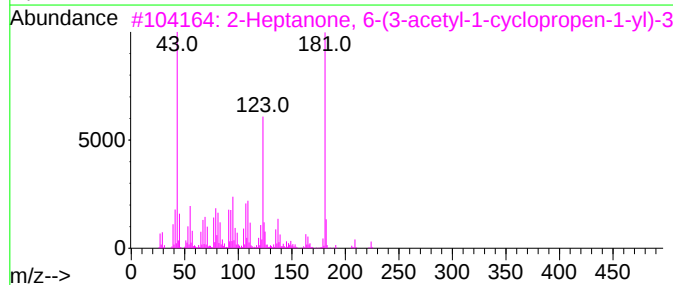
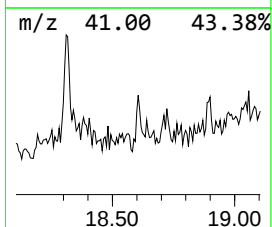
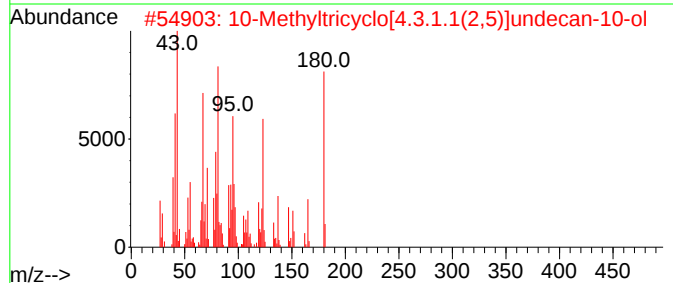
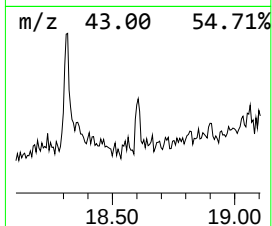
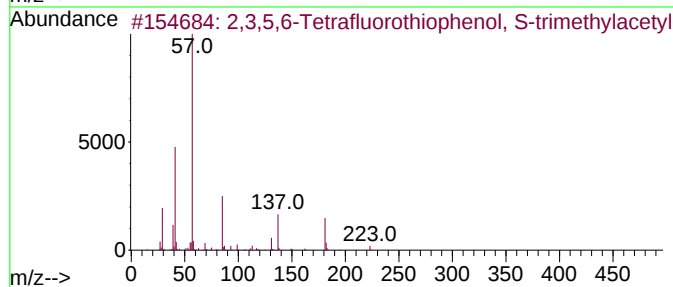
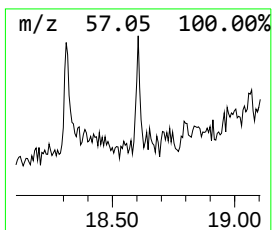
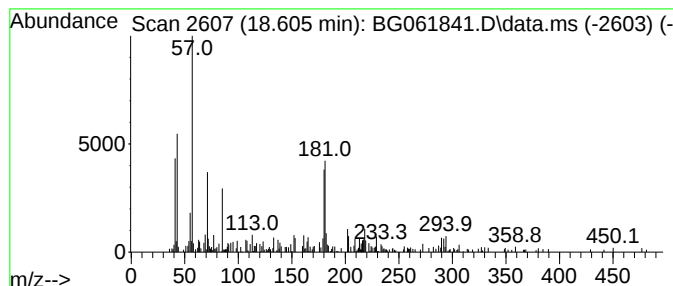
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 unknown-03 Concentration Rank 21

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.605 | 2.36 ng/ul | 139841 | Phenanthrene-d10 | 17.430 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 2,3,5,6-Tetrafluorothiophenol, S... | 266 | C11H10F4OS   | 1000470-48-7 | 32      |      |      |
| 2    | 10-Methyltricyclo[4.3.1.1(2,5)]u... | 180 | C12H20O      | 1000187-06-1 | 22      |      |      |
| 3    | 2-Heptanone, 6-(3-acetyl-1-cyclo... | 224 | C13H20O3     | 078091-85-7  | 12      |      |      |
| 4    | 2,6-Dichloro-5-fluoropyridin-3-a... | 222 | C7H5Cl2FN2O  | 1000503-14-0 | 12      |      |      |
| 5    | 5-Acetamido-1,3,4-thiadiazole-2-... | 222 | C4H6N4O3S2   | 000059-66-5  | 10      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
Acq On : 2 Jul 2024 11:02  
Operator : MA/JU  
Sample : P2963-07  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
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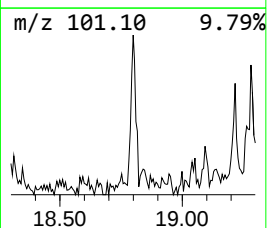
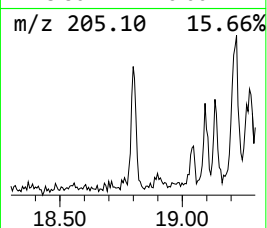
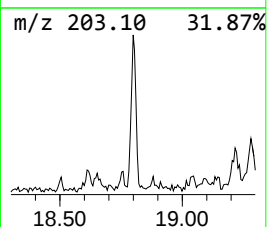
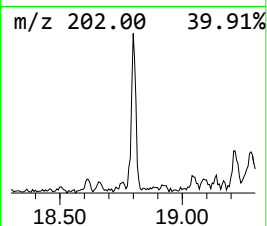
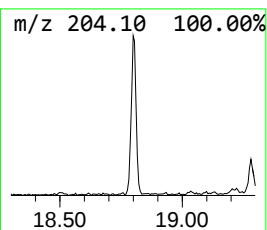
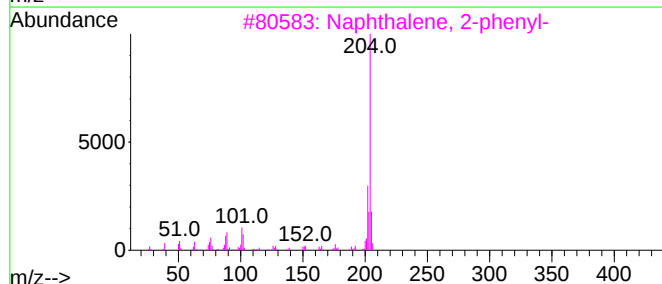
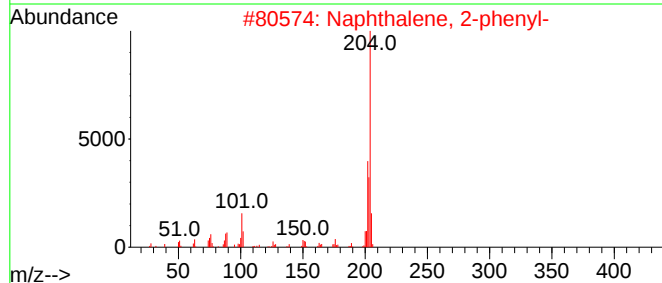
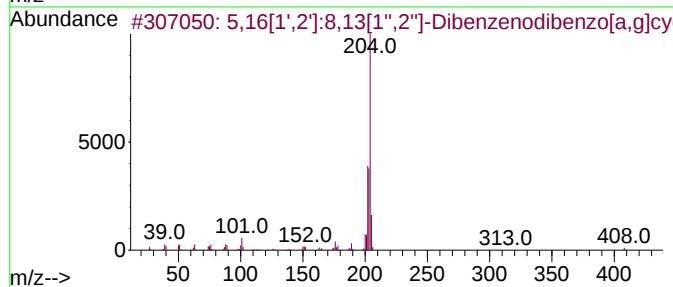
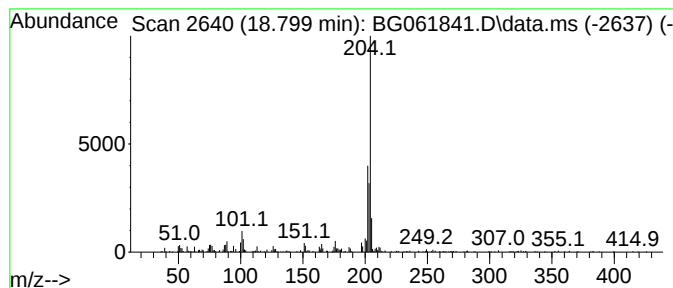
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TIC Integration Parameters: LSCINT.P

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Peak Number 15 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 12

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.799 | 3.76 ng/ul | 223073 | Phenanthrene-d10 | 17.430 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 5,16[1',2']:8,13[1'',2'']-Dibenz... | 408 | C32H24       | 005672-97-9 | 86      |      |      |
| 2    | Naphthalene, 2-phenyl-              | 204 | C16H12       | 000612-94-2 | 70      |      |      |
| 3    | Naphthalene, 2-phenyl-              | 204 | C16H12       | 000612-94-2 | 60      |      |      |
| 4    | Naphthalene, 2-phenyl-              | 204 | C16H12       | 000612-94-2 | 60      |      |      |
| 5    | 6-Cyanophenanthridine               | 204 | C14H8N2      | 002176-28-5 | 55      |      |      |



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Data File : BG061841.D  
Acq On : 2 Jul 2024 11:02  
Operator : MA/JU  
Sample : P2963-07  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

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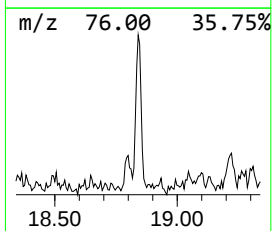
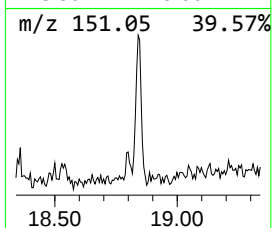
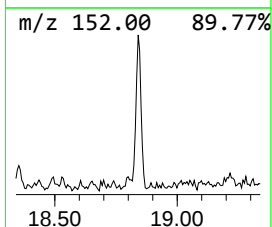
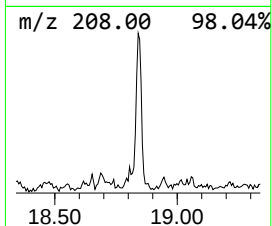
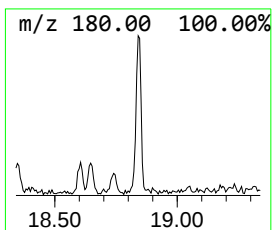
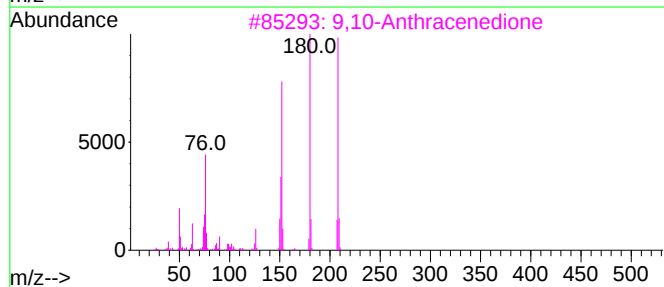
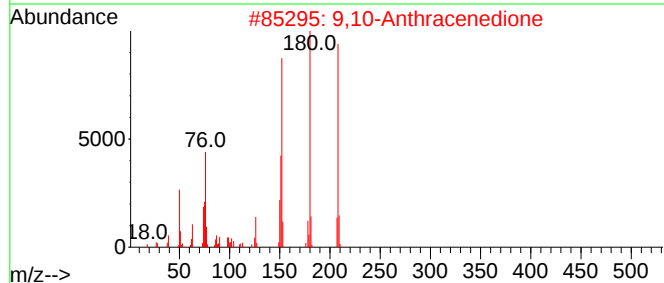
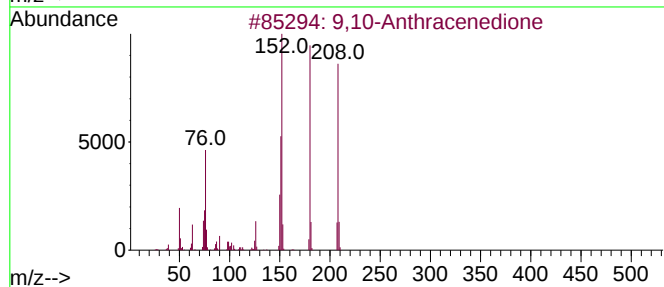
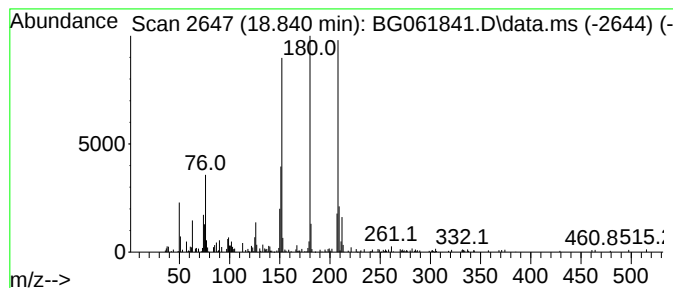
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 9,10-Anthracenedione Concentration Rank 8

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.840 | 4.67 ng/ul | 277261 | Phenanthrene-d10 | 17.430 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | 9,10-Anthracenedione | 208 | C14H8O2 | 000084-65-1 | 93   |
| 2    |    |   | 9,10-Anthracenedione | 208 | C14H8O2 | 000084-65-1 | 93   |
| 3    |    |   | 9,10-Anthracenedione | 208 | C14H8O2 | 000084-65-1 | 83   |
| 4    |    |   | Benzo[c]cinnoline    | 180 | C12H8N2 | 000230-17-1 | 83   |
| 5    |    |   | 9H-Fluoren-9-one     | 180 | C13H8O  | 000486-25-9 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
Acq On : 2 Jul 2024 11:02  
Operator : MA/JU  
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Misc :  
ALS Vial : 38 Sample Multiplier: 1

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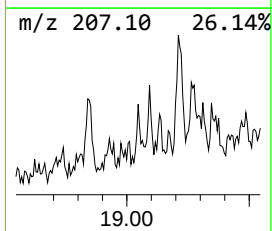
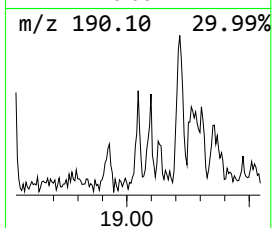
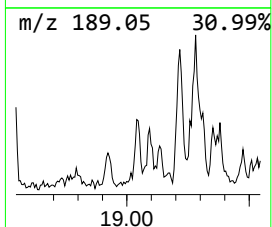
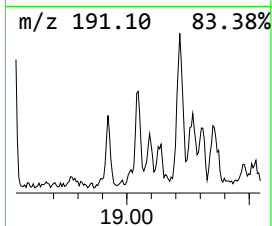
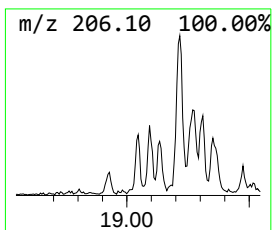
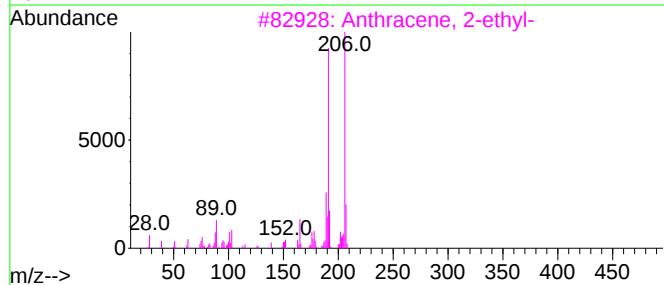
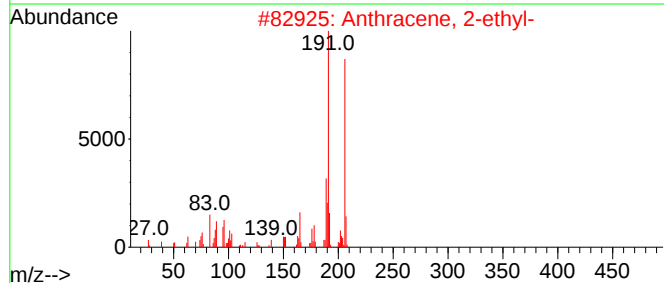
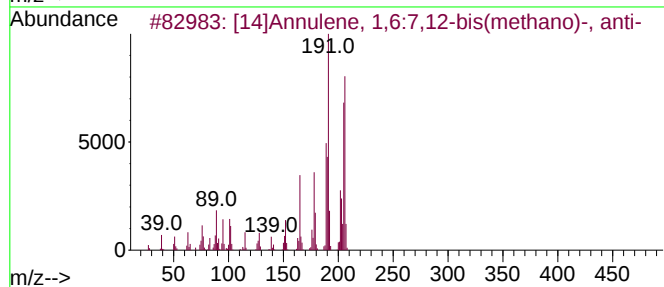
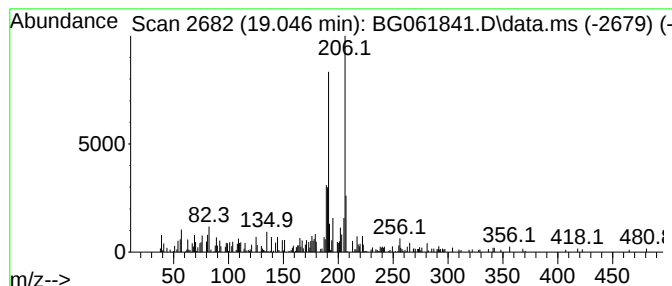
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 [14]Annulene, 1,6:7,12-bis(...) Concentration Rank 22

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 19.046    | 2.06 ng/ul                          | 122112 | Phenanthrene-d10 | 17.430      |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | [14]Annulene, 1,6:7,12-bis(metha... | 206    | C16H14           | 085440-62-6 | 90   |
| 2         | Anthracene, 2-ethyl-                | 206    | C16H14           | 052251-71-5 | 83   |
| 3         | Anthracene, 2-ethyl-                | 206    | C16H14           | 052251-71-5 | 83   |
| 4         | 1-Methyl-3-phenyl-1H-indene         | 206    | C16H14           | 022360-62-9 | 80   |
| 5         | Anthracene, 2-ethyl-                | 206    | C16H14           | 052251-71-5 | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
Acq On : 2 Jul 2024 11:02  
Operator : MA/JU  
Sample : P2963-07  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
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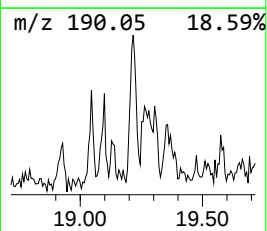
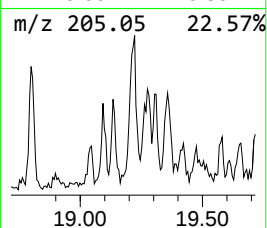
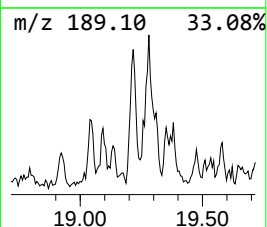
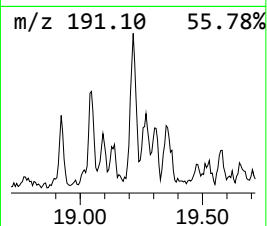
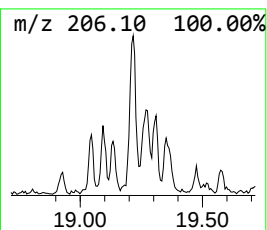
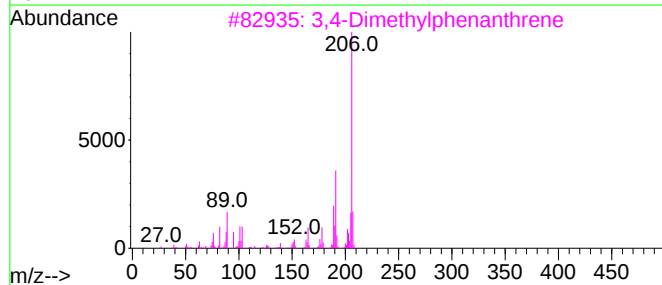
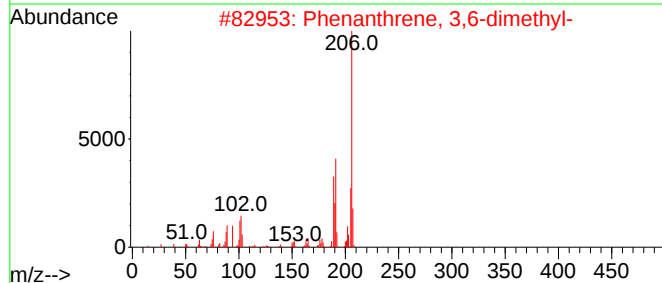
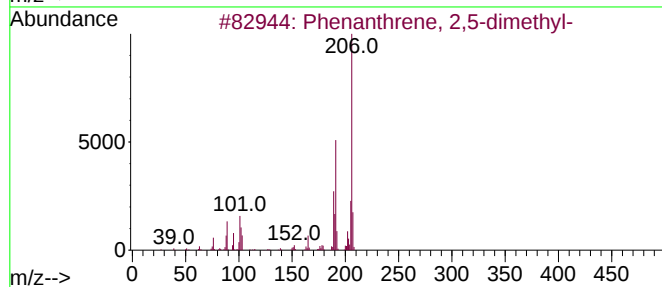
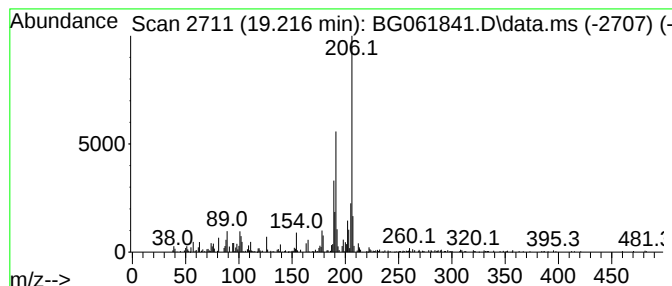
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Phenanthrene, 2,5-dimethyl- Concentration Rank 5

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 19.216 | 6.67 ng/ul | 395844 | Phenanthrene-d10 | 17.430 |

| Hit# | of 5 | Tentative ID                | MW  | MolForm | CAS#         | Qual |
|------|------|-----------------------------|-----|---------|--------------|------|
| 1    |      | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6  | 95   |
| 2    |      | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6  | 93   |
| 3    |      | 3,4-Dimethylphenanthrene    | 206 | C16H14  | 1000452-24-5 | 93   |
| 4    |      | Anthracene, 1,4-dimethyl-   | 206 | C16H14  | 000781-92-0  | 93   |
| 5    |      | Phenanthrene, 2,7-dimethyl- | 206 | C16H14  | 001576-69-8  | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
Acq On : 2 Jul 2024 11:02  
Operator : MA/JU  
Sample : P2963-07  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

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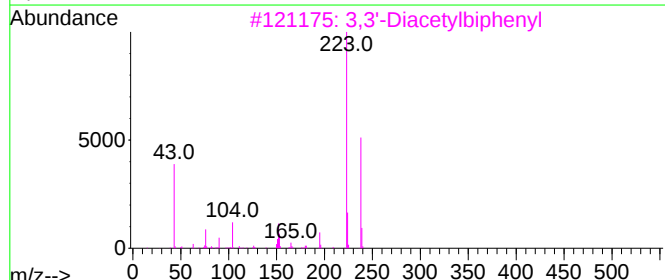
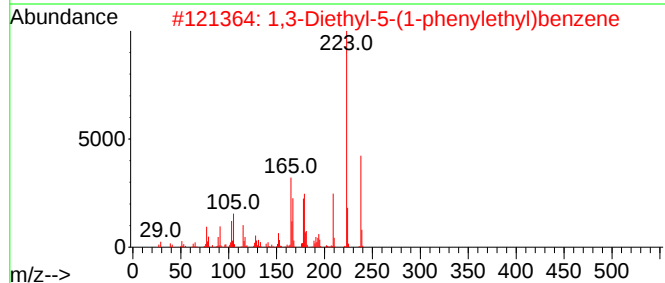
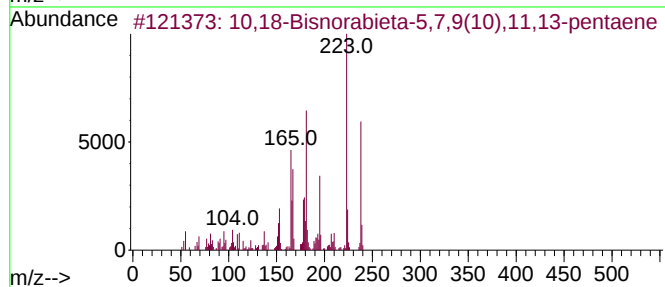
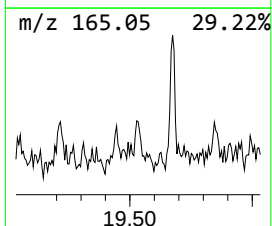
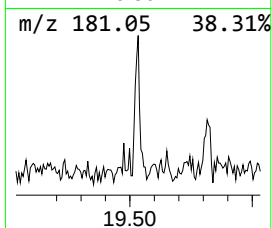
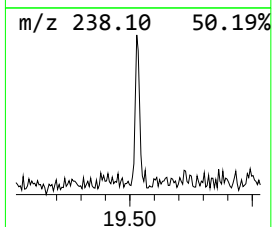
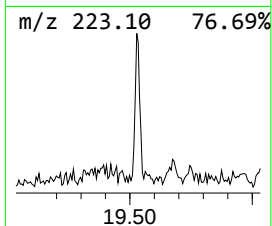
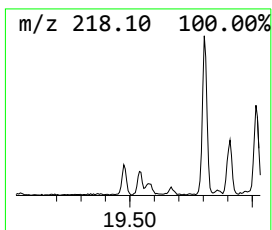
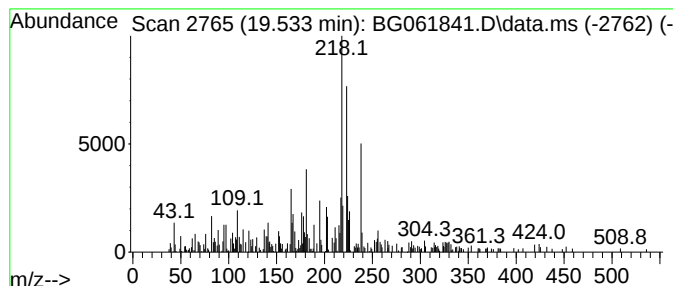
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 unknown-04 Concentration Rank 20

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 19.534 | 2.56 ng/ul | 151961 | Phenanthrene-d10 | 17.430 |

| Hit# | of 5                                | Tentative ID | MW       | MolForm      | CAS# | Qual |
|------|-------------------------------------|--------------|----------|--------------|------|------|
| 1    | 10,18-Bisnorabieta-5,7,9(10),11,... | 238          | C18H22   | 006566-19-4  | 49   |      |
| 2    | 1,3-Diethyl-5-(1-phenylethyl)ben... | 238          | C18H22   | 1000482-32-6 | 41   |      |
| 3    | 3,3'-Diacetylbiophenyl              | 238          | C16H14O2 | 094113-07-2  | 41   |      |
| 4    | Anthraquinone, 2-amino-             | 223          | C14H9NO2 | 000117-79-3  | 35   |      |
| 5    | .alpha.-Amyrone                     | 424          | C30H48O  | 000638-96-0  | 30   |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
Acq On : 2 Jul 2024 11:02  
Operator : MA/JU  
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ALS Vial : 38 Sample Multiplier: 1

Instrument :  
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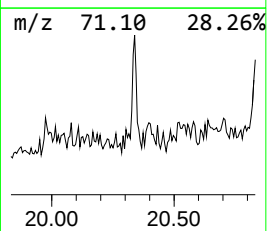
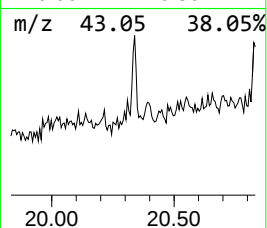
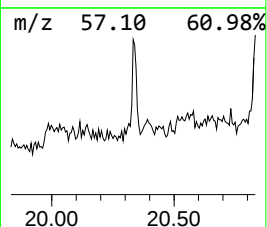
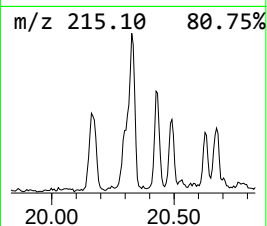
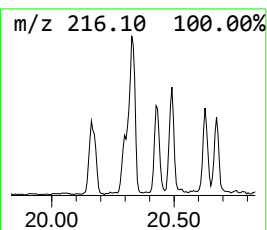
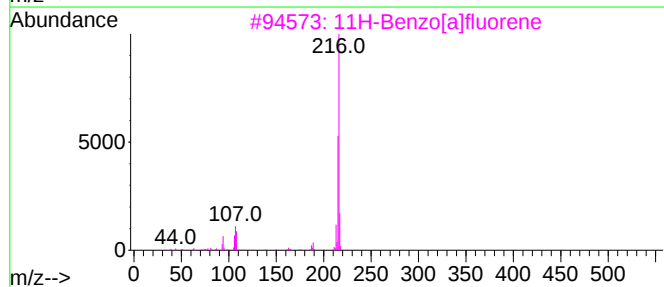
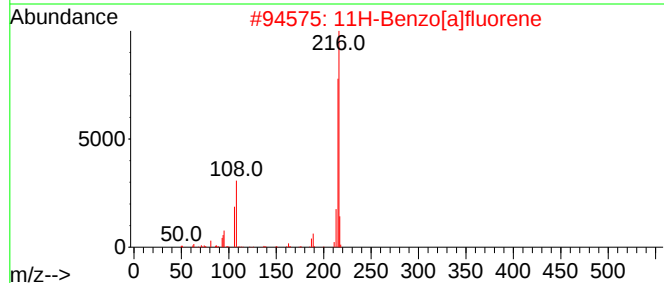
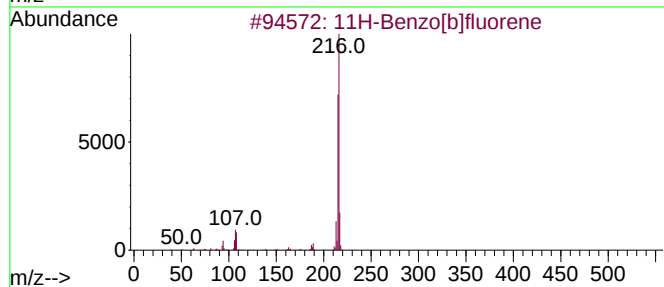
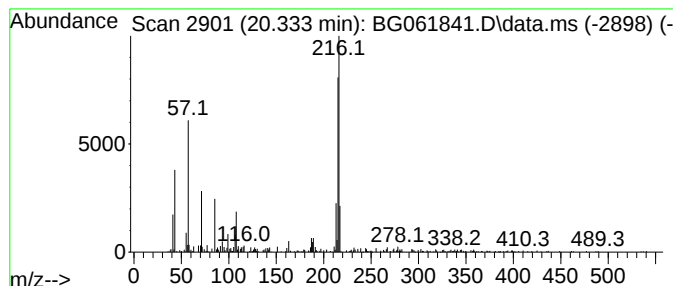
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 11H-Benzo[b]fluorene Concentration Rank 10

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 20.333 | 4.14 ng/ul | 592864 | Chrysene-d12     | 21.696 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | 11H-Benzo[b]fluorene | 216 | C17H12  | 000243-17-4 | 87   |
| 2    |    |   | 11H-Benzo[a]fluorene | 216 | C17H12  | 000238-84-6 | 83   |
| 3    |    |   | 11H-Benzo[a]fluorene | 216 | C17H12  | 000238-84-6 | 83   |
| 4    |    |   | 11H-Benzo[b]fluorene | 216 | C17H12  | 000243-17-4 | 81   |
| 5    |    |   | Pyrene, 1-methyl-    | 216 | C17H12  | 002381-21-7 | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
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Operator : MA/JU  
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ALS Vial : 38 Sample Multiplier: 1

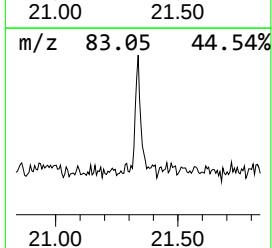
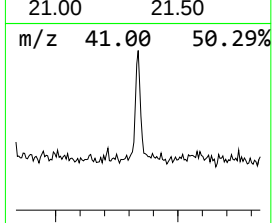
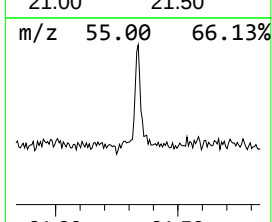
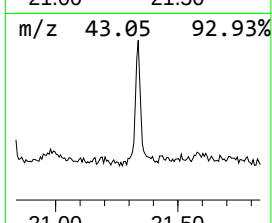
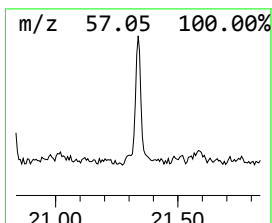
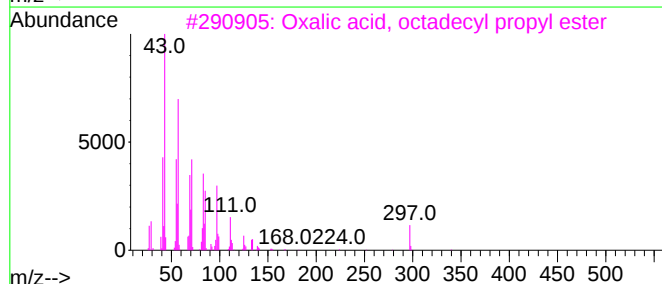
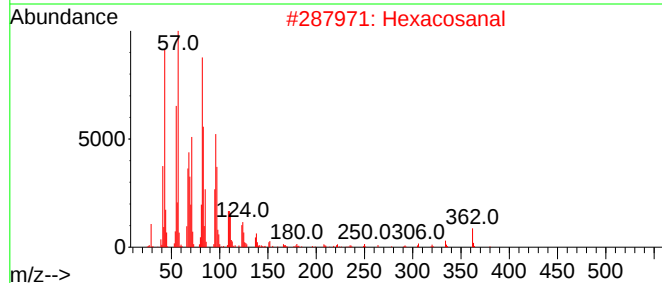
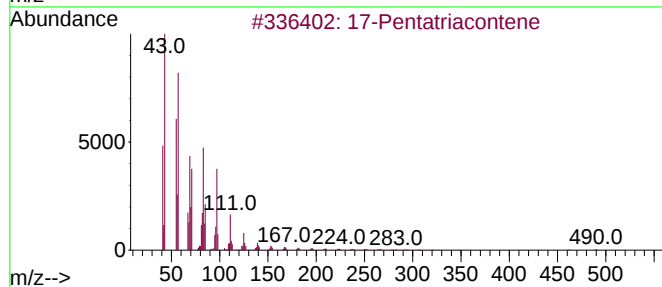
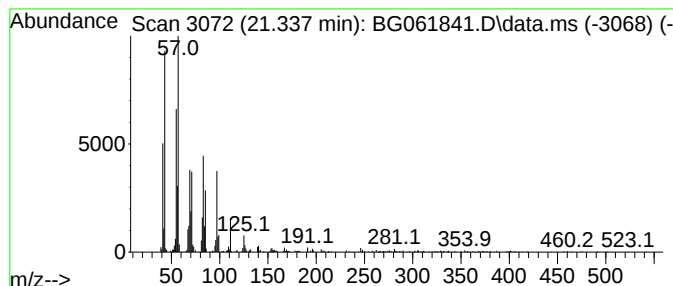
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T.      | EstConc                             | Area   | Relative to ISTD |          | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|----------|--------------|------|
| 21.337    | 4.15 ng/ul                          | 595191 | Chrysene-d12     |          | 21.696       |      |
| Hit# of 5 | Tentative ID                        |        | MW               | MolForm  | CAS#         | Qual |
| 1         | 17-Pentatriacontene                 |        | 491              | C35H70   | 006971-40-0  | 90   |
| 2         | Hexacosanal                         |        | 380              | C26H52O  | 026627-85-0  | 87   |
| 3         | Oxalic acid, octadecyl propyl ester |        | 384              | C23H44O4 | 1000309-27-0 | 87   |
| 4         | Oxalic acid, hexadecyl propyl ester |        | 356              | C21H40O4 | 1000309-26-9 | 83   |
| 5         | n-Tetracosanol-1                    |        | 354              | C24H50O  | 000506-51-4  | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061841.D  
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Operator : MA/JU  
Sample : P2963-07  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

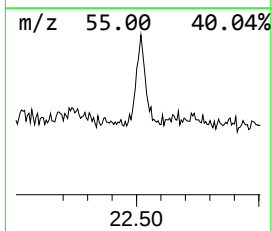
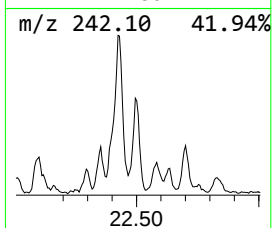
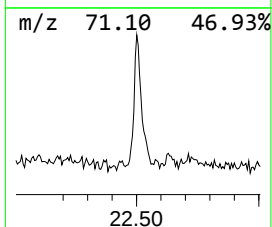
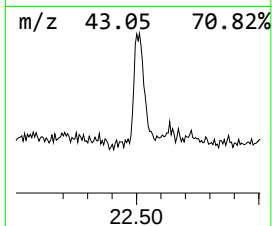
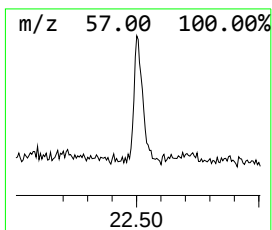
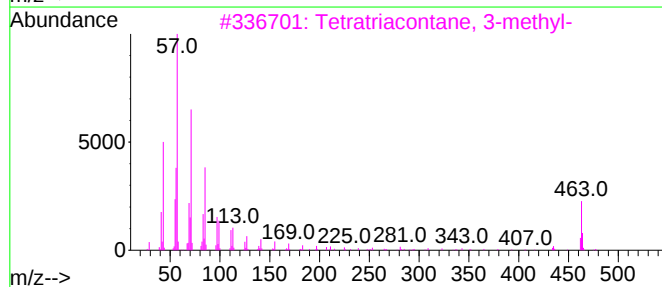
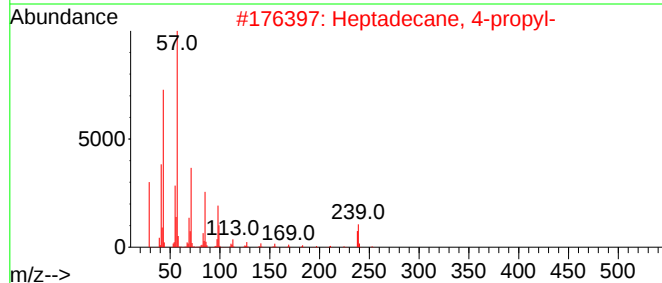
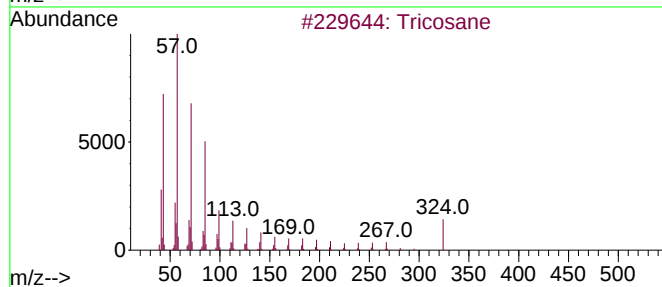
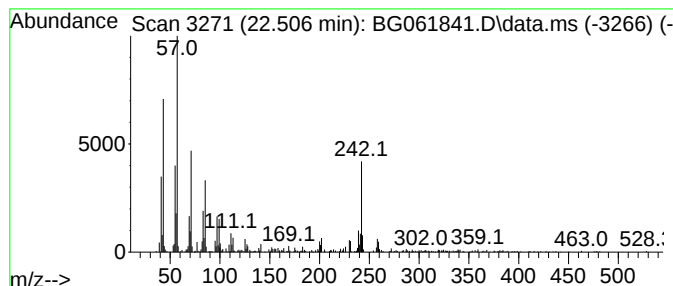
Instrument :  
BNA\_G  
ClientSampleId :  
A4CQ8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062024.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.      | EstConc                     | Area    | Relative to ISTD |         | R.T.         |      |
|-----------|-----------------------------|---------|------------------|---------|--------------|------|
| 22.506    | 8.08 ng/ul                  | 1157970 | Chrysene-d12     |         | 21.696       |      |
| Hit# of 5 | Tentative ID                |         | MW               | MolForm | CAS#         | Qual |
| 1         | Tricosane                   |         | 324              | C23H48  | 000638-67-5  | 46   |
| 2         | Heptadecane, 4-propyl-      |         | 282              | C20H42  | 055044-10-5  | 46   |
| 3         | Tetratriacontane, 3-methyl- |         | 493              | C35H72  | 066309-88-4  | 37   |
| 4         | 3,3-Diethylpentadecane      |         | 268              | C19H40  | 1000360-41-1 | 32   |
| 5         | Nonadecane, 2-methyl-       |         | 282              | C20H42  | 001560-86-7  | 32   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
 Data File : BG061841.D  
 Acq On : 2 Jul 2024 11:02  
 Operator : MA/JU  
 Sample : P2963-07  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 A4CQ8

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062024.MA.M  
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
|                    |        |         |       |          | #                     | RT     | Resp    | Conc |
| Methacrylamide     | 3.987  | 3.9     | ng/ul | 98955    | 1                     | 8.059  | 502537  | 20.0 |
| unknown-01         | 5.133  | 5.2     | ng/ul | 130630   | 1                     | 8.059  | 502537  | 20.0 |
| (DEL) Alkane: S... | 16.402 | 3.0     | ng/ul | 180757   | 4                     | 17.430 | 1187500 | 20.0 |
| 9H-Fluoren-9-one   | 17.083 | 2.6     | ng/ul | 156581   | 4                     | 17.430 | 1187500 | 20.0 |
| unknown-02         | 17.177 | 3.3     | ng/ul | 195658   | 4                     | 17.430 | 1187500 | 20.0 |
| Dibenzothiophene   | 17.248 | 3.3     | ng/ul | 193300   | 4                     | 17.430 | 1187500 | 20.0 |
| 1,1'-Biphenyl, ... | 18.018 | 3.6     | ng/ul | 213844   | 4                     | 17.430 | 1187500 | 20.0 |
| 1H-Cyclopropa[1... | 18.311 | 10.1    | ng/ul | 601386   | 4                     | 17.430 | 1187500 | 20.0 |
| Phenanthrene, 1... | 18.352 | 10.7    | ng/ul | 633925   | 4                     | 17.430 | 1187500 | 20.0 |
| Phenanthrene, 2... | 18.435 | 3.0     | ng/ul | 176886   | 4                     | 17.430 | 1187500 | 20.0 |
| Benzaldehyde, 3... | 18.499 | 13.1    | ng/ul | 776113   | 4                     | 17.430 | 1187500 | 20.0 |
| Naphtho[2,3-b]n... | 18.535 | 5.3     | ng/ul | 314646   | 4                     | 17.430 | 1187500 | 20.0 |
| unknown-03         | 18.605 | 2.4     | ng/ul | 139841   | 4                     | 17.430 | 1187500 | 20.0 |
| 5,16[1',2']:8,1... | 18.799 | 3.8     | ng/ul | 223073   | 4                     | 17.430 | 1187500 | 20.0 |
| 9,10-Anthracene... | 18.840 | 4.7     | ng/ul | 277261   | 4                     | 17.430 | 1187500 | 20.0 |
| [14]Annulene, 1... | 19.046 | 2.1     | ng/ul | 122112   | 4                     | 17.430 | 1187500 | 20.0 |
| Phenanthrene, 2... | 19.216 | 6.7     | ng/ul | 395844   | 4                     | 17.430 | 1187500 | 20.0 |
| unknown-04         | 19.534 | 2.6     | ng/ul | 151961   | 4                     | 17.430 | 1187500 | 20.0 |
| 11H-Benzo[b]flu... | 20.333 | 4.1     | ng/ul | 592864   | 5                     | 21.696 | 2866020 | 20.0 |
| (DEL) Alkane: S... | 21.337 | 4.2     | ng/ul | 595191   | 5                     | 21.696 | 2866020 | 20.0 |
| (DEL) Alkane: S... | 22.506 | 8.1     | ng/ul | 1157970  | 5                     | 21.696 | 2866020 | 20.0 |