

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080319\  
 Data File : BG041949.D  
 Acq On : 4 Aug 2019 19:31  
 Operator : HP/JU  
 Sample : K3962-10  
 Misc :  
 ALS Vial : 50 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BEP03

## Integration Parameters: LSCINT.P

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

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080319MA.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         | 3.158       | 7             | 12          | 39           | rVB      | 1399889        | 2431443       | 100.00%         | 8.791%        |
| 2         | 3.422       | 54            | 57          | 65           | rVB4     | 10578          | 19246         | 0.79%           | 0.070%        |
| 3         | 3.951       | 143           | 147         | 152          | rBV2     | 11023          | 16805         | 0.69%           | 0.061%        |
| 4         | 4.086       | 164           | 170         | 176          | rVB      | 10758          | 19935         | 0.82%           | 0.072%        |
| 5         | 7.182       | 688           | 697         | 711          | rBV      | 171068         | 341361        | 14.04%          | 1.234%        |
| 6         | 7.353       | 720           | 726         | 737          | rBV      | 127820         | 224942        | 9.25%           | 0.813%        |
| 7         | 7.564       | 755           | 762         | 775          | rVB      | 190219         | 343442        | 14.13%          | 1.242%        |
| 8         | 8.034       | 835           | 842         | 850          | rBV      | 218307         | 382548        | 15.73%          | 1.383%        |
| 9         | 8.739       | 955           | 962         | 977          | rBV      | 201512         | 381439        | 15.69%          | 1.379%        |
| 10        | 8.869       | 979           | 984         | 991          | rVB5     | 3954           | 8563          | 0.35%           | 0.031%        |
| 11        | 9.204       | 1033          | 1041        | 1051         | rBV      | 168914         | 321929        | 13.24%          | 1.164%        |
| 12        | 9.932       | 1157          | 1165        | 1177         | rBV      | 148282         | 287122        | 11.81%          | 1.038%        |
| 13        | 10.473      | 1250          | 1257        | 1269         | rBV      | 228733         | 454615        | 18.70%          | 1.644%        |
| 14        | 10.860      | 1313          | 1323        | 1331         | rBV      | 306149         | 565853        | 23.27%          | 2.046%        |
| 15        | 10.990      | 1338          | 1345        | 1360         | rBV      | 140929         | 288041        | 11.85%          | 1.041%        |
| 16        | 12.523      | 1601          | 1606        | 1614         | rBV2     | 6233           | 10921         | 0.45%           | 0.039%        |
| 17        | 12.741      | 1637          | 1643        | 1650         | rVB3     | 7226           | 12604         | 0.52%           | 0.046%        |
| 18        | 13.863      | 1829          | 1834        | 1842         | rVB5     | 4818           | 11692         | 0.48%           | 0.042%        |
| 19        | 14.016      | 1853          | 1860        | 1865         | rBV4     | 11448          | 20243         | 0.83%           | 0.073%        |
| 20        | 14.092      | 1865          | 1873        | 1877         | rVV      | 445800         | 667178        | 27.44%          | 2.412%        |
| 21        | 14.139      | 1877          | 1881        | 1892         | rVV      | 219366         | 337319        | 13.87%          | 1.220%        |
| 22        | 14.256      | 1896          | 1901        | 1905         | rVV3     | 6642           | 11505         | 0.47%           | 0.042%        |
| 23        | 14.386      | 1916          | 1923        | 1938         | rVB2     | 510376         | 862180        | 35.46%          | 3.117%        |
| 24        | 14.697      | 1966          | 1976        | 1982         | rBV      | 496109         | 804705        | 33.10%          | 2.909%        |
| 25        | 14.756      | 1982          | 1986        | 1995         | rVB2     | 22784          | 39422         | 1.62%           | 0.143%        |
| 26        | 14.879      | 2000          | 2007        | 2021         | rBV      | 107813         | 233294        | 9.59%           | 0.843%        |
| 27        | 15.097      | 2039          | 2044        | 2050         | rVB4     | 24011          | 38238         | 1.57%           | 0.138%        |
| 28        | 15.173      | 2052          | 2057        | 2060         | rVB3     | 8207           | 13446         | 0.55%           | 0.049%        |
| 29        | 15.314      | 2074          | 2081        | 2085         | rBV5     | 6640           | 11433         | 0.47%           | 0.041%        |
| 30        | 15.367      | 2086          | 2090        | 2094         | rVB2     | 7385           | 9441          | 0.39%           | 0.034%        |
| 31        | 15.484      | 2101          | 2110        | 2118         | rBV8     | 8502           | 24307         | 1.00%           | 0.088%        |
| 32        | 15.690      | 2136          | 2145        | 2151         | rVV      | 700404         | 1082347       | 44.51%          | 3.913%        |
| 33        | 15.743      | 2151          | 2154        | 2159         | rVV3     | 26799          | 40502         | 1.67%           | 0.146%        |
| 34        | 15.802      | 2159          | 2164        | 2172         | rVV      | 120828         | 198202        | 8.15%           | 0.717%        |

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 Max Peaks: 100  
 Peak Location: TOP

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 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080319MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |       |        |         |        |        |
|----|--------|------|------|------|-------|--------|---------|--------|--------|
| 35 | 15.860 | 2172 | 2174 | 2178 | rVV4  | 10169  | 15713   | 0.65%  | 0.057% |
| 36 | 15.907 | 2178 | 2182 | 2184 | rVV4  | 9633   | 15516   | 0.64%  | 0.056% |
| 37 | 15.954 | 2187 | 2190 | 2194 | rVB5  | 9997   | 14820   | 0.61%  | 0.054% |
| 38 | 16.043 | 2200 | 2205 | 2213 | rVB7  | 9757   | 24700   | 1.02%  | 0.089% |
| 39 | 16.266 | 2238 | 2243 | 2246 | rBV5  | 4926   | 9384    | 0.39%  | 0.034% |
| 40 | 16.419 | 2262 | 2269 | 2275 | rVV8  | 17442  | 38130   | 1.57%  | 0.138% |
| 41 | 16.560 | 2288 | 2293 | 2296 | rBV6  | 9042   | 14402   | 0.59%  | 0.052% |
| 42 | 16.748 | 2323 | 2325 | 2330 | rVV4  | 10994  | 17773   | 0.73%  | 0.064% |
| 43 | 16.795 | 2330 | 2333 | 2339 | rVV4  | 11276  | 19496   | 0.80%  | 0.070% |
| 44 | 16.853 | 2340 | 2343 | 2346 | rVV5  | 6615   | 8329    | 0.34%  | 0.030% |
| 45 | 16.906 | 2349 | 2352 | 2356 | rVB5  | 10588  | 14141   | 0.58%  | 0.051% |
| 46 | 16.959 | 2356 | 2361 | 2362 | rBV4  | 10283  | 12910   | 0.53%  | 0.047% |
| 47 | 17.024 | 2367 | 2372 | 2375 | rVB7  | 11367  | 14123   | 0.58%  | 0.051% |
| 48 | 17.100 | 2380 | 2385 | 2392 | rVB5  | 14030  | 24478   | 1.01%  | 0.088% |
| 49 | 17.200 | 2392 | 2402 | 2406 | rBV10 | 12444  | 33991   | 1.40%  | 0.123% |
| 50 | 17.265 | 2409 | 2413 | 2421 | rVB   | 21340  | 33416   | 1.37%  | 0.121% |
| 51 | 17.335 | 2421 | 2425 | 2427 | rBV4  | 9016   | 11631   | 0.48%  | 0.042% |
| 52 | 17.447 | 2435 | 2444 | 2448 | rBV   | 638617 | 986557  | 40.57% | 3.567% |
| 53 | 17.488 | 2448 | 2451 | 2456 | rVV   | 298191 | 438914  | 18.05% | 1.587% |
| 54 | 17.547 | 2456 | 2461 | 2471 | rVV2  | 722381 | 1205729 | 49.59% | 4.359% |
| 55 | 17.635 | 2472 | 2476 | 2482 | rVV3  | 18717  | 41898   | 1.72%  | 0.151% |
| 56 | 17.688 | 2482 | 2485 | 2487 | rVV3  | 8442   | 12171   | 0.50%  | 0.044% |
| 57 | 17.717 | 2487 | 2490 | 2491 | rVV3  | 9971   | 10959   | 0.45%  | 0.040% |
| 58 | 17.752 | 2494 | 2496 | 2501 | rVB5  | 13154  | 18158   | 0.75%  | 0.066% |
| 59 | 17.852 | 2509 | 2513 | 2516 | rBV3  | 8513   | 13378   | 0.55%  | 0.048% |
| 60 | 18.028 | 2539 | 2543 | 2550 | rBV3  | 42307  | 63009   | 2.59%  | 0.228% |
| 61 | 18.175 | 2563 | 2568 | 2572 | rBV   | 19236  | 34589   | 1.42%  | 0.125% |
| 62 | 18.334 | 2589 | 2595 | 2598 | rBV2  | 155829 | 274538  | 11.29% | 0.993% |
| 63 | 18.375 | 2598 | 2602 | 2607 | rVV   | 158958 | 236205  | 9.71%  | 0.854% |
| 64 | 18.457 | 2611 | 2616 | 2621 | rBV   | 47083  | 79398   | 3.27%  | 0.287% |
| 65 | 18.516 | 2621 | 2626 | 2630 | rVV2  | 167401 | 309107  | 12.71% | 1.118% |
| 66 | 18.557 | 2630 | 2633 | 2640 | rVB   | 97627  | 166378  | 6.84%  | 0.602% |
| 67 | 18.640 | 2643 | 2647 | 2650 | rVV6  | 16369  | 20944   | 0.86%  | 0.076% |
| 68 | 18.669 | 2650 | 2652 | 2657 | rVB6  | 14162  | 18187   | 0.75%  | 0.066% |
| 69 | 18.728 | 2657 | 2662 | 2666 | rBV3  | 23699  | 45108   | 1.86%  | 0.163% |
| 70 | 18.822 | 2673 | 2678 | 2681 | rBV2  | 66264  | 98725   | 4.06%  | 0.357% |
| 71 | 18.857 | 2681 | 2684 | 2694 | rVB4  | 53805  | 91147   | 3.75%  | 0.330% |

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Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080319MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |        |        |
|-----|--------|------|------|------|------|---------|---------|--------|--------|
| 72  | 19.068 | 2712 | 2720 | 2725 | rBV3 | 61422   | 124932  | 5.14%  | 0.452% |
| 73  | 19.121 | 2725 | 2729 | 2732 | rVV  | 43647   | 68205   | 2.81%  | 0.247% |
| 74  | 19.157 | 2732 | 2735 | 2739 | rVB2 | 36494   | 49934   | 2.05%  | 0.181% |
| 75  | 19.239 | 2744 | 2749 | 2753 | rBV  | 130470  | 215231  | 8.85%  | 0.778% |
| 76  | 19.374 | 2768 | 2772 | 2780 | rBV2 | 72707   | 150330  | 6.18%  | 0.543% |
| 77  | 19.503 | 2789 | 2794 | 2800 | rVB  | 552792  | 799489  | 32.88% | 2.890% |
| 78  | 19.562 | 2800 | 2804 | 2807 | rBV4 | 46547   | 68344   | 2.81%  | 0.247% |
| 79  | 19.597 | 2808 | 2810 | 2815 | rVB4 | 35072   | 46408   | 1.91%  | 0.168% |
| 80  | 19.744 | 2832 | 2835 | 2838 | rBV2 | 25820   | 32550   | 1.34%  | 0.118% |
| 81  | 19.838 | 2845 | 2851 | 2853 | rBV  | 958633  | 1534842 | 63.12% | 5.549% |
| 82  | 19.862 | 2853 | 2855 | 2864 | rVV  | 717404  | 1014958 | 41.74% | 3.669% |
| 83  | 19.938 | 2864 | 2868 | 2874 | rVB2 | 59468   | 103695  | 4.26%  | 0.375% |
| 84  | 20.191 | 2905 | 2911 | 2917 | rBV2 | 97390   | 246600  | 10.14% | 0.892% |
| 85  | 20.320 | 2930 | 2933 | 2934 | rBV3 | 53553   | 57187   | 2.35%  | 0.207% |
| 86  | 20.355 | 2936 | 2939 | 2945 | rVB  | 150893  | 214437  | 8.82%  | 0.775% |
| 87  | 20.455 | 2952 | 2956 | 2960 | rBV  | 95606   | 125600  | 5.17%  | 0.454% |
| 88  | 20.514 | 2962 | 2966 | 2970 | rVB  | 134150  | 179669  | 7.39%  | 0.650% |
| 89  | 20.655 | 2986 | 2990 | 2993 | rBV  | 107695  | 145595  | 5.99%  | 0.526% |
| 90  | 20.696 | 2994 | 2997 | 3002 | rVB  | 103661  | 141520  | 5.82%  | 0.512% |
| 91  | 21.354 | 3105 | 3109 | 3110 | rBV  | 63557   | 77878   | 3.20%  | 0.282% |
| 92  | 21.595 | 3145 | 3150 | 3151 | rBV2 | 174855  | 246591  | 10.14% | 0.892% |
| 93  | 21.624 | 3151 | 3155 | 3162 | rVV  | 1041695 | 1521027 | 62.56% | 5.499% |
| 94  | 21.730 | 3165 | 3173 | 3178 | rVV2 | 758149  | 1891238 | 77.78% | 6.838% |
| 95  | 21.777 | 3178 | 3181 | 3190 | rVB  | 385990  | 659771  | 27.13% | 2.385% |
| 96  | 23.969 | 3548 | 3554 | 3562 | rBV  | 183771  | 611744  | 25.16% | 2.212% |
| 97  | 24.104 | 3573 | 3577 | 3584 | rVB2 | 90556   | 162093  | 6.67%  | 0.586% |
| 98  | 24.703 | 3674 | 3679 | 3685 | rBV2 | 120989  | 286164  | 11.77% | 1.035% |
| 99  | 24.791 | 3686 | 3694 | 3700 | rBV  | 363657  | 958723  | 39.43% | 3.466% |
| 100 | 25.014 | 3724 | 3732 | 3740 | rBV2 | 347187  | 920556  | 37.86% | 3.328% |

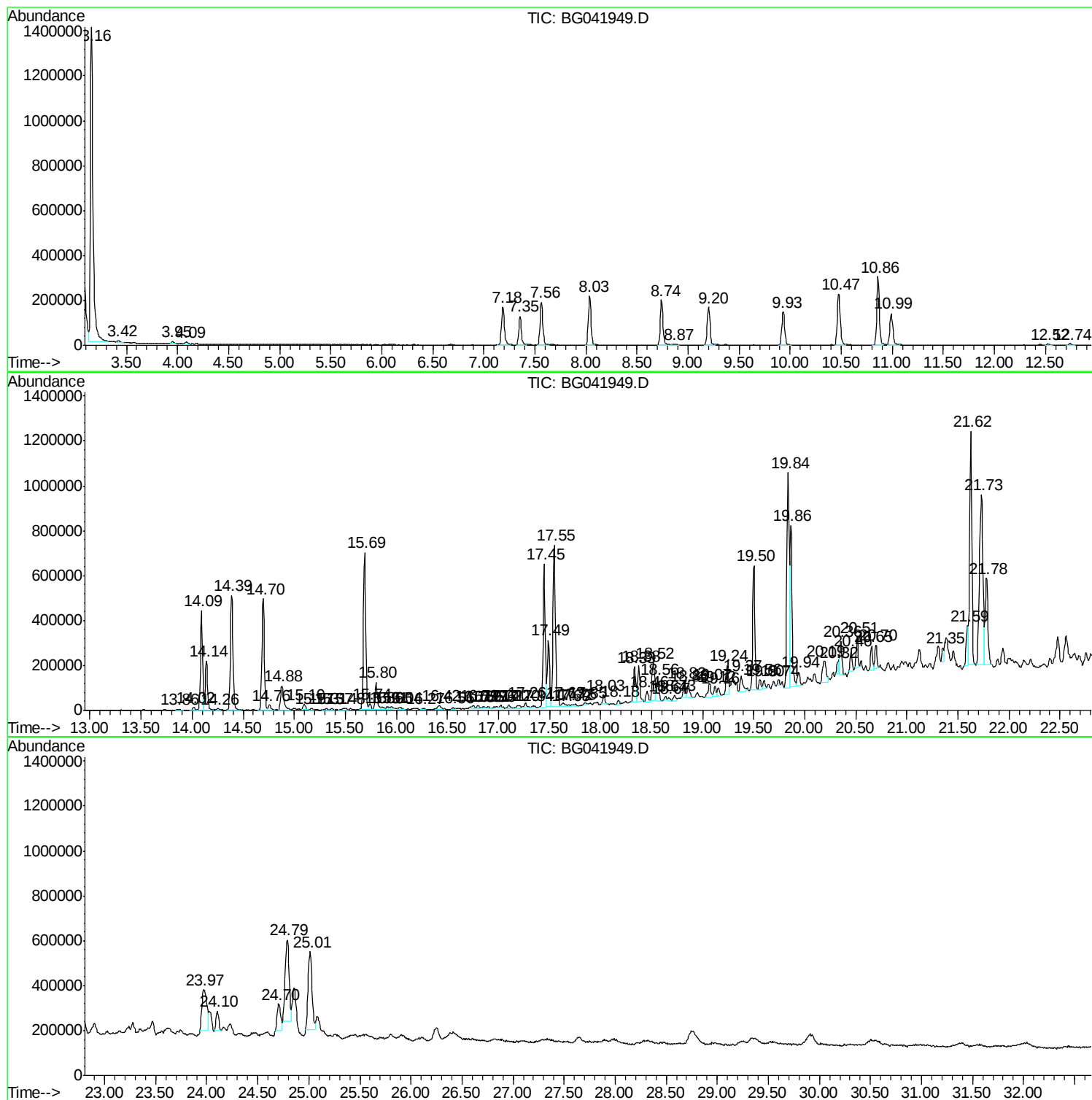
Sum of corrected areas: 27659626

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080319MA.M  
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080319\  
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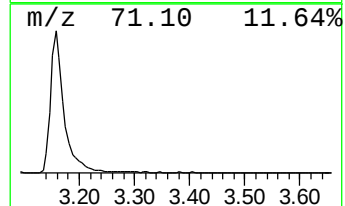
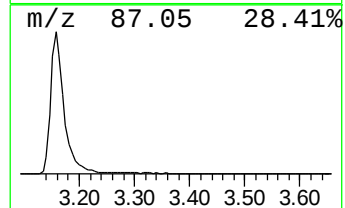
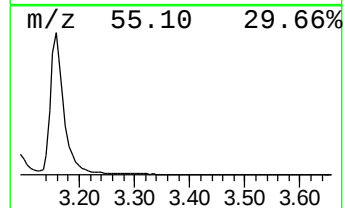
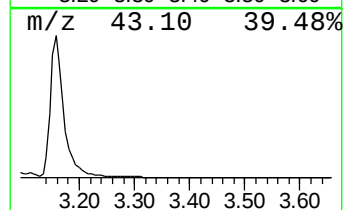
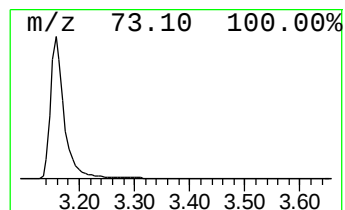
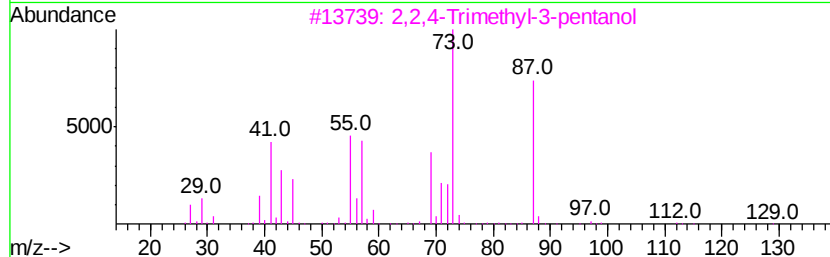
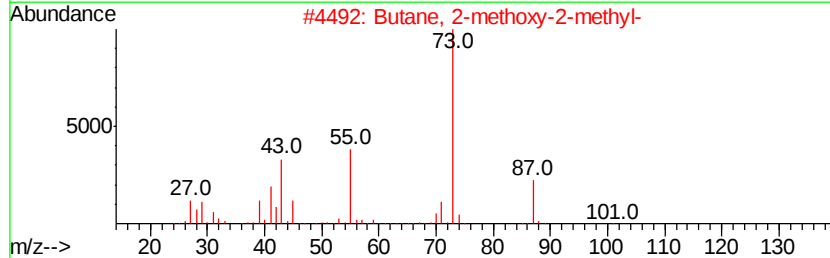
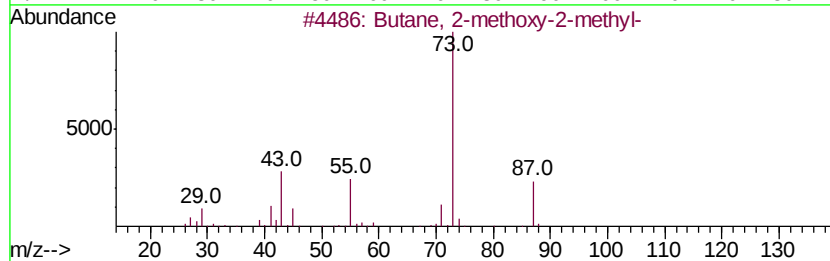
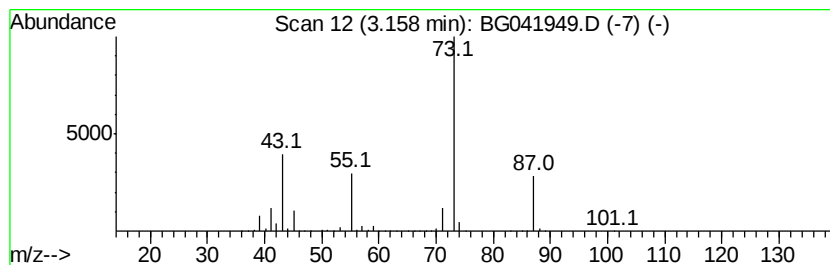
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080319MA.M  
Quant Title : SVOA CALIBRATION

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

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

| R.T.    | EstConc                     | Area         | Relative to ISTD       | R.T.      |
|---------|-----------------------------|--------------|------------------------|-----------|
| 3.16    | 127.12 ng/ul                | 2431440      | 1,4-Dichlorobenzene-d4 | 8.03      |
| Hit# of | 5                           | Tentative ID | MW MolForm             | CAS# Qual |
| 1       | Butane, 2-methoxy-2-methyl- | 102 C6H14O   | 000994-05-8            | 83        |
| 2       | Butane, 2-methoxy-2-methyl- | 102 C6H14O   | 000994-05-8            | 59        |
| 3       | 2,2,4-Trimethyl-3-pentanol  | 130 C8H18O   | 005162-48-1            | 39        |
| 4       | Pentane, 3-methoxy-         | 102 C6H14O   | 036839-67-5            | 12        |
| 5       | 1,3-Dioxolane, 2-methyl-    | 88 C4H8O2    | 000497-26-7            | 9         |



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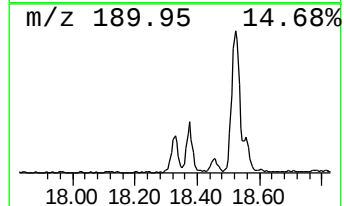
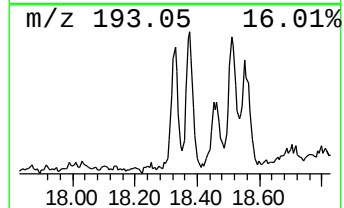
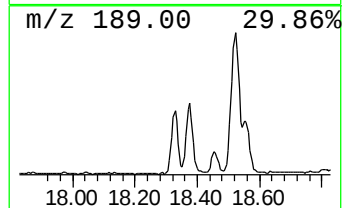
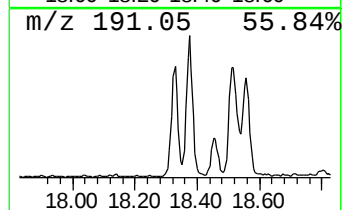
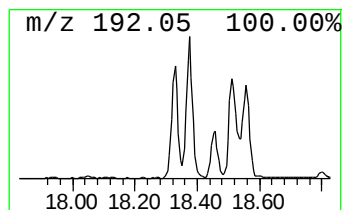
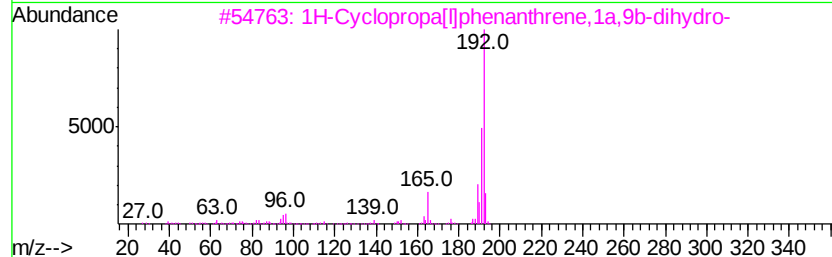
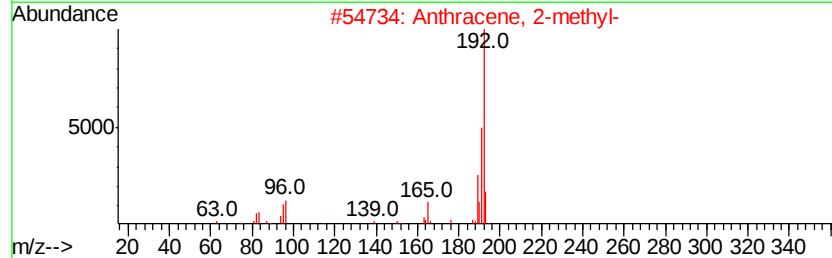
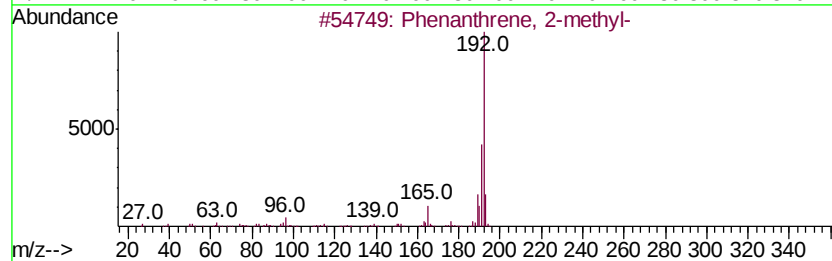
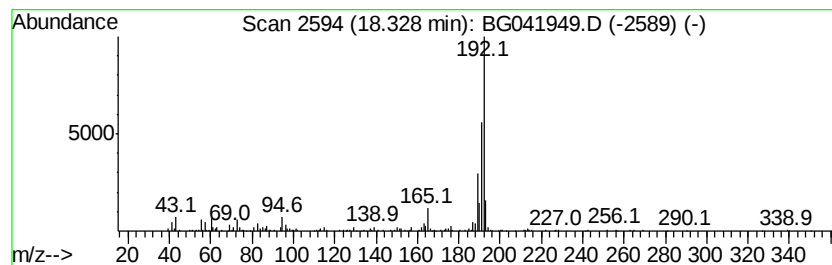
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 Phenanthrene, 2-methyl- Concentration Rank 4

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.33 | 5.57 ng/ul | 274538 | Phenanthrene-d10 | 17.45 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 96   |
| 2    |    | Anthracene, 2-methyl-               | 192 | C15H12  | 000613-12-7 | 93   |
| 3    |    | 1H-Cyclopropa[1]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 92   |
| 4    |    | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 89   |
| 5    |    | 1H-Indene, 2-phenyl-                | 192 | C15H12  | 004505-48-0 | 89   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080319\  
Data File : BG041949.D  
Acq On : 4 Aug 2019 19:31  
Operator : HP/JU  
Sample : K3962-10  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampled :  
BEP03

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Quant Title : SVOA CALIBRATION

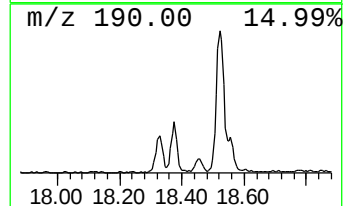
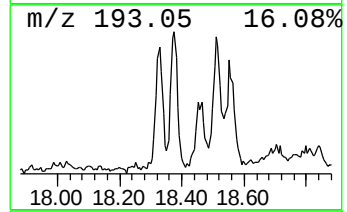
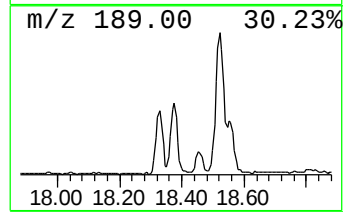
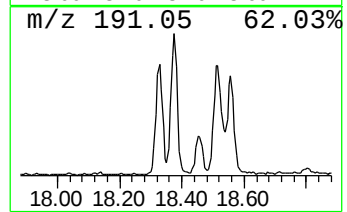
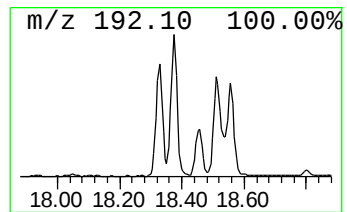
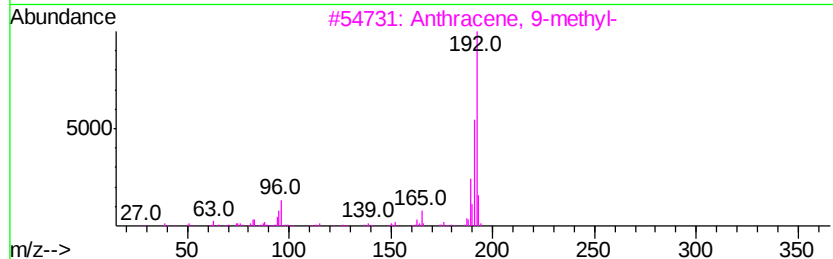
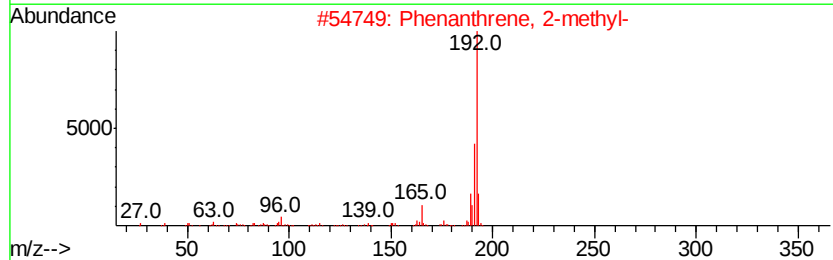
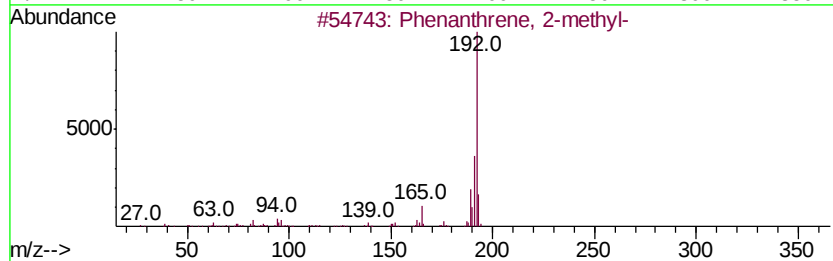
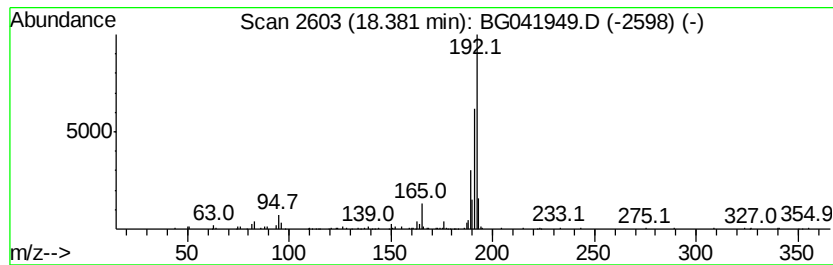
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TIC Integration Parameters: LSCINT.P

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Peak Number 3 Anthracene, 9-methyl- Concentration Rank 5

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.38 | 4.79 ng/ul | 236205 | Phenanthrene-d10 | 17.45 |

| Hit# of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|---------|---|-------------------------|-----|---------|-------------|------|
| 1       |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 95   |
| 2       |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 95   |
| 3       |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 94   |
| 4       |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 93   |
| 5       |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080319\  
Data File : BG041949.D  
Acq On : 4 Aug 2019 19:31  
Operator : HP/JU  
Sample : K3962-10  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

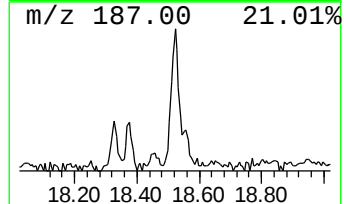
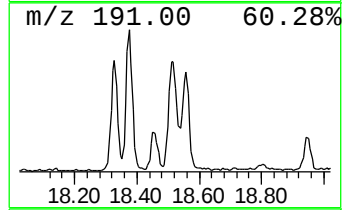
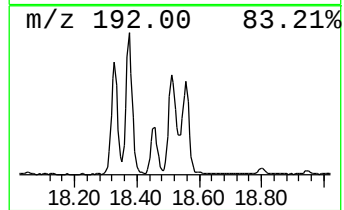
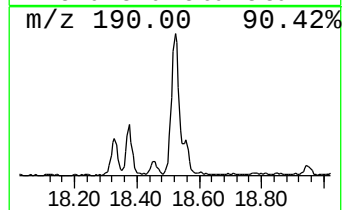
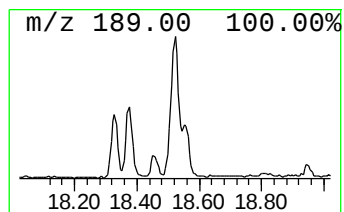
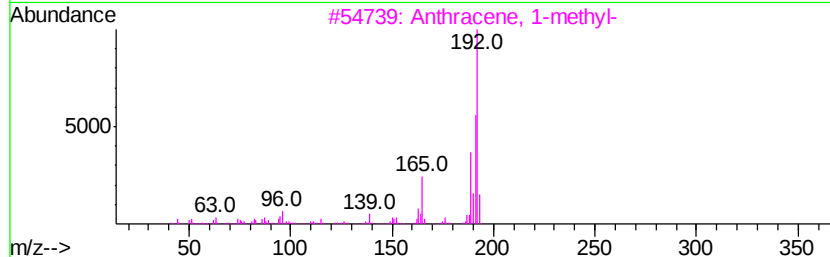
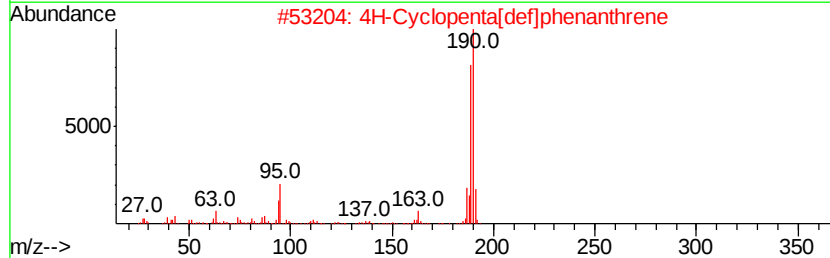
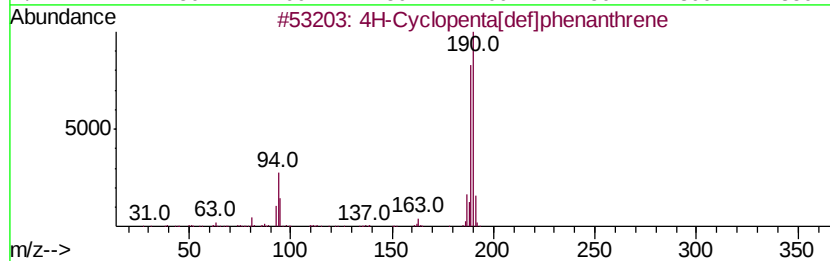
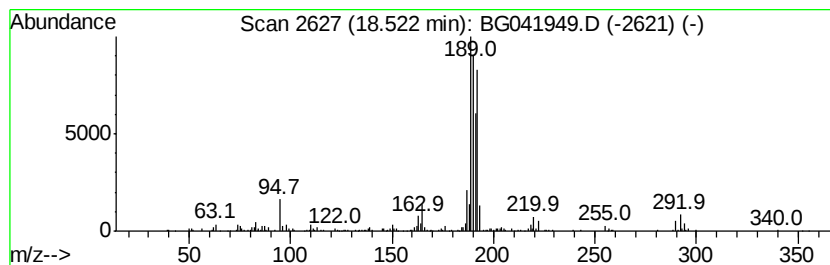
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ClientSampleId :  
BEP03

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Quant Title : SVOA CALIBRATION

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

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Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

| R.T.    | EstConc    | Area                           | Relative to ISTD | R.T.    |             |      |
|---------|------------|--------------------------------|------------------|---------|-------------|------|
| 18.52   | 6.27 ng/ul | 309107                         | Phenanthrene-d10 | 17.45   |             |      |
| Hit# of | 5          | Tentative ID                   | MW               | MolForm | CAS#        | Qual |
| 1       |            | 4H-Cyclopenta[def]phenanthrene | 190              | C15H10  | 000203-64-5 | 50   |
| 2       |            | 4H-Cyclopenta[def]phenanthrene | 190              | C15H10  | 000203-64-5 | 50   |
| 3       |            | Anthracene, 1-methyl-          | 192              | C15H12  | 000610-48-0 | 46   |
| 4       |            | Anthracene, 1-methyl-          | 192              | C15H12  | 000610-48-0 | 46   |
| 5       |            | Anthracene, 1-methyl-          | 192              | C15H12  | 000610-48-0 | 41   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080319\  
Data File : BG041949.D  
Acq On : 4 Aug 2019 19:31  
Operator : HP/JU  
Sample : K3962-10  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BEP03

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080319MA.M  
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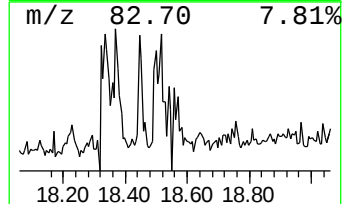
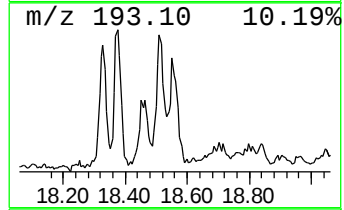
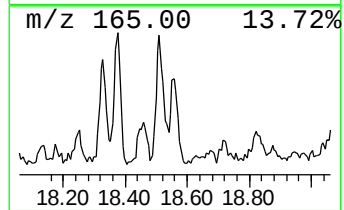
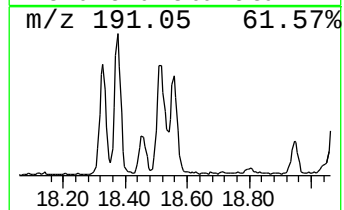
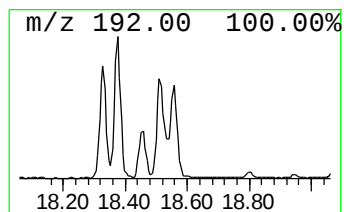
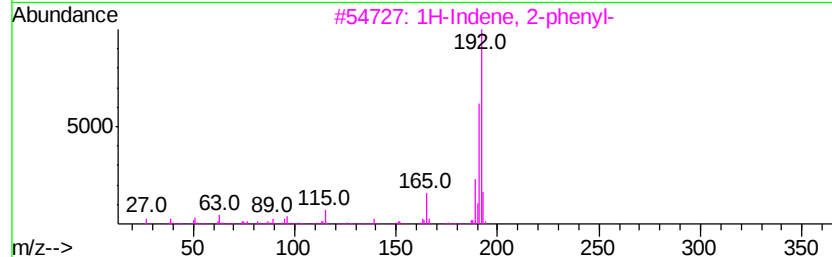
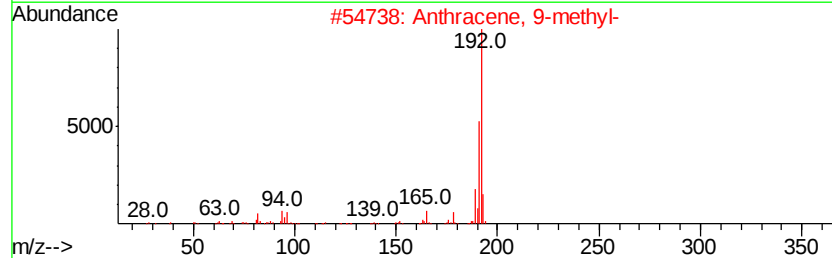
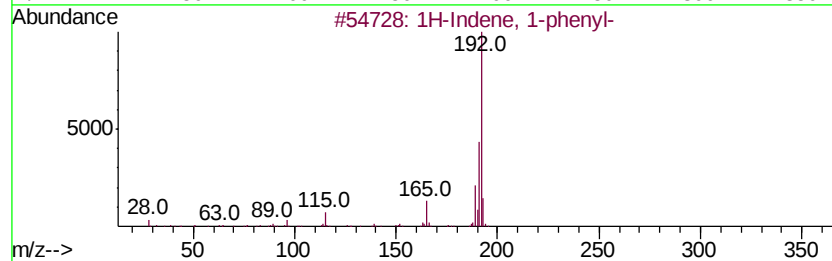
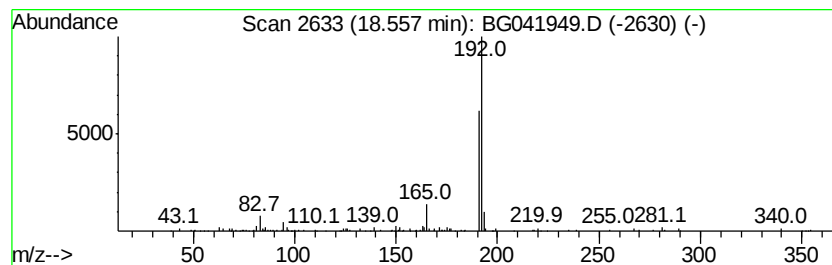
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 1H-Indene, 1-phenyl- Concentration Rank 8

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.56 | 3.37 ng/ul | 166378 | Phenanthrene-d10 | 17.45 |

| Hit# of | 5 | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|---------|---|---------------------------------|-----|---------|-------------|------|
| 1       |   | 1H-Indene, 1-phenyl-            | 192 | C15H12  | 001961-96-2 | 80   |
| 2       |   | Anthracene, 9-methyl-           | 192 | C15H12  | 000779-02-2 | 80   |
| 3       |   | 1H-Indene, 2-phenyl-            | 192 | C15H12  | 004505-48-0 | 80   |
| 4       |   | 3,4-Methylenedioxycinnamic acid | 192 | C10H8O4 | 002373-80-0 | 80   |
| 5       |   | Phenanthrene, 2-methyl-         | 192 | C15H12  | 002531-84-2 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080319\  
Data File : BG041949.D  
Acq On : 4 Aug 2019 19:31  
Operator : HP/JU  
Sample : K3962-10  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

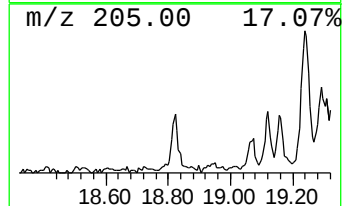
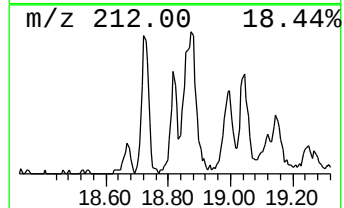
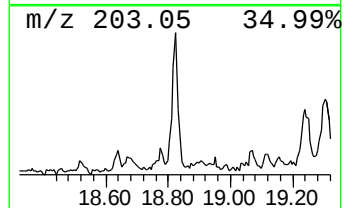
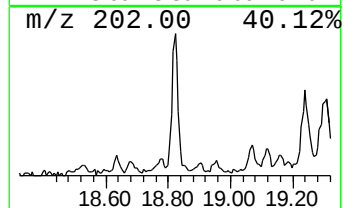
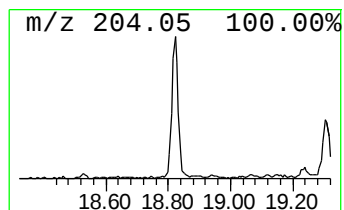
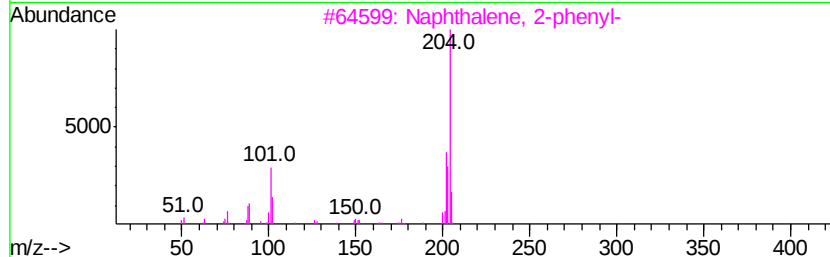
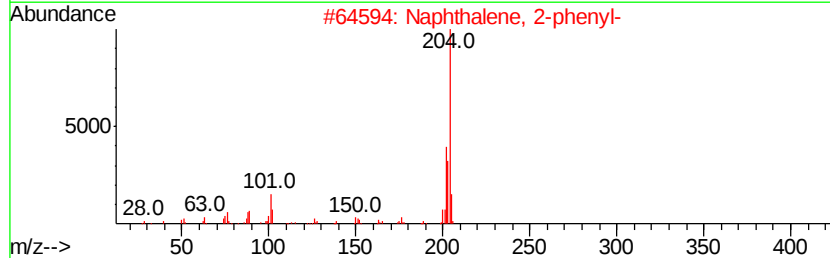
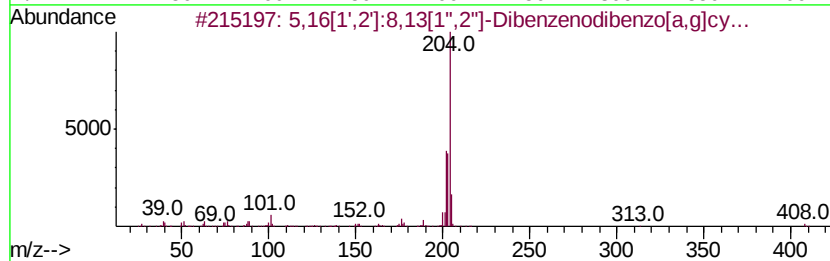
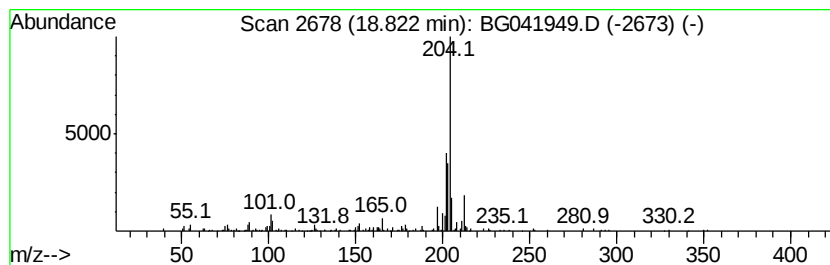
Instrument :  
BNA\_G  
ClientSampleId :  
BEP03

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Quant Title : SVOA CALIBRATION

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

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Peak Number 6 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 13

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.    |             |      |
|---------|-------------------------------------|--------------|------------------|---------|-------------|------|
| 18.82   | 2.00 ng/ul                          | 98725        | Phenanthrene-d10 | 17.45   |             |      |
| Hit# of | 5                                   | Tentative ID | MW               | MolForm | CAS#        | Qual |
| 1       | 5,16[1',2']:8,13[1'',2'']-Dibenz... | 408          | C32H24           |         | 005672-97-9 | 74   |
| 2       | Naphthalene, 2-phenyl-              | 204          | C16H12           |         | 000612-94-2 | 60   |
| 3       | Naphthalene, 2-phenyl-              | 204          | C16H12           |         | 000612-94-2 | 60   |
| 4       | Tricyclo[8.2.2.2(4,7)]hexadeca-2... | 204          | C16H12           |         | 006572-60-7 | 53   |
| 5       | Naphthalene, 2-phenyl-              | 204          | C16H12           |         | 000612-94-2 | 49   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080319\  
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Acq On : 4 Aug 2019 19:31  
Operator : HP/JU  
Sample : K3962-10  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

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

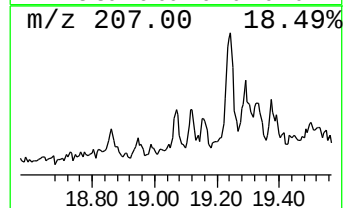
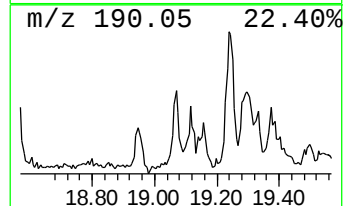
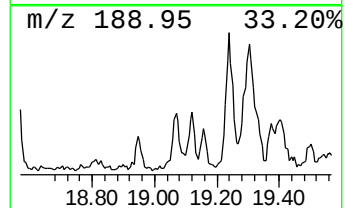
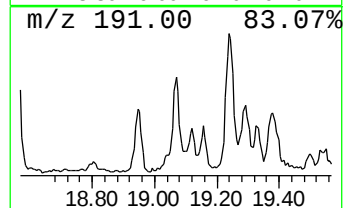
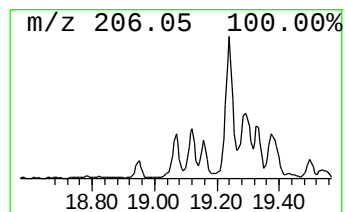
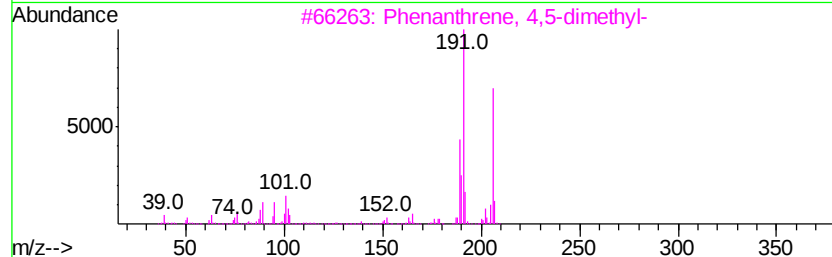
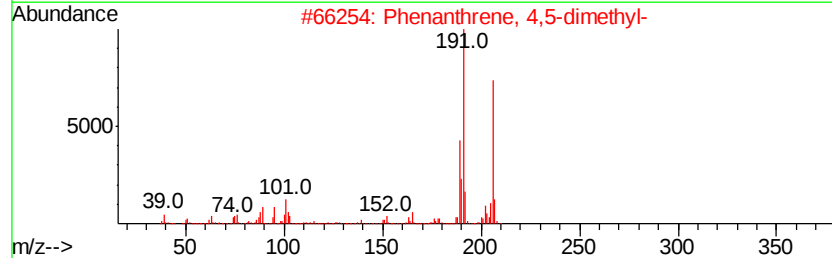
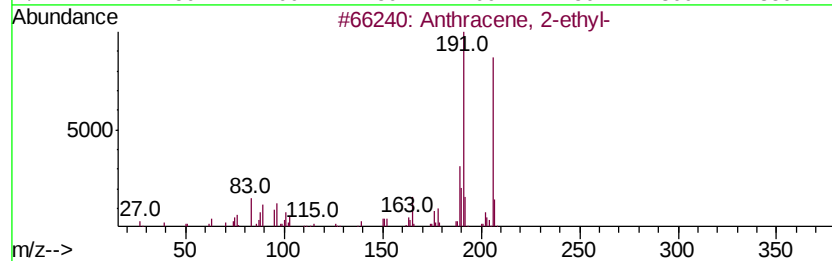
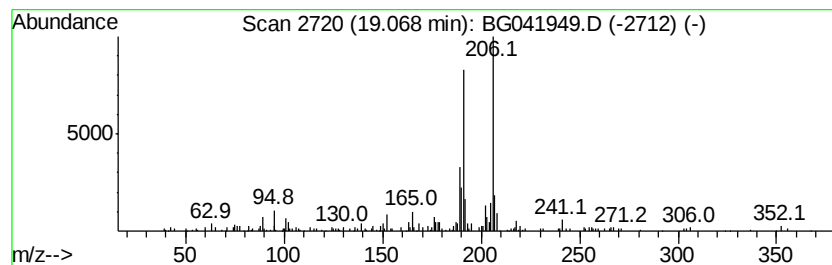
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Anthracene, 2-ethyl- Concentration Rank 11

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.07 | 2.53 ng/ul | 124932 | Phenanthrene-d10 | 17.45 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Anthracene, 2-ethyl-        | 206 | C16H14  | 052251-71-5 | 93   |
| 2    |    | Phenanthrene, 4,5-dimethyl- | 206 | C16H14  | 003674-69-9 | 93   |
| 3    |    | Phenanthrene, 4,5-dimethyl- | 206 | C16H14  | 003674-69-9 | 90   |
| 4    |    | Anthracene, 2-ethyl-        | 206 | C16H14  | 052251-71-5 | 87   |
| 5    |    | Phenanthrene, 4,5-dimethyl- | 206 | C16H14  | 003674-69-9 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080319\  
Data File : BG041949.D  
Acq On : 4 Aug 2019 19:31  
Operator : HP/JU  
Sample : K3962-10  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BEP03

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080319MA.M  
Quant Title : SVOA CALIBRATION

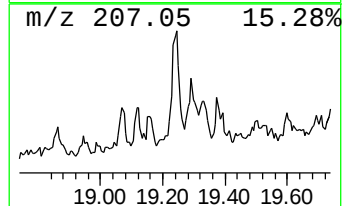
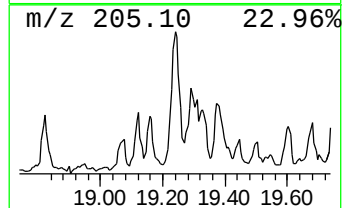
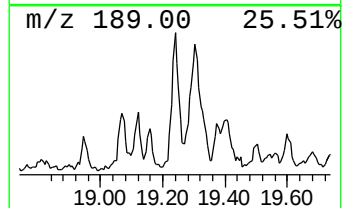
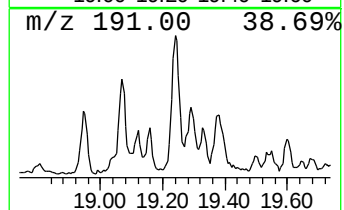
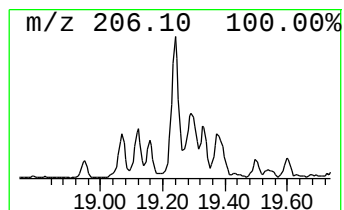
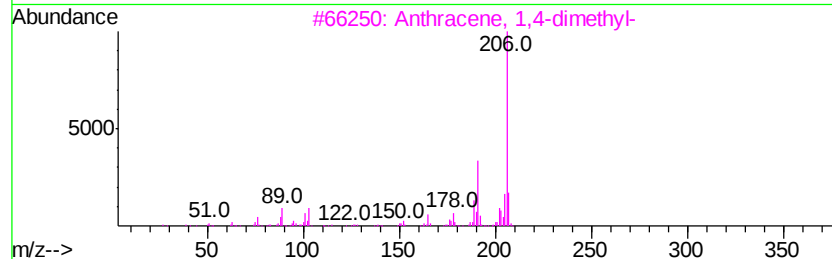
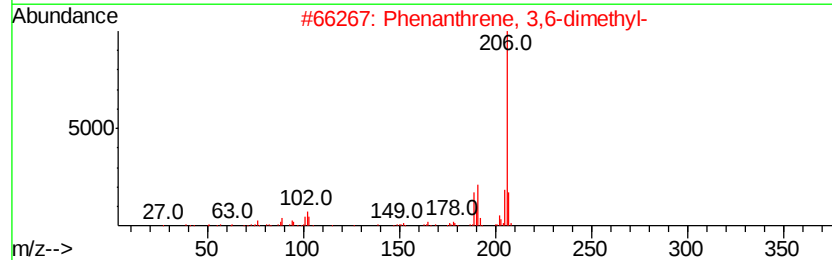
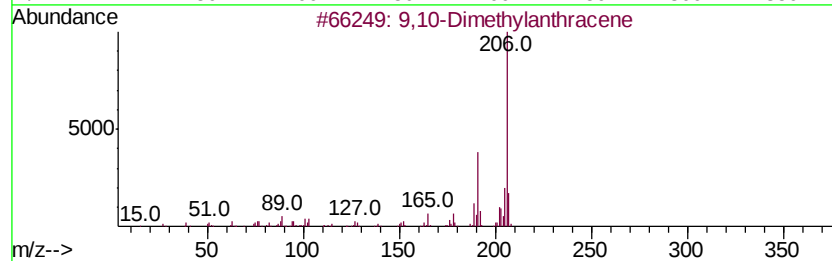
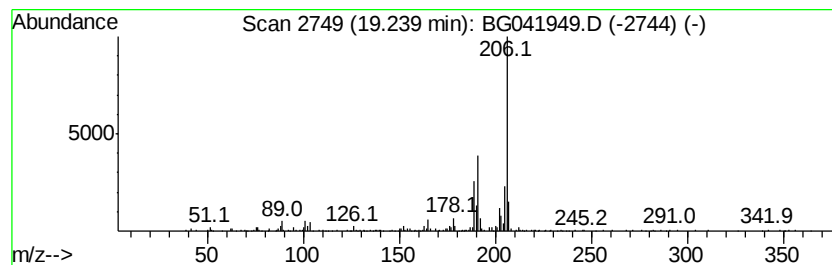
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 9,10-Dimethylantracene Concentration Rank 6

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.24 | 4.36 ng/ul | 215231 | Phenanthrene-d10 | 17.45 |

| Hit# of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|---------|---|-----------------------------|-----|---------|-------------|------|
| 1       |   | 9,10-Dimethylantracene      | 206 | C16H14  | 000781-43-1 | 96   |
| 2       |   | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 94   |
| 3       |   | Anthracene, 1,4-dimethyl-   | 206 | C16H14  | 000781-92-0 | 93   |
| 4       |   | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 93   |
| 5       |   | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080319\  
Data File : BG041949.D  
Acq On : 4 Aug 2019 19:31  
Operator : HP/JU  
Sample : K3962-10  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

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

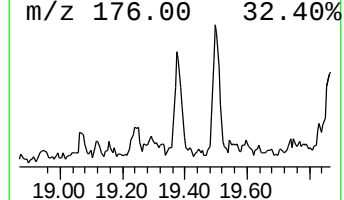
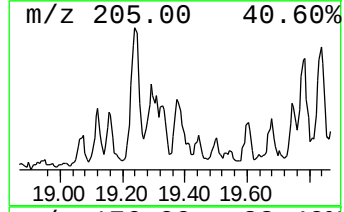
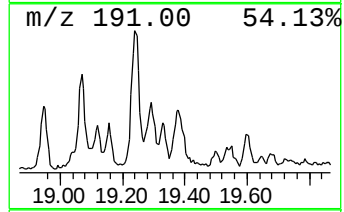
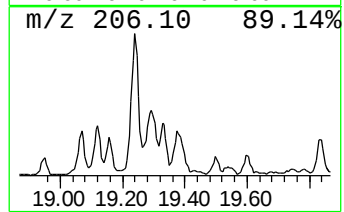
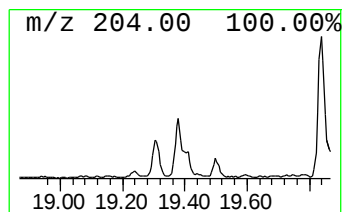
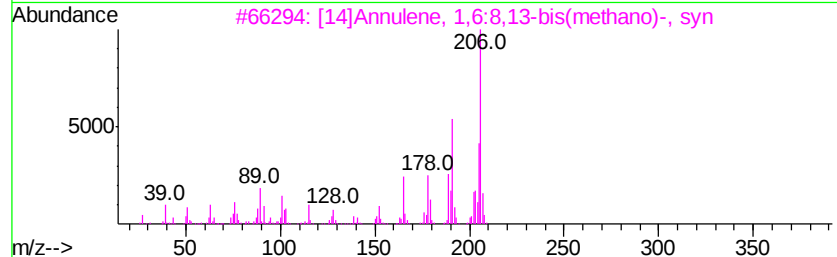
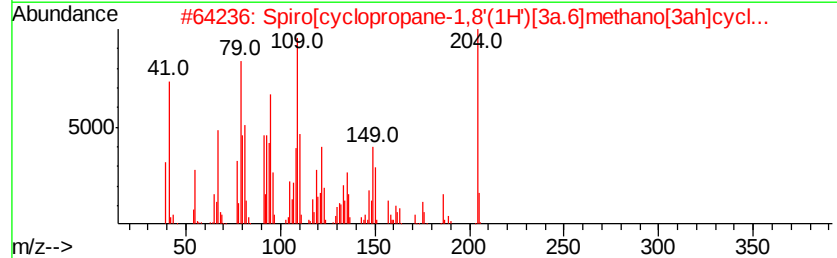
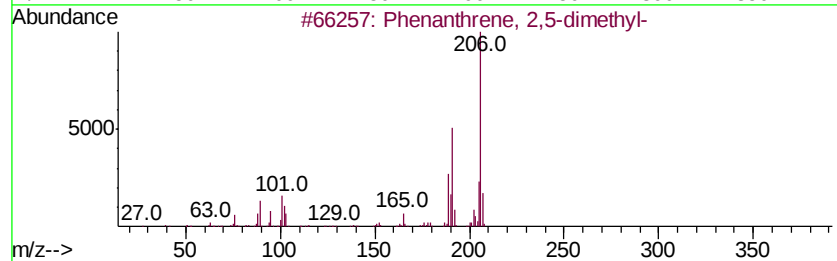
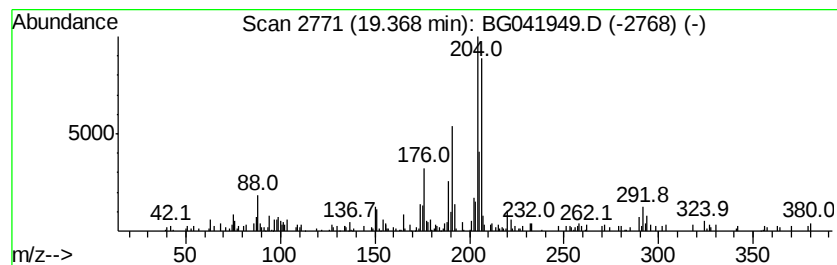
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Phenanthrene, 2,5-dimethyl- Concentration Rank 9

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.37 | 3.05 ng/ul | 150330 | Phenanthrene-d10 | 17.45 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenanthrene, 2,5-dimethyl-         | 206 | C16H14  | 003674-66-6 | 60   |
| 2    |    | Spiro[cyclopropane-1,8'(1H)][3a.... | 204 | C14H20O | 452049-62-6 | 56   |
| 3    |    | [14]Annulene, 1,6:8,13-bis(metha... | 206 | C16H14  | 085385-68-8 | 53   |
| 4    |    | 9,10-Dimethylantracene              | 206 | C16H14  | 000781-43-1 | 49   |
| 5    |    | 1H-Indene, 2-methyl-3-phenyl-       | 206 | C16H14  | 035099-60-6 | 46   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080319\  
Data File : BG041949.D  
Acq On : 4 Aug 2019 19:31  
Operator : HP/JU  
Sample : K3962-10  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BEP03

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080319MA.M  
Quant Title : SVOA CALIBRATION

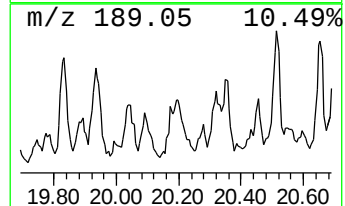
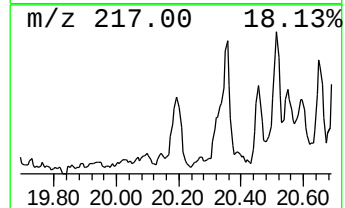
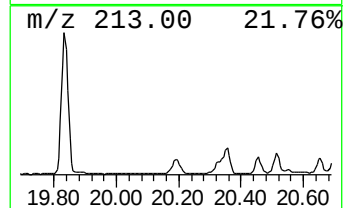
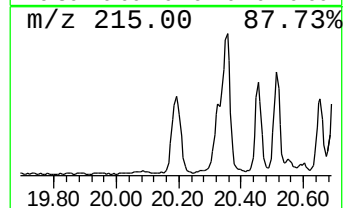
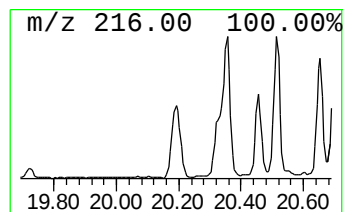
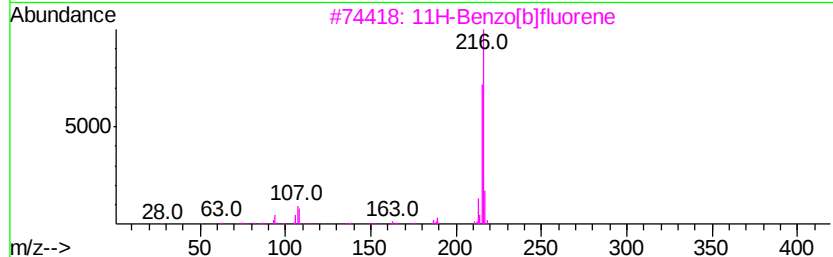
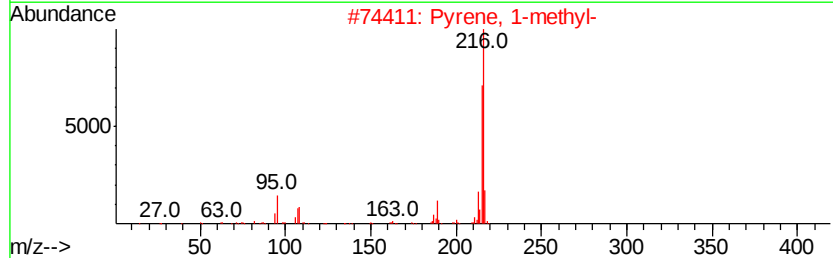
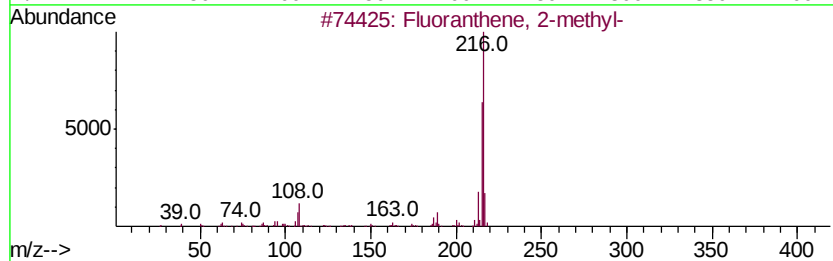
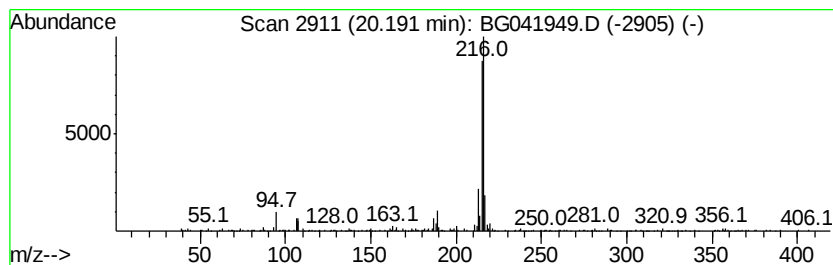
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Fluoranthene, 2-methyl- Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.19 | 2.61 ng/ul | 246600 | Chrysene-d12     | 21.74 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 93   |
| 2    |    |   | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 91   |
| 3    |    |   | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 90   |
| 4    |    |   | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 87   |
| 5    |    |   | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 81   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080319\  
Data File : BG041949.D  
Acq On : 4 Aug 2019 19:31  
Operator : HP/JU  
Sample : K3962-10  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BEP03

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080319MA.M  
Quant Title : SVOA CALIBRATION

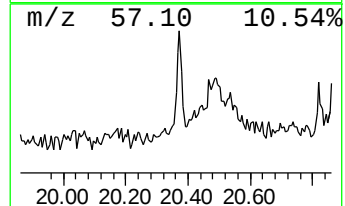
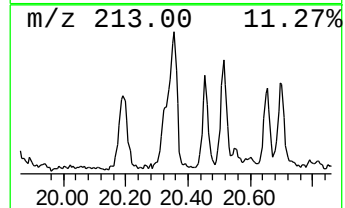
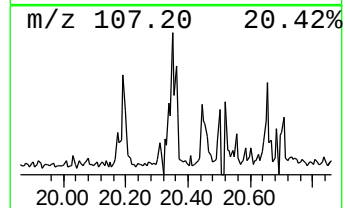
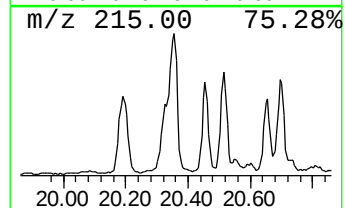
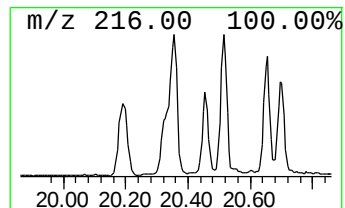
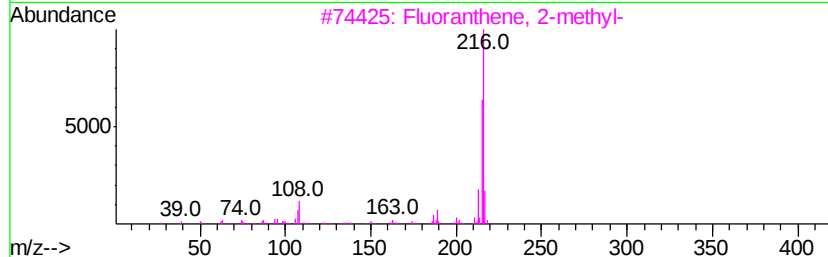
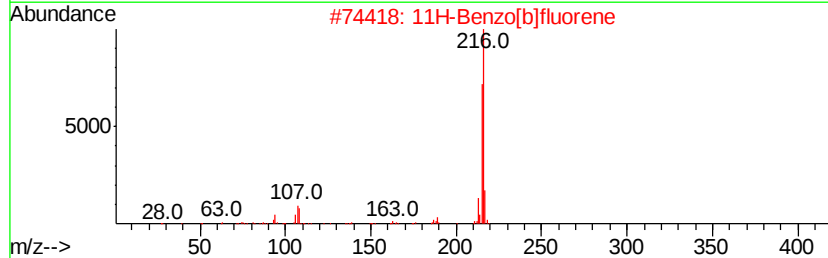
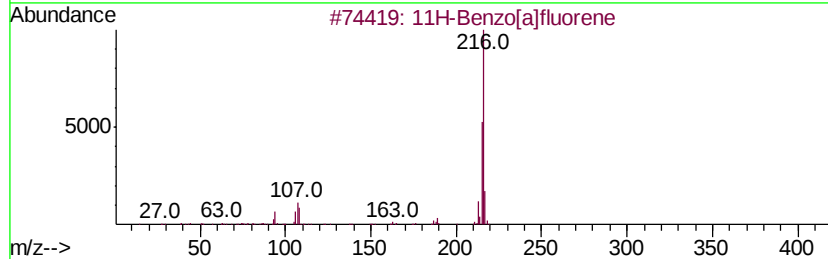
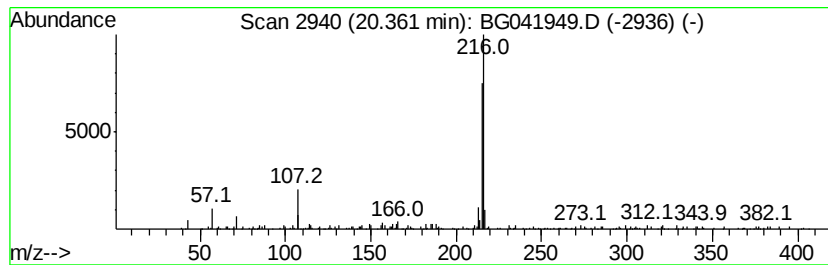
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 11H-Benzo[a]fluorene Concentration Rank 12

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.36 | 2.27 ng/ul | 214437 | Chrysene-d12     | 21.74 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 94   |
| 2    |    | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 90   |
| 3    |    | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 87   |
| 4    |    | Pvrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 83   |
| 5    |    | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080319\  
Data File : BG041949.D  
Acq On : 4 Aug 2019 19:31  
Operator : HP/JU  
Sample : K3962-10  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BEP03

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080319MA.M  
Quant Title : SVOA CALIBRATION

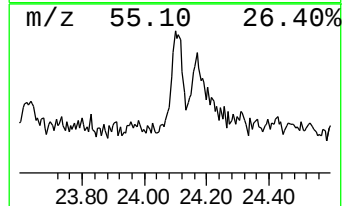
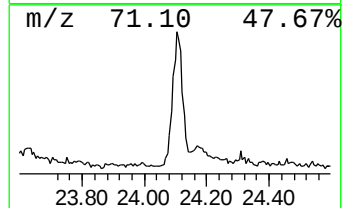
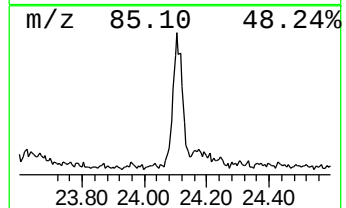
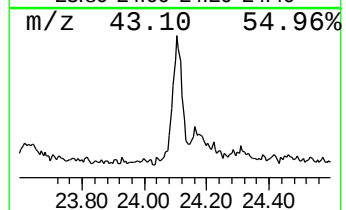
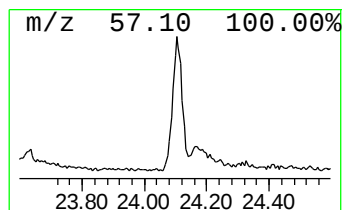
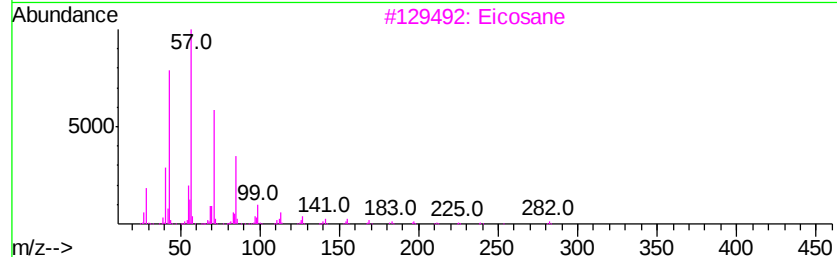
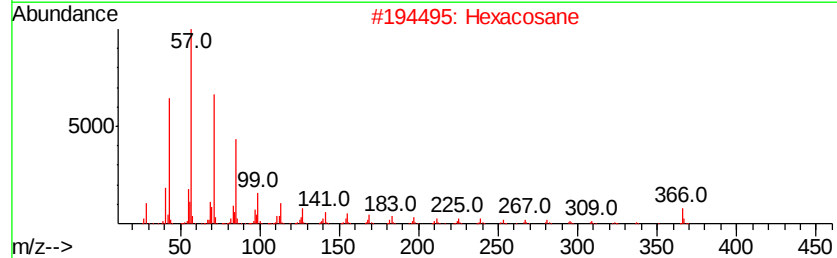
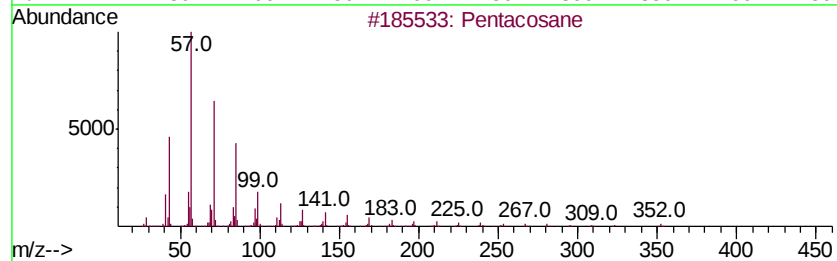
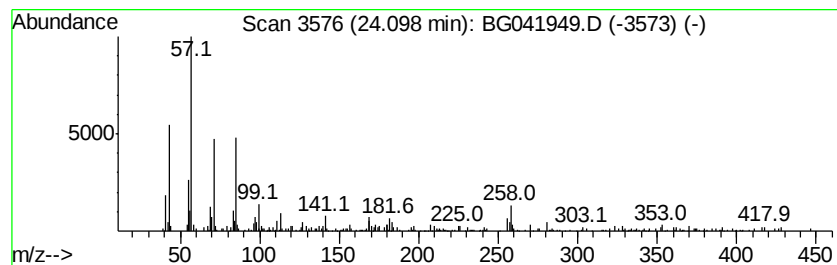
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 24.10 | 3.52 ng/ul | 162093 | Perylene-d12     | 25.01 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Pentacosane  | 352 | C25H52  | 000629-99-2 | 95   |
| 2    |    | Hexacosane   | 366 | C26H54  | 000630-01-3 | 89   |
| 3    |    | Eicosane     | 282 | C20H42  | 000112-95-8 | 87   |
| 4    |    | Heptadecane  | 240 | C17H36  | 000629-78-7 | 76   |
| 5    |    | Heptadecane  | 240 | C17H36  | 000629-78-7 | 76   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080319\  
Data File : BG041949.D  
Acq On : 4 Aug 2019 19:31  
Operator : HP/JU  
Sample : K3962-10  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

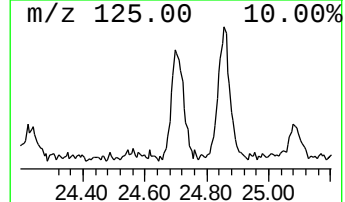
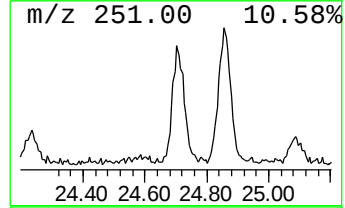
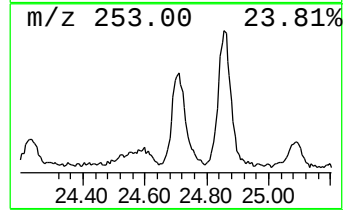
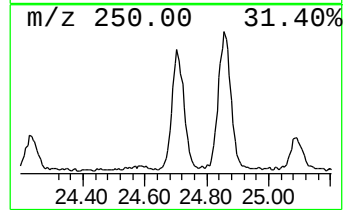
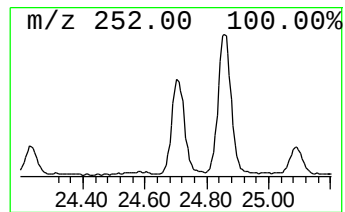
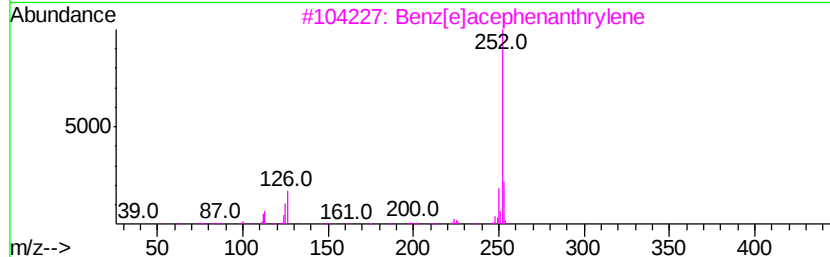
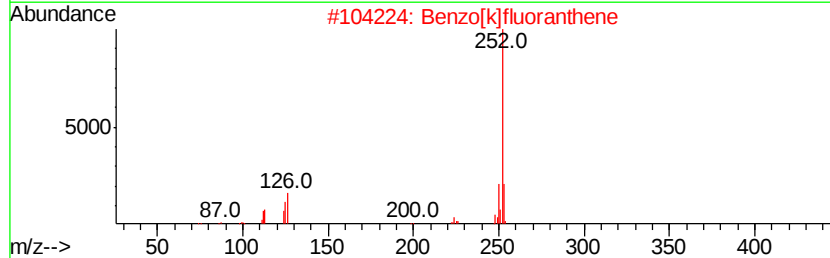
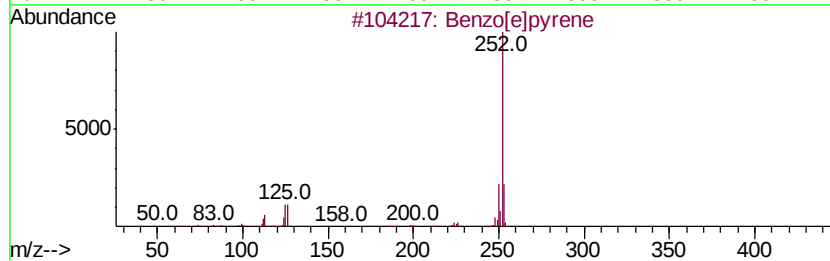
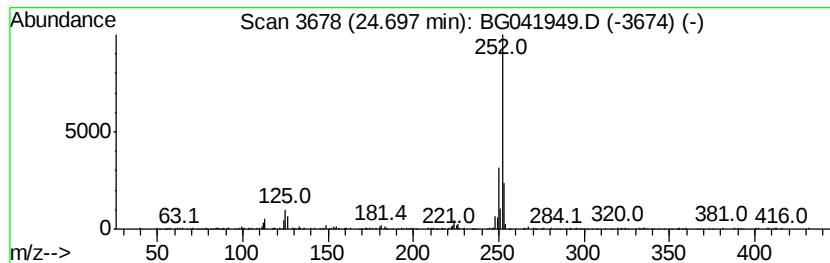
Instrument :  
BNA\_G  
ClientSampleId :  
BEP03

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080319MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 13 Benzo[e]pyrene Concentration Rank 3

| R.T.    | EstConc                   | Area         | Relative to ISTD | R.T.      |
|---------|---------------------------|--------------|------------------|-----------|
| 24.70   | 6.22 ng/ul                | 286164       | Perylene-d12     | 25.01     |
| Hit# of | 5                         | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Benzo[e]pyrene            | 252 C20H12   | 000192-97-2      | 98        |
| 2       | Benzo[k]fluoranthene      | 252 C20H12   | 000207-08-9      | 95        |
| 3       | Benzo[e]acephenanthrylene | 252 C20H12   | 000205-99-2      | 93        |
| 4       | Benzo[a]pyrene            | 252 C20H12   | 000050-32-8      | 93        |
| 5       | Benzo[j]fluoranthene      | 252 C20H12   | 000205-82-3      | 92        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG080319\  
Data File : BG041949.D  
Acq On : 4 Aug 2019 19:31  
Operator : HP/JU  
Sample : K3962-10  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BEP03

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080319MA.M  
Quant Title : SVOA CALIBRATION

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

| TIC Top Hit name      | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|-----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                       |       |         |       |          | #                     | RT    | Resp    | Conc |
| Butane, 2-methoxy...  | 3.16  | 127.1   | ng/ul | 2431440  | 1                     | 8.03  | 382548  | 20.0 |
| Phenanthrene, 2-m...  | 18.33 | 5.6     | ng/ul | 274538   | 4                     | 17.45 | 986557  | 20.0 |
| Anthracene, 9-met...  | 18.38 | 4.8     | ng/ul | 236205   | 4                     | 17.45 | 986557  | 20.0 |
| 4H-Cyclopenta[def...  | 18.52 | 6.3     | ng/ul | 309107   | 4                     | 17.45 | 986557  | 20.0 |
| 1H-Indene, 1-phenyl-  | 18.56 | 3.4     | ng/ul | 166378   | 4                     | 17.45 | 986557  | 20.0 |
| 5,16[1',2']:8,13[...  | 18.82 | 2.0     | ng/ul | 98725    | 4                     | 17.45 | 986557  | 20.0 |
| Anthracene, 2-ethyl-  | 19.07 | 2.5     | ng/ul | 124932   | 4                     | 17.45 | 986557  | 20.0 |
| 9,10-Dimethylanthe... | 19.24 | 4.4     | ng/ul | 215231   | 4                     | 17.45 | 986557  | 20.0 |
| Phenanthrene, 2,5...  | 19.37 | 3.0     | ng/ul | 150330   | 4                     | 17.45 | 986557  | 20.0 |
| Fluoranthene, 2-m...  | 20.19 | 2.6     | ng/ul | 246600   | 5                     | 21.74 | 1891240 | 20.0 |
| 11H-Benzo[a]fluorene  | 20.36 | 2.3     | ng/ul | 214437   | 5                     | 21.74 | 1891240 | 20.0 |
| (DEL) Alkane: Str...  | 24.10 | 3.5     | ng/ul | 162093   | 6                     | 25.01 | 920556  | 20.0 |
| Benzo[e]pyrene        | 24.70 | 6.2     | ng/ul | 286164   | 6                     | 25.01 | 920556  | 20.0 |