

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101619\  
 Data File : BM023203.D  
 Acq On : 16 Oct 2019 22:45  
 Operator : JU  
 Sample : K5262-06  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BFB71

## Integration Parameters: LSCINT.P

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

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101419MA.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         | 3.681       | 113           | 118         | 125          | rVV      | 256159         | 326383        | 2.94%           | 0.237%        |
| 2         | 6.369       | 571           | 575         | 582          | rVB2     | 65744          | 109991        | 0.99%           | 0.080%        |
| 3         | 6.834       | 646           | 654         | 664          | rVV2     | 1002425        | 1922295       | 17.33%          | 1.395%        |
| 4         | 6.998       | 676           | 682         | 690          | rBV      | 720625         | 1153193       | 10.39%          | 0.837%        |
| 5         | 7.198       | 708           | 716         | 722          | rBV      | 985594         | 1571436       | 14.16%          | 1.140%        |
| 6         | 7.322       | 731           | 737         | 742          | rBV3     | 84043          | 140565        | 1.27%           | 0.102%        |
| 7         | 7.551       | 771           | 776         | 783          | rVV3     | 60672          | 109338        | 0.99%           | 0.079%        |
| 8         | 7.663       | 789           | 795         | 806          | rVV      | 1675261        | 2712756       | 24.45%          | 1.968%        |
| 9         | 8.222       | 883           | 890         | 896          | rVB7     | 44941          | 114414        | 1.03%           | 0.083%        |
| 10        | 8.369       | 908           | 915         | 922          | rVV      | 1045270        | 1730080       | 15.59%          | 1.255%        |
| 11        | 8.481       | 928           | 934         | 941          | rVB6     | 43913          | 127891        | 1.15%           | 0.093%        |
| 12        | 8.810       | 977           | 990         | 996          | rBV      | 903867         | 1607663       | 14.49%          | 1.166%        |
| 13        | 9.028       | 1017          | 1027        | 1033         | rVB      | 51622          | 169557        | 1.53%           | 0.123%        |
| 14        | 9.298       | 1067          | 1073        | 1077         | rVV      | 57585          | 104147        | 0.94%           | 0.076%        |
| 15        | 9.527       | 1105          | 1112        | 1120         | rBV      | 738046         | 1228815       | 11.08%          | 0.891%        |
| 16        | 9.933       | 1177          | 1181        | 1192         | rVB5     | 63336          | 107283        | 0.97%           | 0.078%        |
| 17        | 10.069      | 1197          | 1204        | 1213         | rVB      | 1297314        | 2205143       | 19.88%          | 1.600%        |
| 18        | 10.445      | 1261          | 1268        | 1274         | rBV      | 2190312        | 3773050       | 34.01%          | 2.737%        |
| 19        | 10.492      | 1274          | 1276        | 1282         | rVV      | 96750          | 141163        | 1.27%           | 0.102%        |
| 20        | 10.575      | 1283          | 1290        | 1298         | rVV      | 548649         | 1025044       | 9.24%           | 0.744%        |
| 21        | 10.645      | 1298          | 1302        | 1307         | rVB      | 91216          | 133095        | 1.20%           | 0.097%        |
| 22        | 11.510      | 1439          | 1449        | 1454         | rBV2     | 111463         | 226611        | 2.04%           | 0.164%        |
| 23        | 11.798      | 1493          | 1498        | 1508         | rVB2     | 75475          | 168120        | 1.52%           | 0.122%        |
| 24        | 11.916      | 1509          | 1518        | 1523         | rBV5     | 66451          | 171599        | 1.55%           | 0.124%        |
| 25        | 12.116      | 1548          | 1552        | 1563         | rVB3     | 106938         | 218776        | 1.97%           | 0.159%        |
| 26        | 12.339      | 1584          | 1590        | 1597         | rBV      | 73195          | 114737        | 1.03%           | 0.083%        |
| 27        | 12.880      | 1677          | 1682        | 1686         | rBV      | 98348          | 138314        | 1.25%           | 0.100%        |
| 28        | 13.633      | 1805          | 1810        | 1815         | rVV      | 113369         | 208913        | 1.88%           | 0.152%        |
| 29        | 13.686      | 1815          | 1819        | 1821         | rVV      | 90443          | 141741        | 1.28%           | 0.103%        |
| 30        | 13.727      | 1821          | 1826        | 1830         | rVV      | 2118563        | 3157689       | 28.46%          | 2.291%        |
| 31        | 13.768      | 1830          | 1833        | 1838         | rVV      | 791977         | 1112447       | 10.03%          | 0.807%        |
| 32        | 13.833      | 1839          | 1844        | 1847         | rVV3     | 112695         | 200936        | 1.81%           | 0.146%        |
| 33        | 13.868      | 1847          | 1850        | 1860         | rVB3     | 80714          | 159872        | 1.44%           | 0.116%        |
| 34        | 13.998      | 1865          | 1872        | 1887         | rBV      | 2616742        | 4357447       | 39.28%          | 3.161%        |

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Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101419MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 14.310 | 1918 | 1925 | 1932 | rBV  | 3272605 | 4872110 | 43.92% | 3.535% |
| 36 | 14.368 | 1932 | 1935 | 1940 | rVB  | 93606   | 130881  | 1.18%  | 0.095% |
| 37 | 14.498 | 1952 | 1957 | 1968 | rBV  | 757058  | 1135897 | 10.24% | 0.824% |
| 38 | 14.727 | 1988 | 1996 | 2003 | rVV3 | 144868  | 381546  | 3.44%  | 0.277% |
| 39 | 14.792 | 2003 | 2007 | 2013 | rVB  | 82029   | 127262  | 1.15%  | 0.092% |
| 40 | 14.933 | 2027 | 2031 | 2036 | rVV2 | 64474   | 115233  | 1.04%  | 0.084% |
| 41 | 14.986 | 2037 | 2040 | 2046 | rVB2 | 72444   | 103112  | 0.93%  | 0.075% |
| 42 | 15.304 | 2088 | 2094 | 2100 | rBV  | 3190154 | 4561336 | 41.11% | 3.309% |
| 43 | 15.362 | 2100 | 2104 | 2109 | rVB2 | 173340  | 252178  | 2.27%  | 0.183% |
| 44 | 15.421 | 2109 | 2114 | 2122 | rBV  | 548291  | 806148  | 7.27%  | 0.585% |
| 45 | 15.657 | 2151 | 2154 | 2158 | rBV  | 79164   | 98849   | 0.89%  | 0.072% |
| 46 | 15.809 | 2175 | 2180 | 2188 | rVB3 | 103282  | 237686  | 2.14%  | 0.172% |
| 47 | 15.886 | 2188 | 2193 | 2199 | rBV5 | 55680   | 130346  | 1.17%  | 0.095% |
| 48 | 16.092 | 2225 | 2228 | 2234 | rVB2 | 89552   | 123605  | 1.11%  | 0.090% |
| 49 | 16.368 | 2271 | 2275 | 2280 | rVB2 | 127541  | 187323  | 1.69%  | 0.136% |
| 50 | 16.415 | 2280 | 2283 | 2289 | rVB3 | 62941   | 104510  | 0.94%  | 0.076% |
| 51 | 16.633 | 2317 | 2320 | 2324 | rVB2 | 88145   | 129116  | 1.16%  | 0.094% |
| 52 | 16.704 | 2329 | 2332 | 2341 | rVB2 | 149349  | 252700  | 2.28%  | 0.183% |
| 53 | 16.809 | 2344 | 2350 | 2355 | rBV4 | 101539  | 257030  | 2.32%  | 0.186% |
| 54 | 16.868 | 2356 | 2360 | 2364 | rVB2 | 152906  | 238689  | 2.15%  | 0.173% |
| 55 | 17.056 | 2386 | 2392 | 2395 | rBV2 | 3543804 | 5221697 | 47.07% | 3.788% |
| 56 | 17.098 | 2395 | 2399 | 2403 | rVV  | 2920360 | 4061598 | 36.61% | 2.947% |
| 57 | 17.151 | 2403 | 2408 | 2412 | rVV2 | 3421848 | 4935474 | 44.49% | 3.581% |
| 58 | 17.186 | 2412 | 2414 | 2419 | rVV  | 728570  | 825853  | 7.44%  | 0.599% |
| 59 | 17.251 | 2421 | 2425 | 2431 | rVB3 | 102454  | 180107  | 1.62%  | 0.131% |
| 60 | 17.456 | 2456 | 2460 | 2465 | rVB  | 167601  | 278499  | 2.51%  | 0.202% |
| 61 | 17.586 | 2479 | 2482 | 2486 | rBV2 | 67440   | 101031  | 0.91%  | 0.073% |
| 62 | 17.633 | 2487 | 2490 | 2496 | rVB  | 161217  | 217240  | 1.96%  | 0.158% |
| 63 | 17.851 | 2524 | 2527 | 2533 | rBV6 | 62861   | 121926  | 1.10%  | 0.088% |
| 64 | 17.933 | 2536 | 2541 | 2544 | rBV  | 621755  | 871186  | 7.85%  | 0.632% |
| 65 | 17.980 | 2544 | 2549 | 2556 | rVV  | 1325579 | 1886174 | 17.00% | 1.368% |
| 66 | 18.056 | 2557 | 2562 | 2567 | rVV2 | 326995  | 633616  | 5.71%  | 0.460% |
| 67 | 18.121 | 2568 | 2573 | 2577 | rVV2 | 791630  | 1570031 | 14.15% | 1.139% |
| 68 | 18.162 | 2577 | 2580 | 2587 | rVB  | 467680  | 689134  | 6.21%  | 0.500% |
| 69 | 18.433 | 2623 | 2626 | 2629 | rBV  | 340910  | 436262  | 3.93%  | 0.317% |
| 70 | 18.686 | 2666 | 2669 | 2675 | rVB3 | 246894  | 370062  | 3.34%  | 0.268% |
| 71 | 18.745 | 2675 | 2679 | 2682 | rBV2 | 221667  | 393644  | 3.55%  | 0.286% |

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Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101419MA.M  
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|     |        |      |      |      |      |         |          |         |        |
|-----|--------|------|------|------|------|---------|----------|---------|--------|
| 72  | 18.856 | 2693 | 2698 | 2703 | rBV  | 561324  | 952110   | 8.58%   | 0.691% |
| 73  | 18.921 | 2703 | 2709 | 2712 | rVV4 | 276626  | 530087   | 4.78%   | 0.385% |
| 74  | 18.992 | 2717 | 2721 | 2729 | rVB3 | 432825  | 762661   | 6.87%   | 0.553% |
| 75  | 19.115 | 2737 | 2742 | 2748 | rVB  | 4804990 | 6522268  | 58.79%  | 4.732% |
| 76  | 19.192 | 2752 | 2755 | 2758 | rBV  | 195747  | 264958   | 2.39%   | 0.192% |
| 77  | 19.268 | 2764 | 2768 | 2772 | rBV2 | 455825  | 678322   | 6.11%   | 0.492% |
| 78  | 19.450 | 2794 | 2799 | 2801 | rBV2 | 4621222 | 6376334  | 57.47%  | 4.626% |
| 79  | 19.480 | 2801 | 2804 | 2812 | rVB  | 4737195 | 6294241  | 56.73%  | 4.566% |
| 80  | 19.662 | 2832 | 2835 | 2838 | rVV  | 220412  | 260486   | 2.35%   | 0.189% |
| 81  | 19.721 | 2842 | 2845 | 2852 | rVB3 | 407611  | 589211   | 5.31%   | 0.427% |
| 82  | 19.815 | 2857 | 2861 | 2865 | rBV  | 283625  | 544016   | 4.90%   | 0.395% |
| 83  | 19.909 | 2874 | 2877 | 2880 | rBV2 | 318880  | 379768   | 3.42%   | 0.276% |
| 84  | 19.980 | 2886 | 2889 | 2895 | rVB2 | 681598  | 969857   | 8.74%   | 0.704% |
| 85  | 20.080 | 2903 | 2906 | 2910 | rBV2 | 327289  | 392550   | 3.54%   | 0.285% |
| 86  | 20.139 | 2912 | 2916 | 2920 | rVB3 | 578637  | 803048   | 7.24%   | 0.583% |
| 87  | 20.274 | 2936 | 2939 | 2943 | rVB  | 437281  | 534315   | 4.82%   | 0.388% |
| 88  | 20.321 | 2944 | 2947 | 2952 | rVV  | 464917  | 598683   | 5.40%   | 0.434% |
| 89  | 20.727 | 3013 | 3016 | 3022 | rVB  | 568176  | 824303   | 7.43%   | 0.598% |
| 90  | 20.933 | 3048 | 3051 | 3054 | rVB  | 418459  | 428121   | 3.86%   | 0.311% |
| 91  | 20.991 | 3057 | 3061 | 3065 | rBV3 | 1887840 | 2224171  | 20.05%  | 1.614% |
| 92  | 21.197 | 3093 | 3096 | 3098 | rBV2 | 595837  | 661025   | 5.96%   | 0.480% |
| 93  | 21.244 | 3098 | 3104 | 3108 | rVV2 | 5747080 | 11094156 | 100.00% | 8.049% |
| 94  | 21.286 | 3108 | 3111 | 3121 | rVB  | 2763856 | 3775194  | 34.03%  | 2.739% |
| 95  | 21.886 | 3209 | 3213 | 3221 | rVB3 | 1350213 | 2414161  | 21.76%  | 1.751% |
| 96  | 22.803 | 3364 | 3369 | 3373 | rBV  | 2268764 | 4595465  | 41.42%  | 3.334% |
| 97  | 23.274 | 3444 | 3449 | 3452 | rBV  | 1330416 | 2312265  | 20.84%  | 1.678% |
| 98  | 23.321 | 3452 | 3457 | 3461 | rVV  | 3115067 | 5556224  | 50.08%  | 4.031% |
| 99  | 23.362 | 3461 | 3464 | 3470 | rVB  | 1993796 | 3308245  | 29.82%  | 2.400% |
| 100 | 23.462 | 3476 | 3481 | 3486 | rVB  | 3092671 | 5262533  | 47.44%  | 3.818% |

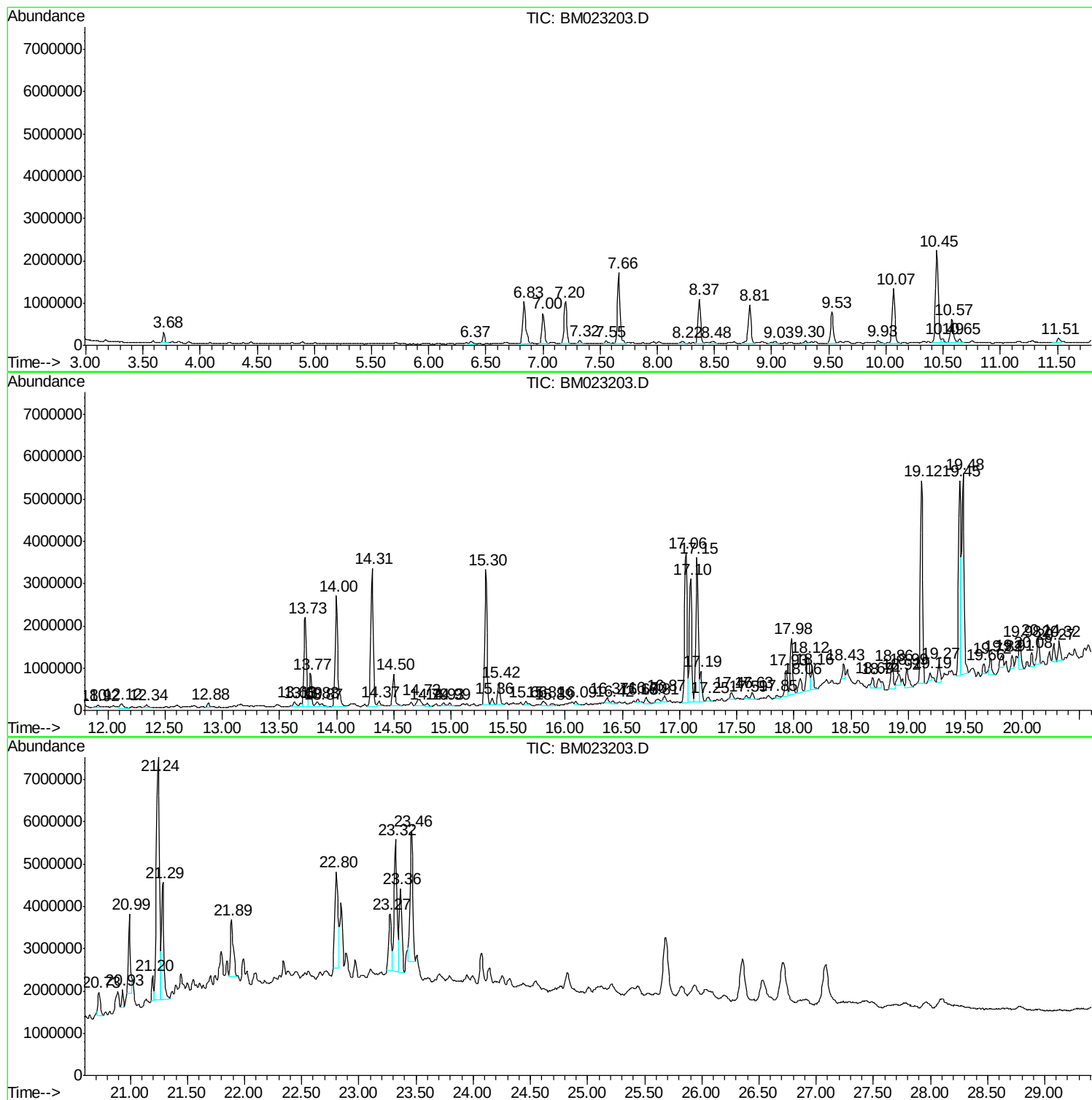
Sum of corrected areas: 137838373

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101419MA.M  
Quant Title : SVOA CALIBRATION

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



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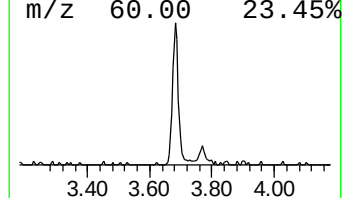
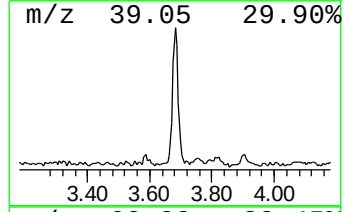
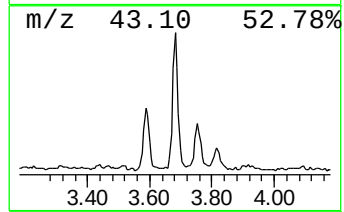
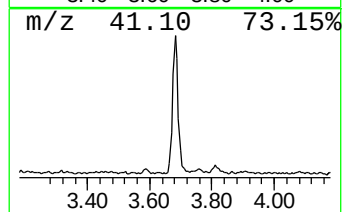
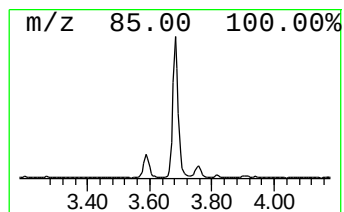
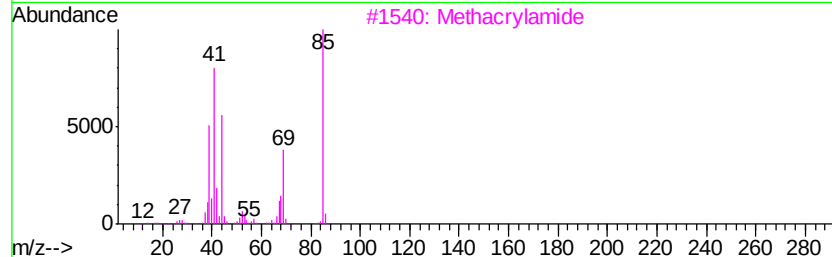
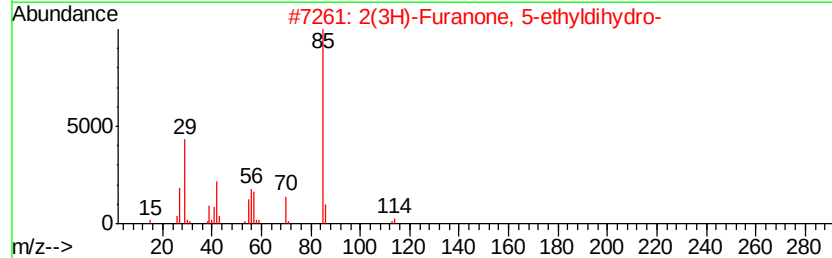
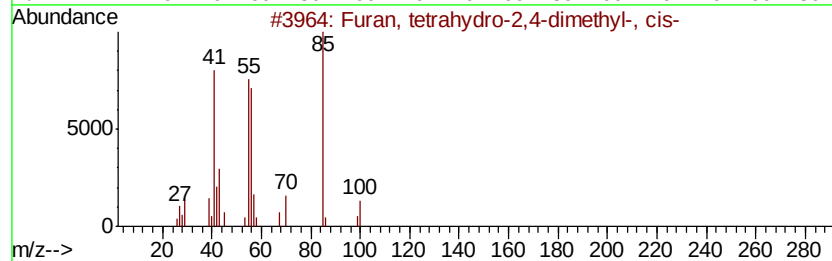
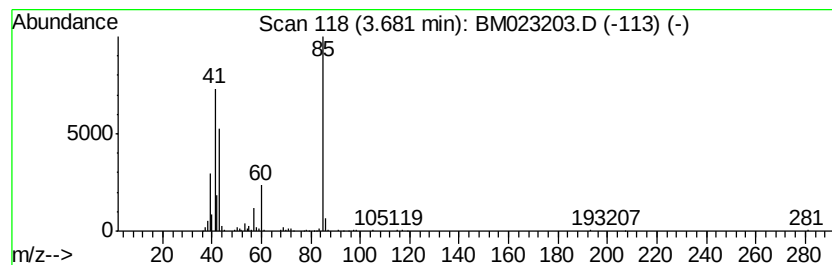
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 Furan, tetrahydro-2,4-dimet... Concentration Rank 10

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 3.68 | 2.41 ng/ul | 326383 | 1,4-Dichlorobenzene-d4 | 7.66 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Furan, tetrahydro-2,4-dimethyl-,... | 100 | C6H12O  | 039168-01-9 | 50   |
| 2    |    |   | 2(3H)-Furanone, 5-ethylidihydro-    | 114 | C6H10O2 | 000695-06-7 | 50   |
| 3    |    |   | Methacrylamide                      | 85  | C4H7NO  | 000079-39-0 | 45   |
| 4    |    |   | 2H-Pyran, tetrahydro-2-methyl-      | 100 | C6H12O  | 010141-72-7 | 36   |
| 5    |    |   | Hexane, 3-iodo-                     | 212 | C6H13I  | 031294-91-4 | 36   |



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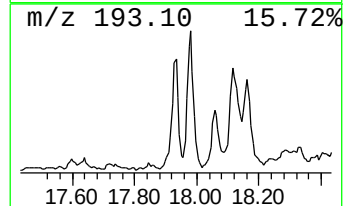
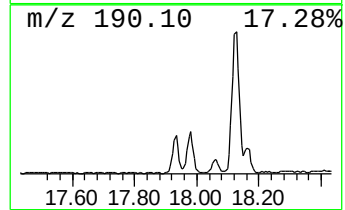
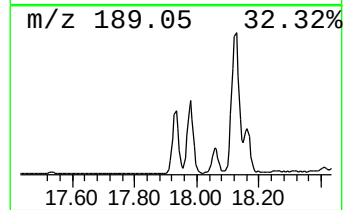
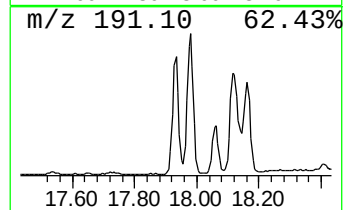
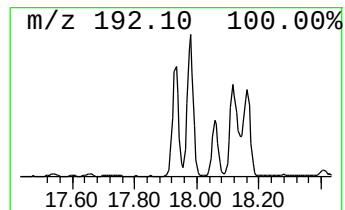
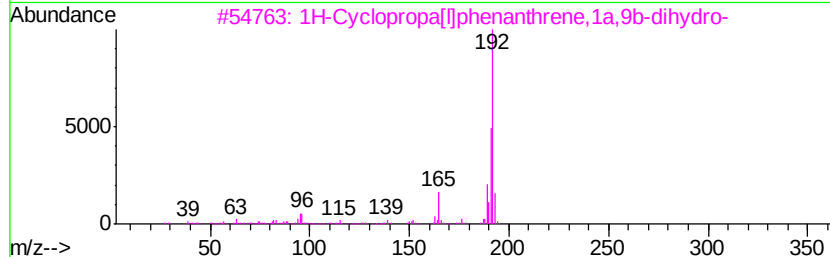
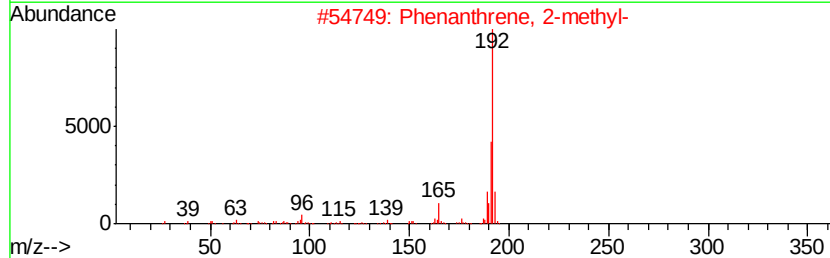
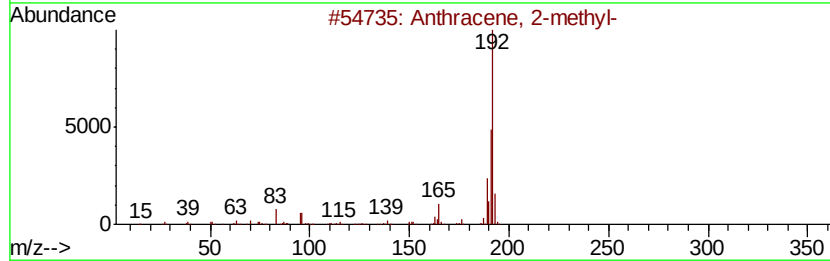
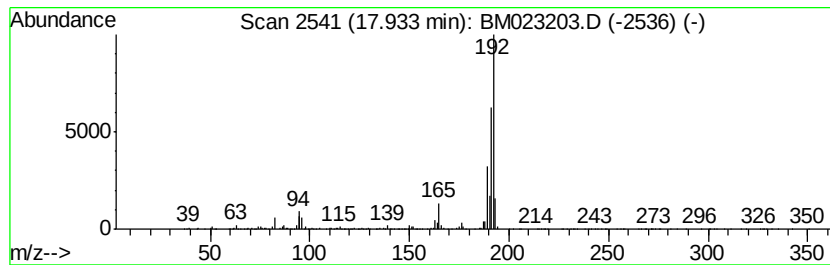
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\*\*\*\*\*  
Peak Number 2 Anthracene, 2-methyl- Concentration Rank 6

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 17.93     | 3.34 ng/ul                          | 871186 | Phenanthrene-d10 | 17.06       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Anthracene, 2-methyl-               | 192    | C15H12           | 000613-12-7 | 96   |
| 2         | Phenanthrene, 2-methyl-             | 192    | C15H12           | 002531-84-2 | 95   |
| 3         | 1H-Cyclopropa[1]phenanthrene,1a,... | 192    | C15H12           | 000949-41-7 | 95   |
| 4         | Phenanthrene, 2-methyl-             | 192    | C15H12           | 002531-84-2 | 94   |
| 5         | Phenanthrene, 1-methyl-             | 192    | C15H12           | 000832-69-9 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101619\  
Data File : BM023203.D  
Acq On : 16 Oct 2019 22:45  
Operator : JU  
Sample : K5262-06  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

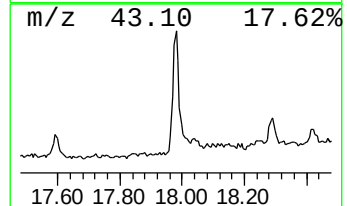
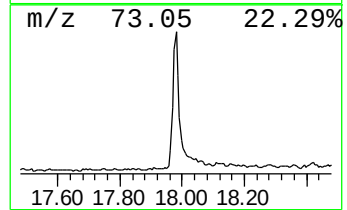
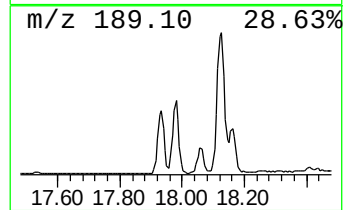
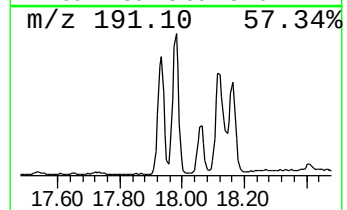
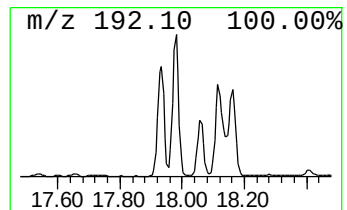
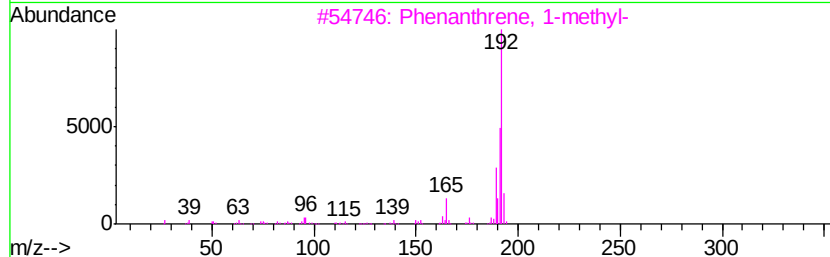
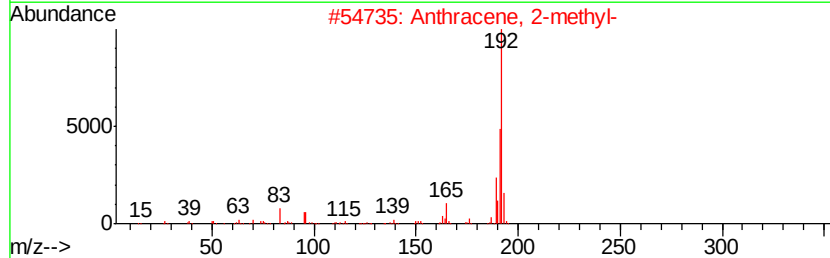
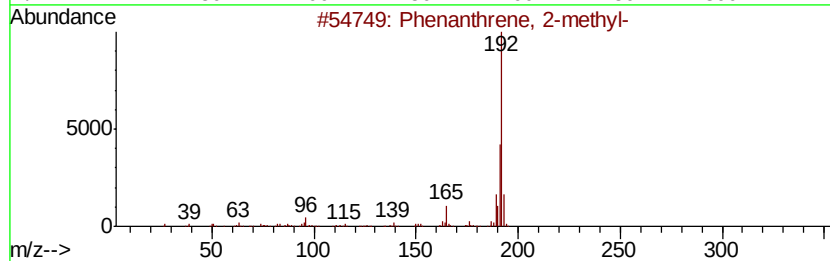
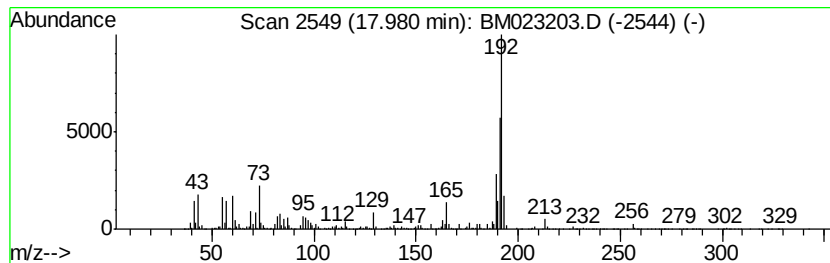
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 1

| R.T.    | EstConc    | Area                                  | Relative to ISTD | R.T.    |             |      |
|---------|------------|---------------------------------------|------------------|---------|-------------|------|
| 17.98   | 7.22 ng/ul | 1886170                               | Phenanthrene-d10 | 17.06   |             |      |
| Hit# of | 5          | Tentative ID                          | MW               | MolForm | CAS#        | Qual |
| 1       |            | Phenanthrene, 2-methyl-               | 192              | C15H12  | 002531-84-2 | 97   |
| 2       |            | Anthracene, 2-methyl-                 | 192              | C15H12  | 000613-12-7 | 97   |
| 3       |            | Phenanthrene, 1-methyl-               | 192              | C15H12  | 000832-69-9 | 97   |
| 4       |            | Phenanthrene, 2-methyl-               | 192              | C15H12  | 002531-84-2 | 97   |
| 5       |            | 1H-Cyclopropa[l]phenanthrene, 1a, ... | 192              | C15H12  | 000949-41-7 | 96   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101619\  
Data File : BM023203.D  
Acq On : 16 Oct 2019 22:45  
Operator : JU  
Sample : K5262-06  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101419MA.M  
Quant Title : SVOA CALIBRATION

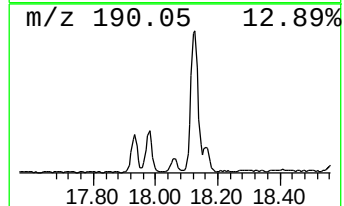
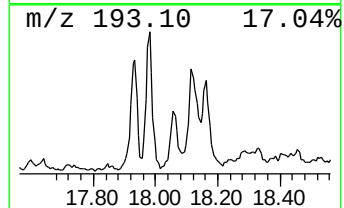
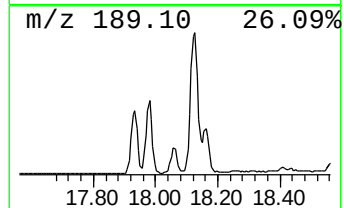
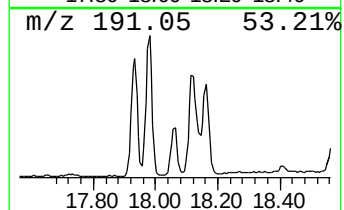
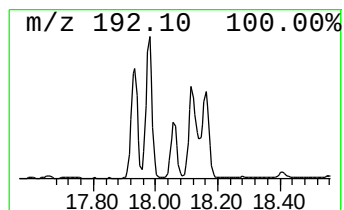
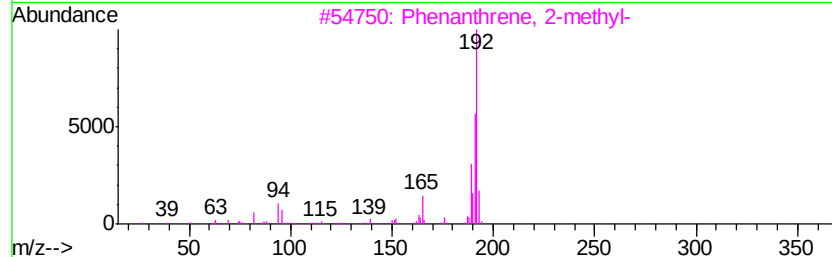
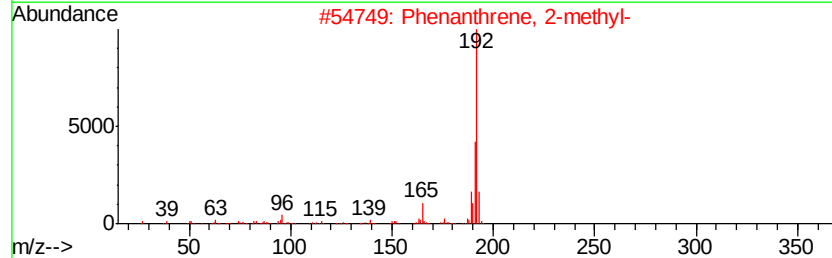
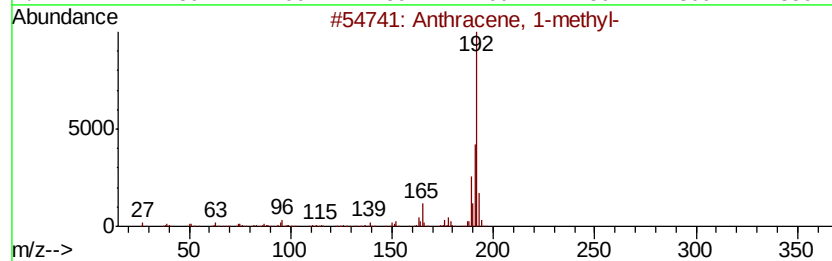
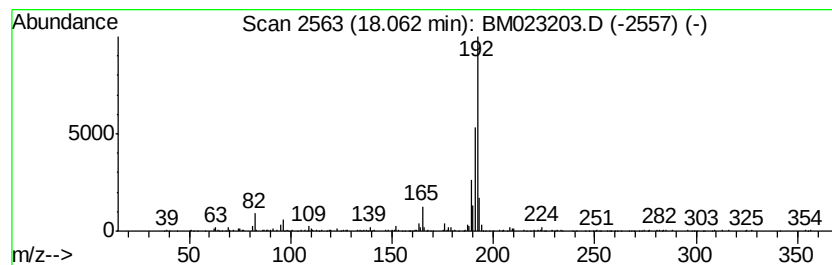
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 Anthracene, 1-methyl- Concentration Rank 9

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.06 | 2.43 ng/ul | 633616 | Phenanthrene-d10 | 17.06 |

| Hit# of | 5 | Tentative ID                           | MW  | MolForm | CAS#        | Qual |
|---------|---|----------------------------------------|-----|---------|-------------|------|
| 1       |   | Anthracene, 1-methyl-                  | 192 | C15H12  | 000610-48-0 | 96   |
| 2       |   | Phenanthrene, 2-methyl-                | 192 | C15H12  | 002531-84-2 | 96   |
| 3       |   | Phenanthrene, 2-methyl-                | 192 | C15H12  | 002531-84-2 | 95   |
| 4       |   | 1H-Cyclopropa[1,1]phenanthrene, 1a,... | 192 | C15H12  | 000949-41-7 | 95   |
| 5       |   | Phenanthrene, 1-methyl-                | 192 | C15H12  | 000832-69-9 | 95   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101619\  
Data File : BM023203.D  
Acq On : 16 Oct 2019 22:45  
Operator : JU  
Sample : K5262-06  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

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

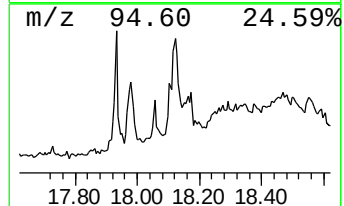
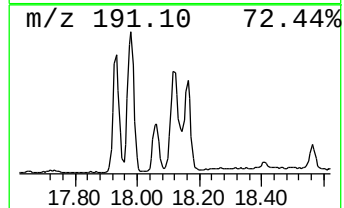
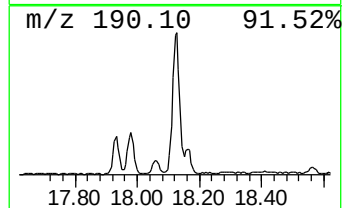
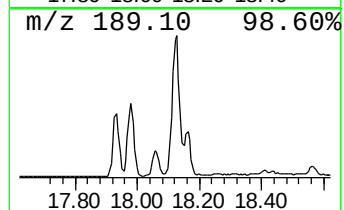
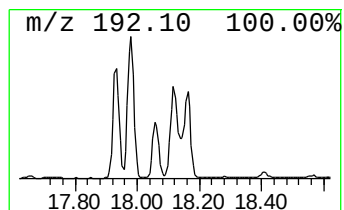
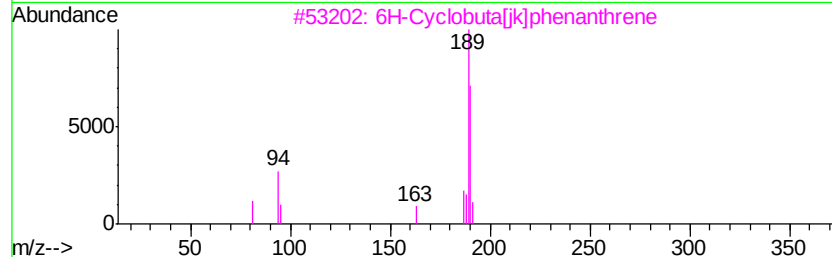
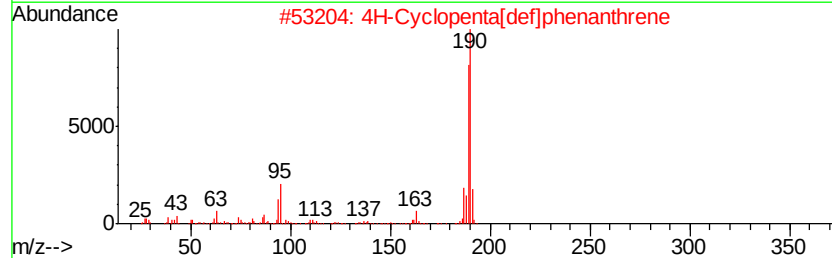
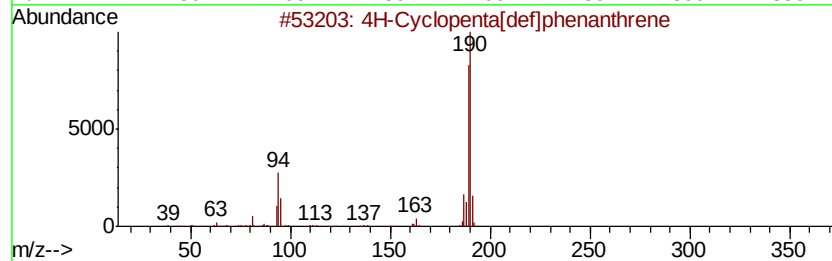
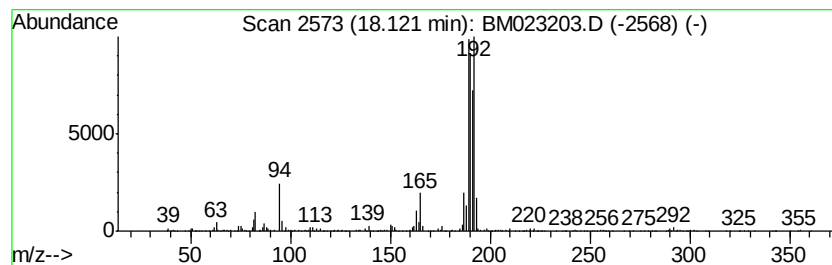
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 18.12 | 6.01 ng/ul | 1570030 | Phenanthrene-d10 | 17.06 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|---------|---|-------------------------------------|-----|---------|-------------|------|
| 1       |   | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10  | 000203-64-5 | 60   |
| 2       |   | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10  | 000203-64-5 | 55   |
| 3       |   | 6H-Cyclobuta[ik]phenanthrene        | 190 | C15H10  | 083469-43-6 | 49   |
| 4       |   | Anthracene, 1-methyl-               | 192 | C15H12  | 000610-48-0 | 46   |
| 5       |   | 1a,9b-Dihydro-1H-cyclopropa[a]an... | 192 | C15H12  | 006109-24-6 | 42   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101619\  
Data File : BM023203.D  
Acq On : 16 Oct 2019 22:45  
Operator : JU  
Sample : K5262-06  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

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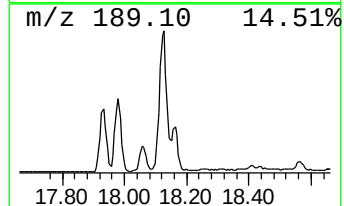
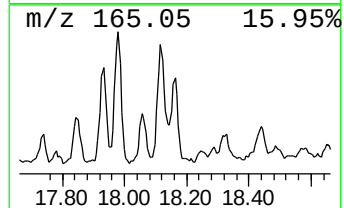
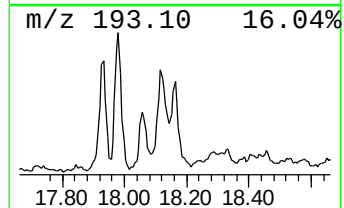
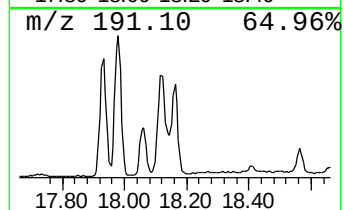
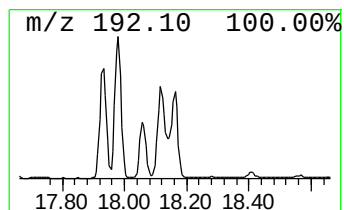
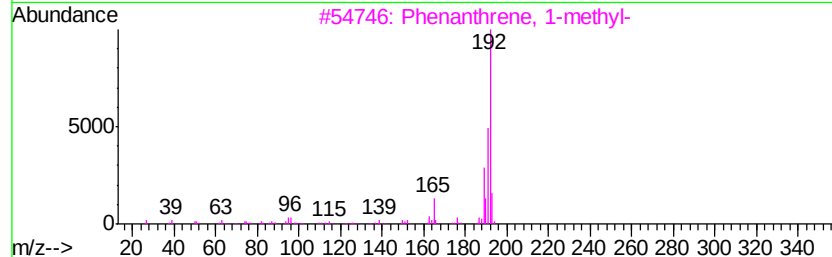
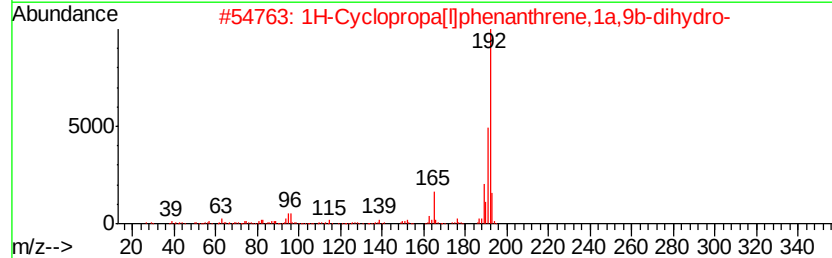
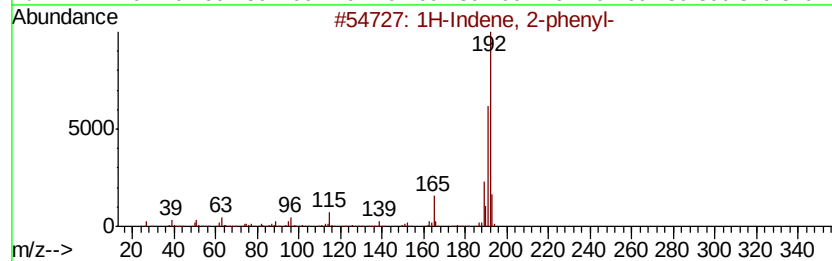
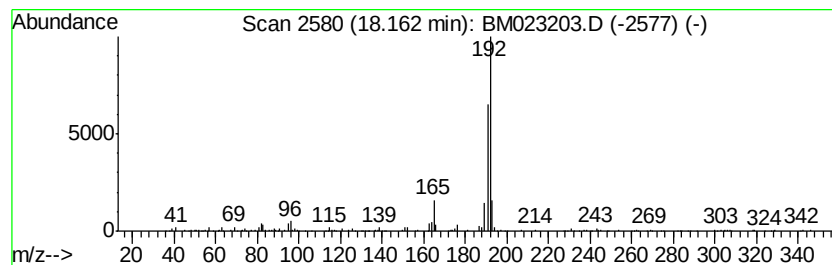
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 1H-Cyclopropa[l]phenanthren... Concentration Rank 8

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.16 | 2.64 ng/ul | 689134 | Phenanthrene-d10 | 17.06 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|---------|---|-------------------------------------|-----|---------|-------------|------|
| 1       |   | 1H-Indene, 2-phenyl-                | 192 | C15H12  | 004505-48-0 | 93   |
| 2       |   | 1H-Cyclopropa[l]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 93   |
| 3       |   | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 93   |
| 4       |   | 1H-Indene, 2-phenyl-                | 192 | C15H12  | 004505-48-0 | 90   |
| 5       |   | Propyne, 1,3-diphenyl-              | 192 | C15H12  | 004980-70-5 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101619\  
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Acq On : 16 Oct 2019 22:45  
Operator : JU  
Sample : K5262-06  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

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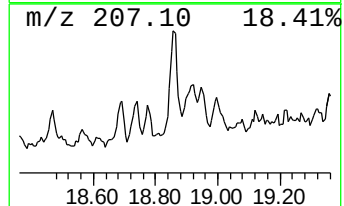
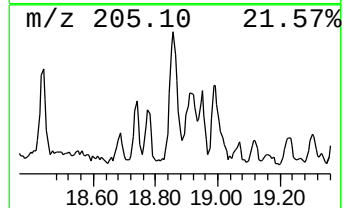
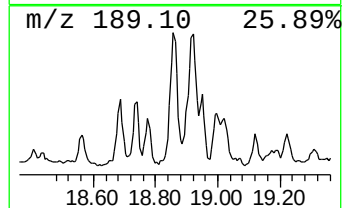
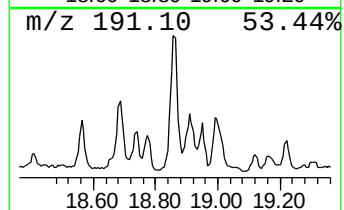
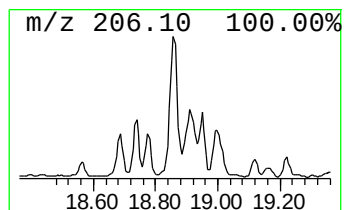
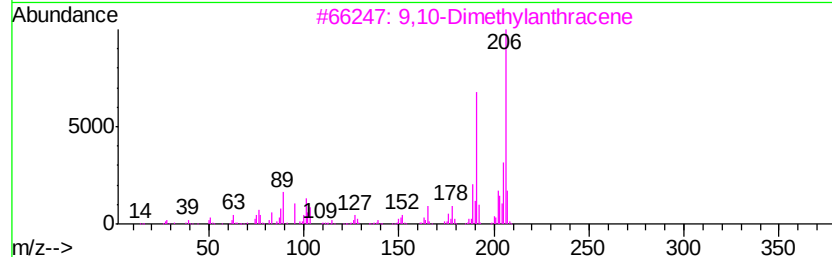
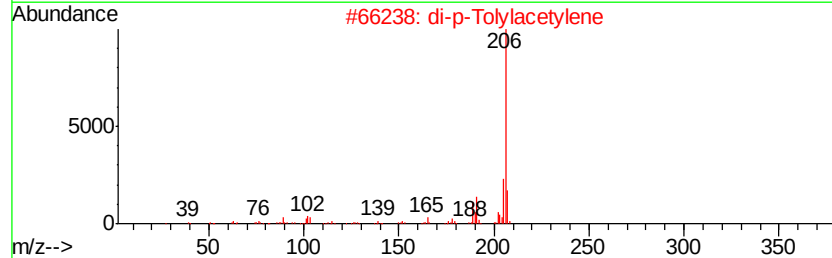
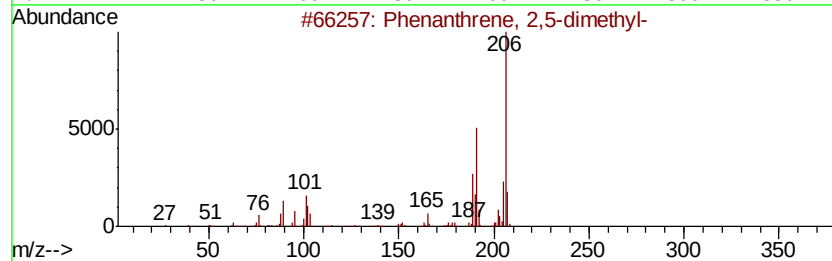
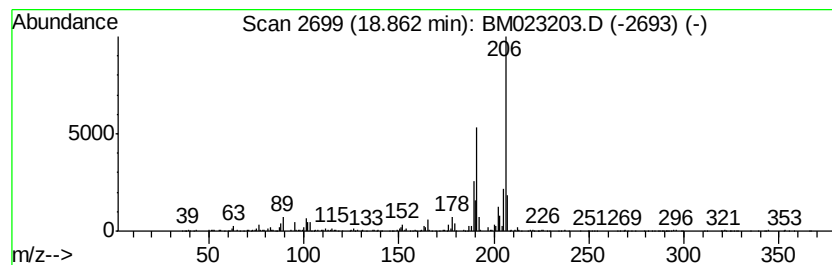
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Phenanthrene, 2,5-dimethyl- Concentration Rank 5

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.86 | 3.65 ng/ul | 952110 | Phenanthrene-d10 | 17.06 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenanthrene, 2,5-dimethyl-   | 206 | C16H14  | 003674-66-6 | 93   |
| 2    |    | di-p-Tolylacetylene           | 206 | C16H14  | 002789-88-0 | 91   |
| 3    |    | 9,10-Dimethylantracene        | 206 | C16H14  | 000781-43-1 | 91   |
| 4    |    | 1H-Indene, 2-methyl-3-phenyl- | 206 | C16H14  | 035099-60-6 | 90   |
| 5    |    | di-p-Tolylacetylene           | 206 | C16H14  | 002789-88-0 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101619\  
Data File : BM023203.D  
Acq On : 16 Oct 2019 22:45  
Operator : JU  
Sample : K5262-06  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

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

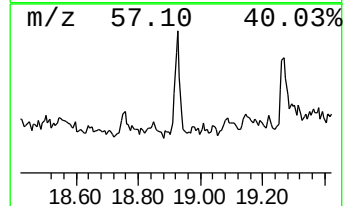
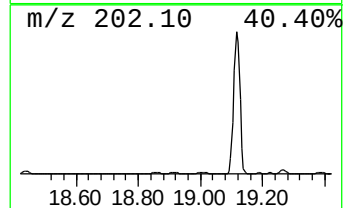
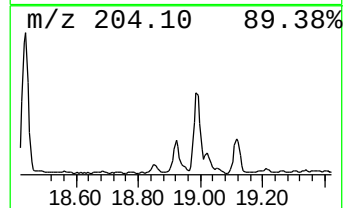
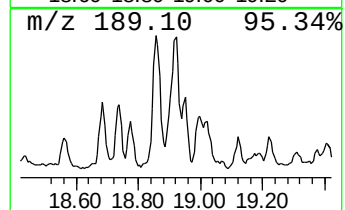
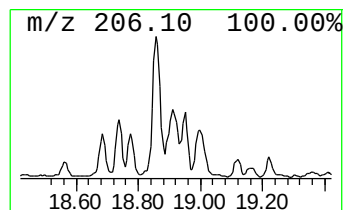
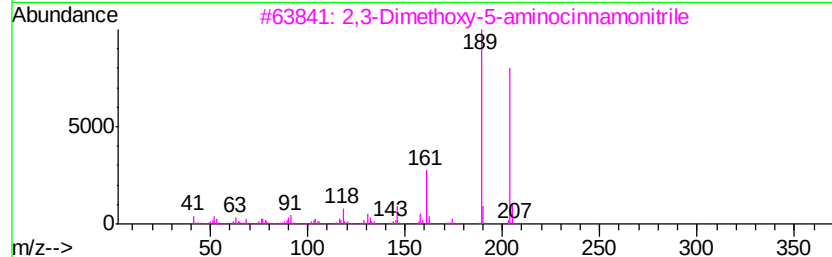
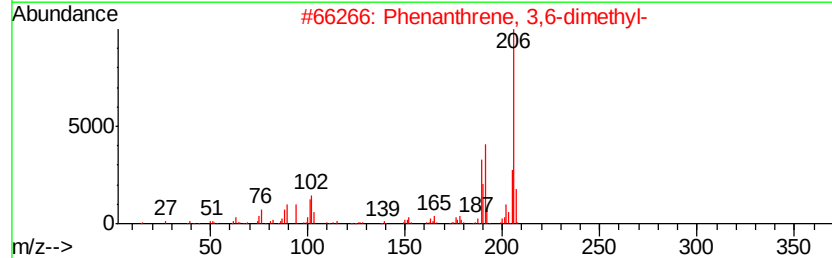
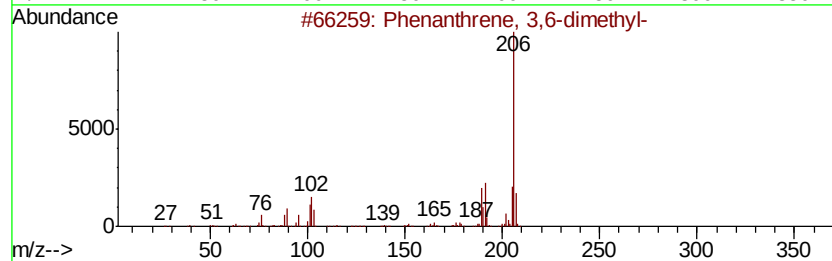
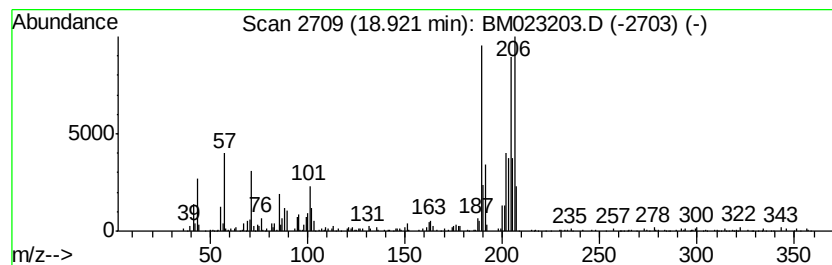
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 unknown-01 Concentration Rank 11

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.92 | 2.03 ng/ul | 530087 | Phenanthrene-d10 | 17.06 |

| Hit# | of | Tentative ID                         | MW  | MolForm    | CAS#         | Qual |
|------|----|--------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Phenanthrene, 3,6-dimethyl-          | 206 | C16H14     | 001576-67-6  | 38   |
| 2    |    | Phenanthrene, 3,6-dimethyl-          | 206 | C16H14     | 001576-67-6  | 38   |
| 3    |    | 2,3-Dimethoxy-5-aminocinnamionitrile | 204 | C11H12N2O2 | 1000214-46-5 | 38   |
| 4    |    | Naphthalene, 2-phenyl-               | 204 | C16H12     | 000612-94-2  | 35   |
| 5    |    | Naphthalene, 2-phenyl-               | 204 | C16H12     | 000612-94-2  | 35   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101619\  
Data File : BM023203.D  
Acq On : 16 Oct 2019 22:45  
Operator : JU  
Sample : K5262-06  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

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

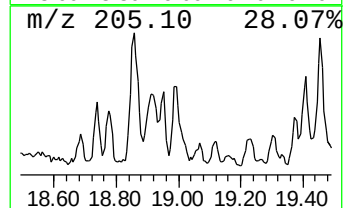
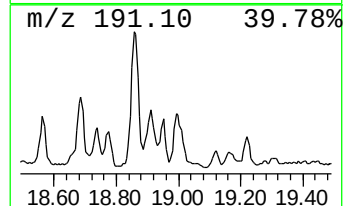
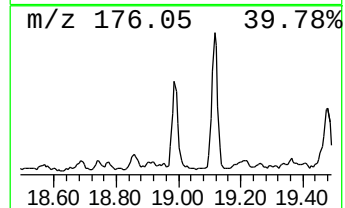
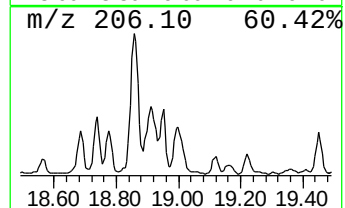
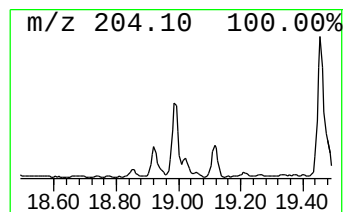
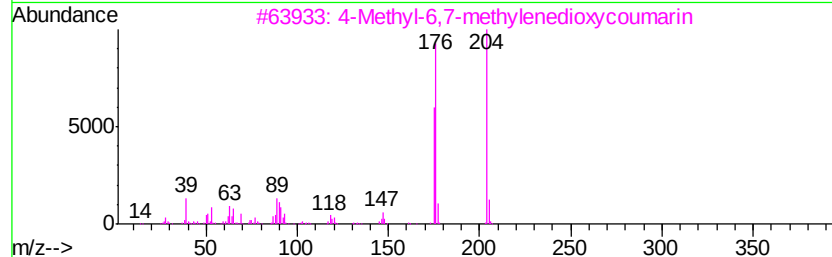
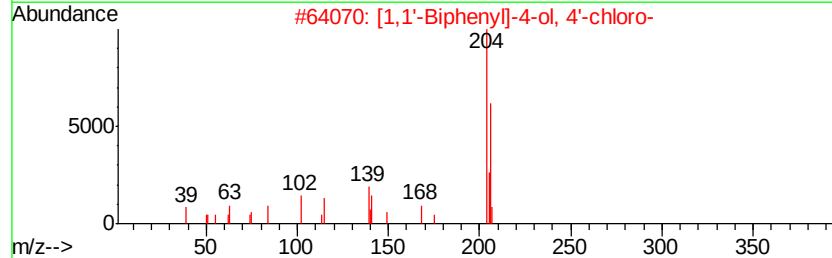
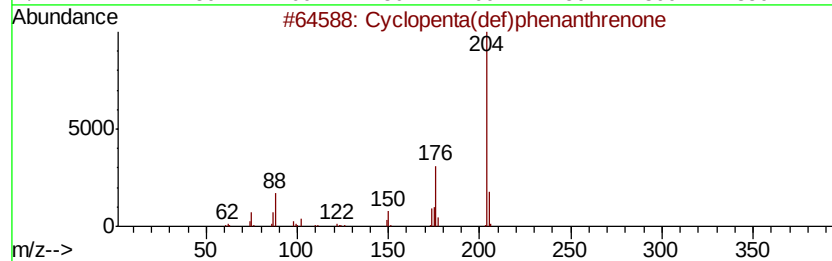
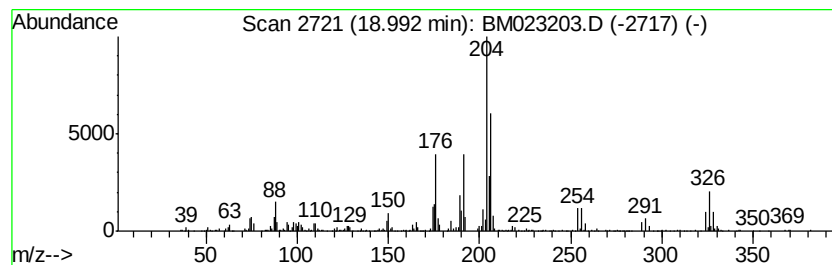
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Cyclopenta(def)phenanthrene Concentration Rank 7

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.99 | 2.92 ng/ul | 762661 | Phenanthrene-d10 | 17.06 |

| Hit# | of | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Cyclopenta(def)phenanthrene          | 204 | C15H8O   | 005737-13-3 | 89   |
| 2    |    | [1,1'-Biphenyl]-4-ol, 4'-chloro-     | 204 | C12H9ClO | 028034-99-3 | 43   |
| 3    |    | 4-Methyl-6,7-methylenedioxy coumarin | 204 | C11H8O4  | 015071-04-2 | 38   |
| 4    |    | 9,10-Dimethylantracene               | 206 | C16H14   | 000781-43-1 | 38   |
| 5    |    | [14]Annulene, 1,6:8,13-bis(metha...  | 206 | C16H14   | 085385-68-8 | 30   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101619\  
Data File : BM023203.D  
Acq On : 16 Oct 2019 22:45  
Operator : JU  
Sample : K5262-06  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BFB71

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101419MA.M  
Quant Title : SVOA CALIBRATION

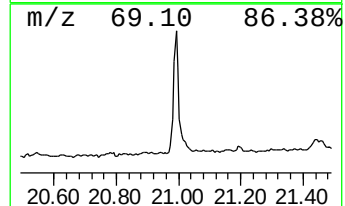
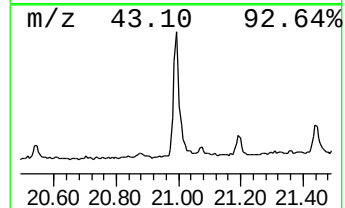
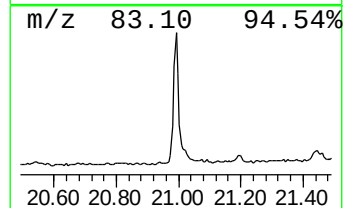
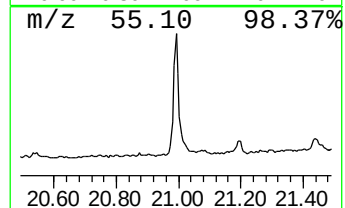
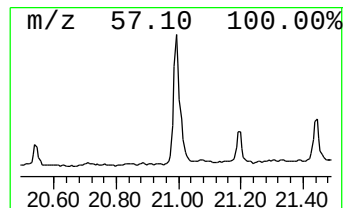
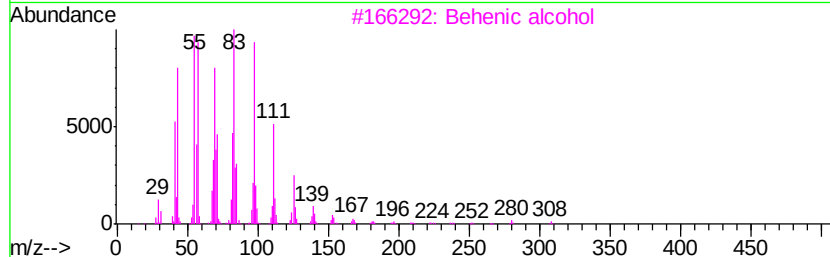
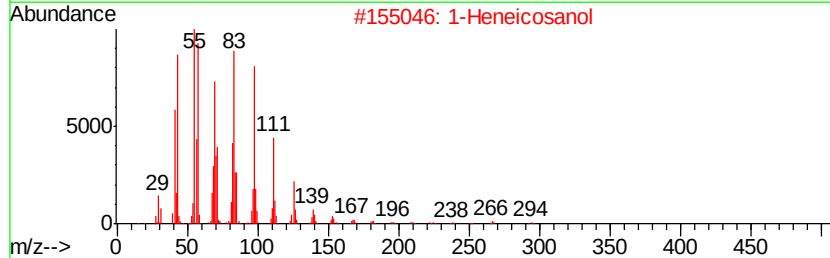
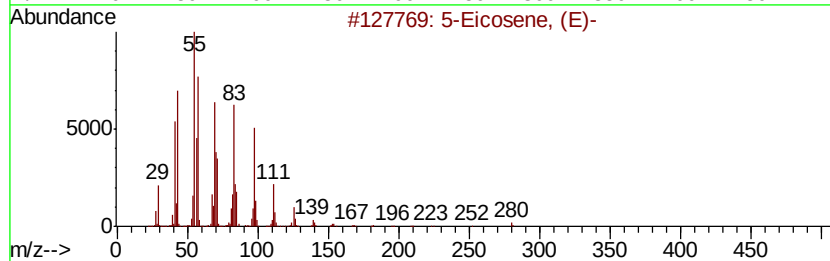
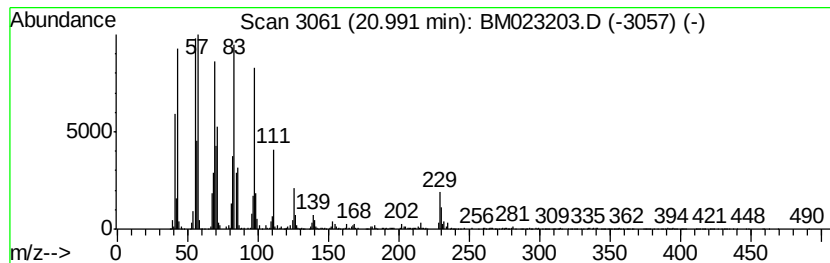
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 (DEL) Alkane: Cyclic20.99 Concentration Rank 4

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 20.99 | 4.01 ng/ul | 2224170 | Chrysene-d12     | 21.24 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|-------------------------------------|-----|-------------|-------------|------|
| 1    | 5  | 5-Eicosene, (E)-                    | 280 | C20H40      | 074685-30-6 | 95   |
| 2    |    | 1-Heneicosanol                      | 312 | C21H44O     | 015594-90-8 | 94   |
| 3    |    | Behenic alcohol                     | 326 | C22H46O     | 000661-19-8 | 94   |
| 4    |    | n-Tetracosanol-1                    | 354 | C24H50O     | 000506-51-4 | 94   |
| 5    |    | Trichloroacetic acid, hexadecyl ... | 386 | C18H33Cl3O2 | 074339-54-1 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101619\  
Data File : BM023203.D  
Acq On : 16 Oct 2019 22:45  
Operator : JU  
Sample : K5262-06  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BFB71

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101419MA.M  
Quant Title : SVOA CALIBRATION

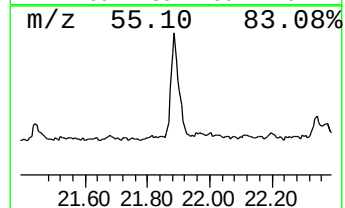
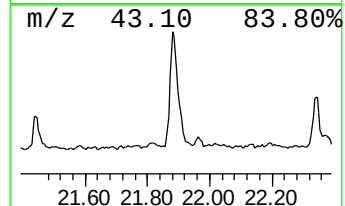
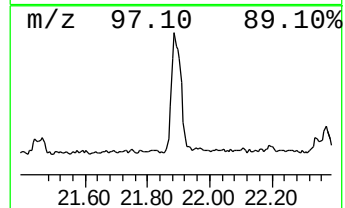
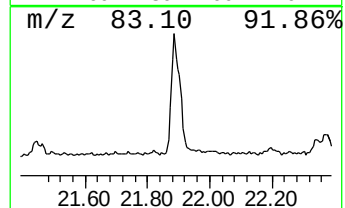
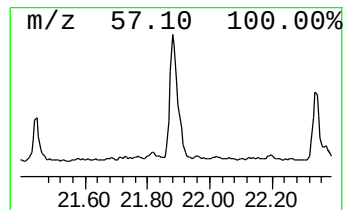
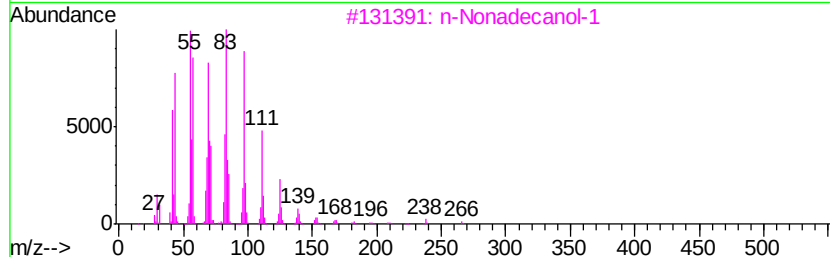
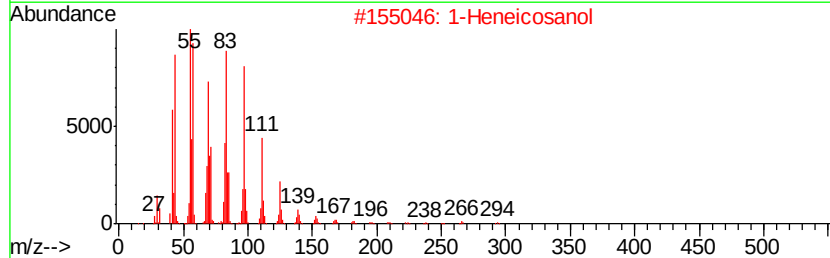
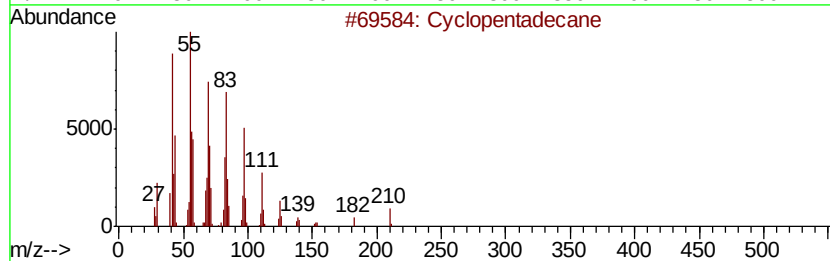
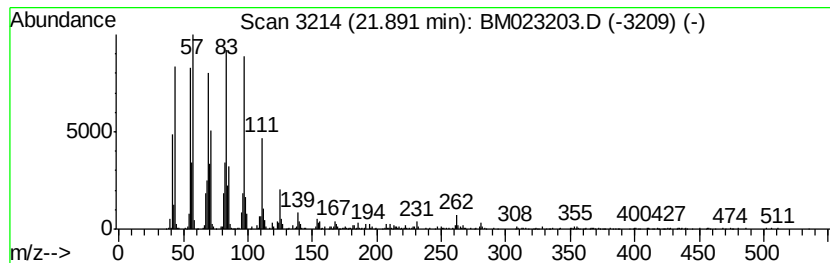
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 (DEL) Alkane: Cyclic21.89 Concentration Rank 3

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 21.89 | 4.35 ng/ul | 2414160 | Chrysene-d12     | 21.24 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | Cyclopentadecane                    | 210 | C15H30     | 000295-48-7 | 93   |
| 2    |    | 1-Heneicosanol                      | 312 | C21H44O    | 015594-90-8 | 93   |
| 3    |    | n-Nonadecanol-1                     | 284 | C19H40O    | 001454-84-8 | 93   |
| 4    |    | Pentafluoropropionic acid, octad... | 416 | C21H37F5O2 | 959261-25-7 | 91   |
| 5    |    | n-Tetracosanol-1                    | 354 | C24H50O    | 000506-51-4 | 91   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM101619\  
 Data File : BM023203.D  
 Acq On : 16 Oct 2019 22:45  
 Operator : JU  
 Sample : K5262-06  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BFB71

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101419MA.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 |
| Furan, tetrahydro... | 3.68  | 2.4     | ng/ul | 326383   | 1                     | 7.66  | 2712760  | 20.0 |
| Anthracene, 2-met... | 17.93 | 3.3     | ng/ul | 871186   | 4                     | 17.06 | 5221700  | 20.0 |
| Phenanthrene, 2-m... | 17.98 | 7.2     | ng/ul | 1886170  | 4                     | 17.06 | 5221700  | 20.0 |
| Anthracene, 1-met... | 18.06 | 2.4     | ng/ul | 633616   | 4                     | 17.06 | 5221700  | 20.0 |
| 4H-Cyclopenta[def... | 18.12 | 6.0     | ng/ul | 1570030  | 4                     | 17.06 | 5221700  | 20.0 |
| 1H-Cyclopropa[l]p... | 18.16 | 2.6     | ng/ul | 689134   | 4                     | 17.06 | 5221700  | 20.0 |
| Phenanthrene, 2,5... | 18.86 | 3.6     | ng/ul | 952110   | 4                     | 17.06 | 5221700  | 20.0 |
| unknown-01           | 18.92 | 2.0     | ng/ul | 530087   | 4                     | 17.06 | 5221700  | 20.0 |
| Cyclopenta(def)ph... | 18.99 | 2.9     | ng/ul | 762661   | 4                     | 17.06 | 5221700  | 20.0 |
| (DEL) Alkane: Cyc... | 20.99 | 4.0     | ng/ul | 2224170  | 5                     | 21.24 | 11094200 | 20.0 |
| (DEL) Alkane: Cyc... | 21.89 | 4.3     | ng/ul | 2414160  | 5                     | 21.24 | 11094200 | 20.0 |