

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM011516\  
 Data File : BM004032.D  
 Acq On : 15 Jan 2016 03:41  
 Operator : SJ/UM  
 Sample : H1099-18  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BC9E4

## Integration Parameters: LSCINT.P

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

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

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

Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010116.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         | 6.845       | 633           | 639         | 648          | rBV      | 121457         | 194352        | 16.40%          | 1.041%        |
| 2         | 6.998       | 660           | 665         | 674          | rBB      | 96872          | 149340        | 12.60%          | 0.800%        |
| 3         | 7.198       | 693           | 699         | 706          | rBV      | 131962         | 200082        | 16.88%          | 1.072%        |
| 4         | 7.663       | 772           | 778         | 785          | rBB      | 167053         | 256092        | 21.61%          | 1.372%        |
| 5         | 8.375       | 894           | 899         | 907          | rBV      | 147012         | 229354        | 19.35%          | 1.229%        |
| 6         | 8.822       | 969           | 975         | 983          | rBB      | 120773         | 190907        | 16.11%          | 1.023%        |
| 7         | 9.539       | 1092          | 1097        | 1104         | rBB      | 101308         | 159682        | 13.48%          | 0.855%        |
| 8         | 10.081      | 1183          | 1189        | 1200         | rBB      | 160780         | 264637        | 22.33%          | 1.418%        |
| 9         | 10.451      | 1245          | 1252        | 1259         | rBV      | 266379         | 429305        | 36.23%          | 2.300%        |
| 10        | 10.592      | 1270          | 1276        | 1286         | rBV      | 73097          | 124135        | 10.48%          | 0.665%        |
| 11        | 13.727      | 1804          | 1809        | 1814         | rBV      | 298093         | 421301        | 35.55%          | 2.257%        |
| 12        | 13.774      | 1814          | 1817        | 1824         | rVB      | 141199         | 193178        | 16.30%          | 1.035%        |
| 13        | 14.010      | 1851          | 1857        | 1871         | rBB      | 359222         | 552805        | 46.65%          | 2.961%        |
| 14        | 14.216      | 1888          | 1892        | 1897         | rBB      | 21158          | 26873         | 2.27%           | 0.144%        |
| 15        | 14.321      | 1903          | 1910        | 1917         | rBB2     | 395859         | 596482        | 50.34%          | 3.195%        |
| 16        | 14.539      | 1942          | 1947        | 1959         | rBV      | 98692          | 160885        | 13.58%          | 0.862%        |
| 17        | 15.174      | 2051          | 2055        | 2061         | rVB      | 37578          | 46391         | 3.91%           | 0.248%        |
| 18        | 15.316      | 2073          | 2079        | 2085         | rVV      | 470745         | 672546        | 56.75%          | 3.603%        |
| 19        | 15.368      | 2085          | 2088        | 2097         | rVB3     | 16637          | 26450         | 2.23%           | 0.142%        |
| 20        | 15.451      | 2097          | 2102        | 2108         | rBV      | 81540          | 110633        | 9.34%           | 0.593%        |
| 21        | 16.039      | 2197          | 2202        | 2209         | rBV2     | 42247          | 79084         | 6.67%           | 0.424%        |
| 22        | 16.374      | 2252          | 2259        | 2263         | rBV4     | 16376          | 30173         | 2.55%           | 0.162%        |
| 23        | 16.827      | 2333          | 2336        | 2340         | rVV      | 18976          | 22615         | 1.91%           | 0.121%        |
| 24        | 16.886      | 2342          | 2346        | 2355         | rVB2     | 16379          | 28082         | 2.37%           | 0.150%        |
| 25        | 17.068      | 2371          | 2377        | 2381         | rBV      | 486332         | 710521        | 59.96%          | 3.806%        |
| 26        | 17.115      | 2381          | 2385        | 2389         | rVV      | 285670         | 390562        | 32.96%          | 2.092%        |
| 27        | 17.168      | 2389          | 2394        | 2399         | rVV2     | 516129         | 738042        | 62.28%          | 3.953%        |
| 28        | 17.204      | 2399          | 2400        | 2406         | rVB      | 54799          | 42686         | 3.60%           | 0.229%        |
| 29        | 17.480      | 2443          | 2447        | 2452         | rBV      | 18889          | 27073         | 2.28%           | 0.145%        |
| 30        | 17.651      | 2469          | 2476        | 2483         | rVB4     | 16817          | 27711         | 2.34%           | 0.148%        |
| 31        | 17.945      | 2518          | 2526        | 2531         | rBV      | 78902          | 145529        | 12.28%          | 0.780%        |
| 32        | 17.992      | 2531          | 2534        | 2540         | rVV      | 101774         | 143998        | 12.15%          | 0.771%        |
| 33        | 18.074      | 2544          | 2548        | 2553         | rVV      | 22233          | 33760         | 2.85%           | 0.181%        |
| 34        | 18.139      | 2553          | 2559        | 2563         | rVV2     | 109861         | 199705        | 16.85%          | 1.070%        |

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Filtering: 5  
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 Max Peaks: 100  
 Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010116.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 35 | 18.180 | 2563 | 2566 | 2571 | rVB  | 64919  | 85030  | 7.18%  | 0.455% |
| 36 | 18.451 | 2608 | 2612 | 2616 | rBV  | 49117  | 63958  | 5.40%  | 0.343% |
| 37 | 18.498 | 2616 | 2620 | 2625 | rVB2 | 49877  | 67172  | 5.67%  | 0.360% |
| 38 | 18.698 | 2651 | 2654 | 2659 | rVB  | 23493  | 30280  | 2.56%  | 0.162% |
| 39 | 18.751 | 2659 | 2663 | 2666 | rBV  | 27039  | 31682  | 2.67%  | 0.170% |
| 40 | 18.792 | 2666 | 2670 | 2674 | rVB2 | 21281  | 28158  | 2.38%  | 0.151% |
| 41 | 18.874 | 2678 | 2684 | 2689 | rBV  | 78743  | 137085 | 11.57% | 0.734% |
| 42 | 18.933 | 2689 | 2694 | 2697 | rVV2 | 34966  | 70484  | 5.95%  | 0.378% |
| 43 | 18.968 | 2697 | 2700 | 2703 | rVB  | 23832  | 29095  | 2.46%  | 0.156% |
| 44 | 19.015 | 2703 | 2708 | 2715 | rBV2 | 42210  | 88266  | 7.45%  | 0.473% |
| 45 | 19.139 | 2721 | 2729 | 2735 | rVB  | 598162 | 806633 | 68.07% | 4.321% |
| 46 | 19.203 | 2735 | 2740 | 2743 | rBV  | 19336  | 25189  | 2.13%  | 0.135% |
| 47 | 19.239 | 2743 | 2746 | 2751 | rVV5 | 24293  | 35693  | 3.01%  | 0.191% |
| 48 | 19.339 | 2758 | 2763 | 2766 | rBV3 | 18493  | 29715  | 2.51%  | 0.159% |
| 49 | 19.386 | 2767 | 2771 | 2774 | rVV2 | 23561  | 34439  | 2.91%  | 0.184% |
| 50 | 19.474 | 2780 | 2786 | 2789 | rBV2 | 633985 | 993463 | 83.84% | 5.322% |
| 51 | 19.503 | 2789 | 2791 | 2799 | rVV  | 602175 | 721703 | 60.90% | 3.866% |
| 52 | 19.580 | 2799 | 2804 | 2809 | rVV2 | 38062  | 56911  | 4.80%  | 0.305% |
| 53 | 19.680 | 2815 | 2821 | 2828 | rBV2 | 21354  | 41553  | 3.51%  | 0.223% |
| 54 | 19.745 | 2828 | 2832 | 2837 | rVB3 | 28852  | 42434  | 3.58%  | 0.227% |
| 55 | 19.833 | 2841 | 2847 | 2853 | rBV4 | 56945  | 140522 | 11.86% | 0.753% |
| 56 | 19.968 | 2864 | 2870 | 2871 | rBV3 | 45984  | 65059  | 5.49%  | 0.348% |
| 57 | 20.003 | 2872 | 2876 | 2882 | rVB2 | 107253 | 164558 | 13.89% | 0.881% |
| 58 | 20.103 | 2889 | 2893 | 2897 | rVV  | 69442  | 90956  | 7.68%  | 0.487% |
| 59 | 20.162 | 2897 | 2903 | 2907 | rVV2 | 90922  | 143233 | 12.09% | 0.767% |
| 60 | 20.203 | 2907 | 2910 | 2914 | rVV3 | 17406  | 23732  | 2.00%  | 0.127% |
| 61 | 20.303 | 2923 | 2927 | 2931 | rBV  | 66829  | 90989  | 7.68%  | 0.487% |
| 62 | 20.345 | 2931 | 2934 | 2938 | rVV  | 66123  | 87364  | 7.37%  | 0.468% |
| 63 | 20.409 | 2942 | 2945 | 2949 | rVB2 | 19131  | 22446  | 1.89%  | 0.120% |
| 64 | 20.521 | 2959 | 2964 | 2967 | rBV7 | 15476  | 21959  | 1.85%  | 0.118% |
| 65 | 20.568 | 2967 | 2972 | 2974 | rBV4 | 16816  | 26898  | 2.27%  | 0.144% |
| 66 | 20.598 | 2974 | 2977 | 2980 | rVV4 | 20844  | 29788  | 2.51%  | 0.160% |
| 67 | 20.715 | 2993 | 2997 | 2999 | rBV2 | 18800  | 27094  | 2.29%  | 0.145% |
| 68 | 20.756 | 3000 | 3004 | 3010 | rVB  | 57408  | 94164  | 7.95%  | 0.504% |
| 69 | 20.921 | 3024 | 3032 | 3035 | rBV3 | 64534  | 126684 | 10.69% | 0.679% |
| 70 | 20.980 | 3039 | 3042 | 3050 | rVV4 | 136441 | 280484 | 23.67% | 1.502% |
| 71 | 21.050 | 3050 | 3054 | 3061 | rVB2 | 54827  | 97263  | 8.21%  | 0.521% |

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Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010116.M  
Title : SVOA CALIBRATION

|     |        |      |      |      |      |        |         |         |        |
|-----|--------|------|------|------|------|--------|---------|---------|--------|
| 72  | 21.274 | 3085 | 3092 | 3096 | rBV2 | 661810 | 1185018 | 100.00% | 6.348% |
| 73  | 21.309 | 3096 | 3098 | 3108 | rVB  | 299147 | 379686  | 32.04%  | 2.034% |
| 74  | 21.392 | 3108 | 3112 | 3115 | rBV2 | 21199  | 29700   | 2.51%   | 0.159% |
| 75  | 21.427 | 3115 | 3118 | 3124 | rBV3 | 50718  | 80572   | 6.80%   | 0.432% |
| 76  | 21.492 | 3124 | 3129 | 3130 | rBV2 | 36199  | 48088   | 4.06%   | 0.258% |
| 77  | 21.592 | 3141 | 3146 | 3150 | rVB3 | 29778  | 50068   | 4.23%   | 0.268% |
| 78  | 21.827 | 3179 | 3186 | 3189 | rBV  | 77709  | 137581  | 11.61%  | 0.737% |
| 79  | 21.874 | 3189 | 3194 | 3201 | rVB2 | 79074  | 155890  | 13.16%  | 0.835% |
| 80  | 22.021 | 3215 | 3219 | 3222 | rBV3 | 51093  | 70207   | 5.92%   | 0.376% |
| 81  | 22.115 | 3231 | 3235 | 3241 | rVB2 | 31466  | 51712   | 4.36%   | 0.277% |
| 82  | 22.309 | 3264 | 3268 | 3273 | rBV5 | 45529  | 76190   | 6.43%   | 0.408% |
| 83  | 22.439 | 3286 | 3290 | 3294 | rBV  | 42266  | 54635   | 4.61%   | 0.293% |
| 84  | 22.586 | 3310 | 3315 | 3321 | rBV3 | 24171  | 53857   | 4.54%   | 0.288% |
| 85  | 22.839 | 3349 | 3358 | 3370 | rBV2 | 202973 | 722929  | 61.01%  | 3.872% |
| 86  | 23.009 | 3383 | 3387 | 3392 | rVB  | 43305  | 69993   | 5.91%   | 0.375% |
| 87  | 23.139 | 3406 | 3409 | 3417 | rBV6 | 13710  | 33131   | 2.80%   | 0.177% |
| 88  | 23.321 | 3434 | 3440 | 3443 | rBV  | 123112 | 223349  | 18.85%  | 1.196% |
| 89  | 23.368 | 3443 | 3448 | 3452 | rVV  | 359490 | 657657  | 55.50%  | 3.523% |
| 90  | 23.415 | 3452 | 3456 | 3463 | rVB  | 202459 | 345884  | 29.19%  | 1.853% |
| 91  | 23.509 | 3465 | 3472 | 3478 | rVV  | 320612 | 634011  | 53.50%  | 3.396% |
| 92  | 23.556 | 3478 | 3480 | 3489 | rVB2 | 55416  | 98305   | 8.30%   | 0.527% |
| 93  | 24.015 | 3551 | 3558 | 3568 | rVB2 | 86310  | 186628  | 15.75%  | 1.000% |
| 94  | 24.750 | 3679 | 3683 | 3690 | rBV6 | 14849  | 31034   | 2.62%   | 0.166% |
| 95  | 25.521 | 3808 | 3814 | 3821 | rVB2 | 15740  | 37967   | 3.20%   | 0.203% |
| 96  | 25.615 | 3824 | 3830 | 3837 | rVB3 | 21492  | 51533   | 4.35%   | 0.276% |
| 97  | 25.756 | 3846 | 3854 | 3865 | rVB  | 94717  | 264727  | 22.34%  | 1.418% |
| 98  | 26.450 | 3963 | 3972 | 3981 | rBV2 | 63898  | 193312  | 16.31%  | 1.035% |
| 99  | 26.815 | 4028 | 4034 | 4049 | rVB2 | 17971  | 64184   | 5.42%   | 0.344% |
| 100 | 27.809 | 4196 | 4203 | 4218 | rVB2 | 28271  | 103560  | 8.74%   | 0.555% |

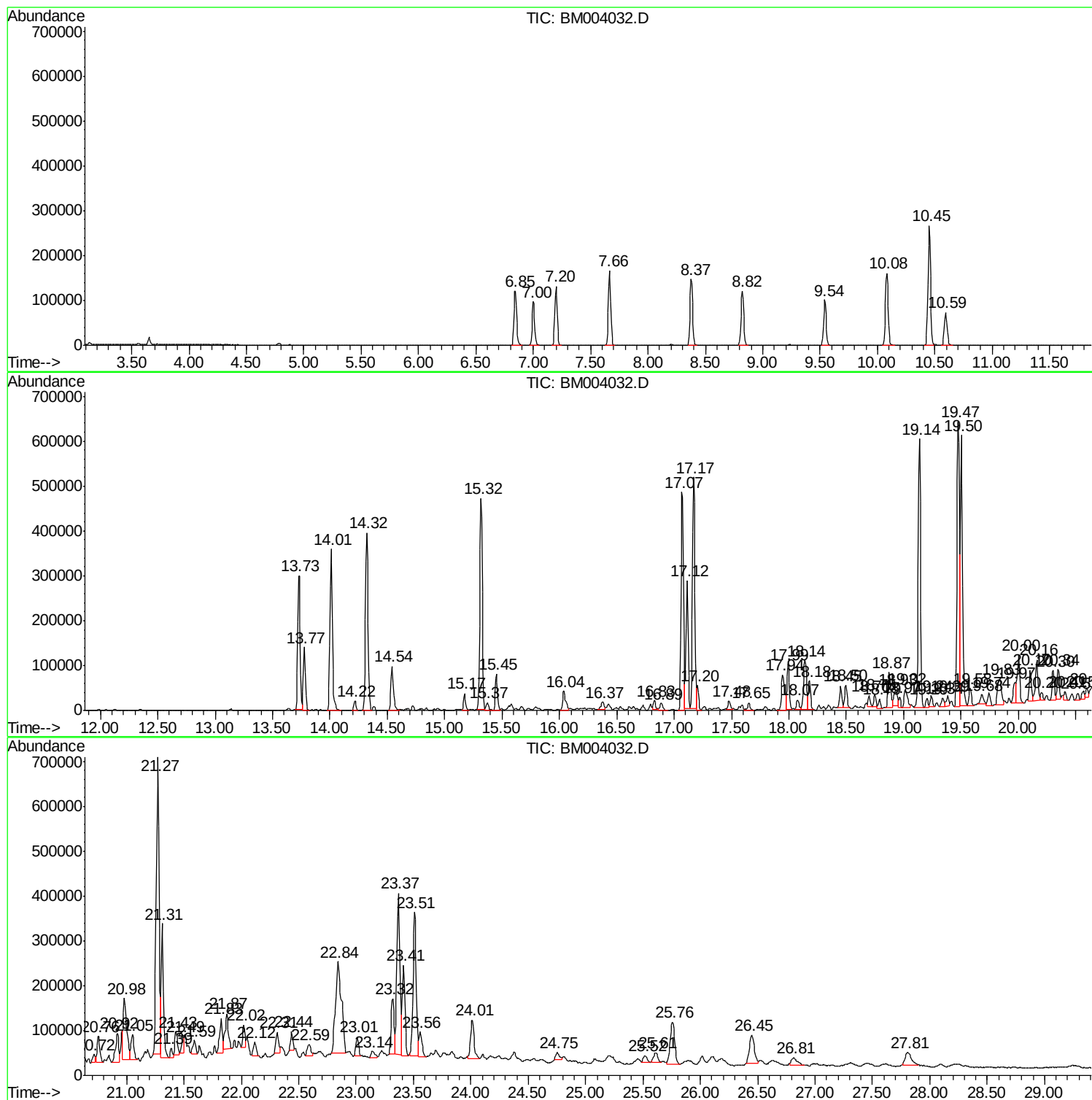
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Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010116.M  
Quant Title : SVOA CALIBRATION

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



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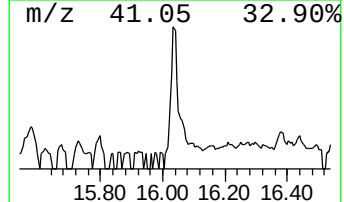
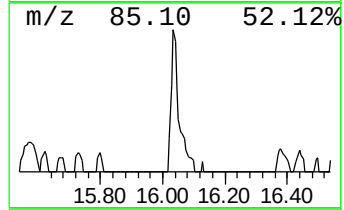
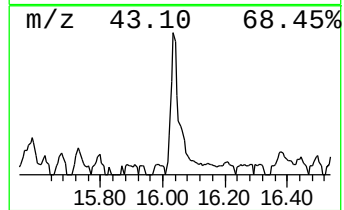
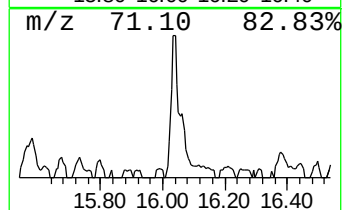
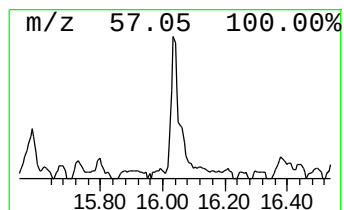
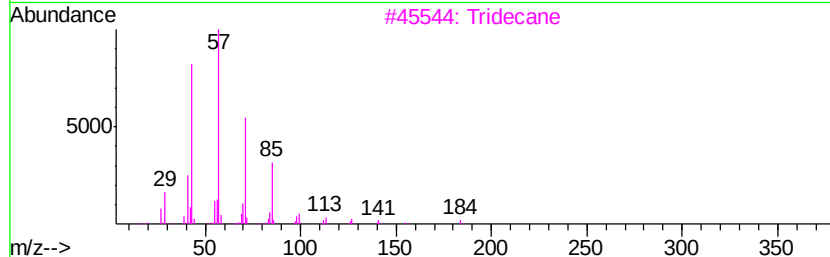
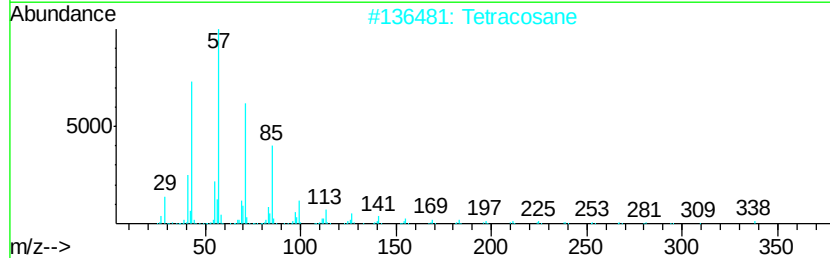
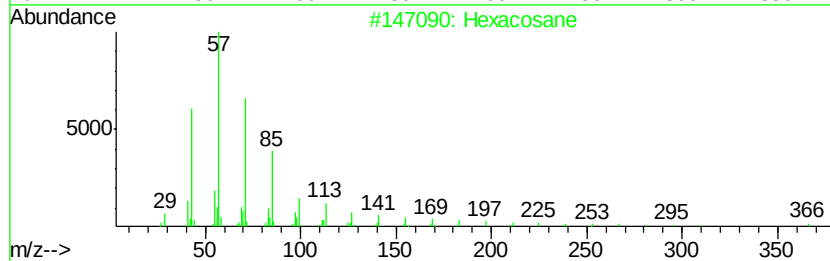
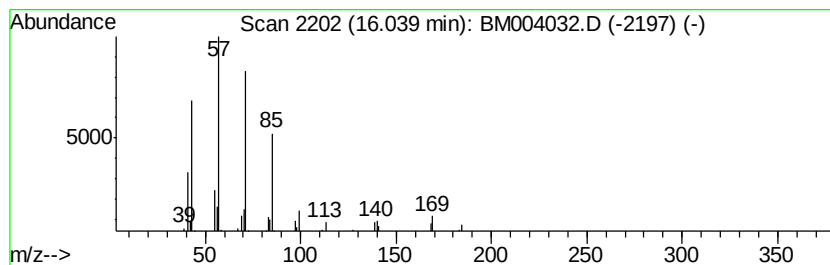
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 16.04 | 2.23 ng/ul | 79084 | Phenanthrene-d10 | 17.07 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------|-----|---------|-------------|------|
| 1    | 5  | Hexacosane          | 366 | C26H54  | 000630-01-3 | 86   |
| 2    |    | Tetracosane         | 338 | C24H50  | 000646-31-1 | 83   |
| 3    |    | Tridecane           | 184 | C13H28  | 000629-50-5 | 81   |
| 4    |    | 10-Methylnonadecane | 282 | C20H42  | 056862-62-5 | 80   |
| 5    |    | Tridecane           | 184 | C13H28  | 000629-50-5 | 74   |



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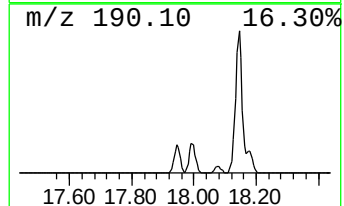
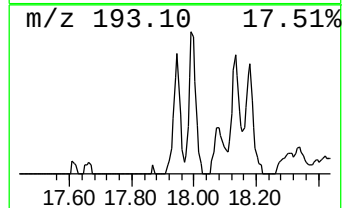
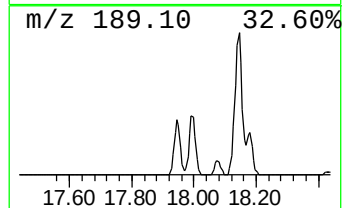
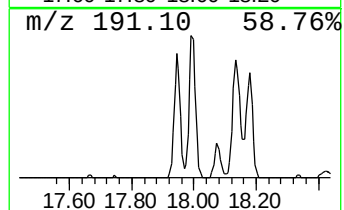
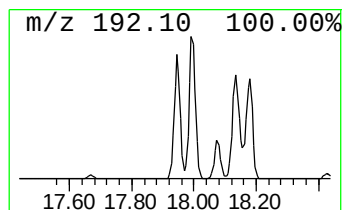
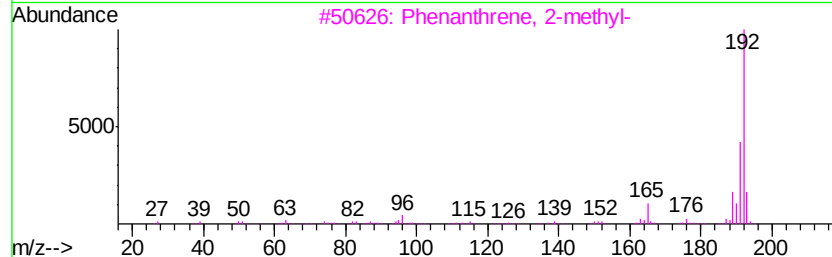
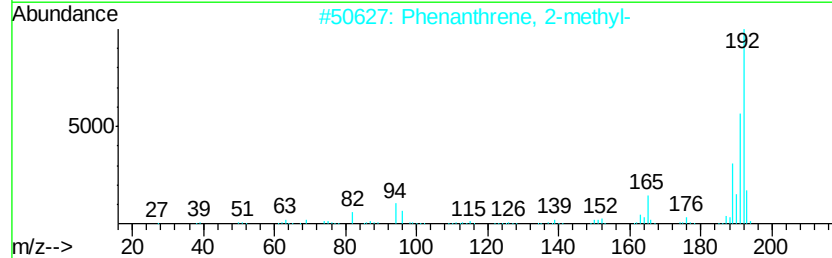
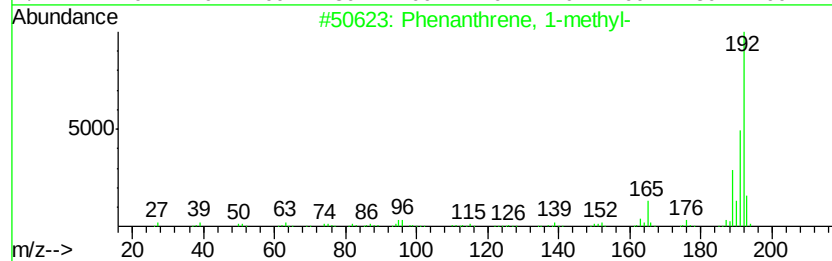
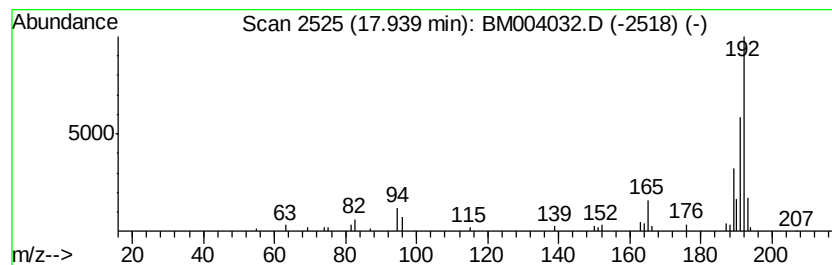
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 Phenanthrene, 1-methyl- Concentration Rank 4

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.94 | 4.10 ng/ul | 145529 | Phenanthrene-d10 | 17.07 |

| Hit# of | 5                                     | Tentative ID | MW     | MolForm     | CAS# | Qual |
|---------|---------------------------------------|--------------|--------|-------------|------|------|
| 1       | Phenanthrene, 1-methyl-               | 192          | C15H12 | 000832-69-9 | 97   |      |
| 2       | Phenanthrene, 2-methyl-               | 192          | C15H12 | 002531-84-2 | 97   |      |
| 3       | Phenanthrene, 2-methyl-               | 192          | C15H12 | 002531-84-2 | 96   |      |
| 4       | Anthracene, 2-methyl-                 | 192          | C15H12 | 000613-12-7 | 96   |      |
| 5       | 1H-Cyclopropa[l]phenanthrene, 1a, ... | 192          | C15H12 | 000949-41-7 | 96   |      |



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Acq On : 15 Jan 2016 03:41  
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Sample : H1099-18  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

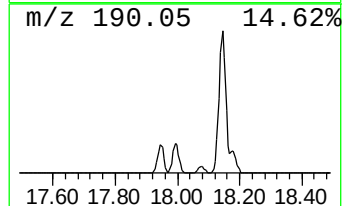
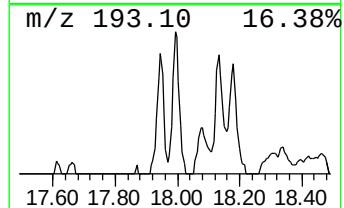
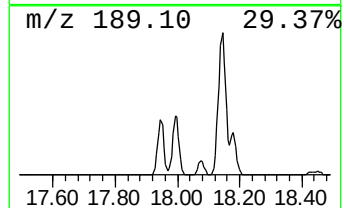
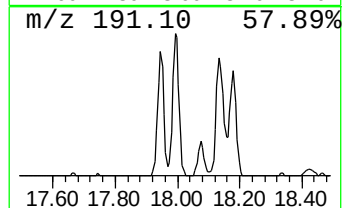
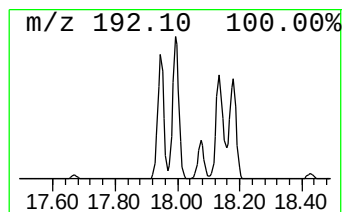
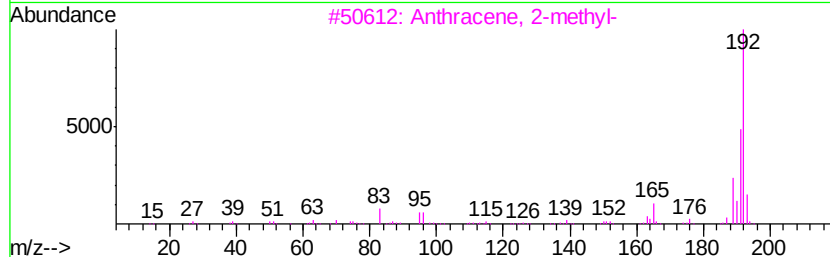
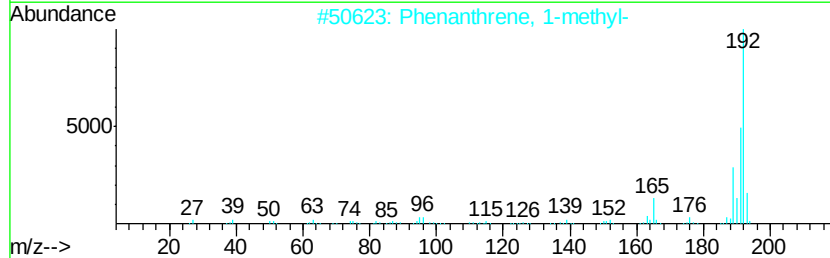
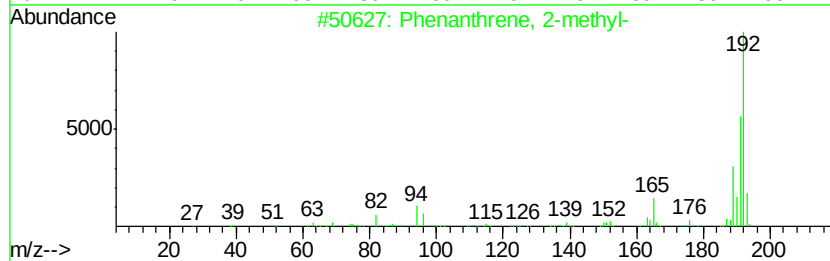
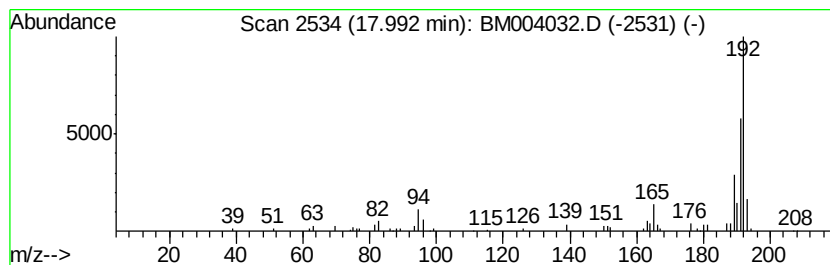
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 5

| R.T.      | EstConc                               | Area   | Relative to ISTD | R.T.        |      |
|-----------|---------------------------------------|--------|------------------|-------------|------|
| 17.99     | 4.05 ng/ul                            | 143998 | Phenanthrene-d10 | 17.07       |      |
| Hit# of 5 | Tentative ID                          | MW     | MolForm          | CAS#        | Qual |
| 1         | Phenanthrene, 2-methyl-               | 192    | C15H12           | 002531-84-2 | 98   |
| 2         | Phenanthrene, 1-methyl-               | 192    | C15H12           | 000832-69-9 | 97   |
| 3         | Anthracene, 2-methyl-                 | 192    | C15H12           | 000613-12-7 | 95   |
| 4         | Phenanthrene, 1-methyl-               | 192    | C15H12           | 000832-69-9 | 95   |
| 5         | 1H-Cyclopropa[l]phenanthrene, 1a, ... | 192    | C15H12           | 000949-41-7 | 95   |





Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM011516\  
Data File : BM004032.D  
Acq On : 15 Jan 2016 03:41  
Operator : SJ/UM  
Sample : H1099-18  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

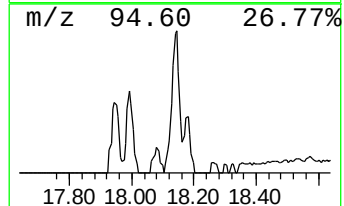
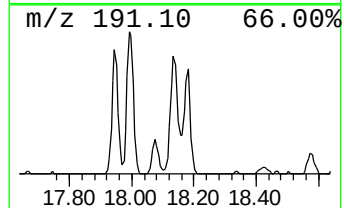
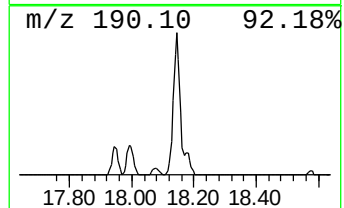
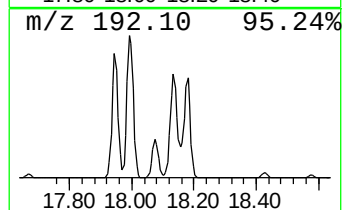
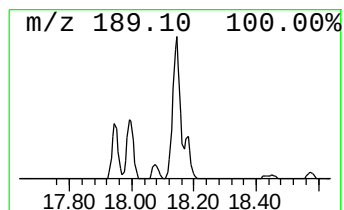
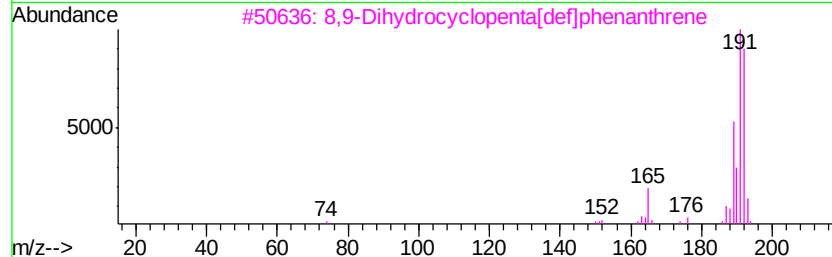
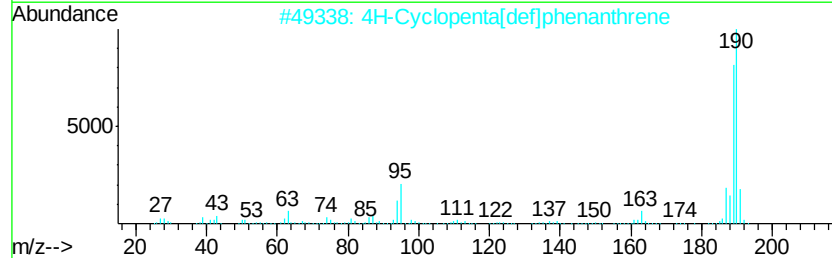
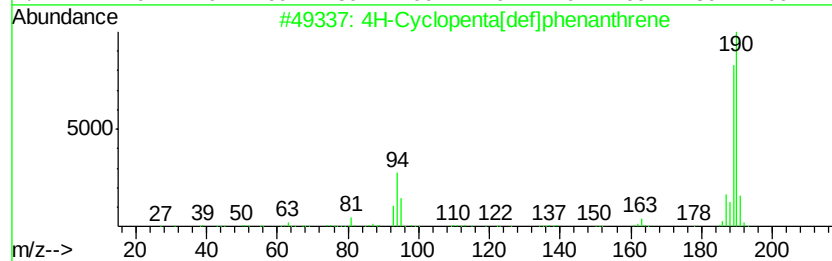
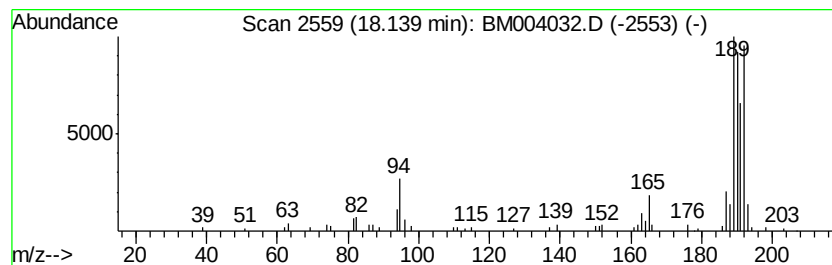
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ClientSampleId :  
BC9E4

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010116.M  
Quant Title : SVOA CALIBRATION

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

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

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 18.14     | 5.62 ng/ul                          | 199705 | Phenanthrene-d10 | 17.07       |      |
| 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         | 8,9-Dihydrocyclopenta[def]phenan... | 192    | C15H12           | 027410-55-5 | 50   |
| 4         | Anthracene, 1-methyl-               | 192    | C15H12           | 000610-48-0 | 49   |
| 5         | 5H-Dibenzo[a,d]cycloheptene         | 192    | C15H12           | 000256-81-5 | 46   |





Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM011516\  
Data File : BM004032.D  
Acq On : 15 Jan 2016 03:41  
Operator : SJ/UM  
Sample : H1099-18  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010116.M  
Quant Title : SVOA CALIBRATION

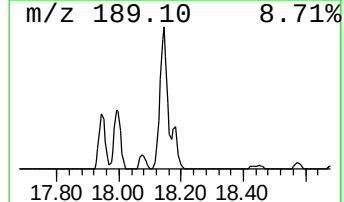
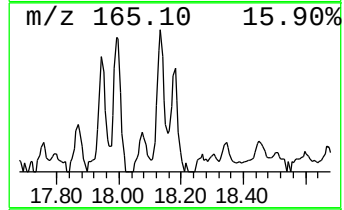
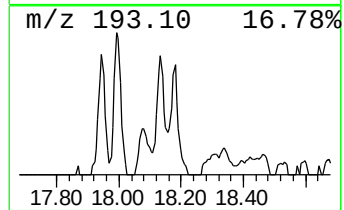
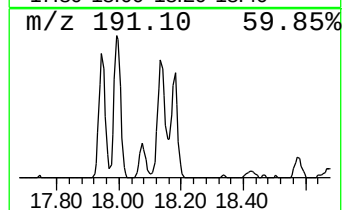
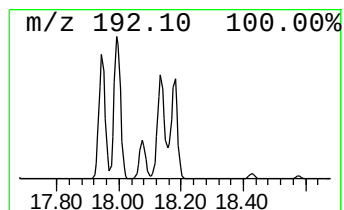
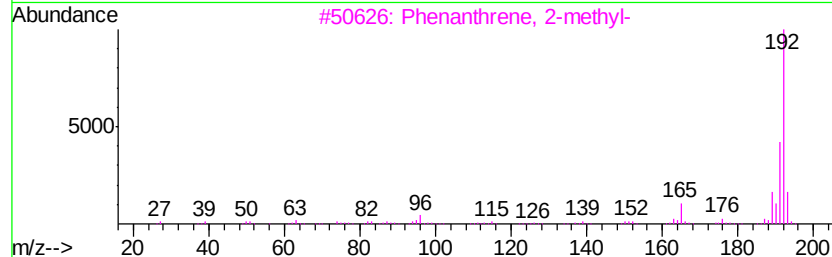
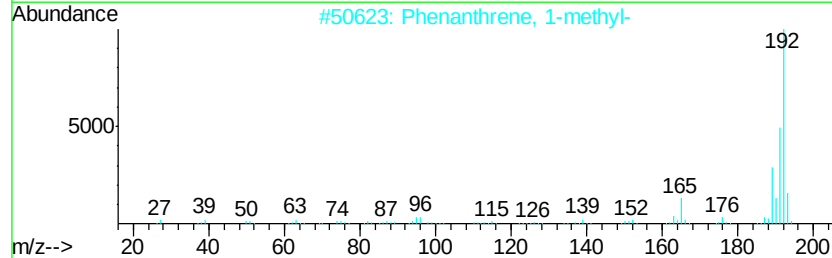
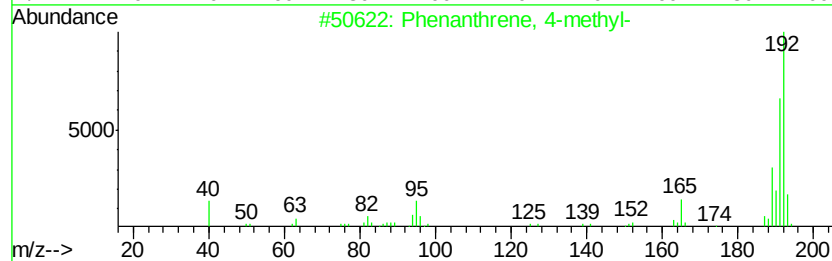
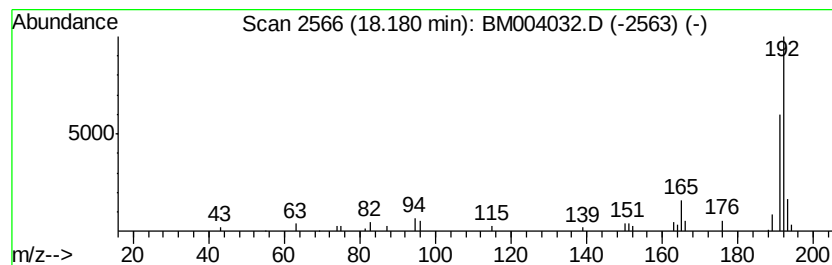
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Phenanthrene, 4-methyl- Concentration Rank 12

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 18.18 | 2.39 ng/ul | 85030 | Phenanthrene-d10 | 17.07 |

| Hit# of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|---------|---|-------------------------|-----|---------|-------------|------|
| 1       |   | Phenanthrene, 4-methyl- | 192 | C15H12  | 000832-64-4 | 93   |
| 2       |   | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 91   |
| 3       |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 91   |
| 4       |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 91   |
| 5       |   | 1H-Indene, 2-phenyl-    | 192 | C15H12  | 004505-48-0 | 90   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM011516\  
Data File : BM004032.D  
Acq On : 15 Jan 2016 03:41  
Operator : SJ/UM  
Sample : H1099-18  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampled :  
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Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010116.M  
Quant Title : SVOA CALIBRATION

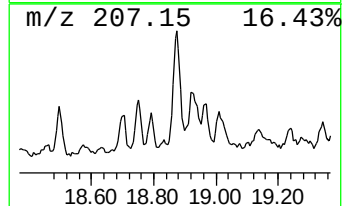
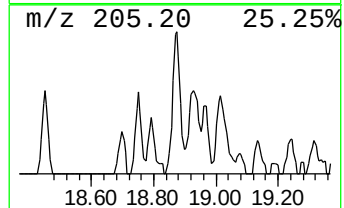
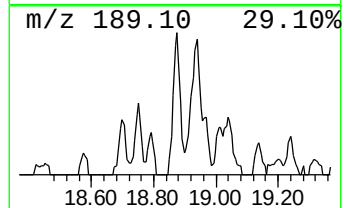
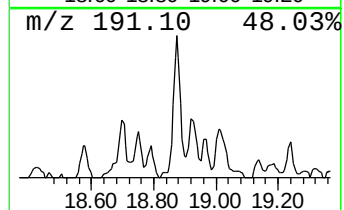
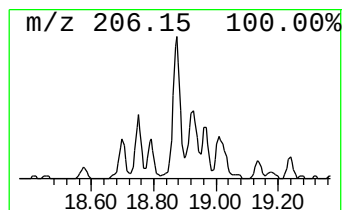
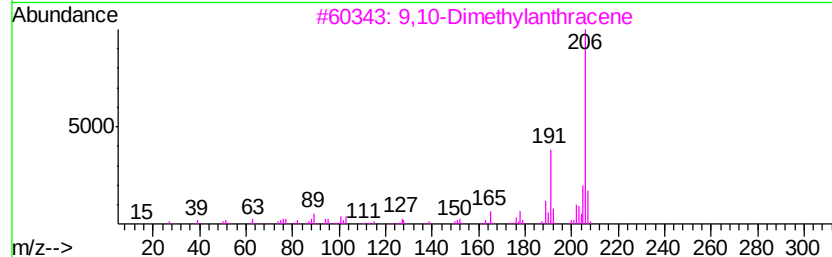
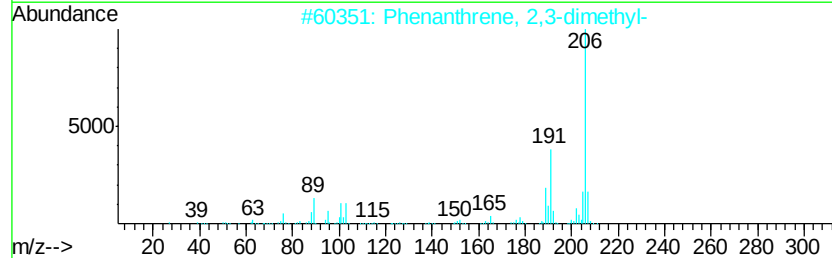
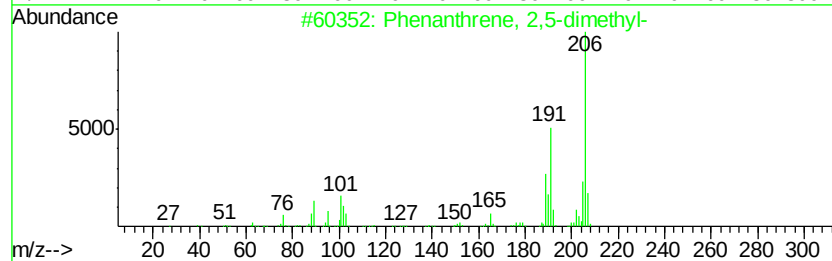
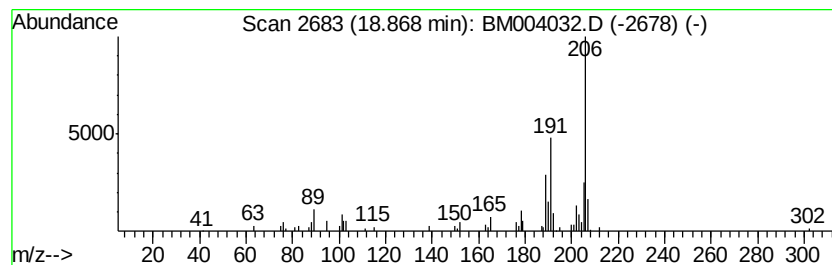
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Phenanthrene, 2,5-dimethyl- Concentration Rank 6

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.87 | 3.86 ng/ul | 137085 | Phenanthrene-d10 | 17.07 |

| Hit# of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|---------|---|-------------------------------|-----|---------|-------------|------|
| 1       |   | Phenanthrene, 2,5-dimethyl-   | 206 | C16H14  | 003674-66-6 | 93   |
| 2       |   | Phenanthrene, 2,3-dimethyl-   | 206 | C16H14  | 003674-65-5 | 93   |
| 3       |   | 9,10-Dimethylantracene        | 206 | C16H14  | 000781-43-1 | 91   |
| 4       |   | 1H-Indene, 2-methyl-3-phenyl- | 206 | C16H14  | 035099-60-6 | 91   |
| 5       |   | di-p-Tolylacetylene           | 206 | C16H14  | 002789-88-0 | 91   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM011516\  
Data File : BM004032.D  
Acq On : 15 Jan 2016 03:41  
Operator : SJ/UM  
Sample : H1099-18  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

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

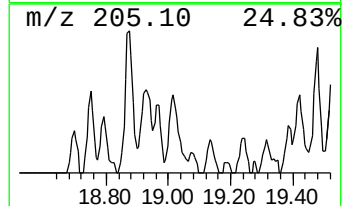
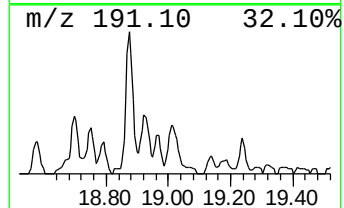
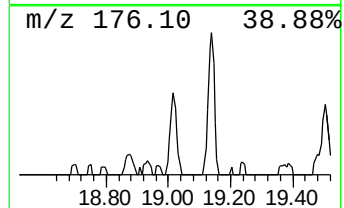
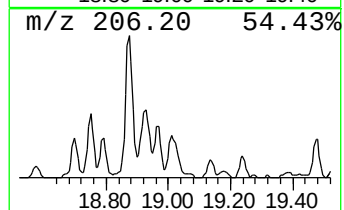
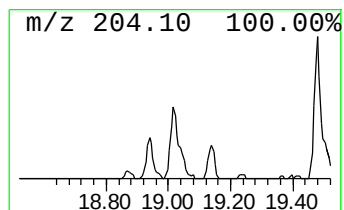
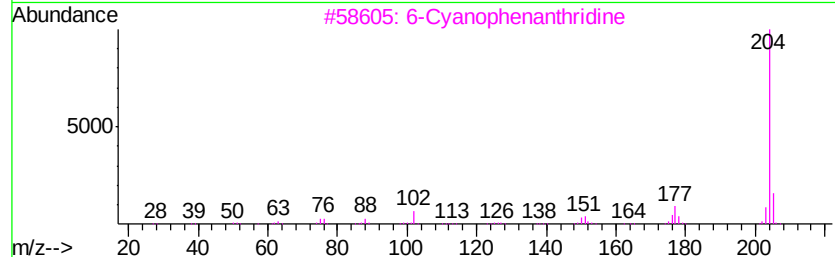
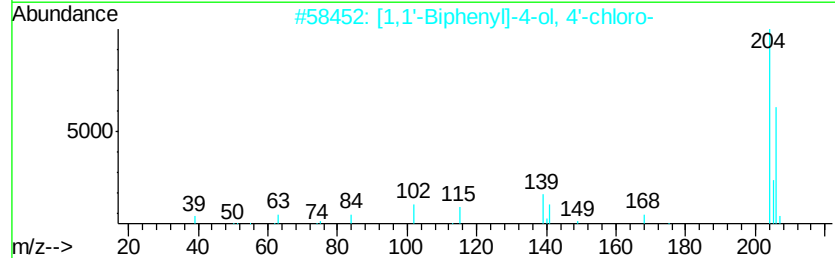
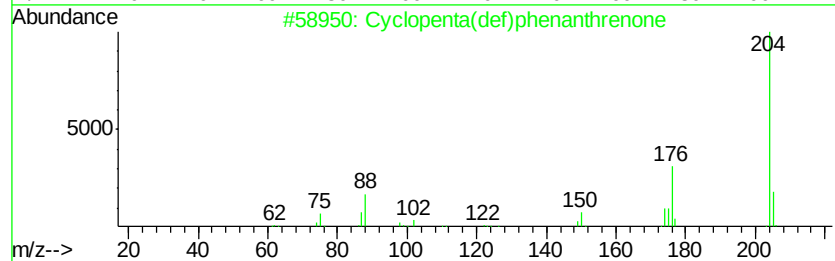
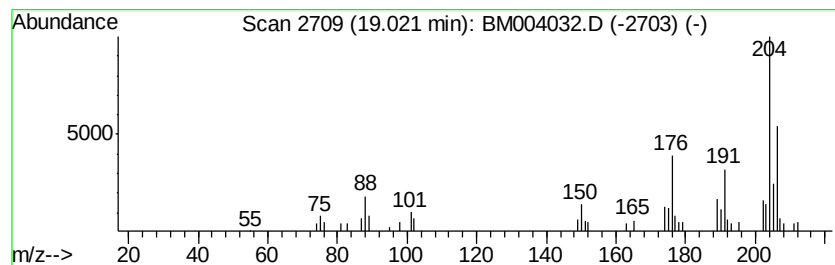
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Cyclopenta(def)phenanthrene Concentration Rank 10

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 19.02 | 2.48 ng/ul | 88266 | Phenanthrene-d10 | 17.07 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Cyclopenta(def)phenanthrene         | 204 | C15H8O   | 005737-13-3 | 86   |
| 2    |    | [1,1'-Biphenyl]-4-ol, 4'-chloro-    | 204 | C12H9ClO | 028034-99-3 | 46   |
| 3    |    | 6-Cyanophenanthridine               | 204 | C14H8N2  | 002176-28-5 | 30   |
| 4    |    | Benzene, (2-methylene-1-phenylcy... | 206 | C16H14   | 025152-47-0 | 30   |
| 5    |    | 9,10-Dimethylantracene              | 206 | C16H14   | 000781-43-1 | 30   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM011516\  
Data File : BM004032.D  
Acq On : 15 Jan 2016 03:41  
Operator : SJ/UM  
Sample : H1099-18  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010116.M  
Quant Title : SVOA CALIBRATION

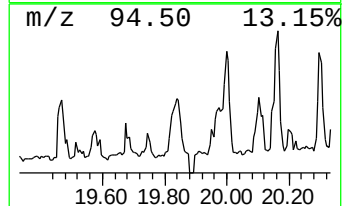
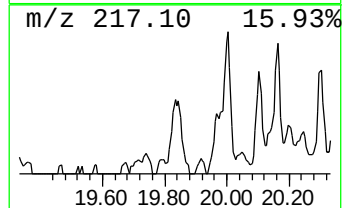
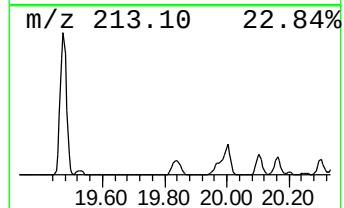
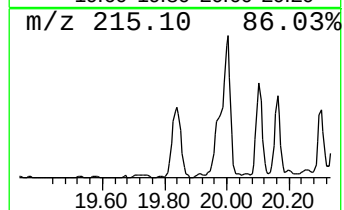
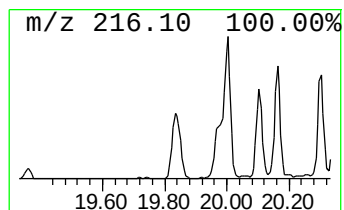
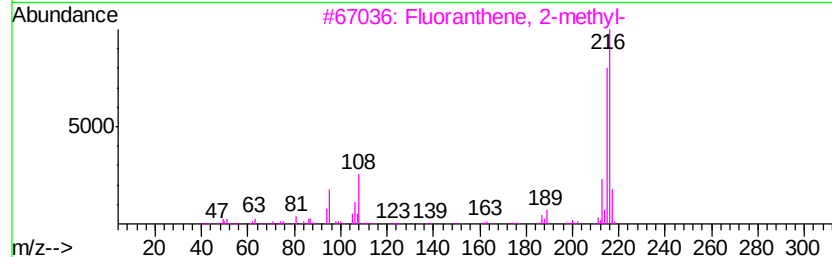
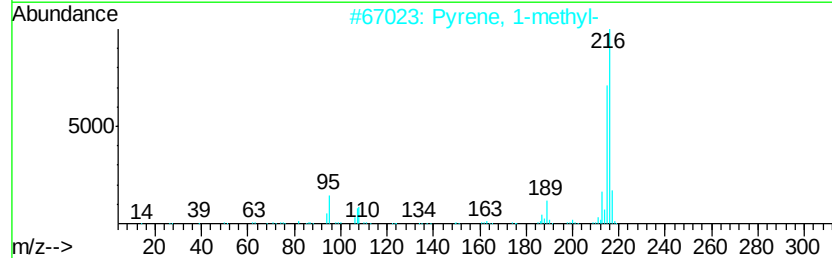
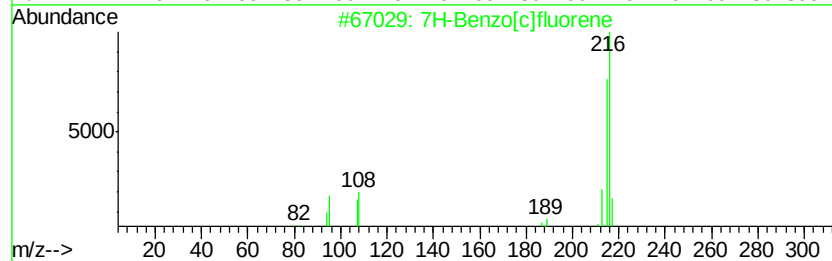
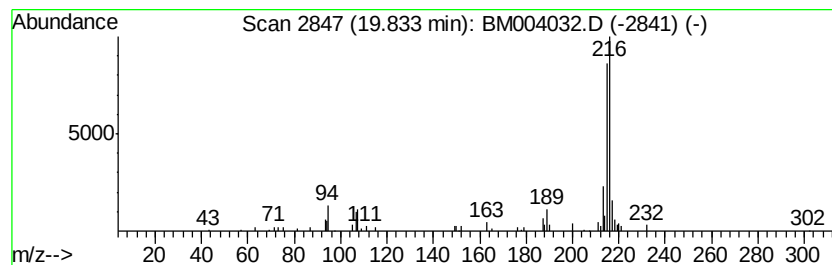
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 7H-Benzo[c]fluorene Concentration Rank 13

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.83 | 2.37 ng/ul | 140522 | Chrysene-d12     | 21.27 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | 7H-Benzo[c]fluorene     | 216 | C17H12  | 000205-12-9 | 87   |
| 2    |    | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 87   |
| 3    |    | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 81   |
| 4    |    | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 81   |
| 5    |    | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 80   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM011516\  
Data File : BM004032.D  
Acq On : 15 Jan 2016 03:41  
Operator : SJ/UM  
Sample : H1099-18  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BC9E4

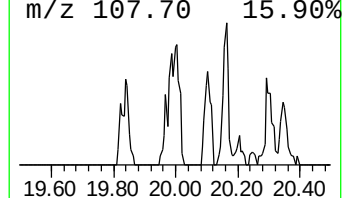
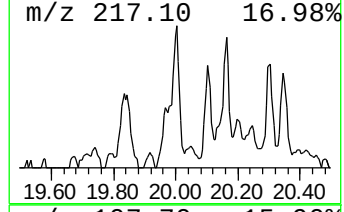
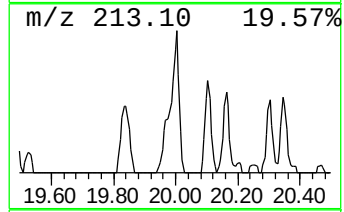
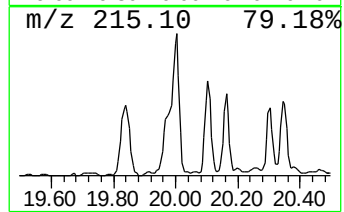
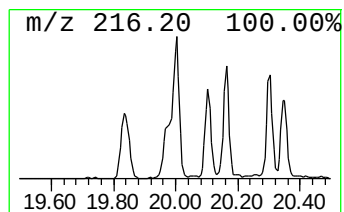
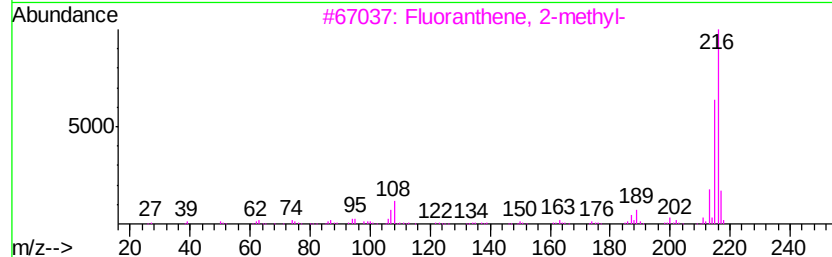
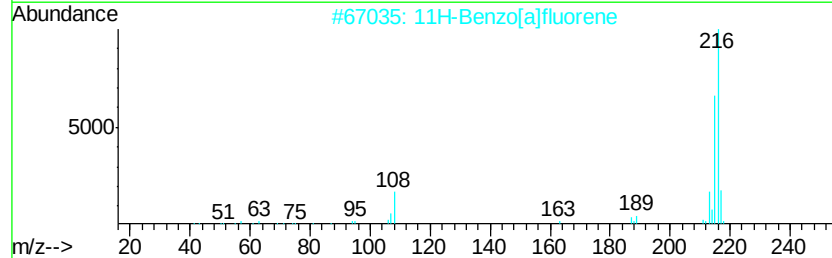
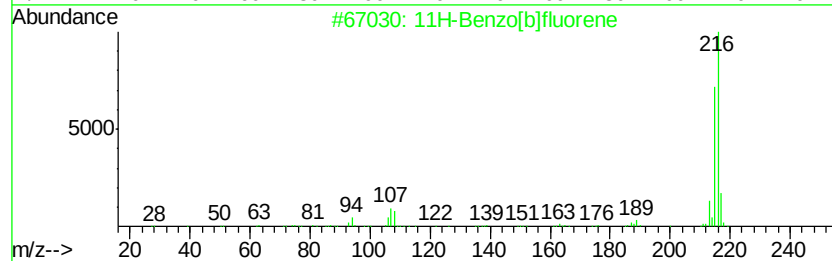
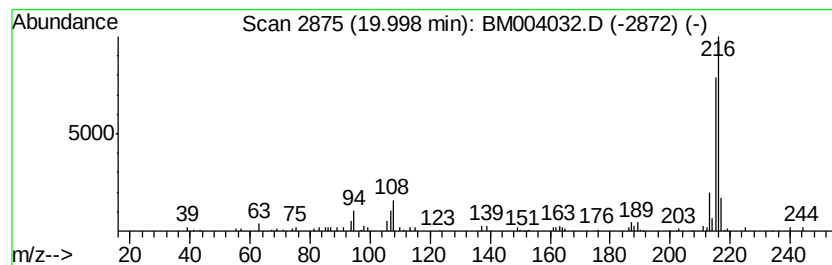
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Quant Title : SVOA CALIBRATION

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

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Peak Number 9 11H-Benzo[b]fluorene Concentration Rank 8

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.00 | 2.78 ng/ul | 164558 | Chrysene-d12     | 21.27 |

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



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM011516\  
Data File : BM004032.D  
Acq On : 15 Jan 2016 03:41  
Operator : SJ/UM  
Sample : H1099-18  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BC9E4

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010116.M  
Quant Title : SVOA CALIBRATION

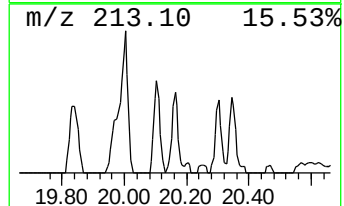
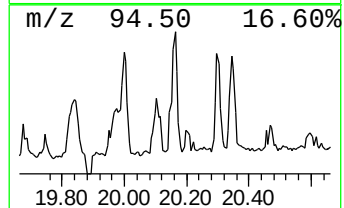
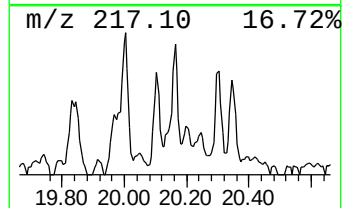
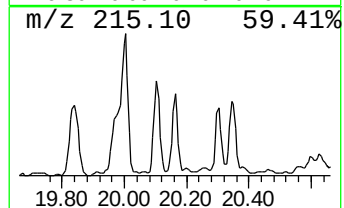
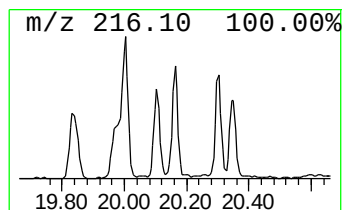
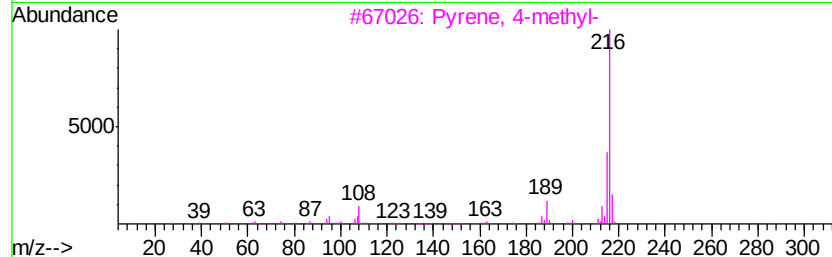
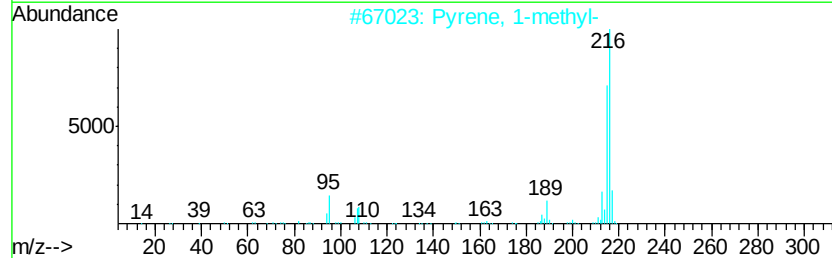
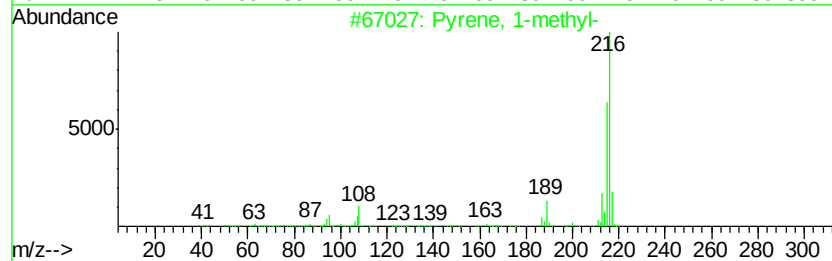
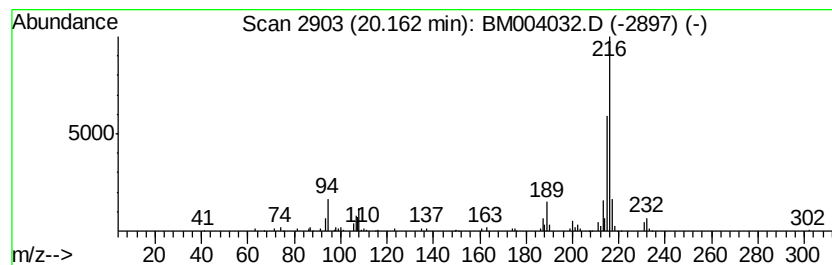
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TIC Integration Parameters: LSCINT.P

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Peak Number 10 Pyrene, 1-methyl- Concentration Rank 11

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.16 | 2.42 ng/ul | 143233 | Chrysene-d12     | 21.27 |

| Hit# | of | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------|-----|---------|-------------|------|
| 1    | 5  | Pyrene, 1-methyl- | 216 | C17H12  | 002381-21-7 | 96   |
| 2    |    | Pyrene, 1-methyl- | 216 | C17H12  | 002381-21-7 | 96   |
| 3    |    | Pyrene, 4-methyl- | 216 | C17H12  | 003353-12-6 | 95   |
| 4    |    | Pyrene, 2-methyl- | 216 | C17H12  | 003442-78-2 | 95   |
| 5    |    | Pyrene, 1-methyl- | 216 | C17H12  | 002381-21-7 | 94   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM011516\  
Data File : BM004032.D  
Acq On : 15 Jan 2016 03:41  
Operator : SJ/UM  
Sample : H1099-18  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BC9E4

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010116.M  
Quant Title : SVOA CALIBRATION

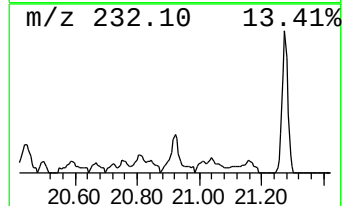
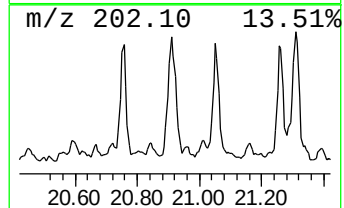
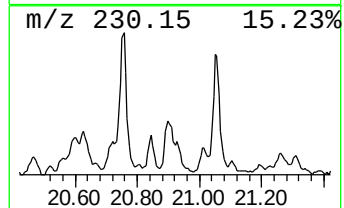
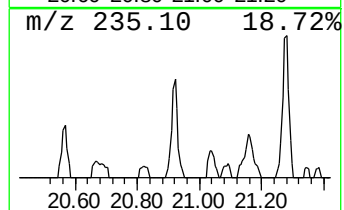
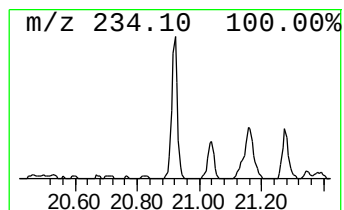
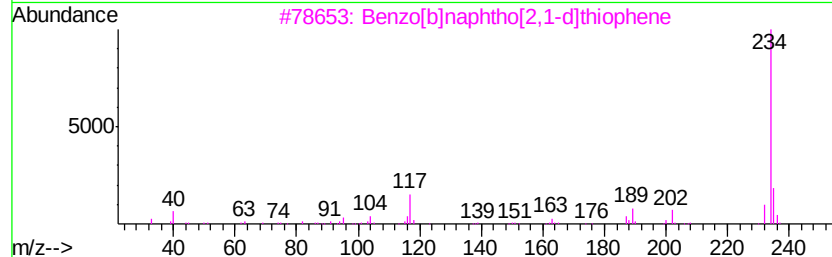
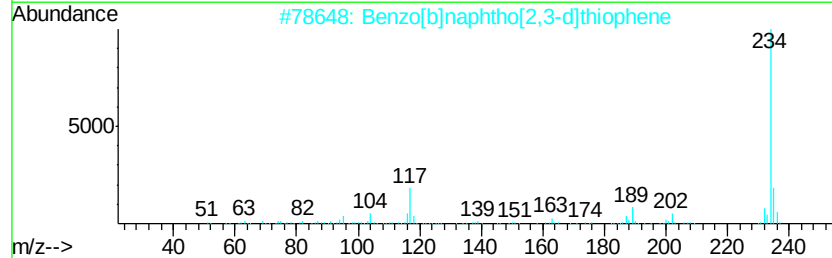
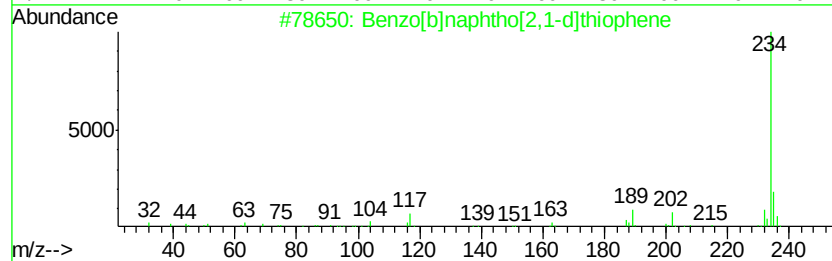
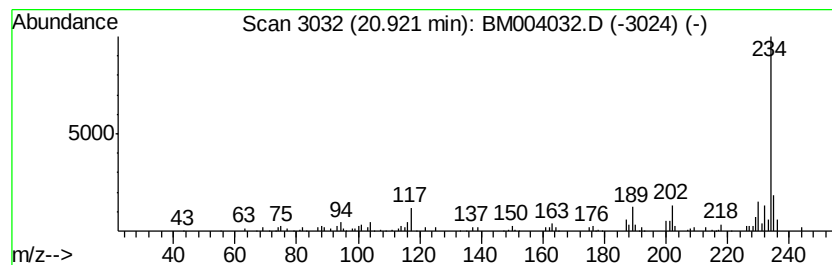
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 17

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.92 | 2.14 ng/ul | 126684 | Chrysene-d12     | 21.27 |

| Hit# | of | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzo[b]naphtho[2,1-d]thiophene | 234 | C16H10S | 000239-35-0 | 96   |
| 2    |    | Benzo[b]naphtho[2,3-d]thiophene | 234 | C16H10S | 000243-46-9 | 93   |
| 3    |    | Benzo[b]naphtho[2,1-d]thiophene | 234 | C16H10S | 000239-35-0 | 91   |
| 4    |    | Benzo[b]naphtho[2,1-d]thiophene | 234 | C16H10S | 000239-35-0 | 90   |
| 5    |    | Benzo[b]naphtho[2,3-d]thiophene | 234 | C16H10S | 000243-46-9 | 87   |





Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM011516\  
Data File : BM004032.D  
Acq On : 15 Jan 2016 03:41  
Operator : SJ/UM  
Sample : H1099-18  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BC9E4

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010116.M  
Quant Title : SVOA CALIBRATION

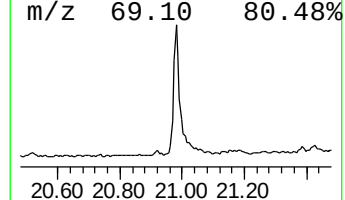
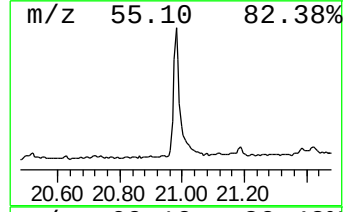
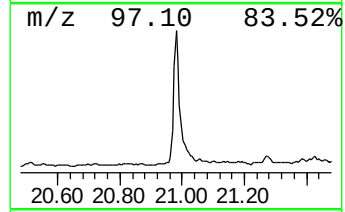
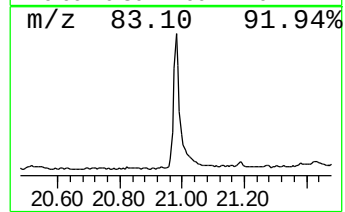
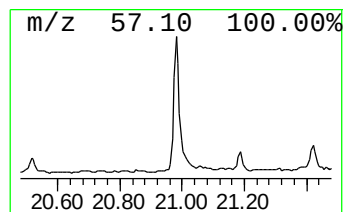
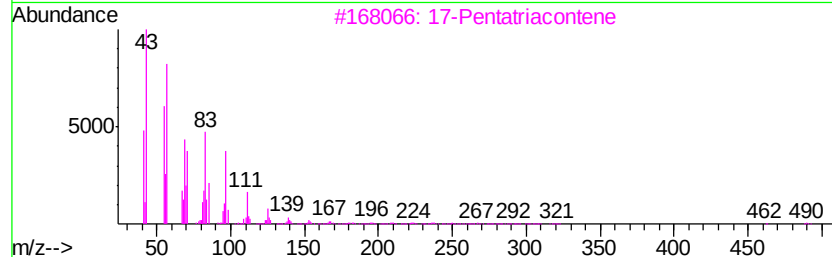
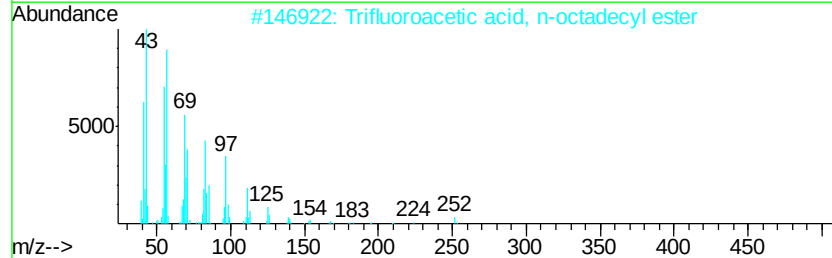
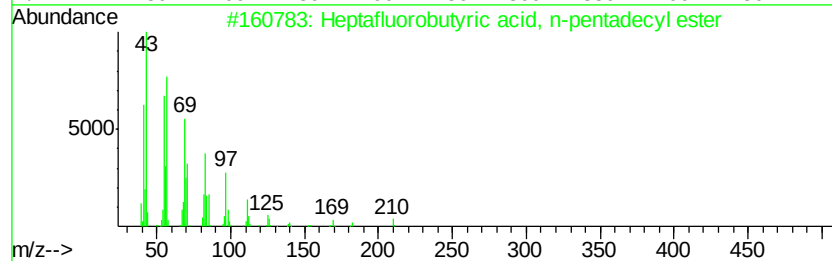
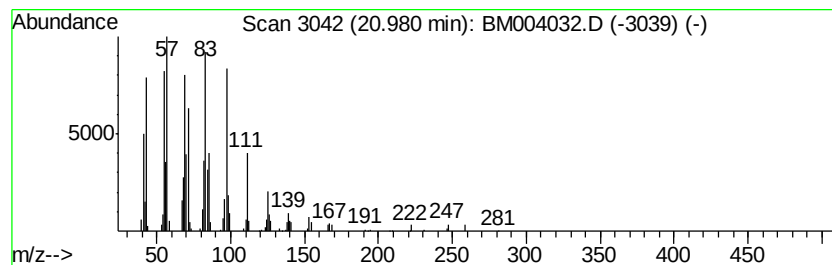
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Heptafluorobutyric acid, n-... Concentration Rank 3

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.98 | 4.73 ng/ul | 280484 | Chrysene-d12     | 21.27 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | Heptafluorobutyric acid, n-penta... | 424 | C19H31F7O2  | 1000216-79-3 | 91   |
| 2    |    | Trifluoroacetic acid, n-octadecy... | 366 | C20H37F3O2  | 079392-43-1  | 91   |
| 3    |    | 17-Pentatriacontene                 | 491 | C35H70      | 006971-40-0  | 91   |
| 4    |    | Heptafluorobutyric acid, n-octad... | 466 | C22H37F7O2  | 000400-57-7  | 91   |
| 5    |    | Trichloroacetic acid, hexadecyl ... | 386 | C18H33Cl3O2 | 074339-54-1  | 91   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM011516\  
Data File : BM004032.D  
Acq On : 15 Jan 2016 03:41  
Operator : SJ/UM  
Sample : H1099-18  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

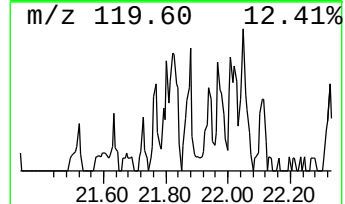
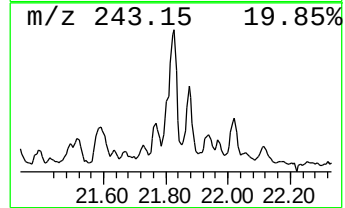
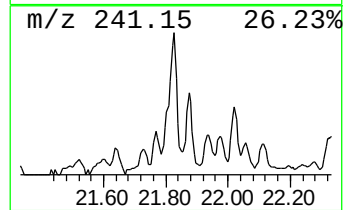
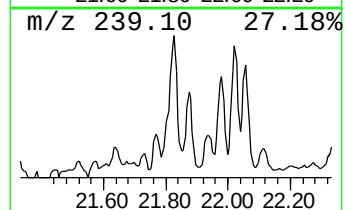
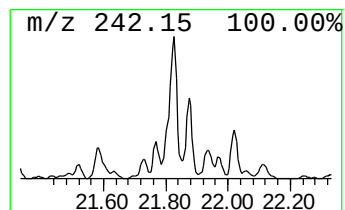
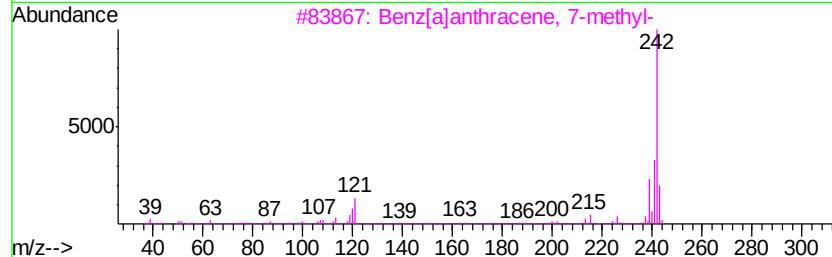
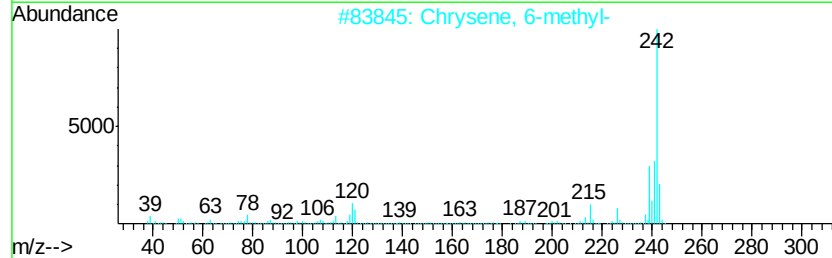
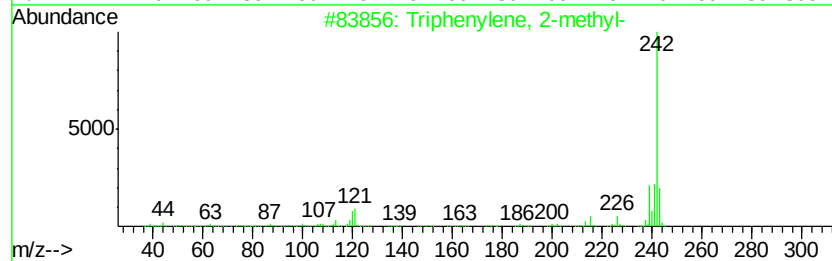
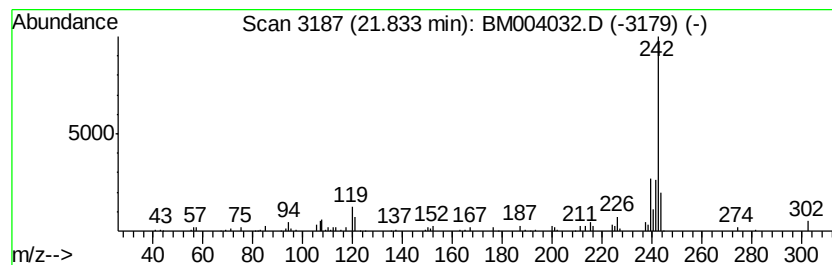
Instrument :  
BNA\_M  
ClientSampleId :  
BC9E4

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010116.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 13 Triphenylene, 2-methyl- Concentration Rank 14

| R.T.    | EstConc    | Area                            | Relative to ISTD | R.T.    |             |      |
|---------|------------|---------------------------------|------------------|---------|-------------|------|
| 21.83   | 2.32 ng/ul | 137581                          | Chrysene-d12     | 21.27   |             |      |
| Hit# of | 5          | Tentative ID                    | MW               | MolForm | CAS#        | Qual |
| 1       |            | Triphenylene, 2-methyl-         | 242              | C19H14  | 001705-84-6 | 93   |
| 2       |            | Chrysene, 6-methyl-             | 242              | C19H14  | 003697-24-3 | 90   |
| 3       |            | Benz[a]anthracene, 7-methyl-    | 242              | C19H14  | 002541-69-7 | 90   |
| 4       |            | Benz[a]anthracene, 8-methyl-    | 242              | C19H14  | 002381-31-9 | 90   |
| 5       |            | Benzo[c]phenanthrene, 6-methyl- | 242              | C19H14  | 002381-34-2 | 89   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM011516\  
Data File : BM004032.D  
Acq On : 15 Jan 2016 03:41  
Operator : SJ/UM  
Sample : H1099-18  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BC9E4

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010116.M  
Quant Title : SVOA CALIBRATION

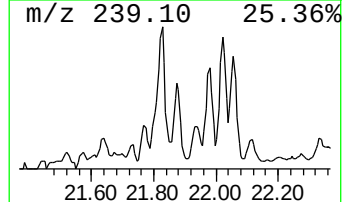
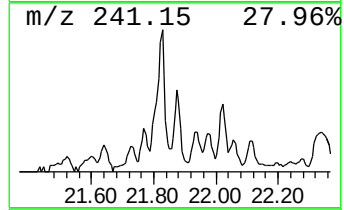
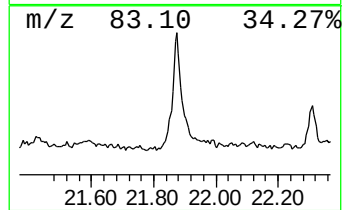
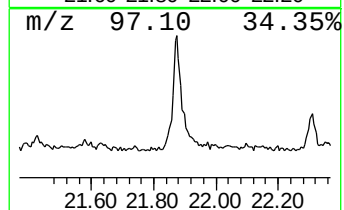
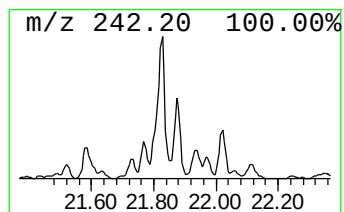
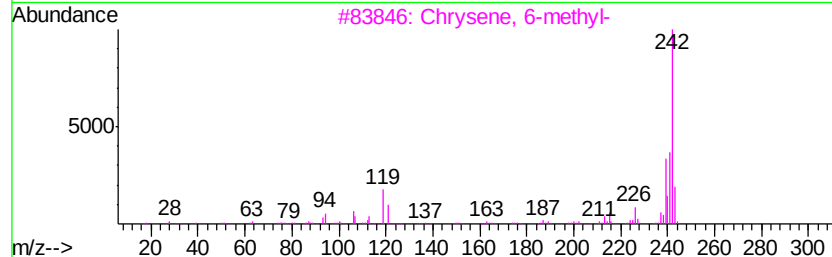
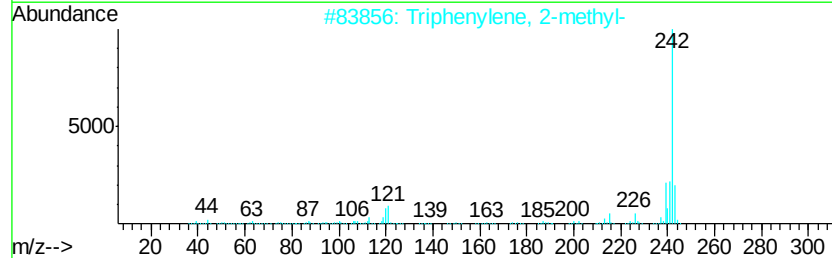
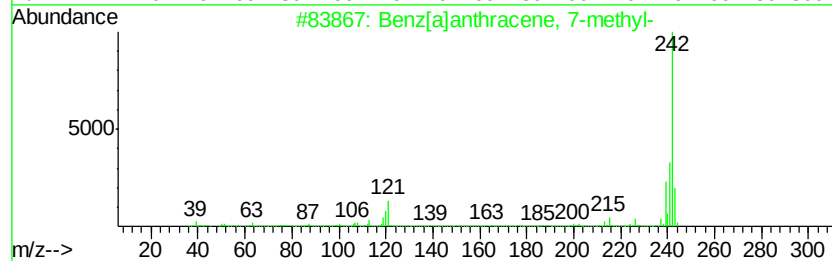
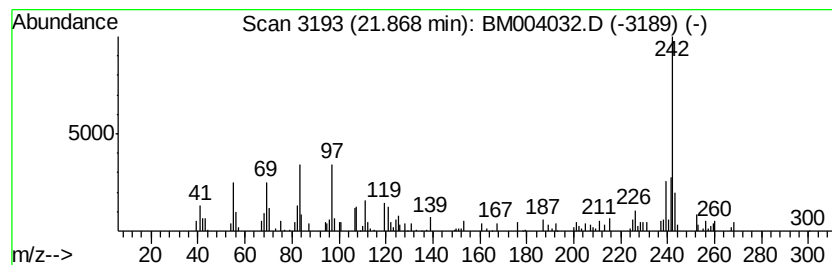
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Benz[a]anthracene, 7-methyl- Concentration Rank 9

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 21.87 | 2.63 ng/ul | 155890 | Chrysene-d12     | 21.27 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benz[a]anthracene, 7-methyl- | 242 | C19H14  | 002541-69-7 | 96   |
| 2    |    | Triphenylene, 2-methyl-      | 242 | C19H14  | 001705-84-6 | 95   |
| 3    |    | Chrysene, 6-methyl-          | 242 | C19H14  | 001705-85-7 | 91   |
| 4    |    | Chrysene, 1-methyl-          | 242 | C19H14  | 003351-28-8 | 86   |
| 5    |    | Benz[a]anthracene, 8-methyl- | 242 | C19H14  | 002381-31-9 | 86   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM011516\  
Data File : BM004032.D  
Acq On : 15 Jan 2016 03:41  
Operator : SJ/UM  
Sample : H1099-18  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BC9E4

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010116.M  
Quant Title : SVOA CALIBRATION

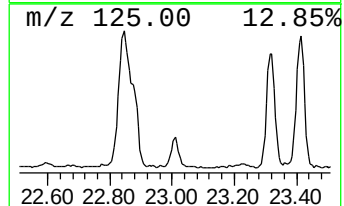
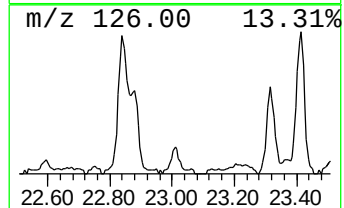
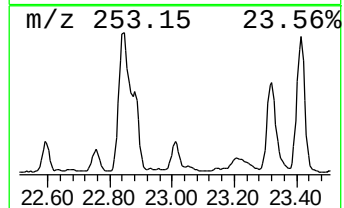
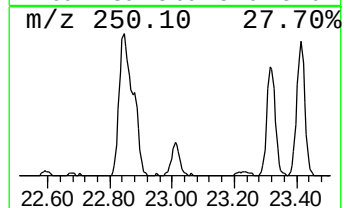
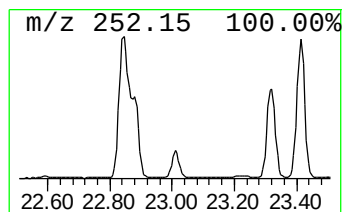
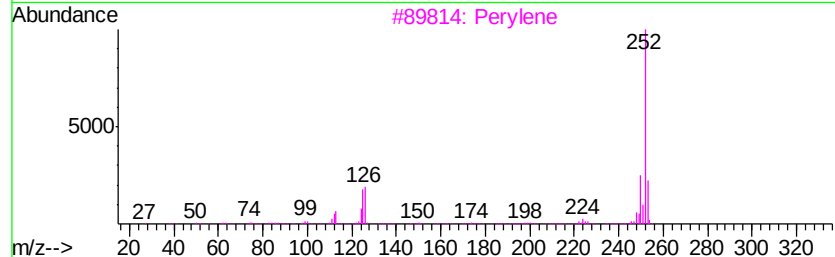
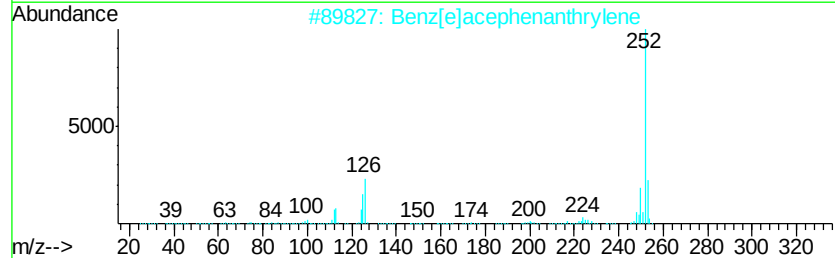
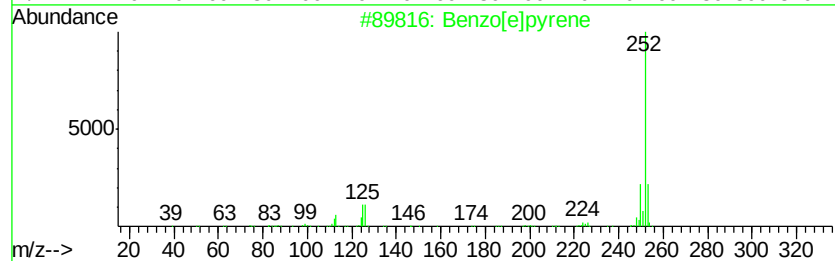
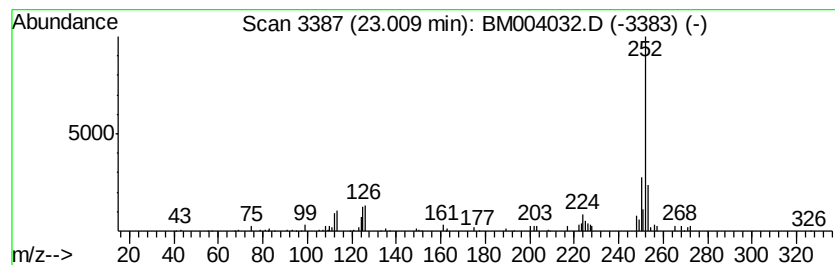
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 Benzo[e]pyrene Concentration Rank 16

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 23.01 | 2.21 ng/ul | 69993 | Perylene-d12     | 23.51 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzo[e]pyrene           | 252 | C20H12  | 000192-97-2 | 95   |
| 2    |    | Benz[e]acephenanthrylene | 252 | C20H12  | 000205-99-2 | 89   |
| 3    |    | Perylene                 | 252 | C20H12  | 000198-55-0 | 87   |
| 4    |    | Perylene                 | 252 | C20H12  | 000198-55-0 | 87   |
| 5    |    | Perylene                 | 252 | C20H12  | 000198-55-0 | 81   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM011516\  
Data File : BM004032.D  
Acq On : 15 Jan 2016 03:41  
Operator : SJ/UM  
Sample : H1099-18  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BC9E4

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010116.M  
Quant Title : SVOA CALIBRATION

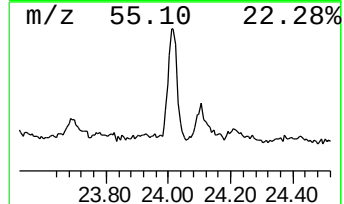
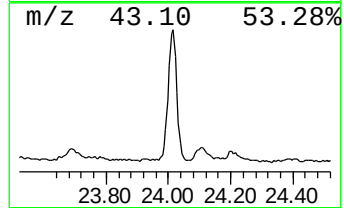
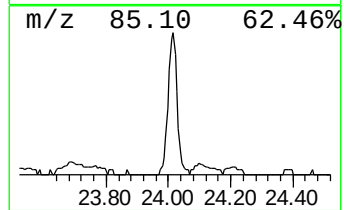
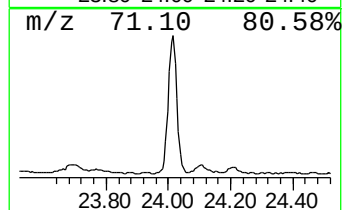
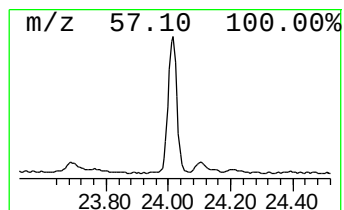
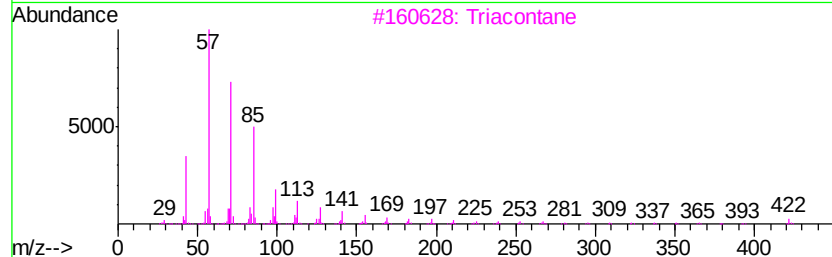
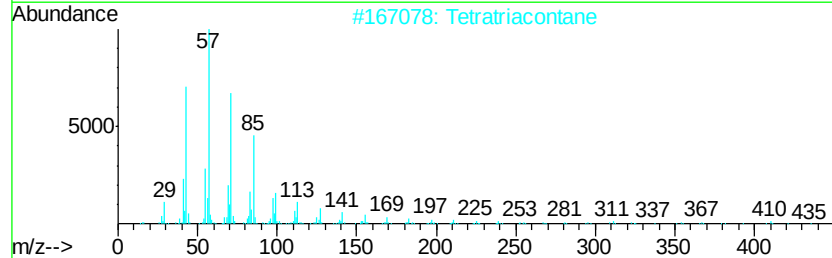
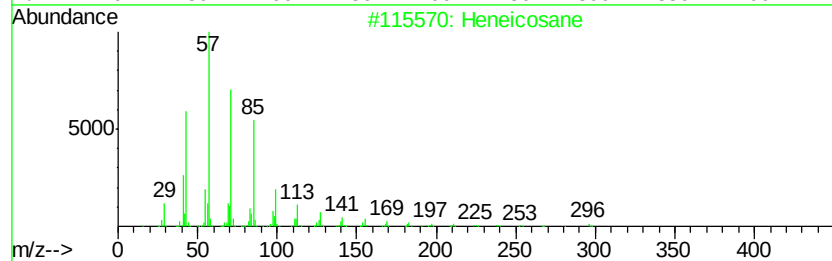
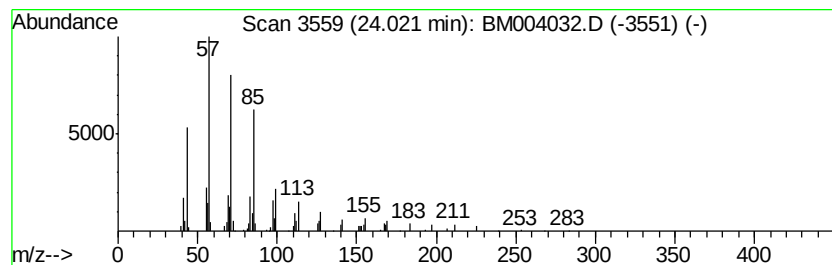
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 1

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 24.02 | 5.89 ng/ul | 186628 | Perylene-d12     | 23.51 |

| Hit# | of | Tentative ID     | MW  | MolForm | CAS#        | Qual |
|------|----|------------------|-----|---------|-------------|------|
| 1    | 5  | Heneicosane      | 296 | C21H44  | 000629-94-7 | 91   |
| 2    |    | Tetratriacontane | 479 | C34H70  | 014167-59-0 | 91   |
| 3    |    | Triacotane       | 422 | C30H62  | 000638-68-6 | 91   |
| 4    |    | Octacosane       | 394 | C28H58  | 000630-02-4 | 91   |
| 5    |    | Pentacosane      | 352 | C25H52  | 000629-99-2 | 91   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM011516\  
 Data File : BM004032.D  
 Acq On : 15 Jan 2016 03:41  
 Operator : SJ/UM  
 Sample : H1099-18  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BC9E4

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010116.M  
 Quant Title : SVOA CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| (DEL) Alkane: Str... | 16.04 | 2.2     | ng/ul | 79084    | 4                     | 17.07 | 710521  | 20.0 |
| Phenanthrene, 1-m... | 17.94 | 4.1     | ng/ul | 145529   | 4                     | 17.07 | 710521  | 20.0 |
| Phenanthrene, 2-m... | 17.99 | 4.0     | ng/ul | 143998   | 4                     | 17.07 | 710521  | 20.0 |
| 4H-Cyclopenta[def... | 18.14 | 5.6     | ng/ul | 199705   | 4                     | 17.07 | 710521  | 20.0 |
| Phenanthrene, 4-m... | 18.18 | 2.4     | ng/ul | 85030    | 4                     | 17.07 | 710521  | 20.0 |
| Phenanthrene, 2,5... | 18.87 | 3.9     | ng/ul | 137085   | 4                     | 17.07 | 710521  | 20.0 |
| Cyclopenta(def)ph... | 19.02 | 2.5     | ng/ul | 88266    | 4                     | 17.07 | 710521  | 20.0 |
| 7H-Benzo[c]fluorene  | 19.83 | 2.4     | ng/ul | 140522   | 5                     | 21.27 | 1185020 | 20.0 |
| 11H-Benzo[b]fluorene | 20.00 | 2.8     | ng/ul | 164558   | 5                     | 21.27 | 1185020 | 20.0 |
| Pyrene, 1-methyl-    | 20.16 | 2.4     | ng/ul | 143233   | 5                     | 21.27 | 1185020 | 20.0 |
| Benzo[b]naphtho[2... | 20.92 | 2.1     | ng/ul | 126684   | 5                     | 21.27 | 1185020 | 20.0 |
| Heptafluorobutyri... | 20.98 | 4.7     | ng/ul | 280484   | 5                     | 21.27 | 1185020 | 20.0 |
| Triphenylene, 2-m... | 21.83 | 2.3     | ng/ul | 137581   | 5                     | 21.27 | 1185020 | 20.0 |
| Benz[a]anthracene... | 21.87 | 2.6     | ng/ul | 155890   | 5                     | 21.27 | 1185020 | 20.0 |
| Benzo[e]pyrene       | 23.01 | 2.2     | ng/ul | 69993    | 6                     | 23.51 | 634011  | 20.0 |
| (DEL) Alkane: Str... | 24.02 | 5.9     | ng/ul | 186628   | 6                     | 23.51 | 634011  | 20.0 |