

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
 Data File : BN005345.D  
 Acq On : 27 Apr 2019 08:00  
 Operator : JU/SJ  
 Sample : K2456-12DL 5X  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 GAHQ0DL

## 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\_N\METHODS\SOM-EPA-BN041919MA.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         | 5.634       | 442           | 450         | 461          | rBV      | 479046         | 750210        | 4.04%           | 0.276%        |
| 2         | 6.775       | 636           | 644         | 652          | rVV      | 253863         | 438580        | 2.36%           | 0.161%        |
| 3         | 7.134       | 696           | 705         | 712          | rVB      | 247805         | 402871        | 2.17%           | 0.148%        |
| 4         | 7.263       | 718           | 727         | 735          | rBV      | 735880         | 1361606       | 7.34%           | 0.500%        |
| 5         | 7.340       | 735           | 740         | 747          | rVB      | 310635         | 479266        | 2.58%           | 0.176%        |
| 6         | 7.593       | 776           | 783         | 791          | rBV      | 1973448        | 3161082       | 17.04%          | 1.162%        |
| 7         | 8.293       | 894           | 902         | 911          | rVB      | 293630         | 518856        | 2.80%           | 0.191%        |
| 8         | 8.734       | 972           | 977         | 984          | rVV      | 231569         | 416821        | 2.25%           | 0.153%        |
| 9         | 8.851       | 991           | 997         | 1009         | rVV      | 363092         | 753095        | 4.06%           | 0.277%        |
| 10        | 9.340       | 1075          | 1080        | 1085         | rBV      | 248219         | 395775        | 2.13%           | 0.145%        |
| 11        | 9.940       | 1173          | 1182        | 1186         | rVV      | 677312         | 1275272       | 6.88%           | 0.469%        |
| 12        | 9.981       | 1186          | 1189        | 1198         | rVB      | 300053         | 497397        | 2.68%           | 0.183%        |
| 13        | 10.357      | 1245          | 1253        | 1271         | rVB      | 2712293        | 4759763       | 25.66%          | 1.749%        |
| 14        | 11.398      | 1425          | 1430        | 1434         | rBV3     | 248735         | 454489        | 2.45%           | 0.167%        |
| 15        | 11.551      | 1450          | 1456        | 1458         | rBV3     | 368334         | 636686        | 3.43%           | 0.234%        |
| 16        | 11.687      | 1469          | 1479        | 1484         | rBV2     | 316294         | 782245        | 4.22%           | 0.288%        |
| 17        | 11.881      | 1504          | 1512        | 1519         | rVB      | 715419         | 1988615       | 10.72%          | 0.731%        |
| 18        | 12.022      | 1531          | 1536        | 1545         | rVB2     | 325730         | 670626        | 3.62%           | 0.246%        |
| 19        | 12.657      | 1640          | 1644        | 1650         | rVB3     | 233119         | 390464        | 2.11%           | 0.144%        |
| 20        | 12.775      | 1660          | 1664        | 1668         | rVB      | 308401         | 404931        | 2.18%           | 0.149%        |
| 21        | 13.069      | 1711          | 1714        | 1722         | rVB      | 334015         | 523844        | 2.82%           | 0.193%        |
| 22        | 13.198      | 1731          | 1736        | 1741         | rBV      | 450904         | 823715        | 4.44%           | 0.303%        |
| 23        | 13.463      | 1776          | 1781        | 1785         | rBV      | 569364         | 903637        | 4.87%           | 0.332%        |
| 24        | 13.545      | 1791          | 1795        | 1800         | rVB      | 286698         | 431865        | 2.33%           | 0.159%        |
| 25        | 13.645      | 1808          | 1812        | 1814         | rBV2     | 896171         | 1228421       | 6.62%           | 0.452%        |
| 26        | 13.681      | 1814          | 1818        | 1824         | rVV      | 4138466        | 6147465       | 33.14%          | 2.260%        |
| 27        | 13.945      | 1854          | 1863        | 1870         | rBV2     | 4435622        | 7714319       | 41.59%          | 2.835%        |
| 28        | 14.157      | 1894          | 1899        | 1902         | rBV      | 518888         | 724955        | 3.91%           | 0.266%        |
| 29        | 14.228      | 1903          | 1911        | 1918         | rBV2     | 3696724        | 6035575       | 32.54%          | 2.218%        |
| 30        | 14.439      | 1943          | 1947        | 1957         | rVB3     | 321743         | 672406        | 3.63%           | 0.247%        |
| 31        | 14.710      | 1991          | 1993        | 1998         | rVB      | 326336         | 399223        | 2.15%           | 0.147%        |
| 32        | 14.775      | 1999          | 2004        | 2008         | rBV2     | 391355         | 674696        | 3.64%           | 0.248%        |
| 33        | 14.839      | 2009          | 2015        | 2020         | rBV      | 2356218        | 3869639       | 20.86%          | 1.422%        |
| 34        | 14.886      | 2020          | 2023        | 2029         | rVV2     | 610537         | 976118        | 5.26%           | 0.359%        |

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
 Data File : BN005345.D  
 Acq On : 27 Apr 2019 08:00  
 Operator : JU/SJ  
 Sample : K2456-12DL 5X  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 GAHQ0DL

## 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\_N\METHODS\SOM-EPA-BN041919MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 14.939 | 2029 | 2032 | 2038 | rVB  | 627150  | 885365  | 4.77%  | 0.325% |
| 36 | 15.004 | 2038 | 2043 | 2051 | rBV2 | 603901  | 1213366 | 6.54%  | 0.446% |
| 37 | 15.110 | 2056 | 2061 | 2066 | rVB2 | 1587833 | 2350114 | 12.67% | 0.864% |
| 38 | 15.228 | 2077 | 2081 | 2086 | rVB  | 1424645 | 1958737 | 10.56% | 0.720% |
| 39 | 15.304 | 2088 | 2094 | 2102 | rVV  | 579055  | 1194822 | 6.44%  | 0.439% |
| 40 | 15.410 | 2108 | 2112 | 2115 | rVB2 | 481769  | 558995  | 3.01%  | 0.205% |
| 41 | 15.510 | 2121 | 2129 | 2136 | rVB3 | 842762  | 2204612 | 11.89% | 0.810% |
| 42 | 15.898 | 2191 | 2195 | 2201 | rBV3 | 263148  | 440093  | 2.37%  | 0.162% |
| 43 | 15.975 | 2202 | 2208 | 2211 | rVV4 | 1044738 | 2033184 | 10.96% | 0.747% |
| 44 | 16.004 | 2211 | 2213 | 2217 | rVV  | 583788  | 658827  | 3.55%  | 0.242% |
| 45 | 16.280 | 2257 | 2260 | 2264 | rBV3 | 246759  | 436700  | 2.35%  | 0.161% |
| 46 | 16.345 | 2268 | 2271 | 2275 | rBV3 | 296396  | 464172  | 2.50%  | 0.171% |
| 47 | 16.563 | 2304 | 2308 | 2313 | rVB  | 1215730 | 1748701 | 9.43%  | 0.643% |
| 48 | 16.775 | 2341 | 2344 | 2347 | rVB  | 535024  | 560490  | 3.02%  | 0.206% |
| 49 | 16.828 | 2350 | 2353 | 2358 | rVB3 | 477094  | 705764  | 3.81%  | 0.259% |
| 50 | 16.986 | 2374 | 2380 | 2384 | rBV  | 4754334 | 6934125 | 37.39% | 2.549% |
| 51 | 17.028 | 2384 | 2387 | 2391 | rVV  | 1546366 | 2044988 | 11.03% | 0.752% |
| 52 | 17.080 | 2392 | 2396 | 2398 | rVV2 | 825587  | 1232748 | 6.65%  | 0.453% |
| 53 | 17.116 | 2398 | 2402 | 2407 | rVB  | 2977554 | 4217731 | 22.74% | 1.550% |
| 54 | 17.510 | 2466 | 2469 | 2474 | rVB  | 1149796 | 1457401 | 7.86%  | 0.536% |
| 55 | 17.863 | 2525 | 2529 | 2533 | rBV  | 1705764 | 2317841 | 12.50% | 0.852% |
| 56 | 17.910 | 2533 | 2537 | 2545 | rVB  | 2364043 | 3458893 | 18.65% | 1.271% |
| 57 | 17.992 | 2547 | 2551 | 2556 | rVB  | 1247644 | 1624635 | 8.76%  | 0.597% |
| 58 | 18.057 | 2557 | 2562 | 2567 | rBV3 | 2409535 | 4849382 | 26.15% | 1.782% |
| 59 | 18.186 | 2582 | 2584 | 2586 | rBV  | 369100  | 402880  | 2.17%  | 0.148% |
| 60 | 18.369 | 2613 | 2615 | 2619 | rVB  | 965504  | 1191053 | 6.42%  | 0.438% |
| 61 | 18.622 | 2655 | 2658 | 2662 | rBV  | 877928  | 1167817 | 6.30%  | 0.429% |
| 62 | 18.674 | 2664 | 2667 | 2671 | rVV  | 1175601 | 1533388 | 8.27%  | 0.564% |
| 63 | 18.716 | 2671 | 2674 | 2678 | rVB  | 733644  | 929890  | 5.01%  | 0.342% |
| 64 | 18.798 | 2683 | 2688 | 2693 | rBV2 | 2902985 | 4533250 | 24.44% | 1.666% |
| 65 | 18.857 | 2693 | 2698 | 2701 | rBV3 | 1726729 | 2784268 | 15.01% | 1.023% |
| 66 | 18.892 | 2701 | 2704 | 2707 | rVB  | 1089761 | 1318114 | 7.11%  | 0.484% |
| 67 | 18.939 | 2707 | 2712 | 2719 | rBV3 | 2273745 | 4243867 | 22.88% | 1.560% |
| 68 | 19.063 | 2728 | 2733 | 2738 | rBV2 | 5703860 | 7942645 | 42.82% | 2.919% |
| 69 | 19.133 | 2743 | 2745 | 2749 | rVB  | 853233  | 988475  | 5.33%  | 0.363% |
| 70 | 19.216 | 2755 | 2759 | 2763 | rVB  | 1411536 | 1855889 | 10.01% | 0.682% |
| 71 | 19.333 | 2775 | 2779 | 2786 | rVB  | 1599421 | 2730943 | 14.72% | 1.004% |

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
 Data File : BN005345.D  
 Acq On : 27 Apr 2019 08:00  
 Operator : JU/SJ  
 Sample : K2456-12DL 5X  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 GAHQ0DL

## 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\_N\METHODS\SOM-EPA-BN041919MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |          |          |         |        |
|-----|--------|------|------|------|------|----------|----------|---------|--------|
| 72  | 19.433 | 2787 | 2796 | 2804 | rBV  | 11399450 | 18547196 | 100.00% | 6.817% |
| 73  | 19.763 | 2848 | 2852 | 2863 | rVB4 | 2568247  | 5523845  | 29.78%  | 2.030% |
| 74  | 19.916 | 2871 | 2878 | 2880 | rBV4 | 2402879  | 5765357  | 31.08%  | 2.119% |
| 75  | 19.986 | 2887 | 2890 | 2894 | rBV3 | 702078   | 1011512  | 5.45%   | 0.372% |
| 76  | 20.033 | 2895 | 2898 | 2904 | rVV  | 1556280  | 2129841  | 11.48%  | 0.783% |
| 77  | 20.098 | 2904 | 2909 | 2913 | rVV  | 3331883  | 4194131  | 22.61%  | 1.542% |
| 78  | 20.239 | 2929 | 2933 | 2936 | rBV  | 3233730  | 4144533  | 22.35%  | 1.523% |
| 79  | 20.280 | 2936 | 2940 | 2946 | rVB  | 3062422  | 4037602  | 21.77%  | 1.484% |
| 80  | 20.492 | 2974 | 2976 | 2979 | rBV3 | 544872   | 514957   | 2.78%   | 0.189% |
| 81  | 20.692 | 3007 | 3010 | 3017 | rVB2 | 1835168  | 3375640  | 18.20%  | 1.241% |
| 82  | 20.833 | 3032 | 3034 | 3038 | rBV4 | 629246   | 915764   | 4.94%   | 0.337% |
| 83  | 20.898 | 3042 | 3045 | 3048 | rVB  | 1379808  | 1435100  | 7.74%   | 0.527% |
| 84  | 20.933 | 3049 | 3051 | 3054 | rBV  | 656445   | 874406   | 4.71%   | 0.321% |
| 85  | 20.992 | 3059 | 3061 | 3069 | rVB2 | 960964   | 1397480  | 7.53%   | 0.514% |
| 86  | 21.210 | 3093 | 3098 | 3104 | rBV2 | 7508378  | 17047275 | 91.91%  | 6.266% |
| 87  | 21.257 | 3104 | 3106 | 3113 | rVB2 | 5911952  | 8124782  | 43.81%  | 2.986% |
| 88  | 21.427 | 3132 | 3135 | 3140 | rVB  | 2306323  | 2796155  | 15.08%  | 1.028% |
| 89  | 21.721 | 3181 | 3185 | 3188 | rBV2 | 1760614  | 2697105  | 14.54%  | 0.991% |
| 90  | 21.774 | 3188 | 3194 | 3198 | rVV  | 3851813  | 6761477  | 36.46%  | 2.485% |
| 91  | 21.968 | 3223 | 3227 | 3229 | rBV3 | 1574623  | 2015306  | 10.87%  | 0.741% |
| 92  | 21.998 | 3229 | 3232 | 3237 | rVB2 | 2099219  | 2734116  | 14.74%  | 1.005% |
| 93  | 22.068 | 3242 | 3244 | 3254 | rVB2 | 1394311  | 2315941  | 12.49%  | 0.851% |
| 94  | 22.780 | 3360 | 3365 | 3383 | rVB2 | 4276446  | 14577157 | 78.59%  | 5.358% |
| 95  | 22.945 | 3389 | 3393 | 3401 | rVB  | 1536639  | 2418317  | 13.04%  | 0.889% |
| 96  | 23.245 | 3439 | 3444 | 3449 | rBV  | 2727291  | 4484971  | 24.18%  | 1.648% |
| 97  | 23.339 | 3454 | 3460 | 3466 | rVB  | 6013557  | 10380467 | 55.97%  | 3.815% |
| 98  | 23.433 | 3471 | 3476 | 3480 | rVV  | 2708521  | 4937846  | 26.62%  | 1.815% |
| 99  | 23.480 | 3481 | 3484 | 3492 | rVB4 | 1183584  | 2731296  | 14.73%  | 1.004% |
| 100 | 25.615 | 3842 | 3847 | 3858 | rVB2 | 1872076  | 5337819  | 28.78%  | 1.962% |

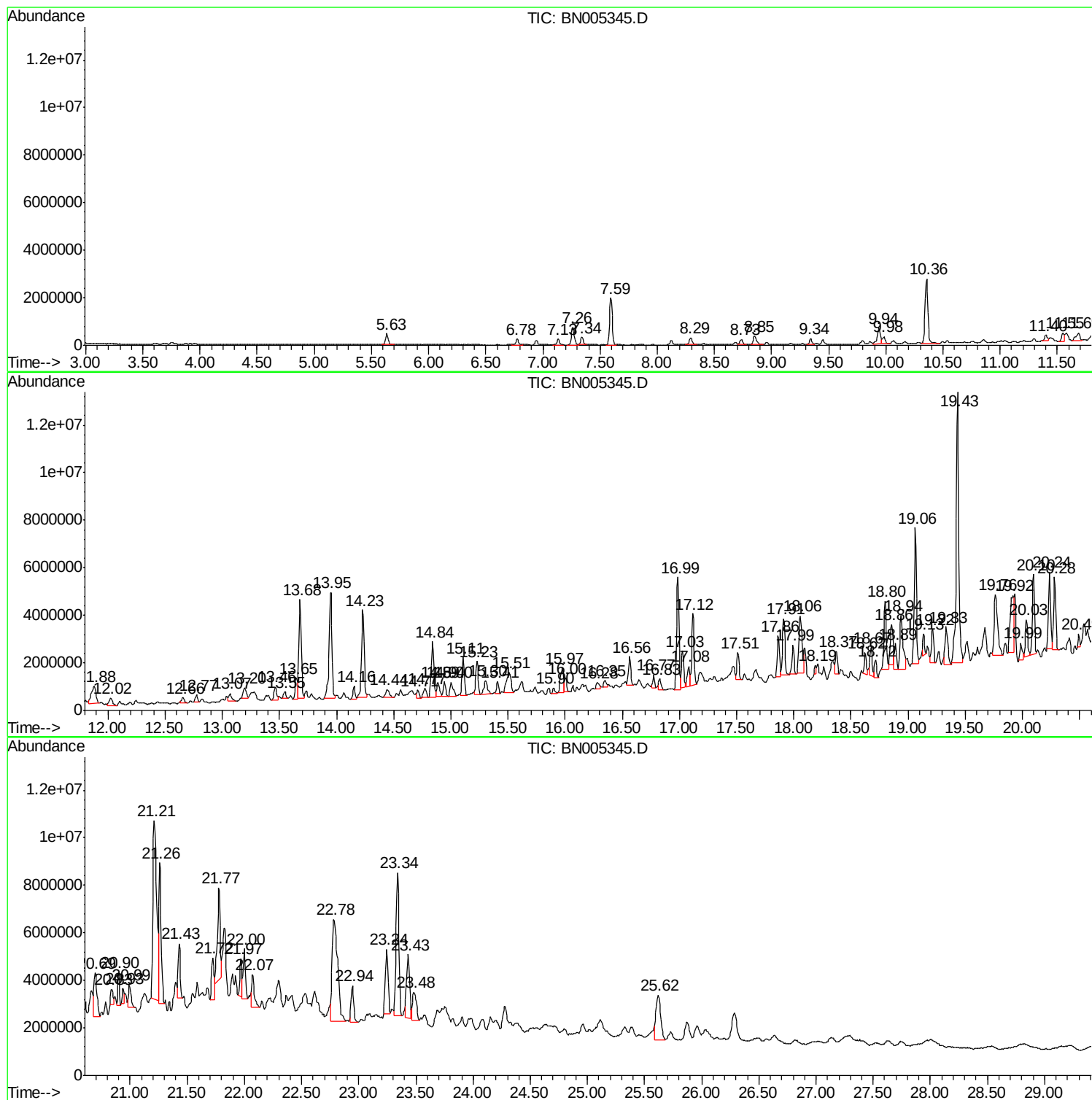
Sum of corrected areas: 272066087

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

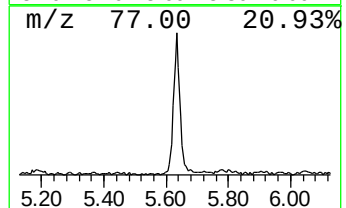
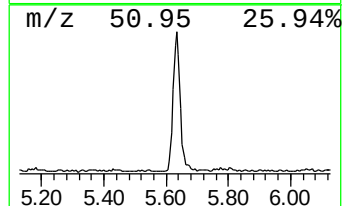
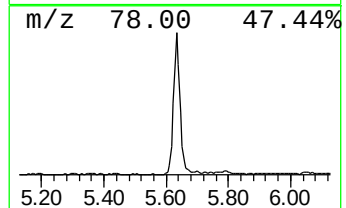
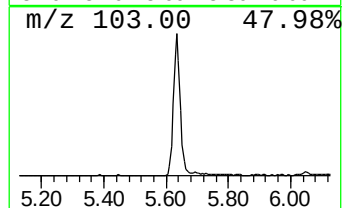
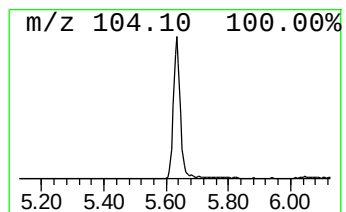
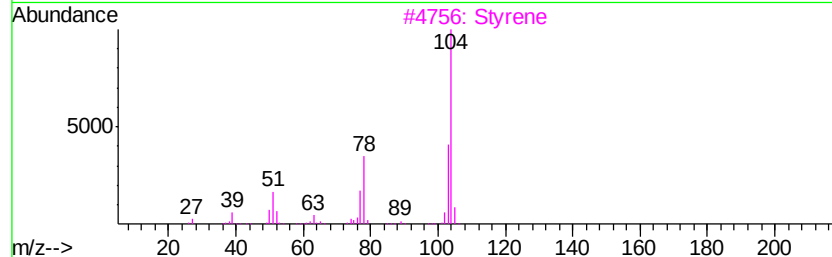
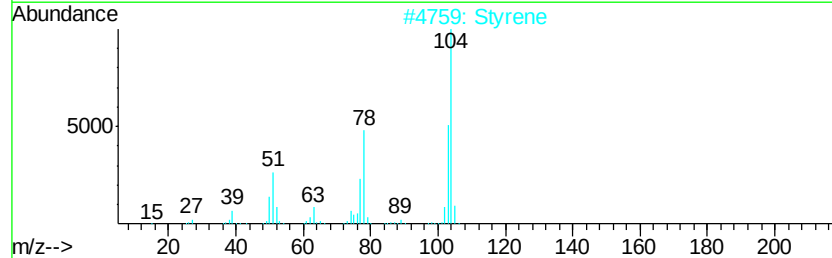
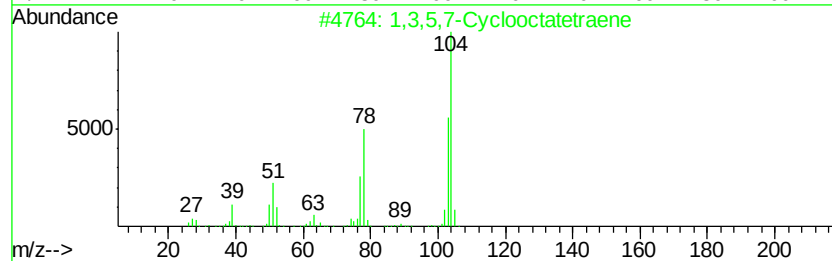
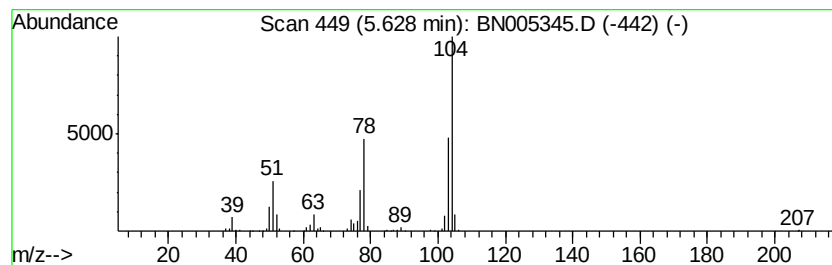
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 1,3,5,7-Cyclooctatetraene Concentration Rank 21

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 5.63 | 4.75 ng/ul | 750210 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1,3,5,7-Cyclooctatetraene       | 104 | C8H8    | 000629-20-9 | 96   |
| 2    |    |   | Styrene                         | 104 | C8H8    | 000100-42-5 | 96   |
| 3    |    |   | Styrene                         | 104 | C8H8    | 000100-42-5 | 95   |
| 4    |    |   | Bicyclo[4.2.0]octa-1,3,5-triene | 104 | C8H8    | 000694-87-1 | 95   |
| 5    |    |   | Styrene                         | 104 | C8H8    | 000100-42-5 | 95   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

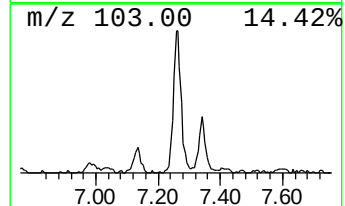
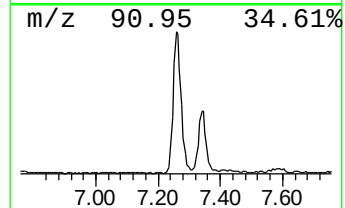
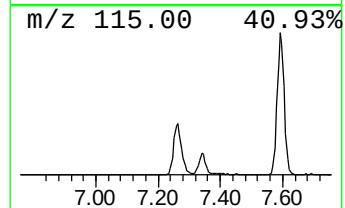
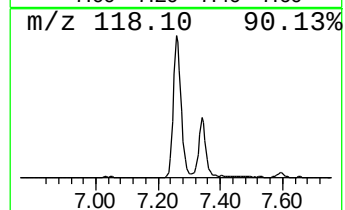
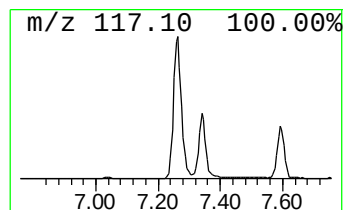
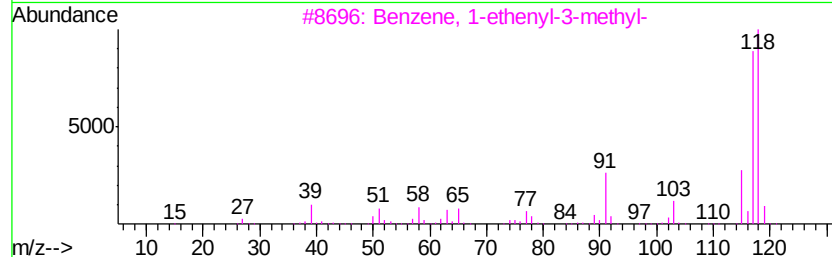
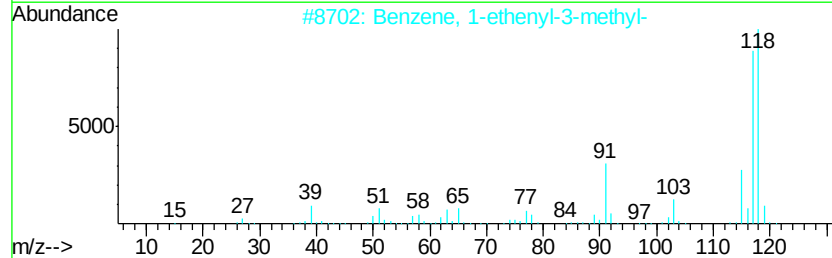
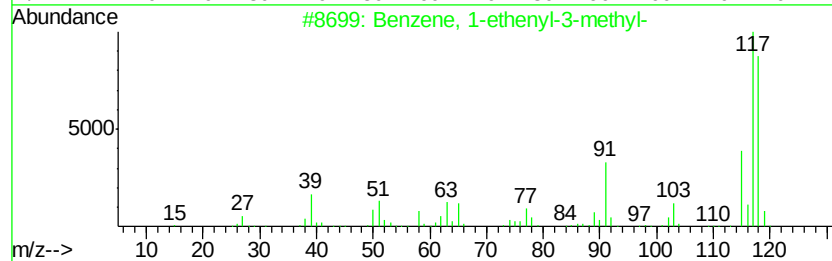
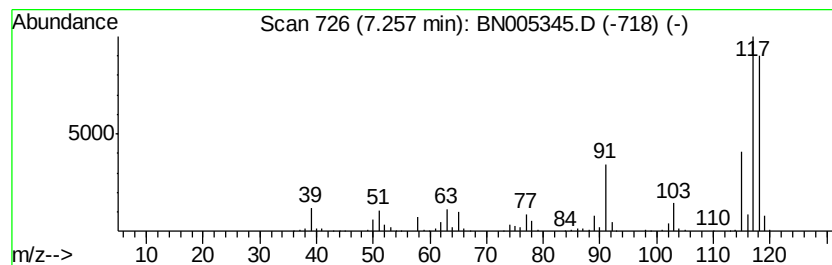
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 Benzene, 1-ethenyl-3-methyl- Concentration Rank 7

| R.T. | EstConc    | Area    | Relative to ISTD       | R.T. |
|------|------------|---------|------------------------|------|
| 7.26 | 8.61 ng/ul | 1361610 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethenyl-3-methyl- | 118 | C9H10   | 000100-80-1 | 96   |
| 2    |    | Benzene, 1-ethenyl-3-methyl- | 118 | C9H10   | 000100-80-1 | 95   |
| 3    |    | Benzene, 1-ethenyl-3-methyl- | 118 | C9H10   | 000100-80-1 | 95   |
| 4    |    | Benzene, 1-propenyl-         | 118 | C9H10   | 000637-50-3 | 94   |
| 5    |    | Benzene, cyclopropyl-        | 118 | C9H10   | 000873-49-4 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

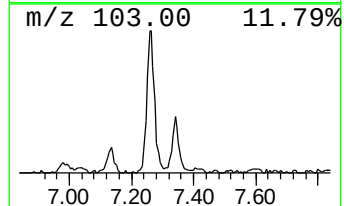
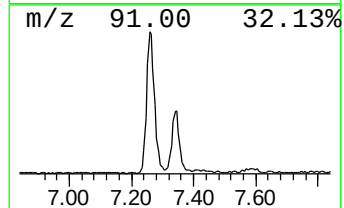
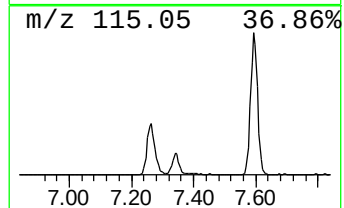
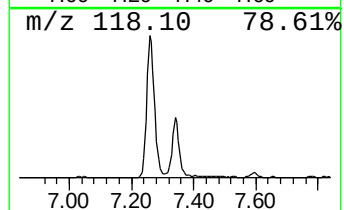
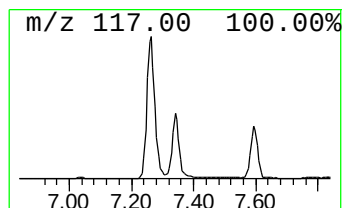
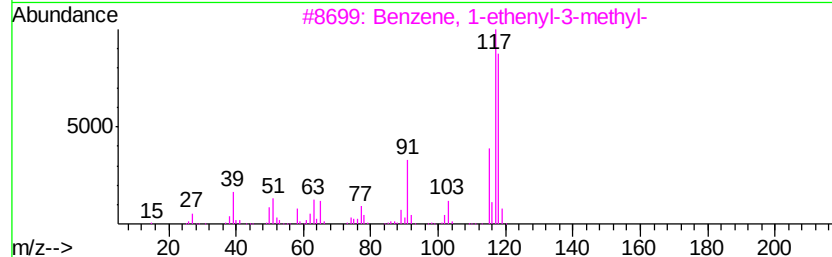
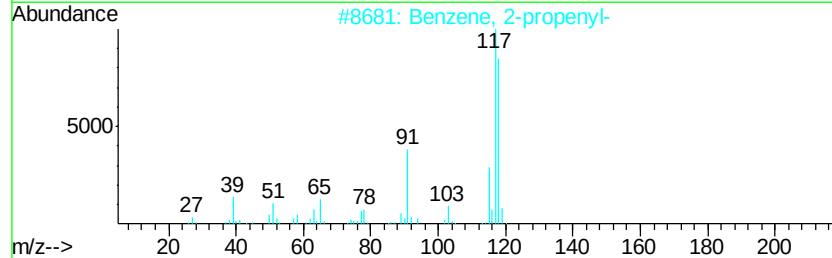
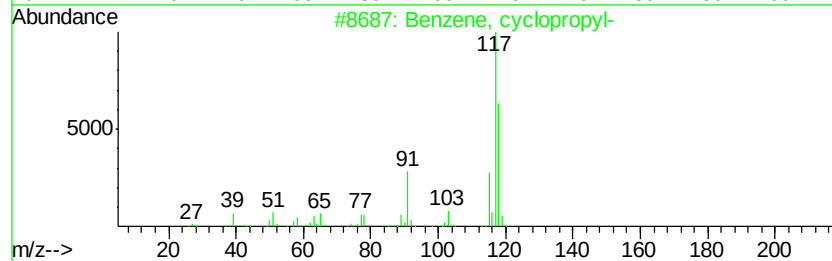
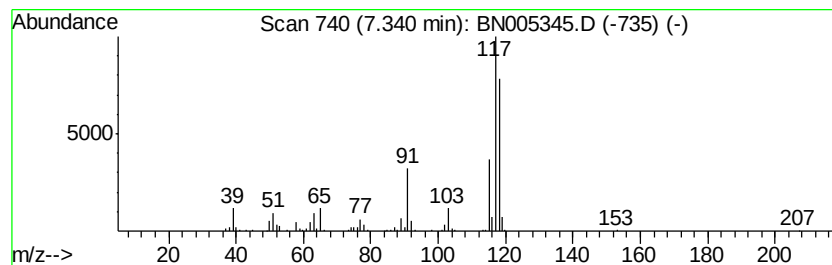
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Benzene, cyclopropyl- Concentration Rank 37

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.34 | 3.03 ng/ul | 479266 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, cyclopropyl-        | 118 | C9H10   | 000873-49-4 | 94   |
| 2    |    | Benzene, 2-propenyl-         | 118 | C9H10   | 000300-57-2 | 94   |
| 3    |    | Benzene, 1-ethenyl-3-methyl- | 118 | C9H10   | 000100-80-1 | 93   |
| 4    |    | Benzene, 1-ethenyl-2-methyl- | 118 | C9H10   | 000611-15-4 | 93   |
| 5    |    | Benzene, 1-ethenyl-3-methyl- | 118 | C9H10   | 000100-80-1 | 92   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

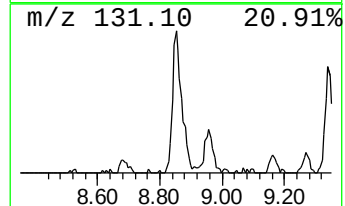
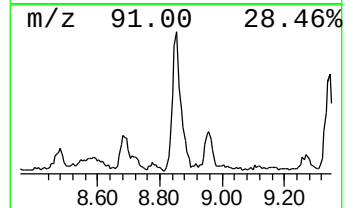
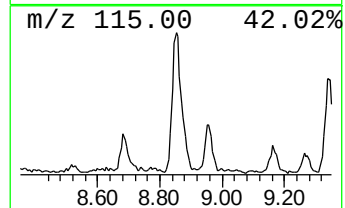
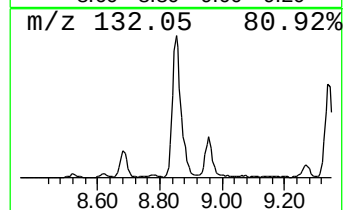
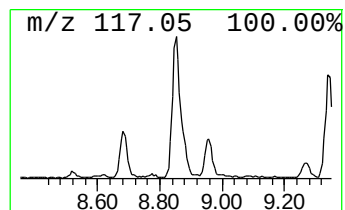
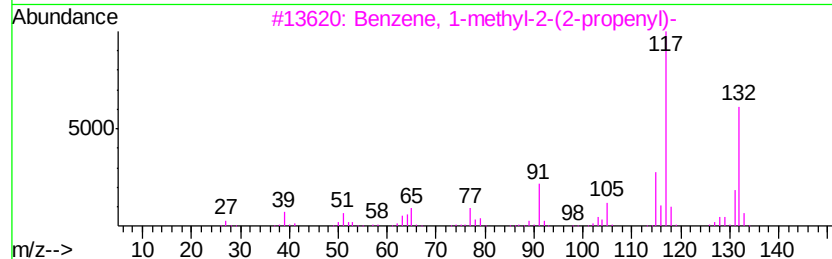
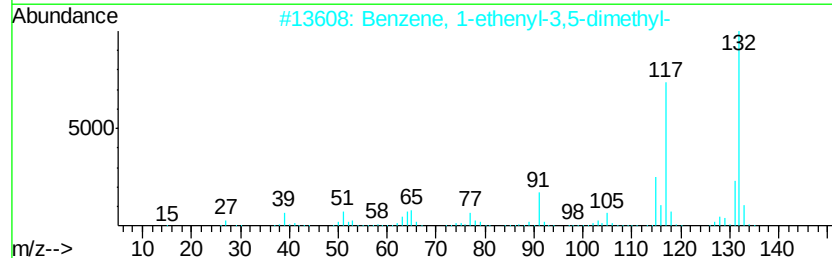
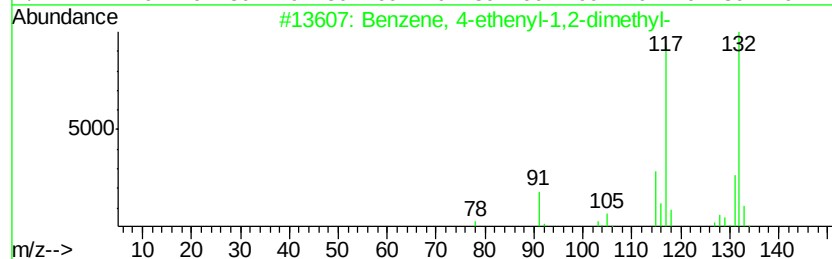
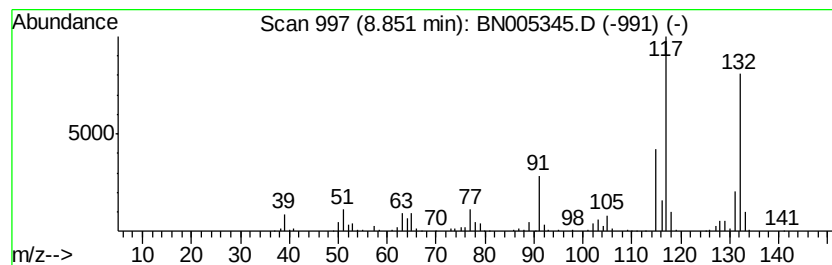
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 Benzene, 4-ethenyl-1,2-dime... Concentration Rank 20

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.85 | 4.76 ng/ul | 753095 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# | of | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 4-ethenyl-1,2-dimethyl-  | 132 | C10H12  | 027831-13-6 | 97   |
| 2    |    | Benzene, 1-ethenyl-3,5-dimethyl-  | 132 | C10H12  | 005379-20-4 | 96   |
| 3    |    | Benzene, 1-methyl-2-(2-propenyl)- | 132 | C10H12  | 001587-04-8 | 95   |
| 4    |    | Benzene, (2-methyl-2-propenyl)-   | 132 | C10H12  | 003290-53-7 | 95   |
| 5    |    | Benzene, (2-methyl-1-propenyl)-   | 132 | C10H12  | 000768-49-0 | 95   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

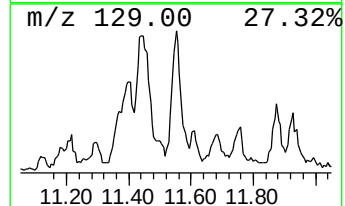
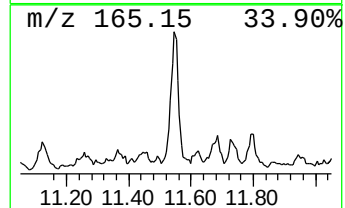
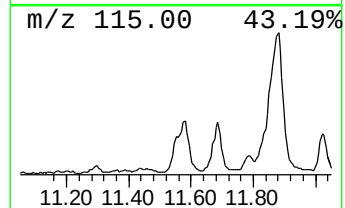
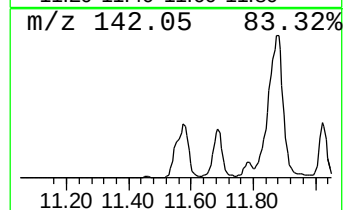
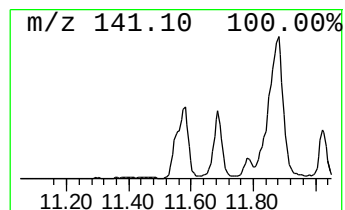
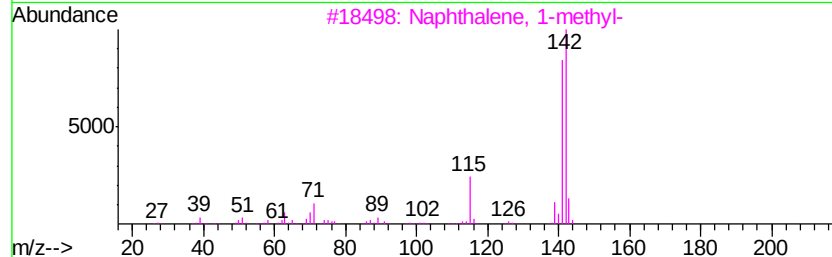
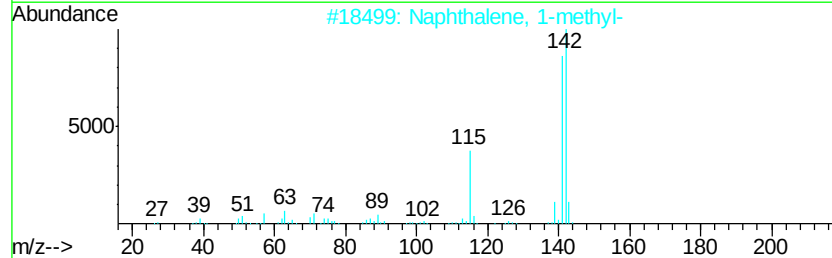
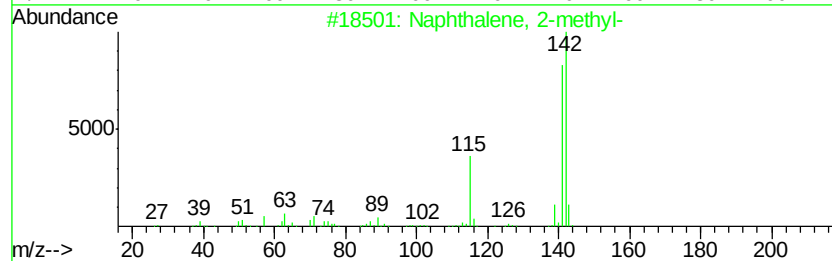
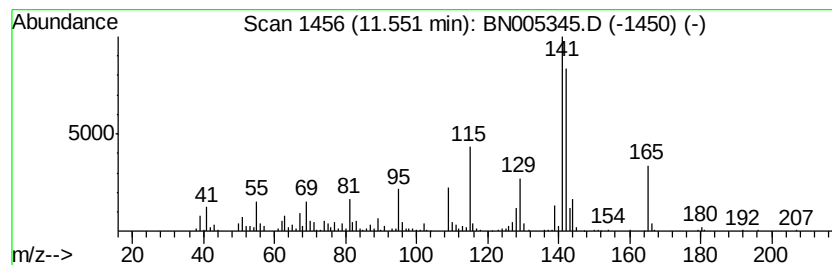
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 1,4-Methanonaphthalene, 1,4... Concentration Rank 43

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.55 | 2.68 ng/ul | 636686 | Naphthalene-d8   | 10.36 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2-methyl-              | 142 | C11H10  | 000091-57-6 | 95   |
| 2    |    | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 95   |
| 3    |    | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 86   |
| 4    |    | 1,4-Methanonaphthalene, 1,4-dihy... | 142 | C11H10  | 004453-90-1 | 86   |
| 5    |    | Bicyclo[4.4.1]undeca-1,3,5,7,9-p... | 142 | C11H10  | 002443-46-1 | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

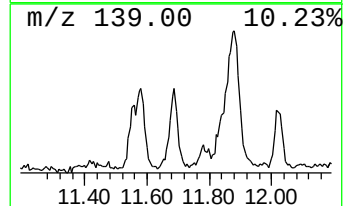
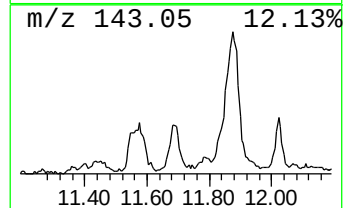
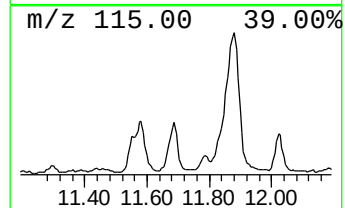
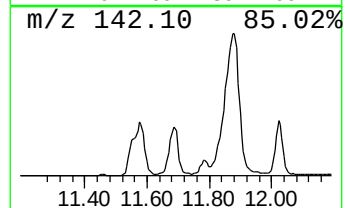
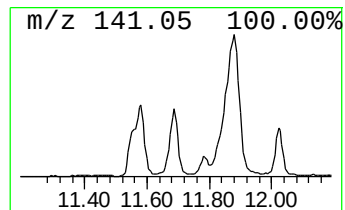
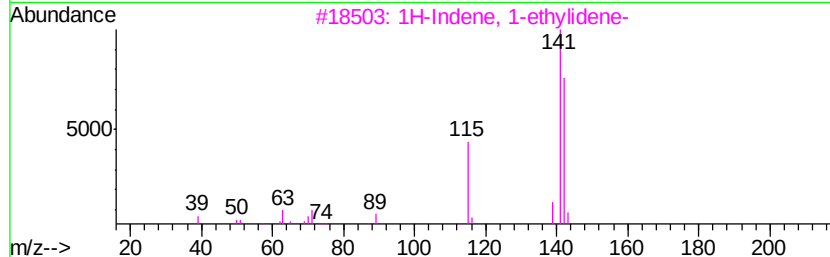
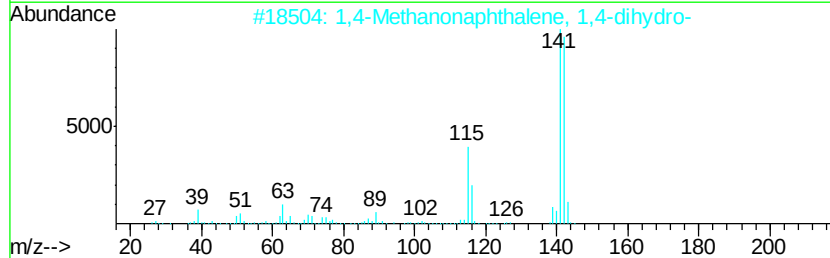
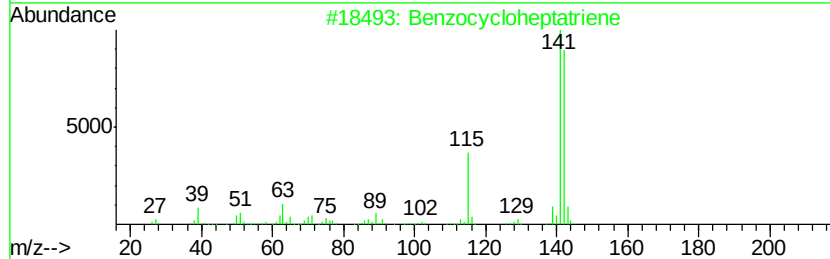
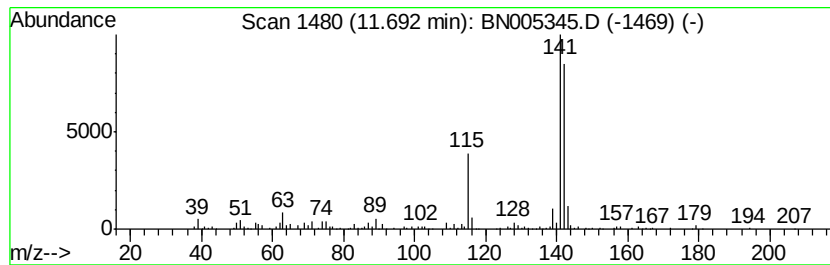
Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 6 1H-Indene, 1-ethylidene- Concentration Rank 31

| R.T.    | EstConc    | Area                                | Relative to ISTD | R.T.           |
|---------|------------|-------------------------------------|------------------|----------------|
| 11.69   | 3.29 ng/ul | 782245                              | Naphthalene-d8   | 10.36          |
| Hit# of | 5          | Tentative ID                        | MW MolForm       | CAS# Qual      |
| 1       |            | Benzocycloheptatriene               | 142 C11H10       | 000264-09-5 94 |
| 2       |            | 1,4-Methanonaphthalene, 1,4-dihy... | 142 C11H10       | 004453-90-1 94 |
| 3       |            | 1H-Indene, 1-ethylidene-            | 142 C11H10       | 002471-83-2 94 |
| 4       |            | Bicyclo[4.4.1]undeca-1,3,5,7,9-p... | 142 C11H10       | 002443-46-1 94 |
| 5       |            | 1H-Indene, 1-ethylidene-            | 142 C11H10       | 002471-83-2 91 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

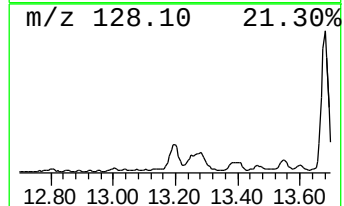
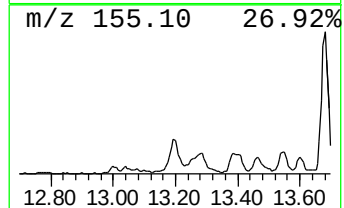
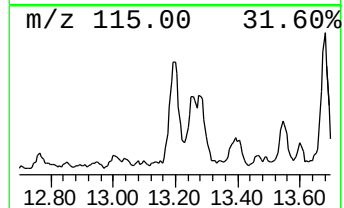
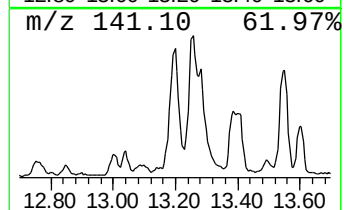
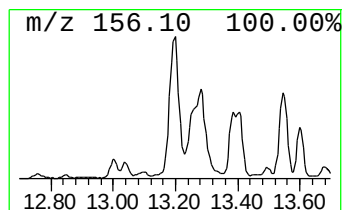
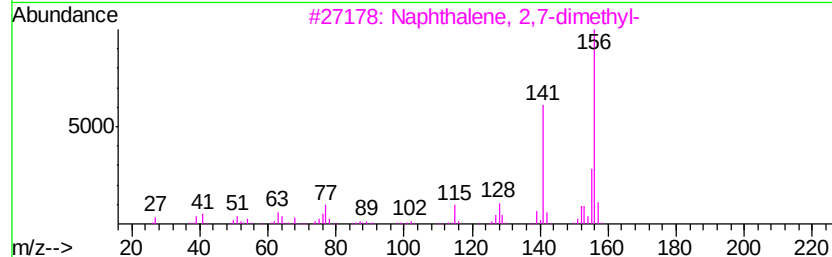
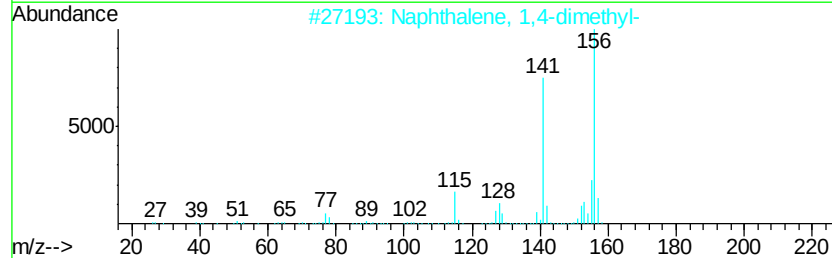
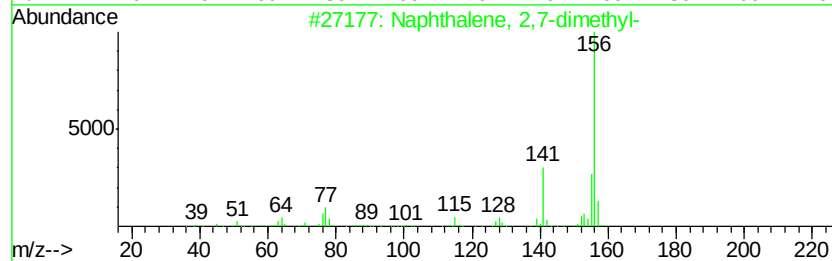
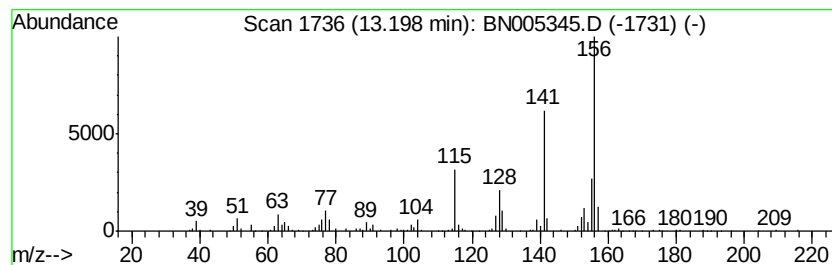
Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 8 Naphthalene, 2,7-dimethyl- Concentration Rank 40

| R.T.    | EstConc    | Area                       | Relative to ISTD | R.T.           |
|---------|------------|----------------------------|------------------|----------------|
| 13.20   | 2.73 ng/ul | 823715                     | Acenaphthene-d10 | 14.23          |
| Hit# of | 5          | Tentative ID               | MW MolForm       | CAS# Qual      |
| 1       |            | Naphthalene, 2,7-dimethyl- | 156 C12H12       | 000582-16-1 95 |
| 2       |            | Naphthalene, 1,4-dimethyl- | 156 C12H12       | 000571-58-4 95 |
| 3       |            | Naphthalene, 2,7-dimethyl- | 156 C12H12       | 000582-16-1 94 |
| 4       |            | Naphthalene, 1,7-dimethyl- | 156 C12H12       | 000575-37-1 94 |
| 5       |            | Naphthalene, 2,7-dimethyl- | 156 C12H12       | 000582-16-1 94 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleID :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

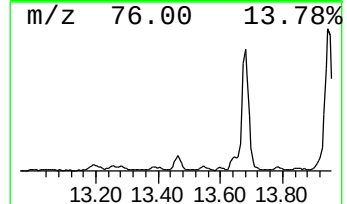
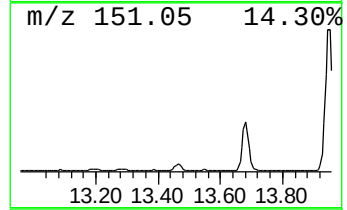
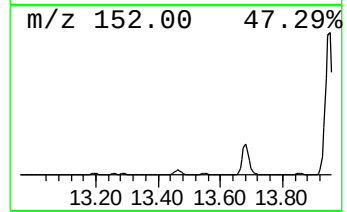
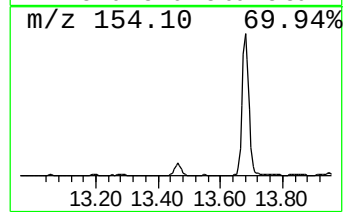
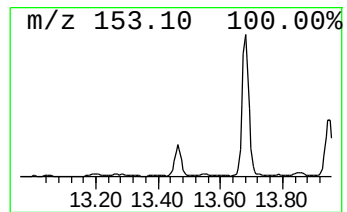
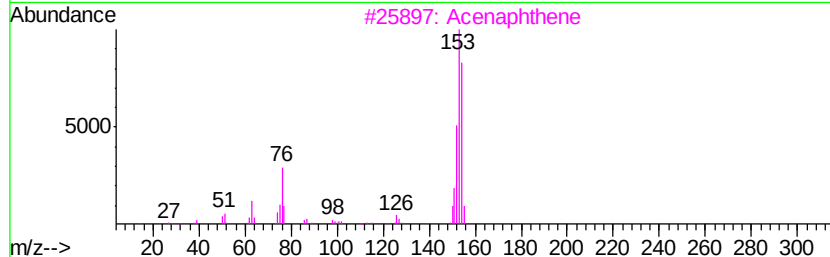
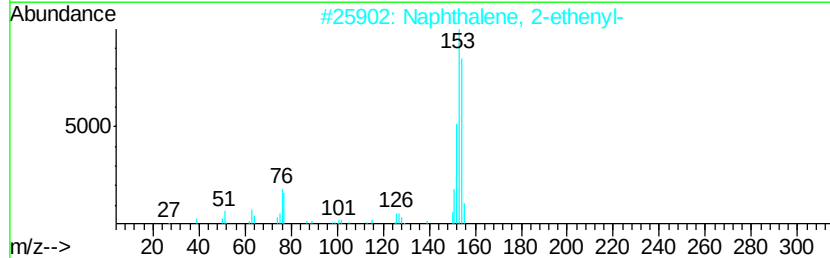
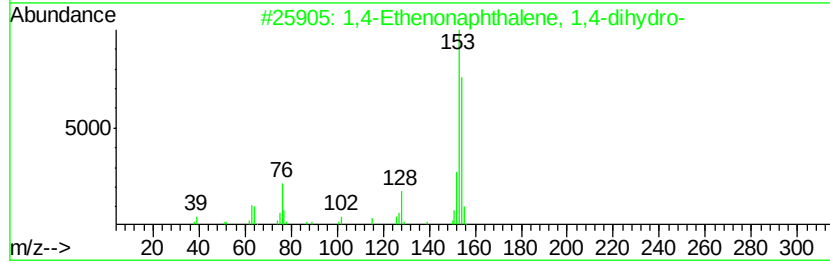
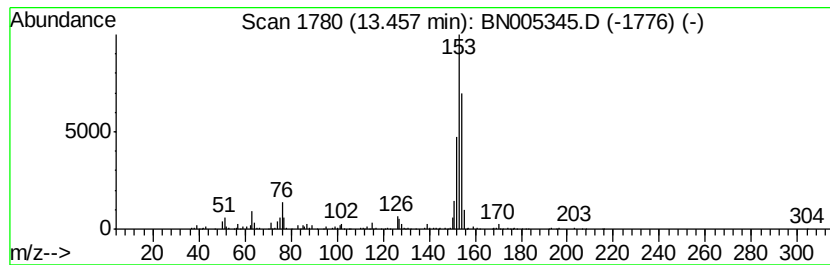
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 1,4-Ethenonaphthalene, 1,4-... Concentration Rank 38

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.46 | 2.99 ng/ul | 903637 | Acenaphthene-d10 | 14.23 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1,4-Ethenonaphthalene, 1,4-dihydro- | 154 | C12H10  | 007322-47-6 | 68   |
| 2    |    | Naphthalene, 2-ethenyl-             | 154 | C12H10  | 000827-54-3 | 55   |
| 3    |    | Acenaphthene                        | 154 | C12H10  | 000083-32-9 | 53   |
| 4    |    | Naphthalene, 2-ethenyl-             | 154 | C12H10  | 000827-54-3 | 50   |
| 5    |    | Naphthalene, 2-ethenyl-             | 154 | C12H10  | 000827-54-3 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

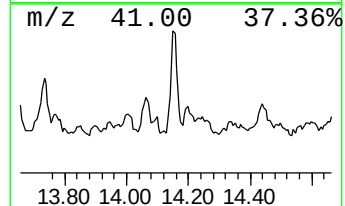
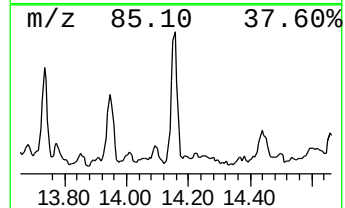
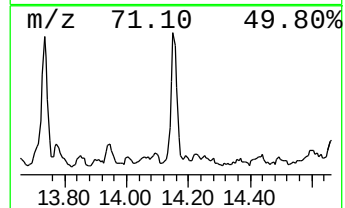
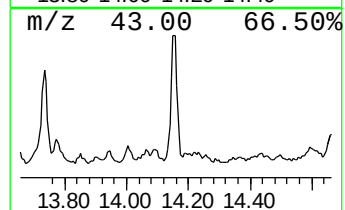
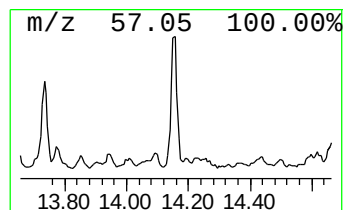
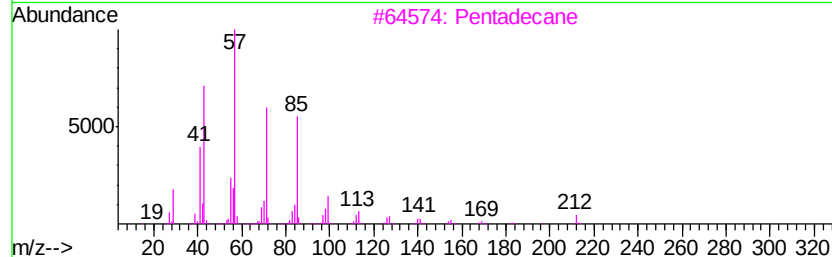
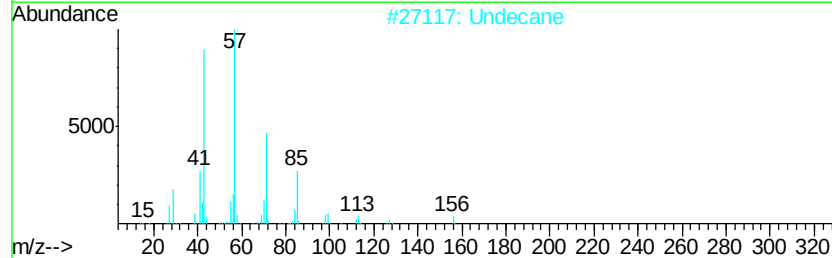
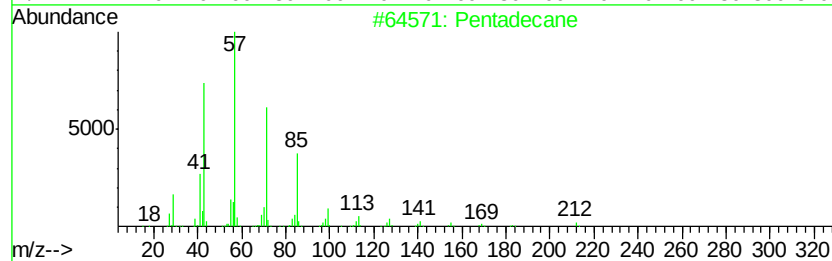
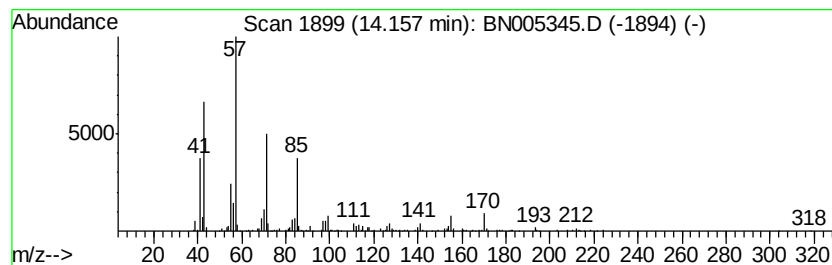
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 45

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.16 | 2.40 ng/ul | 724955 | Acenaphthene-d10 | 14.23 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentadecane            | 212 | C15H32  | 000629-62-9 | 94   |
| 2    |    |   | Undecane               | 156 | C11H24  | 001120-21-4 | 87   |
| 3    |    |   | Pentadecane            | 212 | C15H32  | 000629-62-9 | 86   |
| 4    |    |   | Tetradecane, 2-methyl- | 212 | C15H32  | 001560-95-8 | 83   |
| 5    |    |   | Eicosane               | 282 | C20H42  | 000112-95-8 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

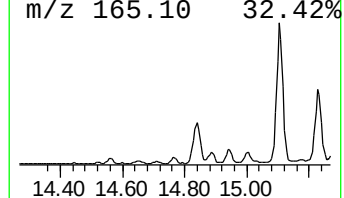
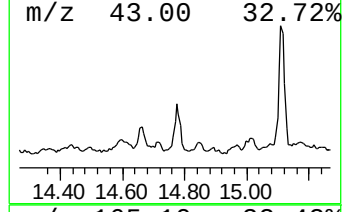
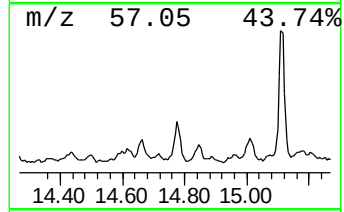
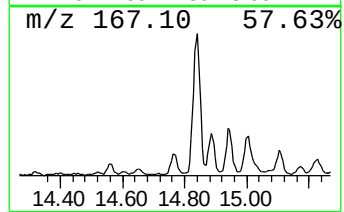
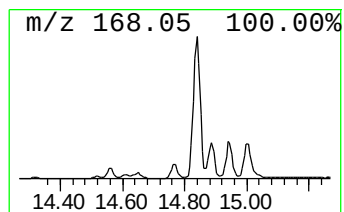
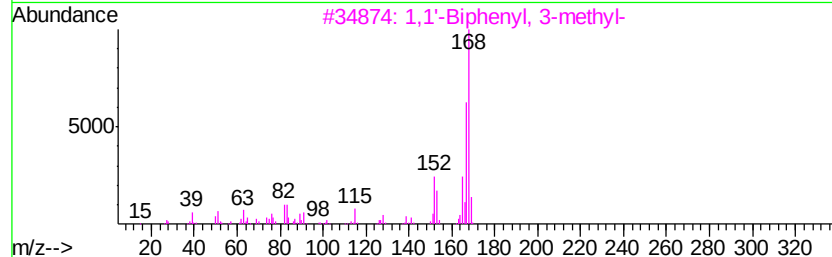
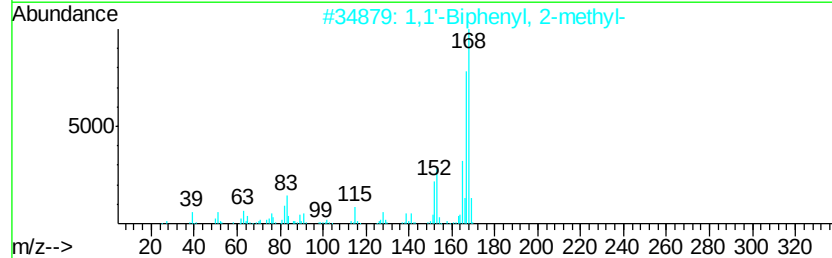
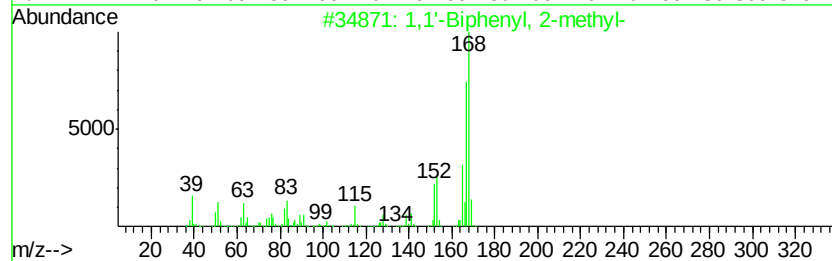
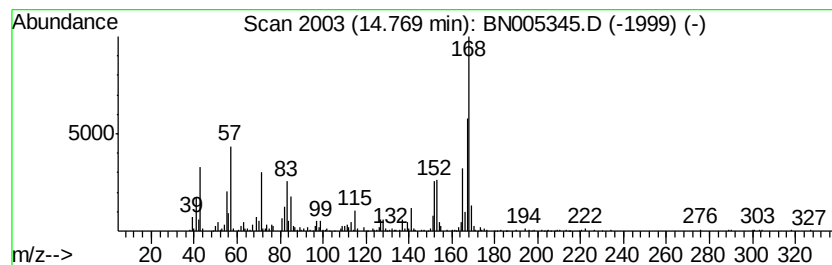
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 1,1'-Biphenyl, 2-methyl- Concentration Rank 47

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.77 | 2.24 ng/ul | 674696 | Acenaphthene-d10 | 14.23 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | 1,1'-Biphenyl, 2-methyl- | 168 | C13H12  | 000643-58-3 | 55   |
| 2    |    | 1,1'-Biphenyl, 2-methyl- | 168 | C13H12  | 000643-58-3 | 55   |
| 3    |    | 1,1'-Biphenyl, 3-methyl- | 168 | C13H12  | 000643-93-6 | 50   |
| 4    |    | 1,1'-Biphenyl, 4-methyl- | 168 | C13H12  | 000644-08-6 | 46   |
| 5    |    | 1,1'-Biphenyl, 3-methyl- | 168 | C13H12  | 000643-93-6 | 45   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

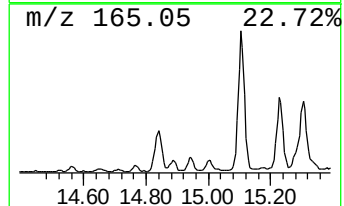
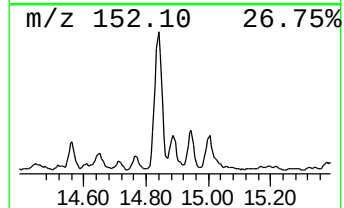
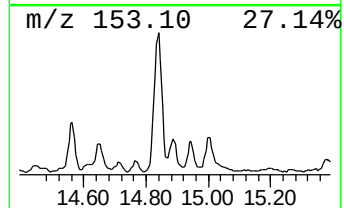
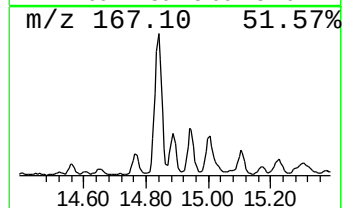
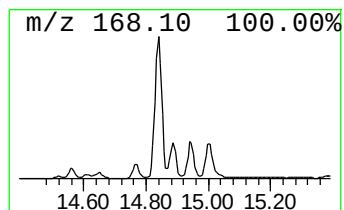
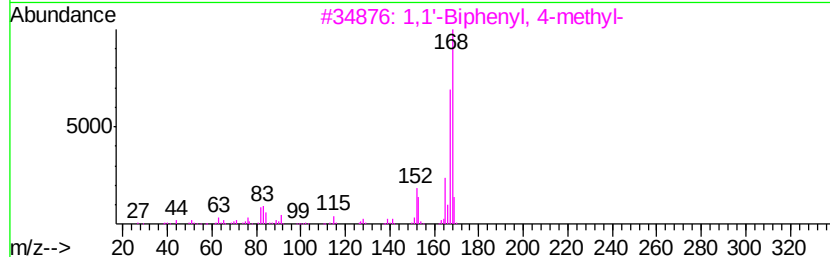
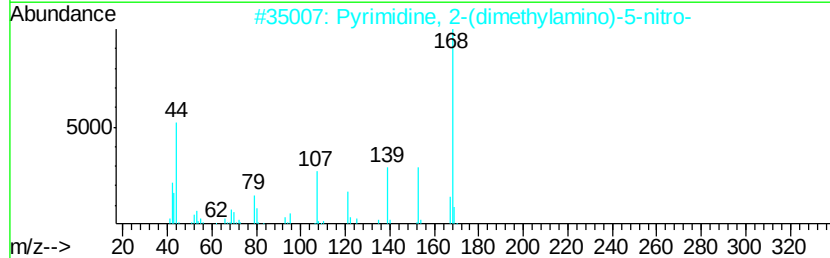
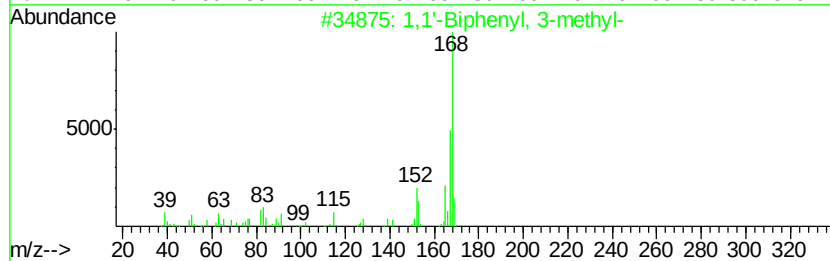
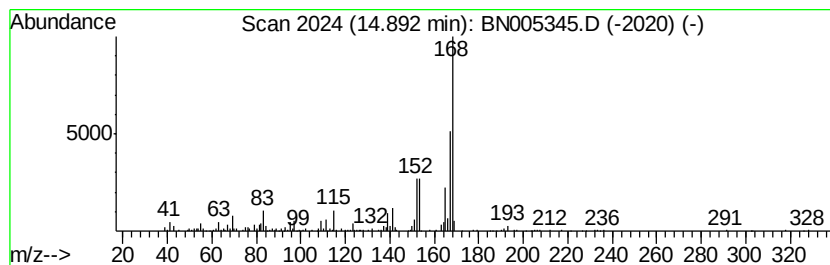
Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 13 unknown-01 Concentration Rank 33

| R.T.    | EstConc    | Area                                | Relative to ISTD | R.T.           |
|---------|------------|-------------------------------------|------------------|----------------|
| 14.89   | 3.23 ng/ul | 976118                              | Acenaphthene-d10 | 14.23          |
| Hit# of | 5          | Tentative ID                        | MW MolForm       | CAS# Qual      |
| 1       |            | 1,1'-Biphenyl, 3-methyl-            | 168 C13H12       | 000643-93-6 68 |
| 2       |            | Pyrimidine, 2-(dimethylamino)-5-... | 168 C6H8N4O2     | 014233-44-4 47 |
| 3       |            | 1,1'-Biphenyl, 4-methyl-            | 168 C13H12       | 000644-08-6 43 |
| 4       |            | 1,1'-Biphenyl, 3-methyl-            | 168 C13H12       | 000643-93-6 42 |
| 5       |            | 2,4-Dimethoxy-5-pyrimidine carbo... | 168 C7H8N2O3     | 052606-02-7 40 |





Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

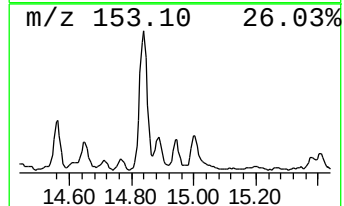
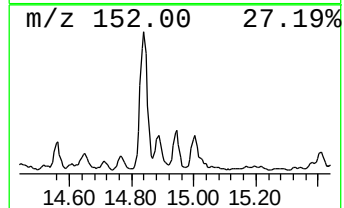
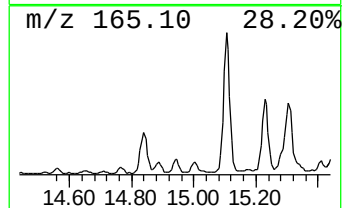
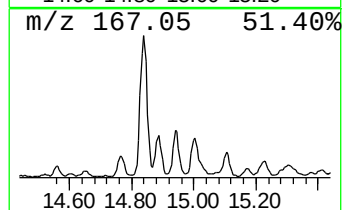
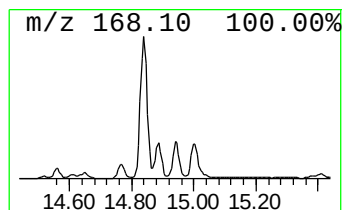
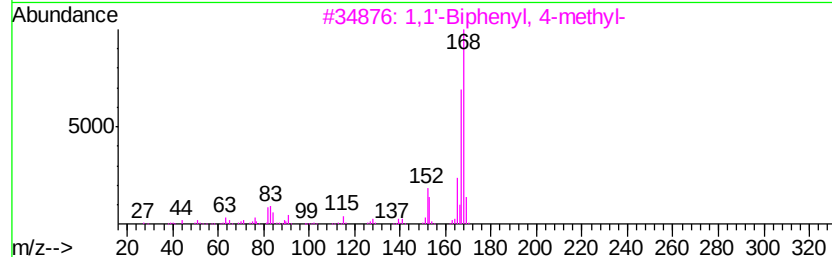
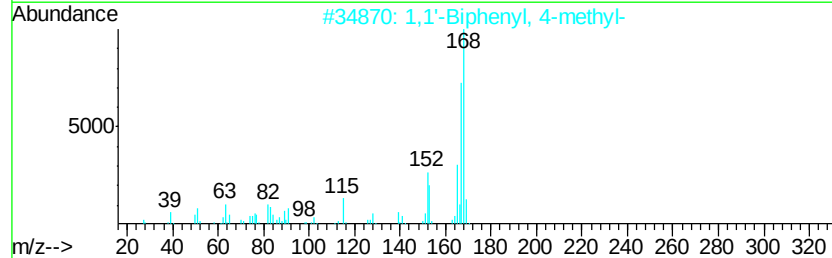
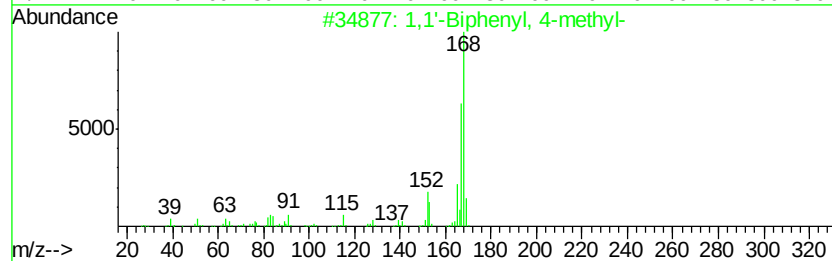
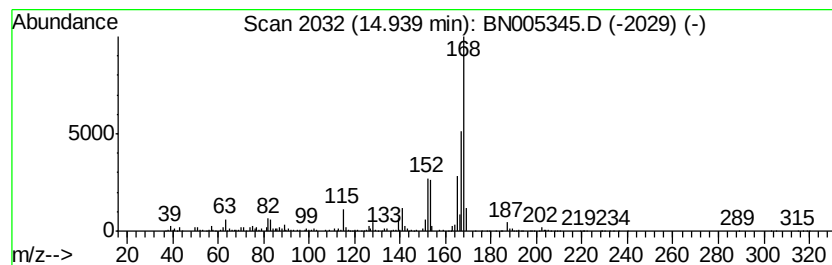
Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 14 1,1'-Biphenyl, 4-methyl- Concentration Rank 39

| R.T.    | EstConc                  | Area         | Relative to ISTD | R.T.      |
|---------|--------------------------|--------------|------------------|-----------|
| 14.94   | 2.93 ng/ul               | 885365       | Acenaphthene-d10 | 14.23     |
| Hit# of | 5                        | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | 1,1'-Biphenyl, 4-methyl- | 168 C13H12   | 000644-08-6      | 90        |
| 2       | 1,1'-Biphenyl, 4-methyl- | 168 C13H12   | 000644-08-6      | 81        |
| 3       | 1,1'-Biphenyl, 4-methyl- | 168 C13H12   | 000644-08-6      | 81        |
| 4       | 1,1'-Biphenyl, 3-methyl- | 168 C13H12   | 000643-93-6      | 68        |
| 5       | 1,1'-Biphenyl, 2-methyl- | 168 C13H12   | 000643-58-3      | 64        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

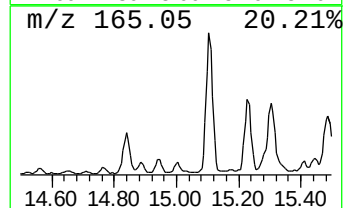
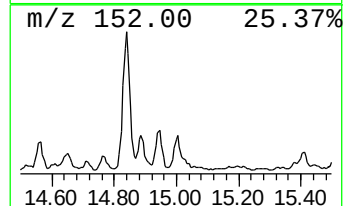
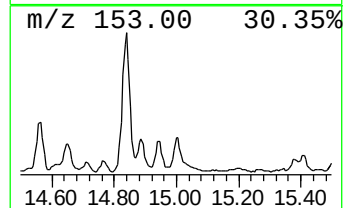
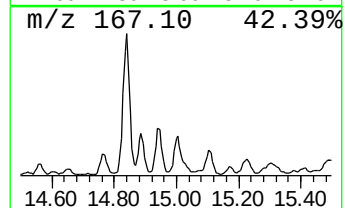
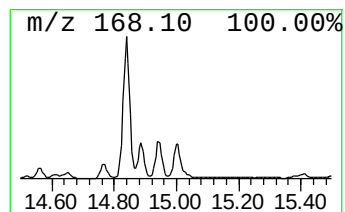
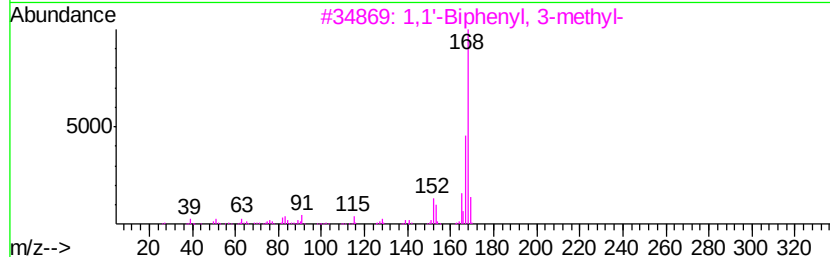
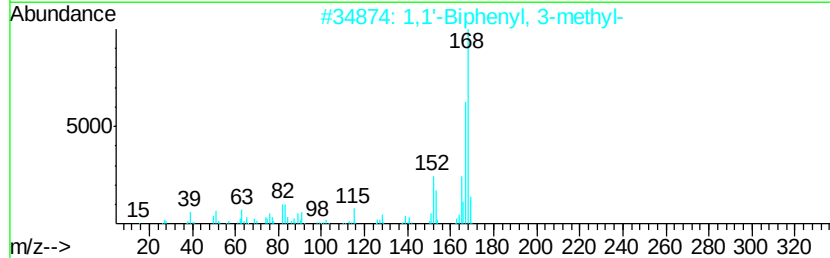
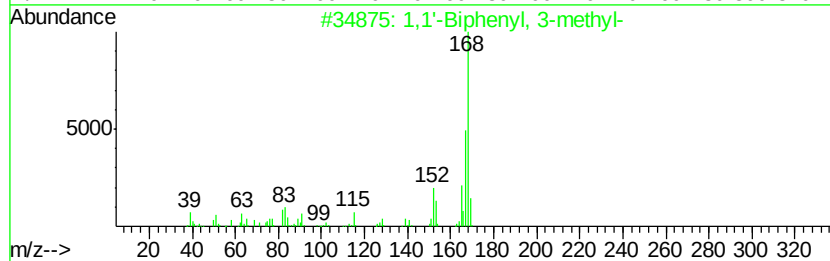
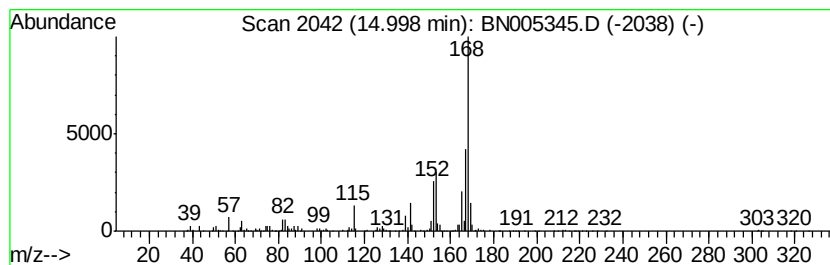
Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 15 1,1'-Biphenyl, 3-methyl- Concentration Rank 26

| R.T.      | EstConc                           | Area    | Relative to ISTD | R.T.        |      |
|-----------|-----------------------------------|---------|------------------|-------------|------|
| 15.00     | 4.02 ng/ul                        | 1213370 | Acenaphthene-d10 | 14.23       |      |
| Hit# of 5 | Tentative ID                      | MW      | MolForm          | CAS#        | Qual |
| 1         | 1,1'-Biphenyl, 3-methyl-          | 168     | C13H12           | 000643-93-6 | 76   |
| 2         | 1,1'-Biphenyl, 3-methyl-          | 168     | C13H12           | 000643-93-6 | 74   |
| 3         | 1,1'-Biphenyl, 3-methyl-          | 168     | C13H12           | 000643-93-6 | 72   |
| 4         | Naphthalene, 2-(1-methylethenyl)- | 168     | C13H12           | 003710-23-4 | 68   |
| 5         | 1,1'-Biphenyl, 3-methyl-          | 168     | C13H12           | 000643-93-6 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

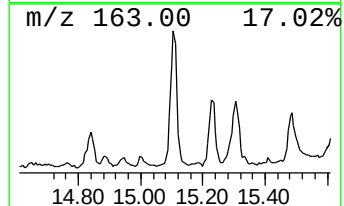
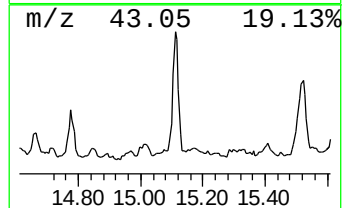
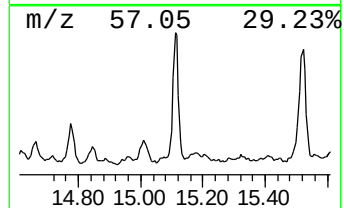
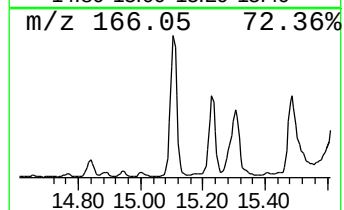
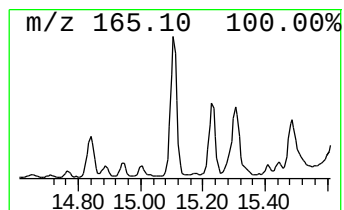
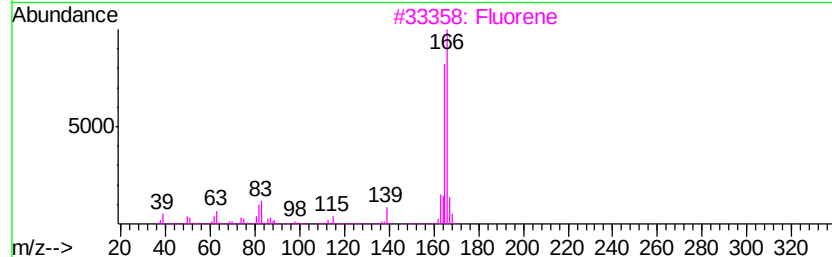
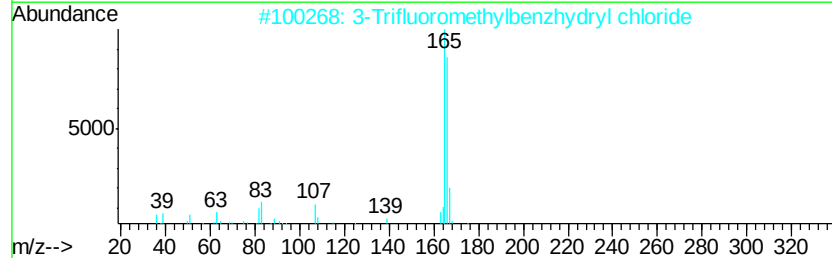
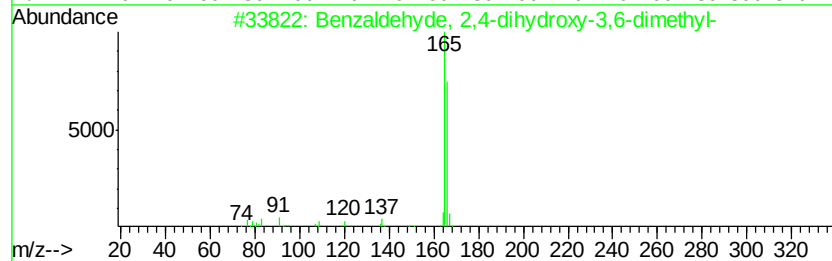
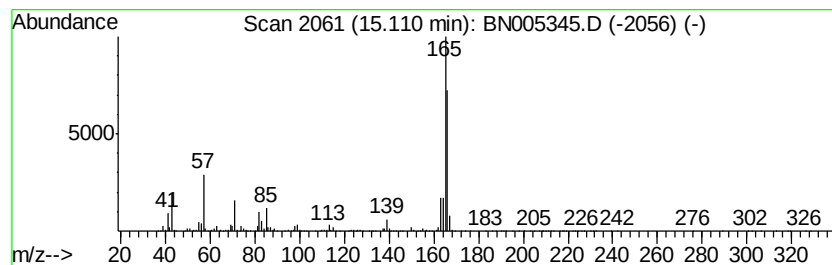
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Benzaldehyde, 2,4-dihydroxy... Concentration Rank 11

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 15.11 | 7.79 ng/ul | 2350110 | Acenaphthene-d10 | 14.23 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | Benzaldehyde, 2,4-dihydroxy-3,6-... | 166 | C9H10O3    | 034883-14-2 | 59   |
| 2    |    | 3-Trifluoromethylbenzhydryl chlo... | 270 | C14H10ClF3 | 067240-79-3 | 50   |
| 3    |    | Fluorene                            | 166 | C13H10     | 000086-73-7 | 47   |
| 4    |    | Fluorene-9-methanol                 | 196 | C14H12O    | 024324-17-2 | 45   |
| 5    |    | Benzene, 1,1'-(diazomethylene)bis-  | 194 | C13H10N2   | 000883-40-9 | 45   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

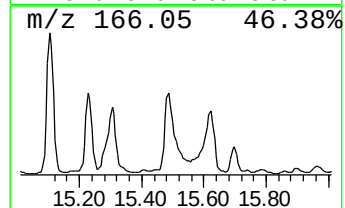
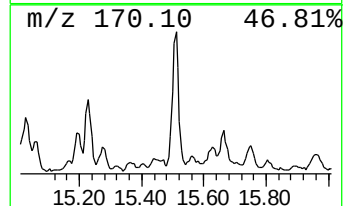
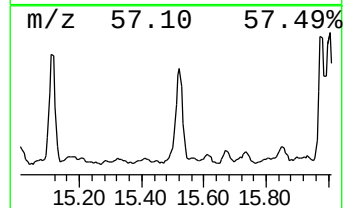
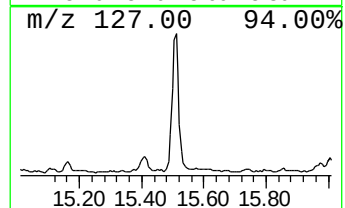
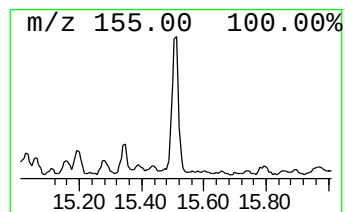
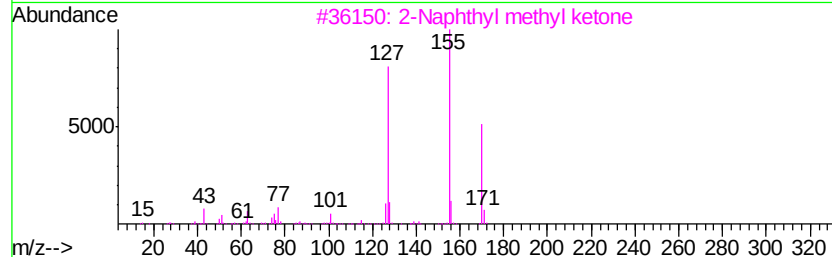
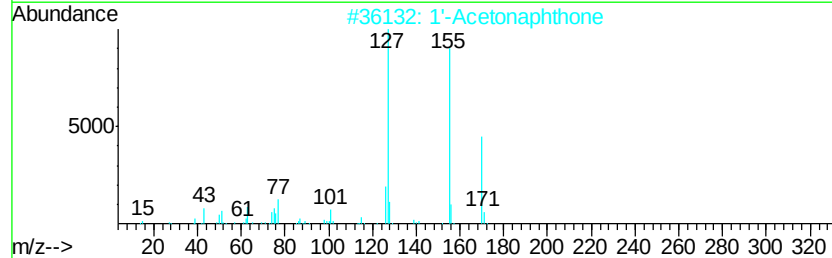
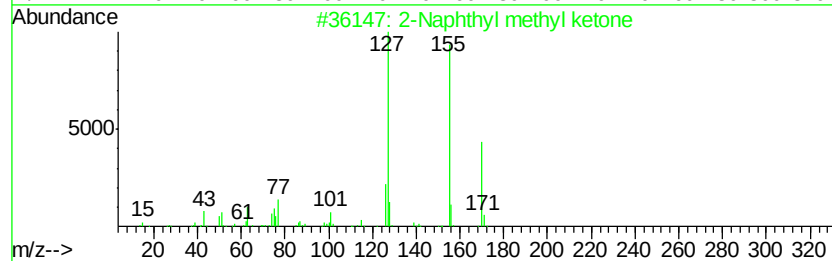
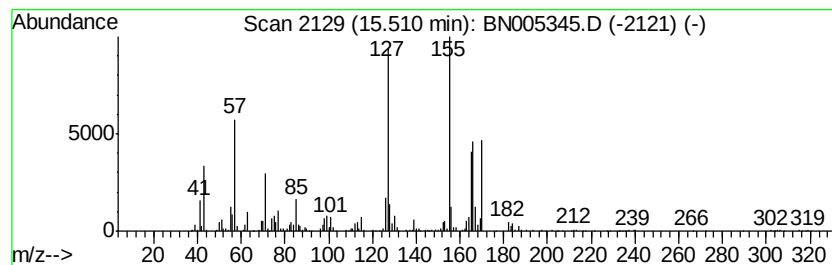
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 2-Naphthyl methyl ketone Concentration Rank 12

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 15.51 | 7.31 ng/ul | 2204610 | Acenaphthene-d10 | 14.23 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | 2-Naphthyl methyl ketone | 170 | C12H10O | 000093-08-3 | 96   |
| 2    |    |   | 1'-Acetonaphthone        | 170 | C12H10O | 000941-98-0 | 95   |
| 3    |    |   | 2-Naphthyl methyl ketone | 170 | C12H10O | 000093-08-3 | 92   |
| 4    |    |   | 2-Naphthyl methyl ketone | 170 | C12H10O | 000093-08-3 | 70   |
| 5    |    |   | 1'-Acetonaphthone        | 170 | C12H10O | 000941-98-0 | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleID :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

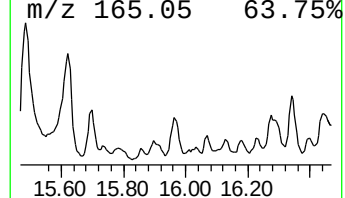
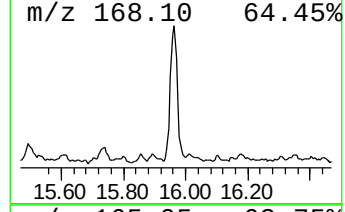
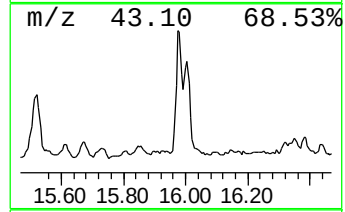
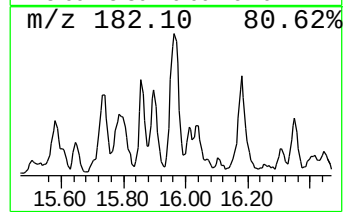
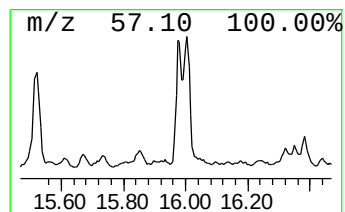
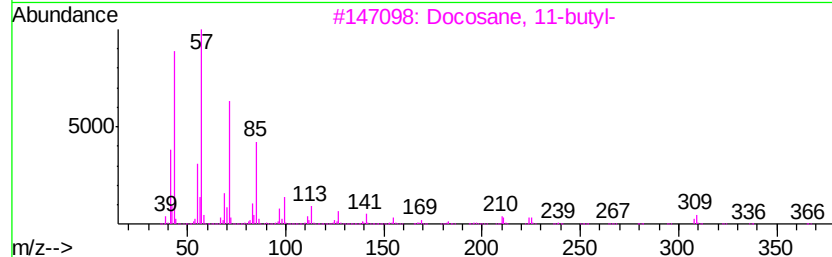
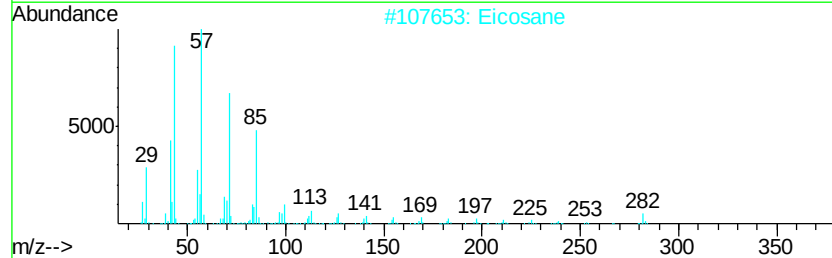
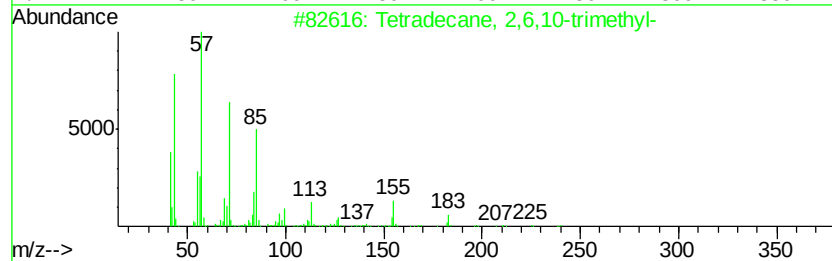
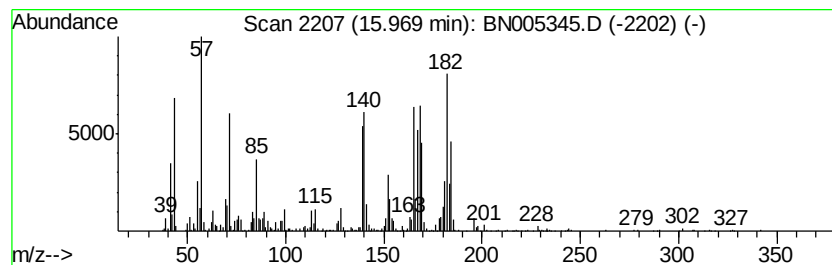
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 16

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 15.97 | 5.86 ng/ul | 2033180 | Phenanthrene-d10 | 16.99 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Tetradecane, 2,6,10-trimethyl- | 240 | C17H36  | 014905-56-7 | 46   |
| 2    |    |   | Eicosane                       | 282 | C20H42  | 000112-95-8 | 42   |
| 3    |    |   | Docosane, 11-butyl-            | 366 | C26H54  | 013475-76-8 | 42   |
| 4    |    |   | Eicosane                       | 282 | C20H42  | 000112-95-8 | 42   |
| 5    |    |   | Docosane, 9-butyl-             | 366 | C26H54  | 055282-14-9 | 41   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

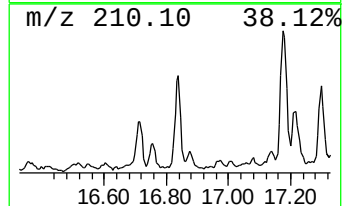
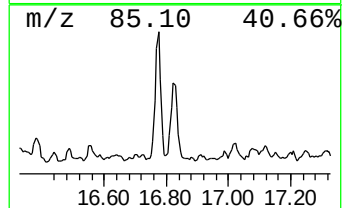
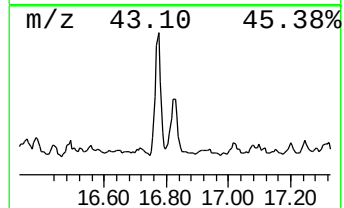
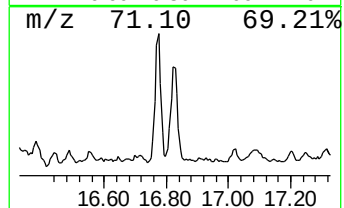
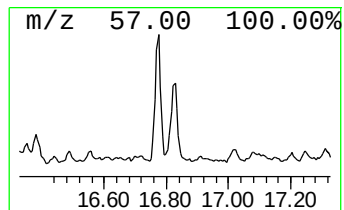
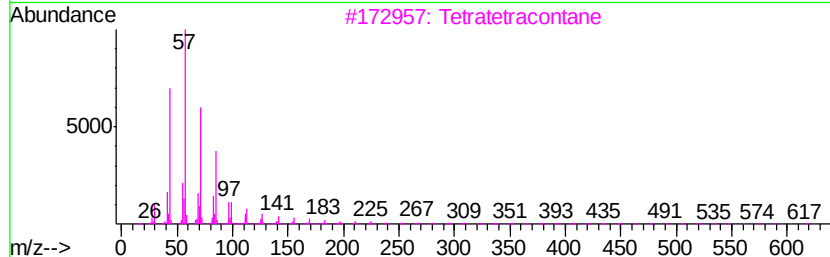
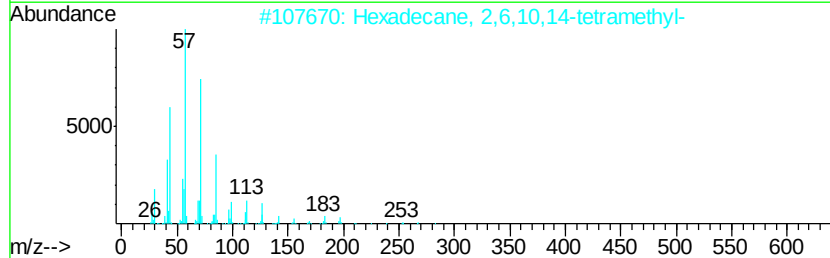
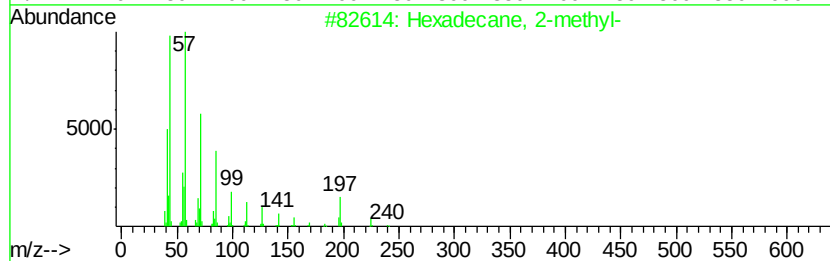
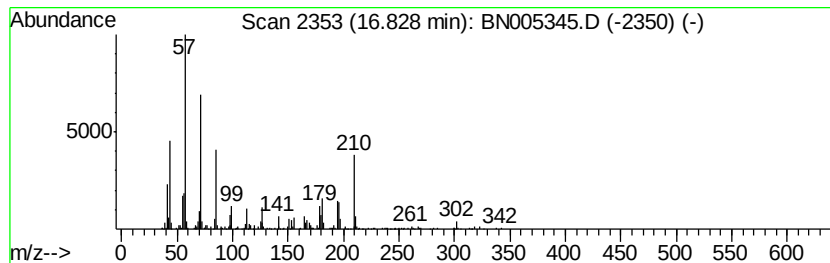
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 49

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.83 | 2.04 ng/ul | 705764 | Phenanthrene-d10 | 16.99 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane, 2-methyl-              | 240 | C17H36  | 001560-92-5 | 55   |
| 2    |    |   | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8 | 49   |
| 3    |    |   | Tetratetracontane                  | 619 | C44H90  | 007098-22-8 | 46   |
| 4    |    |   | Decane, 3,6-dimethyl-              | 170 | C12H26  | 017312-53-7 | 46   |
| 5    |    |   | Eicosane, 2-methyl-                | 296 | C21H44  | 001560-84-5 | 46   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleID :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

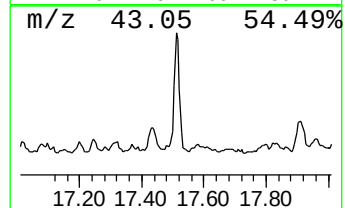
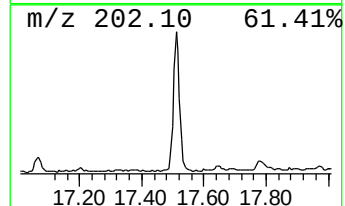
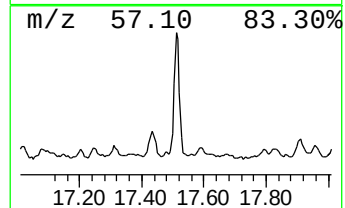
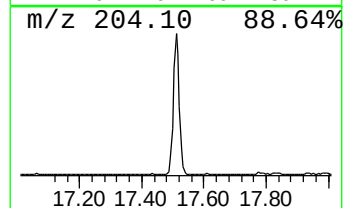
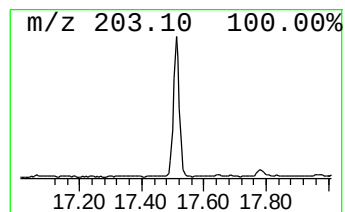
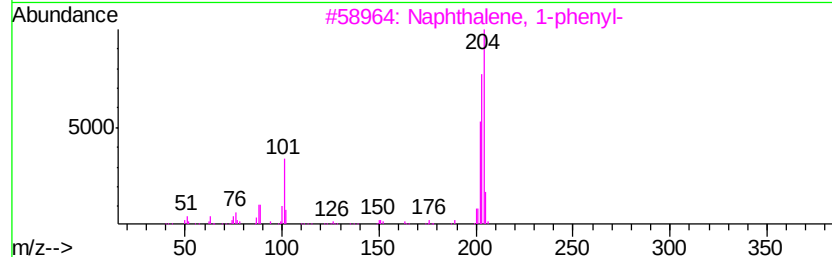
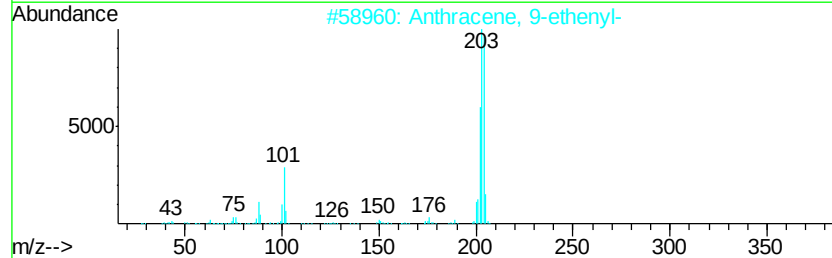
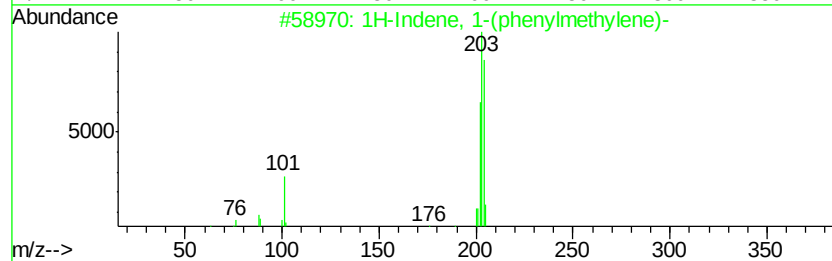
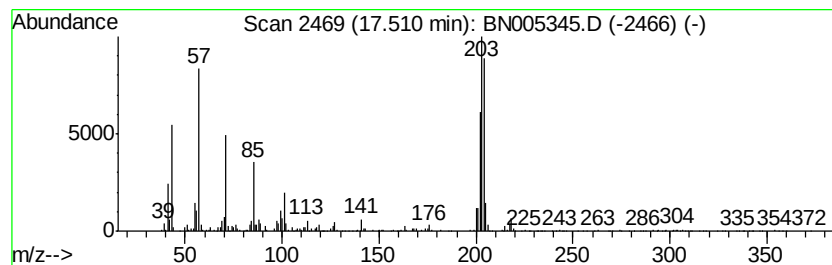
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 1H-Indene, 1-(phenylmethyle... Concentration Rank 25

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 17.51 | 4.20 ng/ul | 1457400 | Phenanthrene-d10 | 16.99 |

| Hit# | of | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1H-Indene, 1-(phenylmethylene)- | 204 | C16H12  | 005394-86-5 | 78   |
| 2    |    | Anthracene, 9-ethenyl-          | 204 | C16H12  | 002444-68-0 | 60   |
| 3    |    | Naphthalene, 1-phenyl-          | 204 | C16H12  | 000605-02-7 | 50   |
| 4    |    | Dibenzo[a,e]cyclooctene         | 204 | C16H12  | 000262-89-5 | 49   |
| 5    |    | Anthracene, 9-ethenyl-          | 204 | C16H12  | 002444-68-0 | 49   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

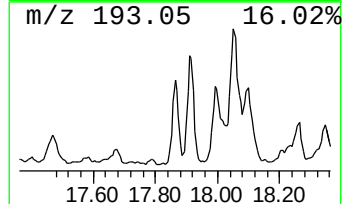
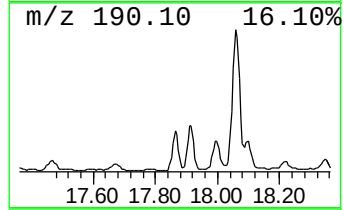
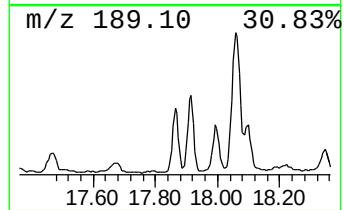
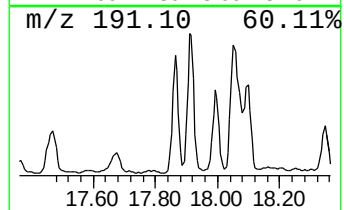
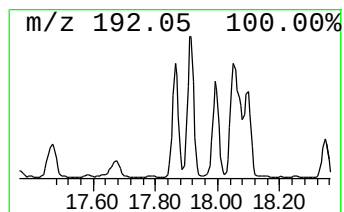
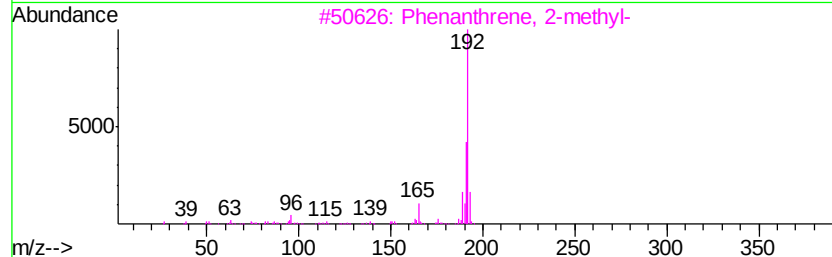
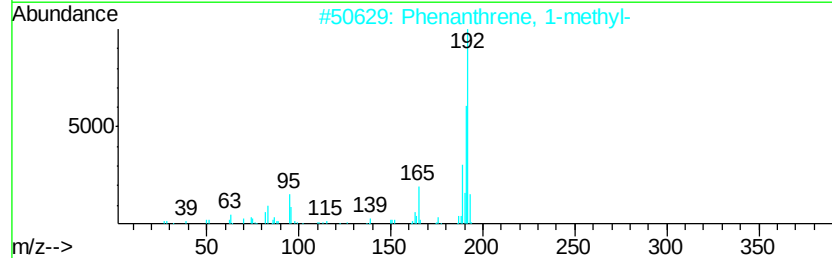
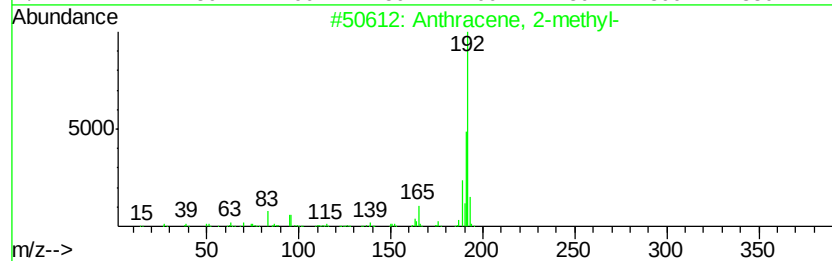
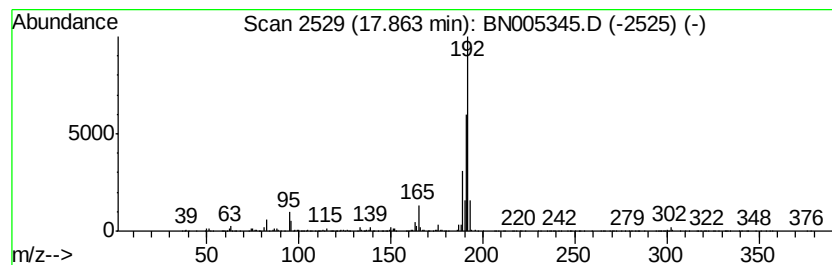
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 Anthracene, 2-methyl- Concentration Rank 14

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 17.86 | 6.69 ng/ul | 2317840 | Phenanthrene-d10 | 16.99 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 95   |
| 2    |    |   | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 94   |
| 3    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 93   |
| 4    |    |   | Anthracene, 1-methyl-   | 192 | C15H12  | 000610-48-0 | 93   |
| 5    |    |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 93   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleID :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

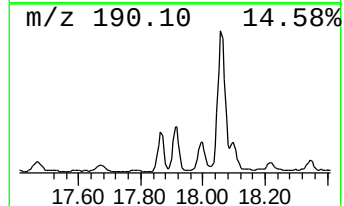
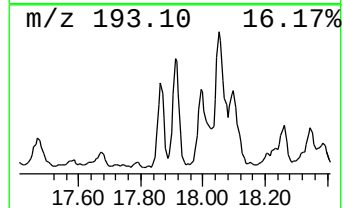
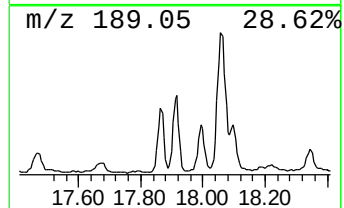
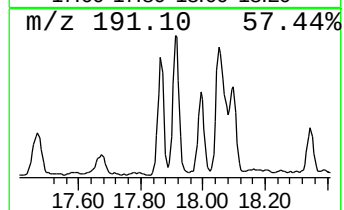
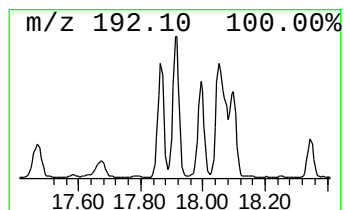
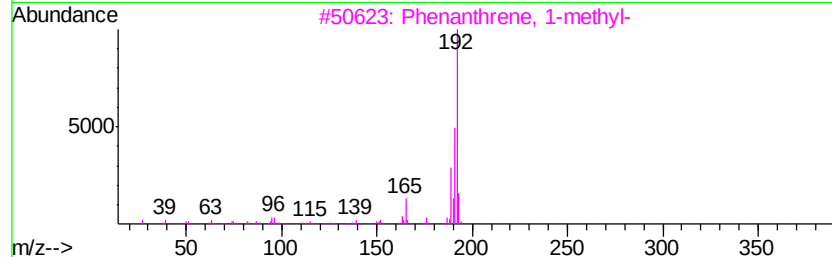
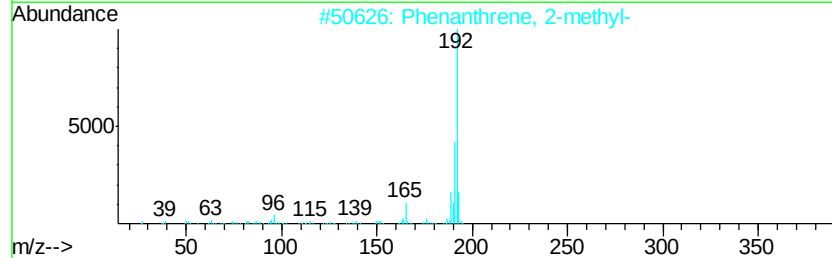
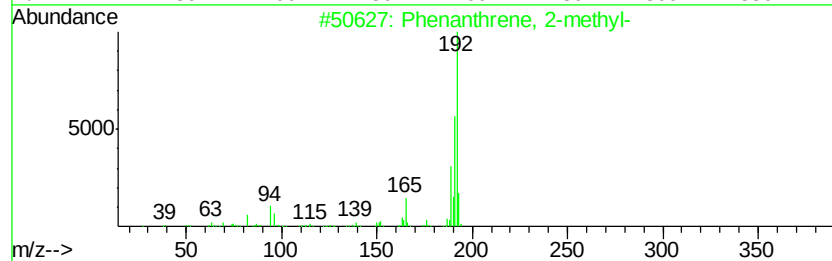
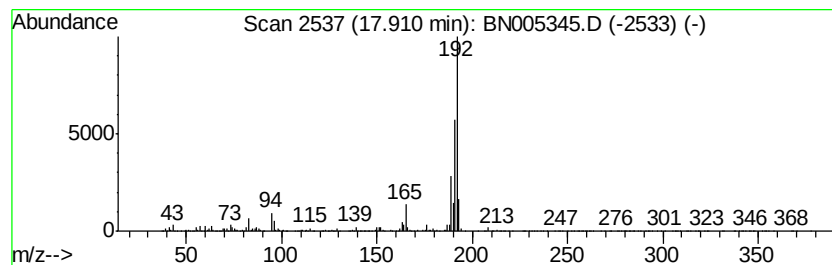
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 23 Phenanthrene, 2-methyl- Concentration Rank 5

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 17.91 | 9.98 ng/ul | 3458890 | Phenanthrene-d10 | 16.99 |

| Hit# | of | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenanthrene, 2-methyl-               | 192 | C15H12  | 002531-84-2 | 98   |
| 2    |    | Phenanthrene, 2-methyl-               | 192 | C15H12  | 002531-84-2 | 97   |
| 3    |    | Phenanthrene, 1-methyl-               | 192 | C15H12  | 000832-69-9 | 97   |
| 4    |    | 1H-Cyclopropa[l]phenanthrene, 1a, ... | 192 | C15H12  | 000949-41-7 | 96   |
| 5    |    | Anthracene, 2-methyl-                 | 192 | C15H12  | 000613-12-7 | 96   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

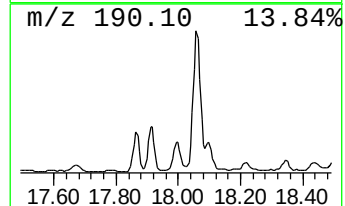
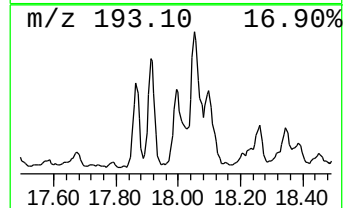
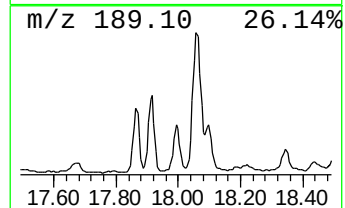
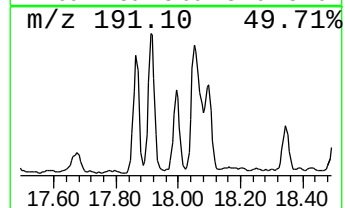
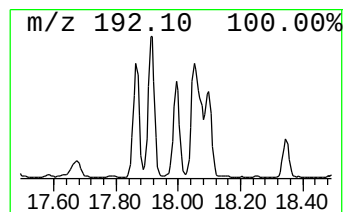
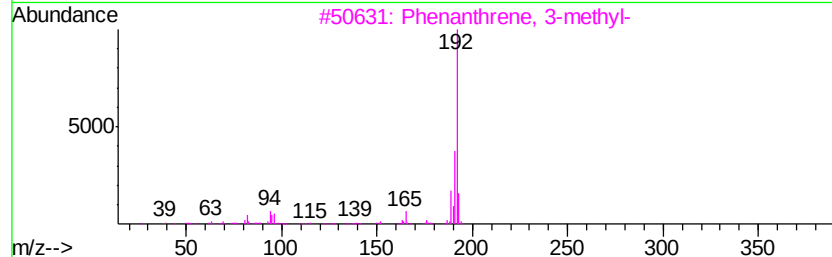
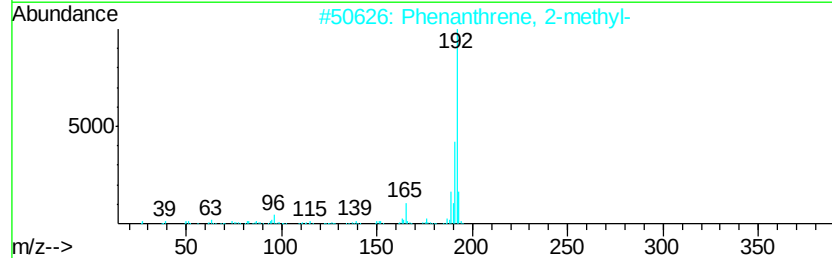
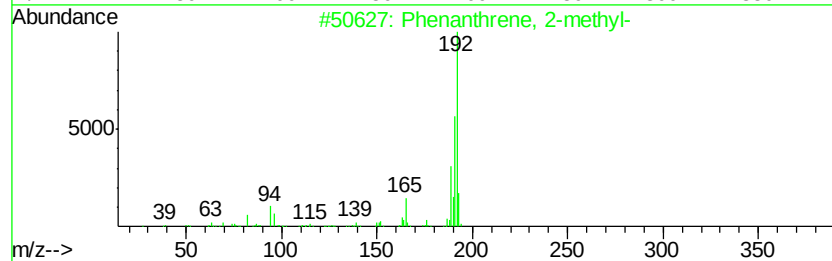
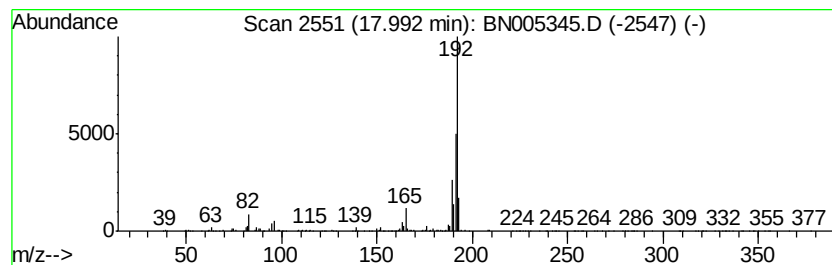
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 24 Phenanthrene, 3-methyl- Concentration Rank 23

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 17.99 | 4.69 ng/ul | 1624640 | Phenanthrene-d10 | 16.99 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 96   |
| 2    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 94   |
| 3    |    |   | Phenanthrene, 3-methyl- | 192 | C15H12  | 000832-71-3 | 93   |
| 4    |    |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 93   |
| 5    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

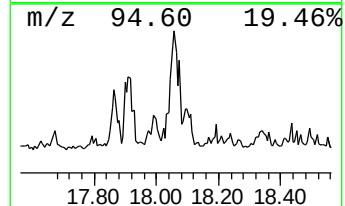
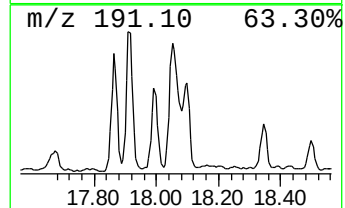
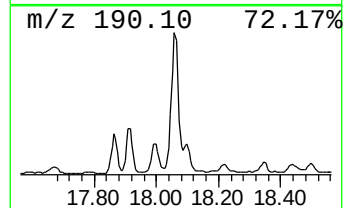
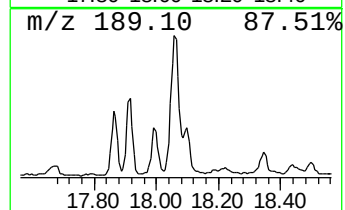
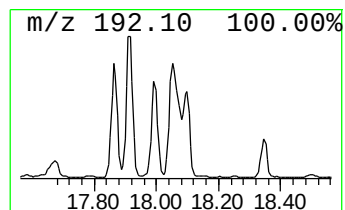
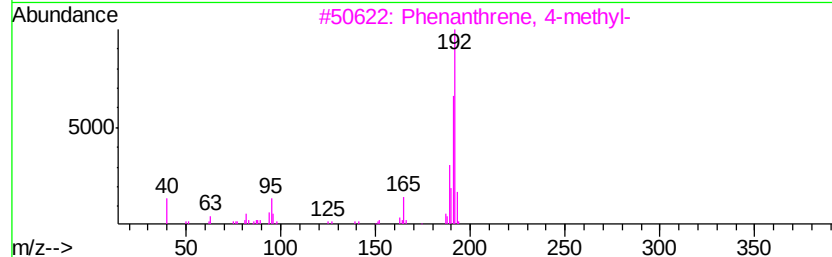
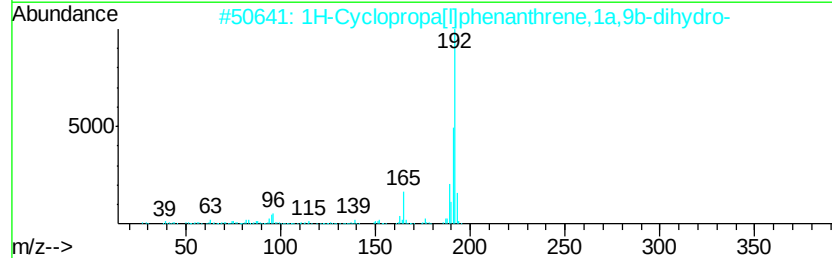
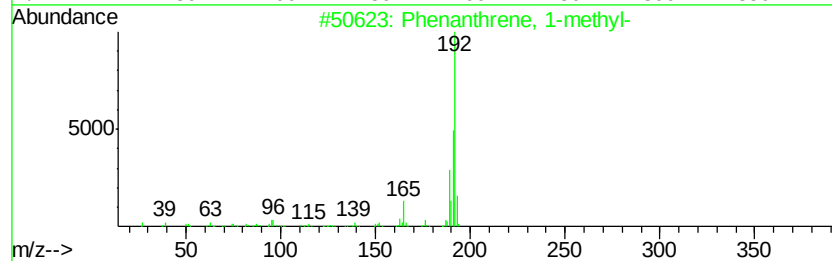
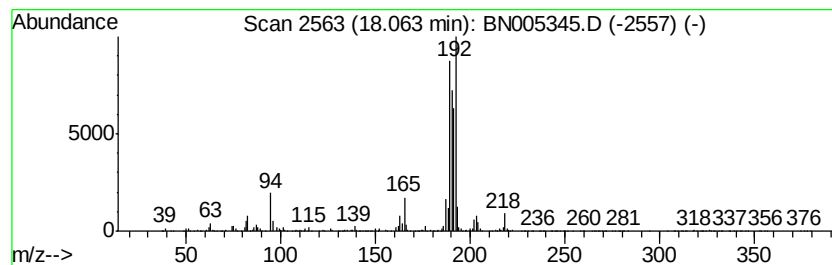
Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 25 Phenanthrene, 4-methyl- Concentration Rank 1

| R.T.    | EstConc     | Area                                | Relative to ISTD | R.T.    |             |      |
|---------|-------------|-------------------------------------|------------------|---------|-------------|------|
| 18.06   | 13.99 ng/ul | 4849380                             | Phenanthrene-d10 | 16.99   |             |      |
| Hit# of | 5           | Tentative ID                        | MW               | MolForm | CAS#        | Qual |
| 1       |             | Phenanthrene, 1-methyl-             | 192              | C15H12  | 000832-69-9 | 90   |
| 2       |             | 1H-Cyclopropa[l]phenanthrene,1a,... | 192              | C15H12  | 000949-41-7 | 60   |
| 3       |             | Phenanthrene, 4-methyl-             | 192              | C15H12  | 000832-64-4 | 60   |
| 4       |             | Phenanthrene, 1-methyl-             | 192              | C15H12  | 000832-69-9 | 60   |
| 5       |             | Anthracene, 1-methyl-               | 192              | C15H12  | 000610-48-0 | 60   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

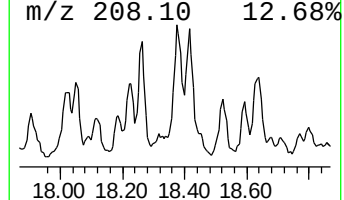
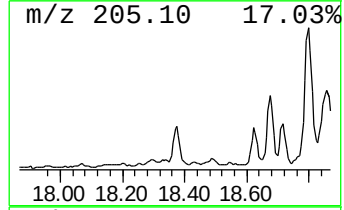
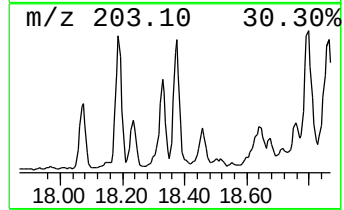
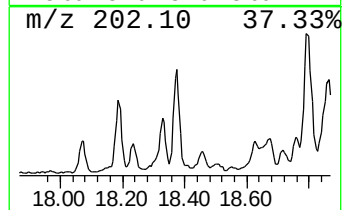
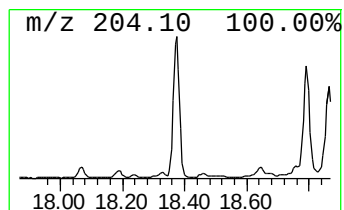
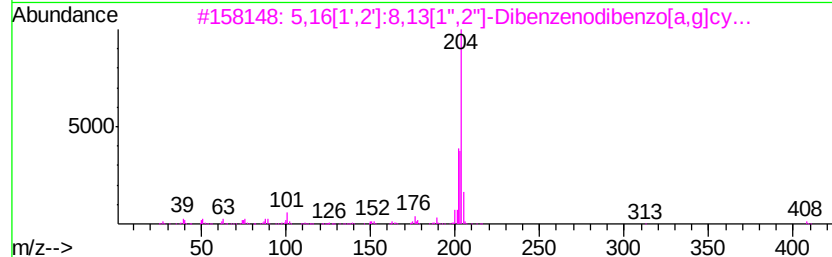
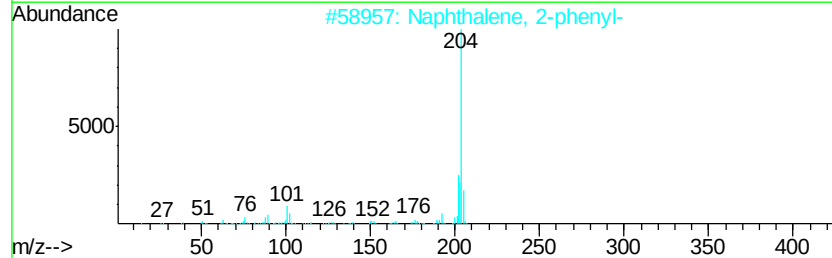
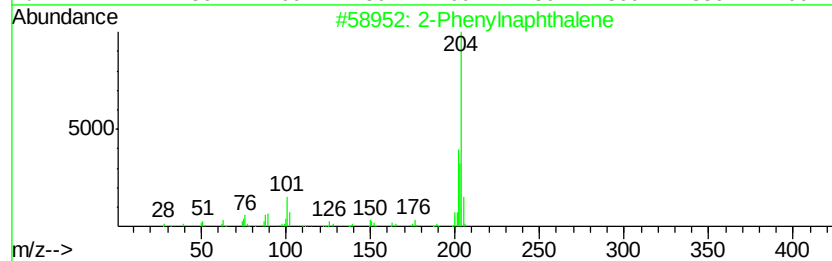
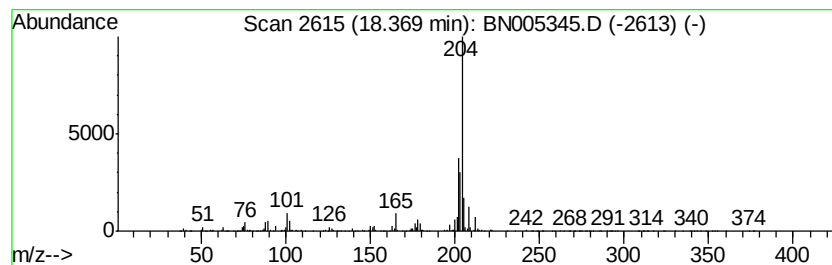
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 26 2-Phenylanthracene Concentration Rank 29

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 18.37 | 3.44 ng/ul | 1191050 | Phenanthrene-d10 | 16.99 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | 2-Phenylanthracene                  | 204 | C16H12    | 035465-71-5 | 94   |
| 2    |    | Naphthalene, 2-phenyl-              | 204 | C16H12    | 000612-94-2 | 93   |
| 3    |    | 5,16[1',2']:8,13[1'',2'']-Dibenz... | 408 | C32H24    | 005672-97-9 | 83   |
| 4    |    | 9,10-di(Chloromethyl)anthracene     | 274 | C16H12Cl2 | 010387-13-0 | 72   |
| 5    |    | Naphthalene, 2-phenyl-              | 204 | C16H12    | 000612-94-2 | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

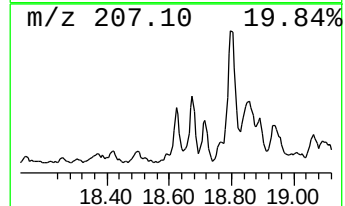
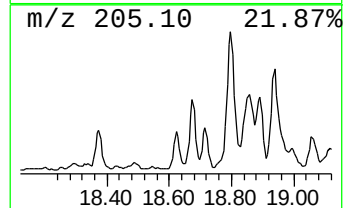
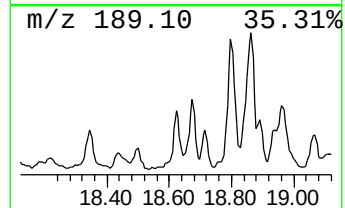
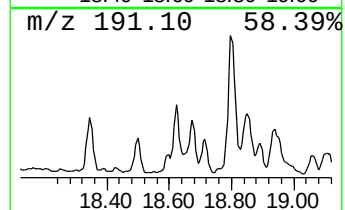
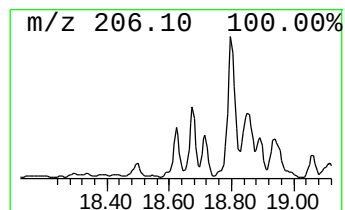
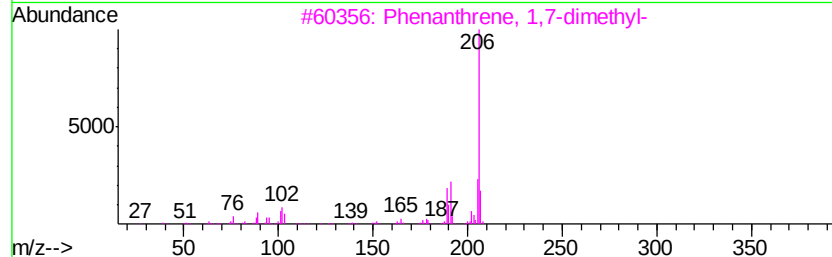
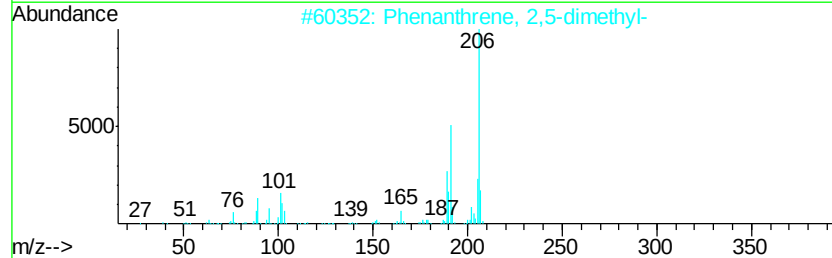
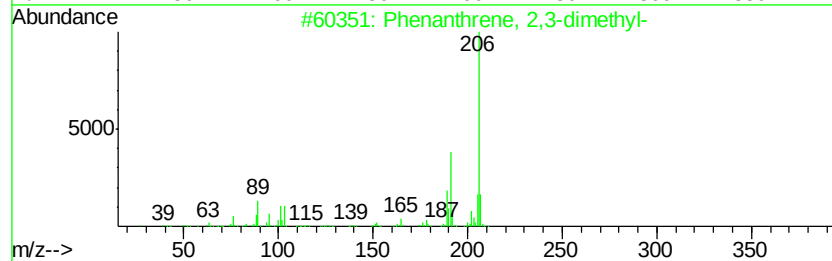
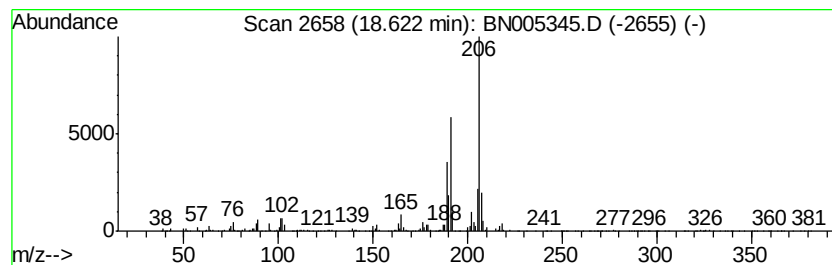
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 27 Phenanthrene, 2,3-dimethyl- Concentration Rank 30

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 18.62 | 3.37 ng/ul | 1167820 | Phenanthrene-d10 | 16.99 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenanthrene, 2,3-dimethyl-   | 206 | C16H14  | 003674-65-5 | 93   |
| 2    |    | Phenanthrene, 2,5-dimethyl-   | 206 | C16H14  | 003674-66-6 | 93   |
| 3    |    | Phenanthrene, 1,7-dimethyl-   | 206 | C16H14  | 000483-87-4 | 90   |
| 4    |    | 1H-Indene, 2-methyl-3-phenyl- | 206 | C16H14  | 035099-60-6 | 83   |
| 5    |    | 9,10-Dimethylantracene        | 206 | C16H14  | 000781-43-1 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

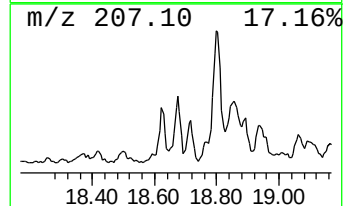
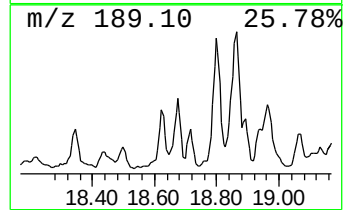
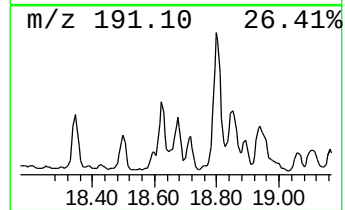
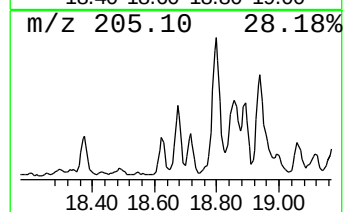
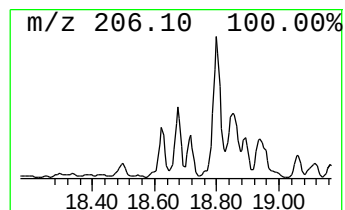
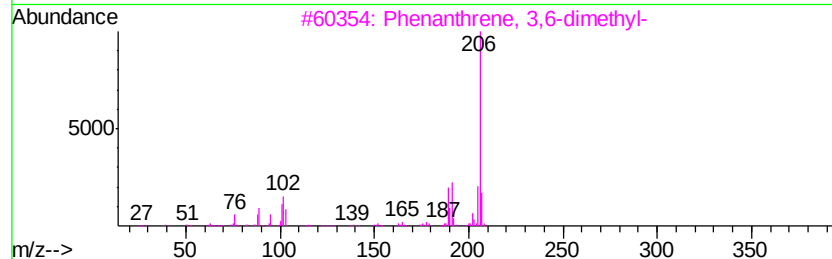
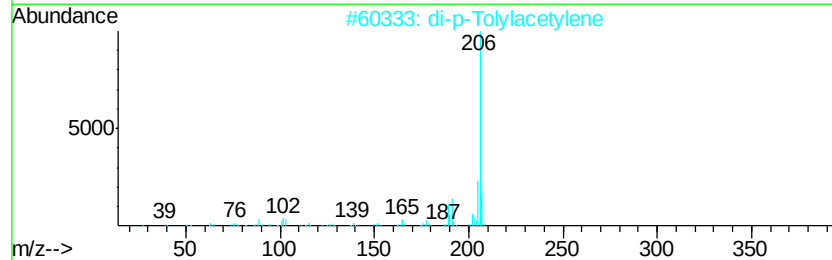
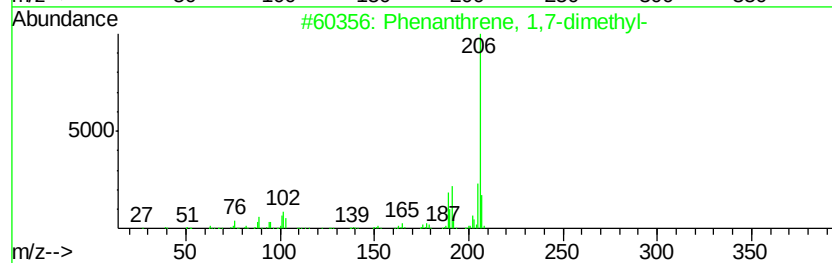
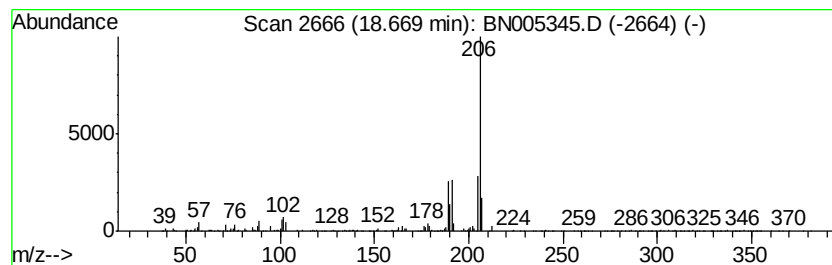
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 28 Phenanthrene, 1,7-dimethyl- Concentration Rank 24

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 18.67 | 4.42 ng/ul | 1533390 | Phenanthrene-d10 | 16.99 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenanthrene, 1,7-dimethyl- | 206 | C16H14  | 000483-87-4 | 94   |
| 2    |    | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 91   |
| 3    |    | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 91   |
| 4    |    | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 87   |
| 5    |    | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
Data File : BN005345.D  
Acq On : 27 Apr 2019 08:00  
Operator : JU/SJ  
Sample : K2456-12DL 5X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION

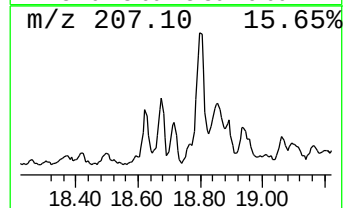
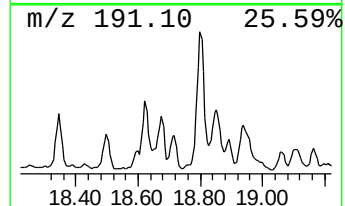
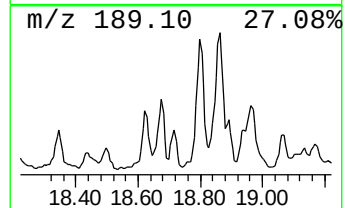
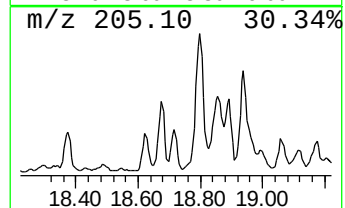
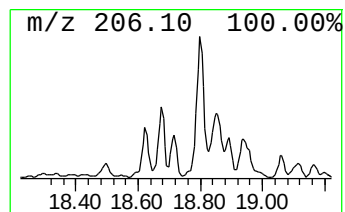
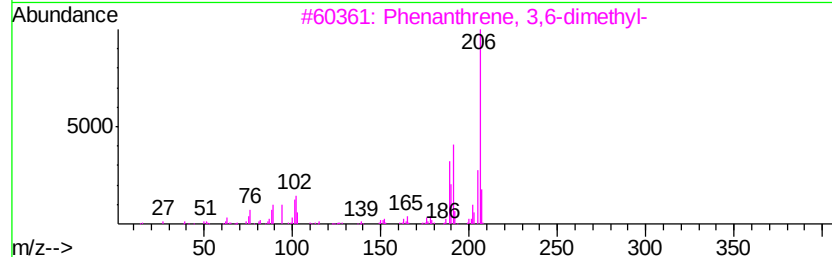
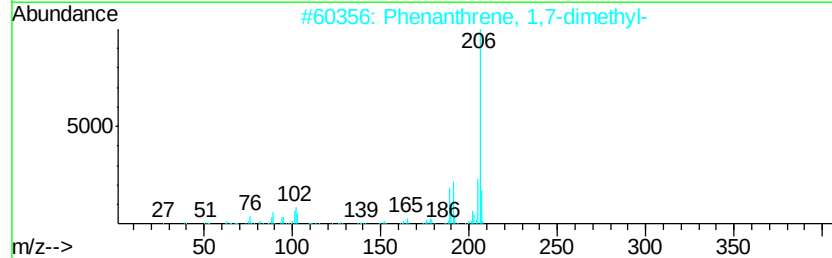
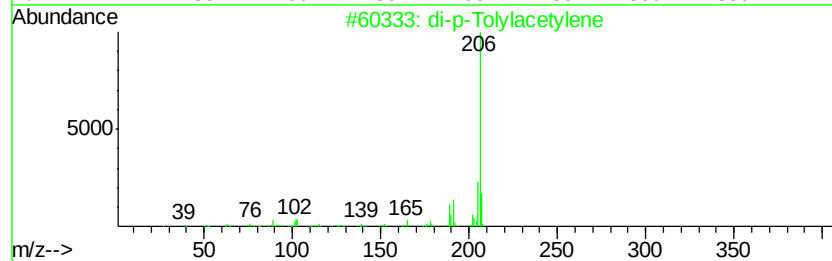
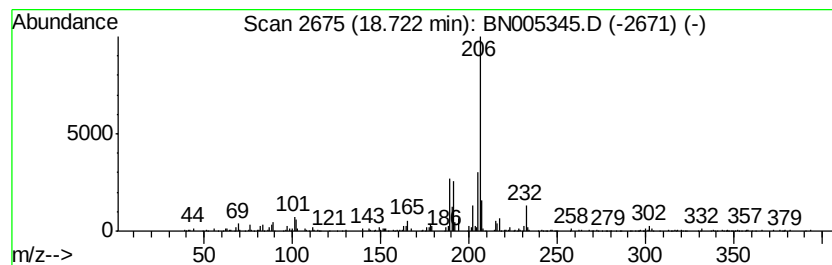
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 29 di-p-Tolylacetylene Concentration Rank 42

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.72 | 2.68 ng/ul | 929890 | Phenanthrene-d10 | 16.99 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 94   |
| 2    |    |   | Phenanthrene, 1,7-dimethyl- | 206 | C16H14  | 000483-87-4 | 93   |
| 3    |    |   | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 93   |
| 4    |    |   | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 93   |
| 5    |    |   | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042619\  
 Data File : BN005345.D  
 Acq On : 27 Apr 2019 08:00  
 Operator : JU/SJ  
 Sample : K2456-12DL 5X  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 GAHQ0DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.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 |
| 1,3,5,7-Cycloocta... | 5.63  | 4.8     | ng/ul | 750210   | 1                     | 7.59  | 3161080 | 20.0 |
| Benzene, 1-etheny... | 7.26  | 8.6     | ng/ul | 1361610  | 1                     | 7.59  | 3161080 | 20.0 |
| Benzene, cyclopro... | 7.34  | 3.0     | ng/ul | 479266   | 1                     | 7.59  | 3161080 | 20.0 |
| Benzene, 4-etheny... | 8.85  | 4.8     | ng/ul | 753095   | 1                     | 7.59  | 3161080 | 20.0 |
| 1,4-Methanonaphth... | 11.55 | 2.7     | ng/ul | 636686   | 2                     | 10.36 | 4759760 | 20.0 |
| 1H-Indene, 1-ethy... | 11.69 | 3.3     | ng/ul | 782245   | 2                     | 10.36 | 4759760 | 20.0 |
| Naphthalene, 2,7-... | 13.20 | 2.7     | ng/ul | 823715   | 3                     | 14.23 | 6035580 | 20.0 |
| 1,4-Ethenonaphtha... | 13.46 | 3.0     | ng/ul | 903637   | 3                     | 14.23 | 6035580 | 20.0 |
| (DEL) Alkane: Str... | 14.16 | 2.4     | ng/ul | 724955   | 3                     | 14.23 | 6035580 | 20.0 |
| 1,1'-Biphenyl, 2-... | 14.77 | 2.2     | ng/ul | 674696   | 3                     | 14.23 | 6035580 | 20.0 |
| unknown-01           | 14.89 | 3.2     | ng/ul | 976118   | 3                     | 14.23 | 6035580 | 20.0 |
| 1,1'-Biphenyl, 4-... | 14.94 | 2.9     | ng/ul | 885365   | 3                     | 14.23 | 6035580 | 20.0 |
| 1,1'-Biphenyl, 3-... | 15.00 | 4.0     | ng/ul | 1213370  | 3                     | 14.23 | 6035580 | 20.0 |
| Benzaldehyde, 2,4... | 15.11 | 7.8     | ng/ul | 2350110  | 3                     | 14.23 | 6035580 | 20.0 |
| 2-Naphthyl methyl... | 15.51 | 7.3     | ng/ul | 2204610  | 3                     | 14.23 | 6035580 | 20.0 |
| (DEL) Alkane: Str... | 15.97 | 5.9     | ng/ul | 2033180  | 4                     | 16.99 | 6934130 | 20.0 |
| (DEL) Alkane: Str... | 16.83 | 2.0     | ng/ul | 705764   | 4                     | 16.99 | 6934130 | 20.0 |
| 1H-Indene, 1-(phe... | 17.51 | 4.2     | ng/ul | 1457400  | 4                     | 16.99 | 6934130 | 20.0 |
| Anthracene, 2-met... | 17.86 | 6.7     | ng/ul | 2317840  | 4                     | 16.99 | 6934130 | 20.0 |
| Phenanthrene, 2-m... | 17.91 | 10.0    | ng/ul | 3458890  | 4                     | 16.99 | 6934130 | 20.0 |
| Phenanthrene, 3-m... | 17.99 | 4.7     | ng/ul | 1624640  | 4                     | 16.99 | 6934130 | 20.0 |
| Phenanthrene, 4-m... | 18.06 | 14.0    | ng/ul | 4849380  | 4                     | 16.99 | 6934130 | 20.0 |
| 2-Phenyl naphthalene | 18.37 | 3.4     | ng/ul | 1191050  | 4                     | 16.99 | 6934130 | 20.0 |
| Phenanthrene, 2,3... | 18.62 | 3.4     | ng/ul | 1167820  | 4                     | 16.99 | 6934130 | 20.0 |
| Phenanthrene, 1,7... | 18.67 | 4.4     | ng/ul | 1533390  | 4                     | 16.99 | 6934130 | 20.0 |
| di-p-Tolylacetylene  | 18.72 | 2.7     | ng/ul | 929890   | 4                     | 16.99 | 6934130 | 20.0 |