

Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
 Data File : BG032588.D  
 Acq On : 7 Feb 2018 7:26  
 Operator : SJ/JU  
 Sample : J1455-08  
 Misc : MDL-W-8 ppm  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 LW-03-OW

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG012918.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 3.364       | 8             | 13          | 22           | rBV2     | 17272          | 39767         | 1.81%           | 0.240%        |
| 2         | 3.458       | 24            | 29          | 40           | rVB      | 65937          | 116340        | 5.29%           | 0.701%        |
| 3         | 6.143       | 480           | 486         | 503          | rBV      | 100411         | 217307        | 9.89%           | 1.310%        |
| 4         | 7.824       | 765           | 772         | 778          | rBV2     | 61082          | 125554        | 5.71%           | 0.757%        |
| 5         | 7.883       | 778           | 782         | 793          | rVB      | 41886          | 77723         | 3.54%           | 0.469%        |
| 6         | 8.065       | 806           | 813         | 821          | rBV      | 21457          | 36137         | 1.64%           | 0.218%        |
| 7         | 8.223       | 832           | 840         | 862          | rBV      | 264630         | 547968        | 24.93%          | 3.304%        |
| 8         | 8.711       | 914           | 923         | 934          | rBV3     | 50259          | 120521        | 5.48%           | 0.727%        |
| 9         | 9.040       | 966           | 979         | 984          | rBV2     | 292641         | 592073        | 26.94%          | 3.570%        |
| 10        | 9.093       | 984           | 988         | 997          | rVV      | 112529         | 207995        | 9.46%           | 1.254%        |
| 11        | 9.199       | 1000          | 1006        | 1012         | rVB3     | 20498          | 35511         | 1.62%           | 0.214%        |
| 12        | 9.340       | 1024          | 1030        | 1036         | rBV3     | 28473          | 54262         | 2.47%           | 0.327%        |
| 13        | 9.404       | 1036          | 1041        | 1049         | rVB3     | 18729          | 32785         | 1.49%           | 0.198%        |
| 14        | 9.522       | 1053          | 1061        | 1068         | rBV2     | 14028          | 24909         | 1.13%           | 0.150%        |
| 15        | 9.833       | 1108          | 1114        | 1120         | rBV3     | 66026          | 129228        | 5.88%           | 0.779%        |
| 16        | 9.910       | 1120          | 1127        | 1137         | rVB2     | 233439         | 538395        | 24.50%          | 3.246%        |
| 17        | 10.186      | 1167          | 1174        | 1178         | rBV4     | 14030          | 31853         | 1.45%           | 0.192%        |
| 18        | 10.368      | 1199          | 1205        | 1210         | rBV      | 78209          | 138276        | 6.29%           | 0.834%        |
| 19        | 10.427      | 1210          | 1215        | 1228         | rVB2     | 112766         | 224313        | 10.21%          | 1.352%        |
| 20        | 10.744      | 1263          | 1269        | 1275         | rBV5     | 28768          | 63526         | 2.89%           | 0.383%        |
| 21        | 10.803      | 1275          | 1279        | 1292         | rVB5     | 19519          | 48210         | 2.19%           | 0.291%        |
| 22        | 10.961      | 1294          | 1306        | 1312         | rBV      | 233539         | 456852        | 20.79%          | 2.754%        |
| 23        | 11.014      | 1313          | 1315        | 1322         | rVB2     | 16573          | 22782         | 1.04%           | 0.137%        |
| 24        | 11.202      | 1339          | 1347        | 1359         | rBV3     | 32592          | 72540         | 3.30%           | 0.437%        |
| 25        | 11.437      | 1378          | 1387        | 1391         | rBV4     | 15242          | 32266         | 1.47%           | 0.195%        |
| 26        | 11.549      | 1391          | 1406        | 1408         | rBV2     | 62550          | 130455        | 5.94%           | 0.786%        |
| 27        | 11.590      | 1408          | 1413        | 1418         | rBV2     | 64720          | 120903        | 5.50%           | 0.729%        |
| 28        | 11.684      | 1424          | 1429        | 1437         | rVB2     | 65231          | 129596        | 5.90%           | 0.781%        |
| 29        | 11.772      | 1437          | 1444        | 1453         | rVB4     | 26037          | 48081         | 2.19%           | 0.290%        |
| 30        | 11.937      | 1466          | 1472        | 1484         | rVB7     | 14623          | 41772         | 1.90%           | 0.252%        |
| 31        | 12.042      | 1484          | 1490        | 1497         | rBV2     | 31567          | 63588         | 2.89%           | 0.383%        |
| 32        | 12.154      | 1501          | 1509        | 1512         | rBV3     | 33326          | 63166         | 2.87%           | 0.381%        |
| 33        | 12.225      | 1517          | 1521        | 1534         | rVB5     | 27457          | 78174         | 3.56%           | 0.471%        |
| 34        | 12.471      | 1557          | 1563        | 1573         | rVB      | 65940          | 115672        | 5.26%           | 0.697%        |

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 Misc : MDL-W-8 ppm  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 LW-03-OW

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG012918.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |         |        |        |
|----|--------|------|------|------|------|--------|---------|--------|--------|
| 35 | 12.659 | 1586 | 1595 | 1603 | rBV2 | 55748  | 125340  | 5.70%  | 0.756% |
| 36 | 12.742 | 1604 | 1609 | 1615 | rVV4 | 21083  | 44313   | 2.02%  | 0.267% |
| 37 | 12.859 | 1626 | 1629 | 1635 | rVB  | 23690  | 39989   | 1.82%  | 0.241% |
| 38 | 12.930 | 1635 | 1641 | 1649 | rBV4 | 41309  | 103418  | 4.71%  | 0.623% |
| 39 | 13.018 | 1650 | 1656 | 1660 | rVV3 | 20060  | 48079   | 2.19%  | 0.290% |
| 40 | 13.088 | 1663 | 1668 | 1675 | rVV3 | 68956  | 131973  | 6.00%  | 0.796% |
| 41 | 13.153 | 1675 | 1679 | 1682 | rVV2 | 20509  | 28290   | 1.29%  | 0.171% |
| 42 | 13.417 | 1713 | 1724 | 1732 | rBV  | 121294 | 296228  | 13.48% | 1.786% |
| 43 | 13.482 | 1732 | 1735 | 1745 | rVB5 | 27496  | 74103   | 3.37%  | 0.447% |
| 44 | 13.693 | 1767 | 1771 | 1778 | rVB3 | 23402  | 47518   | 2.16%  | 0.286% |
| 45 | 13.970 | 1810 | 1818 | 1831 | rBV  | 787867 | 1347510 | 61.31% | 8.124% |
| 46 | 14.122 | 1840 | 1844 | 1850 | rBV9 | 10946  | 23963   | 1.09%  | 0.144% |
| 47 | 14.193 | 1854 | 1856 | 1861 | rVV6 | 14743  | 25436   | 1.16%  | 0.153% |
| 48 | 14.275 | 1864 | 1870 | 1876 | rVV3 | 37169  | 72452   | 3.30%  | 0.437% |
| 49 | 14.404 | 1886 | 1892 | 1898 | rVB2 | 35685  | 66595   | 3.03%  | 0.401% |
| 50 | 14.516 | 1900 | 1911 | 1922 | rBV9 | 77407  | 278898  | 12.69% | 1.681% |
| 51 | 14.663 | 1929 | 1936 | 1940 | rBV  | 100739 | 191535  | 8.71%  | 1.155% |
| 52 | 14.716 | 1940 | 1945 | 1949 | rVV3 | 85801  | 159256  | 7.25%  | 0.960% |
| 53 | 14.763 | 1951 | 1953 | 1961 | rVB2 | 31265  | 56963   | 2.59%  | 0.343% |
| 54 | 14.892 | 1970 | 1975 | 1978 | rBV2 | 60746  | 100727  | 4.58%  | 0.607% |
| 55 | 14.939 | 1978 | 1983 | 1985 | rVV4 | 35263  | 59194   | 2.69%  | 0.357% |
| 56 | 15.062 | 1993 | 2004 | 2009 | rBV3 | 57944  | 141839  | 6.45%  | 0.855% |
| 57 | 15.133 | 2013 | 2016 | 2021 | rVV6 | 27504  | 51447   | 2.34%  | 0.310% |
| 58 | 15.339 | 2040 | 2051 | 2058 | rBV2 | 148499 | 372623  | 16.95% | 2.246% |
| 59 | 15.403 | 2058 | 2062 | 2067 | rBV3 | 54878  | 106728  | 4.86%  | 0.643% |
| 60 | 15.456 | 2067 | 2071 | 2075 | rVB5 | 23241  | 34813   | 1.58%  | 0.210% |
| 61 | 15.521 | 2075 | 2082 | 2085 | rBV4 | 59781  | 143076  | 6.51%  | 0.863% |
| 62 | 15.638 | 2098 | 2102 | 2106 | rVB5 | 35596  | 43880   | 2.00%  | 0.265% |
| 63 | 15.720 | 2108 | 2116 | 2123 | rVV4 | 51172  | 169397  | 7.71%  | 1.021% |
| 64 | 15.785 | 2123 | 2127 | 2131 | rVV3 | 29979  | 63823   | 2.90%  | 0.385% |
| 65 | 15.832 | 2131 | 2135 | 2141 | rVV4 | 54523  | 120309  | 5.47%  | 0.725% |
| 66 | 15.903 | 2143 | 2147 | 2149 | rVV4 | 31823  | 54454   | 2.48%  | 0.328% |
| 67 | 15.944 | 2149 | 2154 | 2159 | rVV4 | 86418  | 175487  | 7.98%  | 1.058% |
| 68 | 15.985 | 2159 | 2161 | 2168 | rVB6 | 14076  | 24437   | 1.11%  | 0.147% |
| 69 | 16.096 | 2175 | 2180 | 2181 | rBV2 | 34736  | 49727   | 2.26%  | 0.300% |
| 70 | 16.155 | 2185 | 2190 | 2198 | rVV5 | 91963  | 206488  | 9.40%  | 1.245% |
| 71 | 16.226 | 2199 | 2202 | 2205 | rVV  | 19780  | 27969   | 1.27%  | 0.169% |

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 ALS Vial : 22 Sample Multiplier: 1

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 BNA\_G  
 ClientSampleId :  
 LW-03-OW

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0  
 Filtering: 5  
 Min Area: 3 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG012918.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |       |         |         |         |         |
|----|--------|------|------|------|-------|---------|---------|---------|---------|
| 72 | 16.373 | 2221 | 2227 | 2231 | rBV4  | 47218   | 108572  | 4.94%   | 0.655%  |
| 73 | 16.508 | 2245 | 2250 | 2256 | rVB   | 81709   | 152412  | 6.93%   | 0.919%  |
| 74 | 16.578 | 2256 | 2262 | 2268 | rVB8  | 38097   | 94469   | 4.30%   | 0.570%  |
| 75 | 16.708 | 2274 | 2284 | 2294 | rBV2  | 354870  | 818854  | 37.26%  | 4.937%  |
| 76 | 16.819 | 2297 | 2303 | 2311 | rBV   | 644447  | 1099307 | 50.02%  | 6.628%  |
| 77 | 17.054 | 2337 | 2343 | 2350 | rBV8  | 35536   | 81868   | 3.72%   | 0.494%  |
| 78 | 17.178 | 2360 | 2364 | 2370 | rVB9  | 21094   | 43219   | 1.97%   | 0.261%  |
| 79 | 17.348 | 2385 | 2393 | 2394 | rBV8  | 20765   | 35742   | 1.63%   | 0.215%  |
| 80 | 17.424 | 2401 | 2406 | 2417 | rBV5  | 42395   | 107303  | 4.88%   | 0.647%  |
| 81 | 17.812 | 2466 | 2472 | 2473 | rBV6  | 16117   | 24474   | 1.11%   | 0.148%  |
| 82 | 17.888 | 2481 | 2485 | 2486 | rBV4  | 19081   | 22923   | 1.04%   | 0.138%  |
| 83 | 18.076 | 2511 | 2517 | 2526 | rVB   | 230651  | 407174  | 18.53%  | 2.455%  |
| 84 | 18.400 | 2567 | 2572 | 2573 | rBV5  | 17149   | 26374   | 1.20%   | 0.159%  |
| 85 | 18.693 | 2619 | 2622 | 2631 | rVB2  | 54892   | 96167   | 4.38%   | 0.580%  |
| 86 | 18.899 | 2653 | 2657 | 2664 | rBV7  | 56271   | 91731   | 4.17%   | 0.553%  |
| 87 | 19.539 | 2760 | 2766 | 2767 | rBV6  | 21210   | 31571   | 1.44%   | 0.190%  |
| 88 | 19.816 | 2807 | 2813 | 2818 | rBV10 | 23577   | 57515   | 2.62%   | 0.347%  |
| 89 | 20.133 | 2863 | 2867 | 2871 | rVB6  | 38764   | 56378   | 2.57%   | 0.340%  |
| 90 | 20.344 | 2899 | 2903 | 2905 | rBV5  | 15050   | 22836   | 1.04%   | 0.138%  |
| 91 | 20.615 | 2943 | 2949 | 2957 | rVB   | 1682247 | 2197800 | 100.00% | 13.250% |
| 92 | 21.619 | 3117 | 3120 | 3124 | rBV6  | 12873   | 22478   | 1.02%   | 0.136%  |
| 93 | 21.943 | 3172 | 3175 | 3180 | rVB7  | 45054   | 72256   | 3.29%   | 0.436%  |
| 94 | 22.401 | 3246 | 3253 | 3263 | rVB   | 272478  | 510539  | 23.23%  | 3.078%  |
| 95 | 26.073 | 3865 | 3878 | 3890 | rBV   | 153196  | 518210  | 23.58%  | 3.124%  |

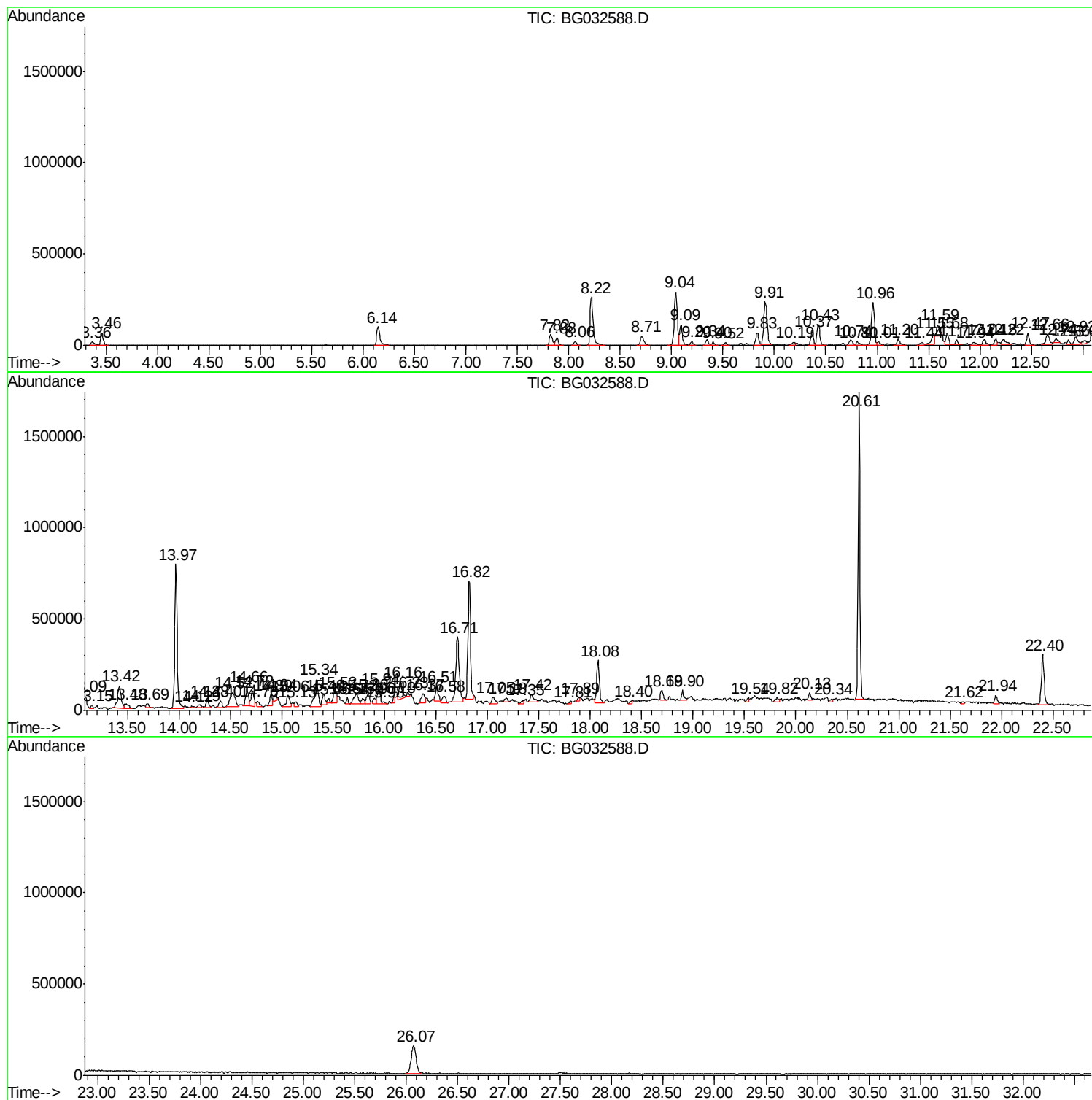
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ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-03-OW

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
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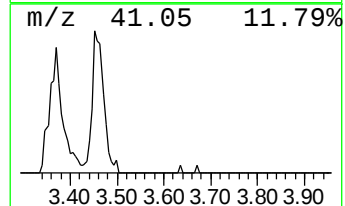
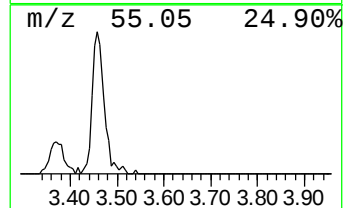
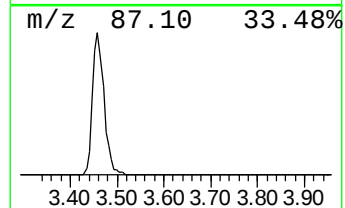
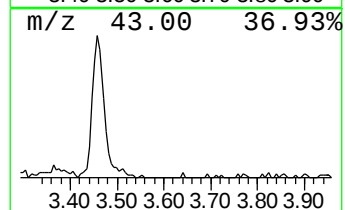
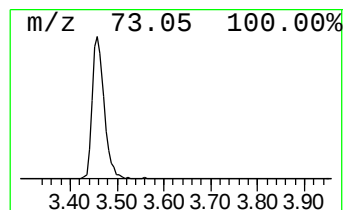
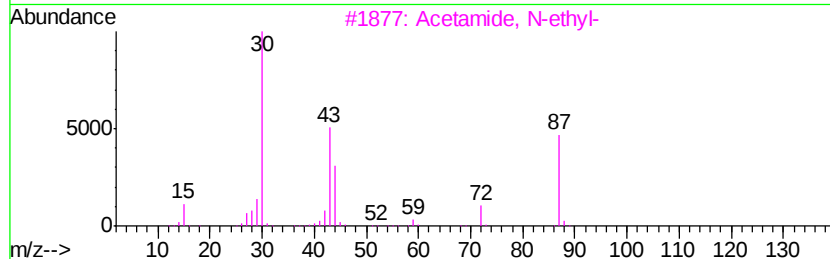
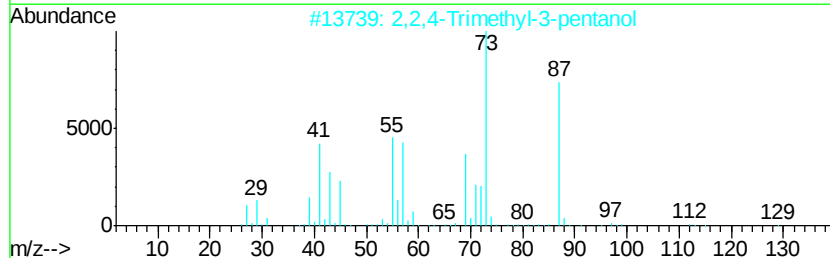
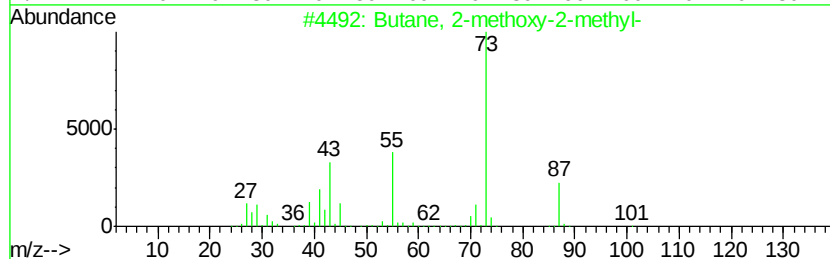
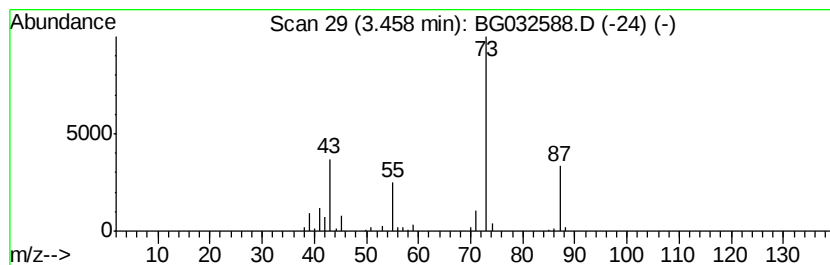
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 13

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 3.46 | 19.31 ng | 116340 | 1,4-Dichlorobenzene-d4 | 8.71 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 78   |
| 2    |    | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 40   |
| 3    |    | Acetamide, N-ethyl-         | 87  | C4H9NO  | 000625-50-3 | 16   |
| 4    |    | 2-Methyl-5-hexen-3-ol       | 114 | C7H14O  | 032815-70-6 | 12   |
| 5    |    | N-Ethylformamide            | 73  | C3H7NO  | 000627-45-2 | 9    |



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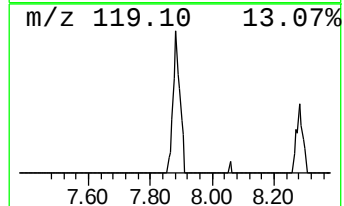
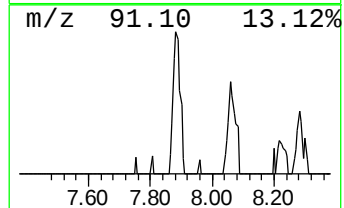
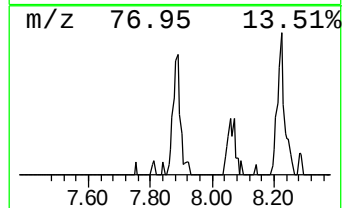
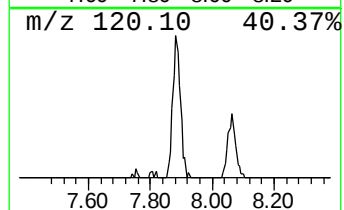
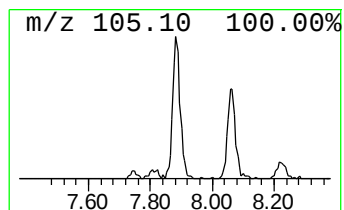
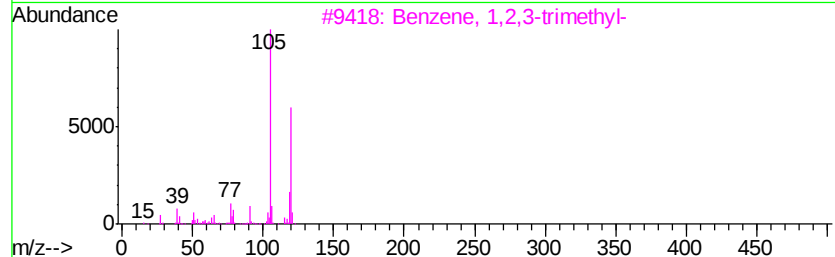
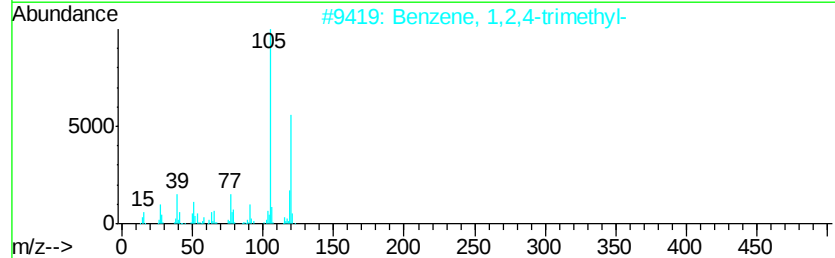
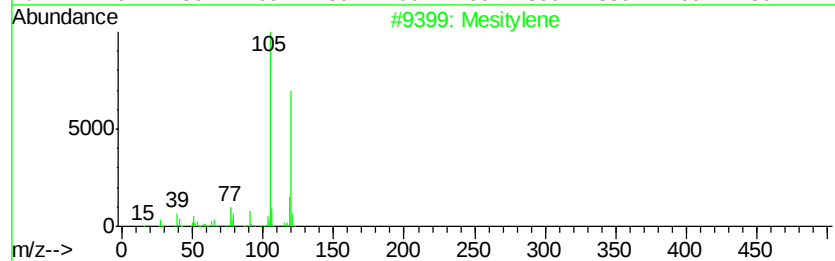
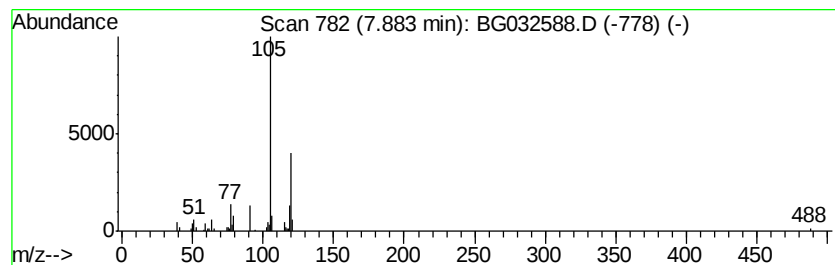
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 Mesitylene Concentration Rank 18

| R.T. | EstConc  | Area  | Relative to ISTD       | R.T. |
|------|----------|-------|------------------------|------|
| 7.88 | 12.90 ng | 77723 | 1,4-Dichlorobenzene-d4 | 8.71 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 91   |
| 2    |    | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 91   |
| 3    |    | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |
| 4    |    | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 87   |
| 5    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 87   |



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ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-03-OW

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

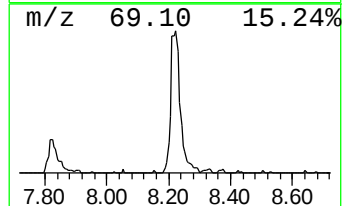
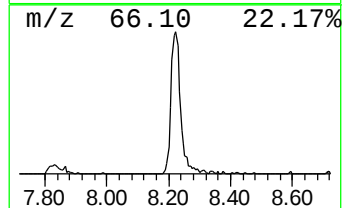
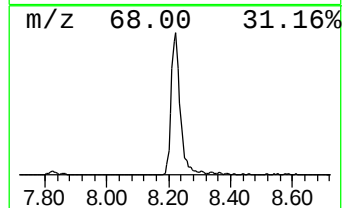
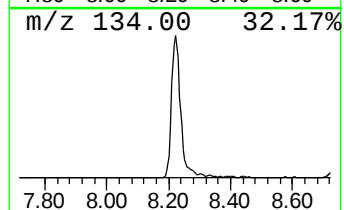
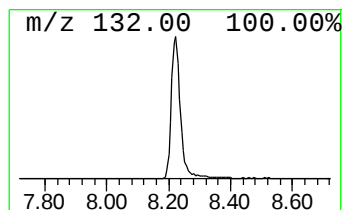
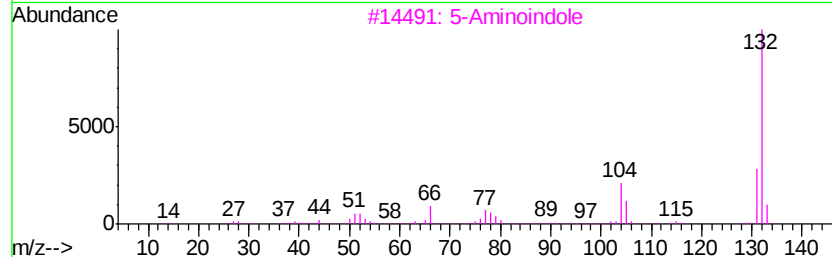
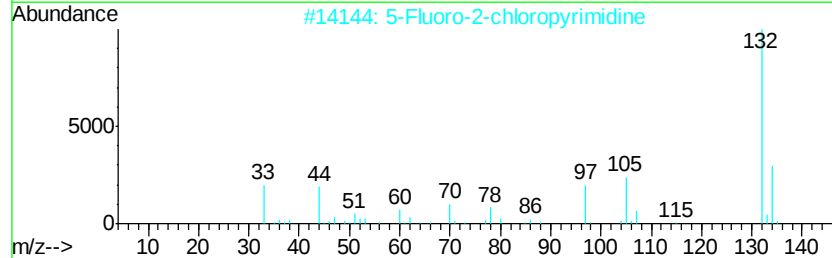
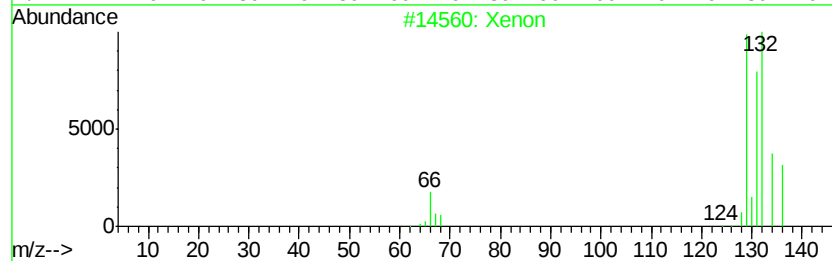
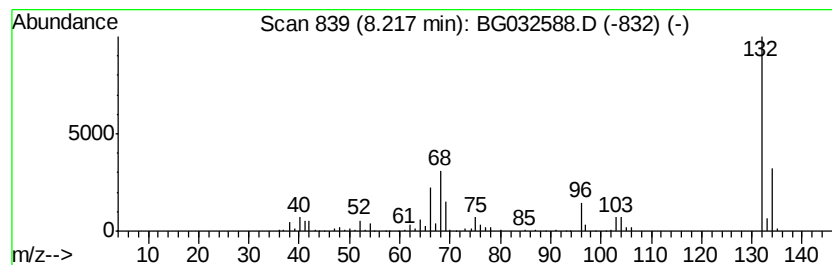
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Xenon Concentration Rank 2

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 8.22 | 90.93 ng | 547968 | 1,4-Dichlorobenzene-d4 | 8.71 |

| Hit# of | 5                                | Tentative ID | MW  | MolForm   | CAS#         | Qual |
|---------|----------------------------------|--------------|-----|-----------|--------------|------|
| 1       | Xenon                            |              | 132 | Xe        | 007440-63-3  | 50   |
| 2       | 5-Fluoro-2-chloropyrimidine      |              | 132 | C4H2ClFN2 | 062802-42-0  | 17   |
| 3       | 5-Aminoindole                    |              | 132 | C8H8N2    | 005192-03-0  | 16   |
| 4       | (E)-3-Chloro-2-methyl-2-pentenal |              | 132 | C6H9ClO   | 031357-76-3  | 12   |
| 5       | Tranylcypromine-propionyl        |              | 189 | C12H15NO  | 1000123-86-3 | 12   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
Data File : BG032588.D  
Acq On : 7 Feb 2018 7:26  
Operator : SJ/JU  
Sample : J1455-08  
Misc : MDL-W-8 ppm  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-03-OW

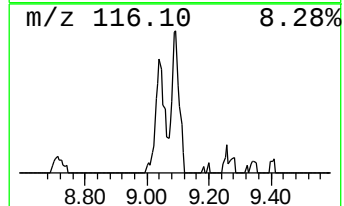
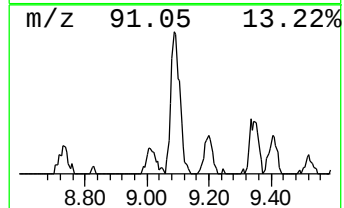
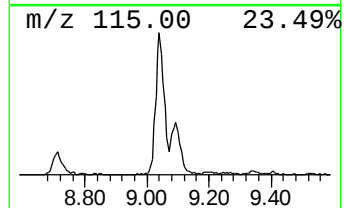
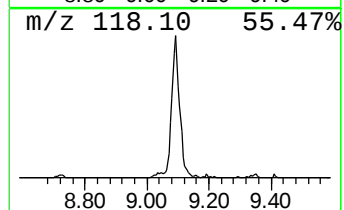
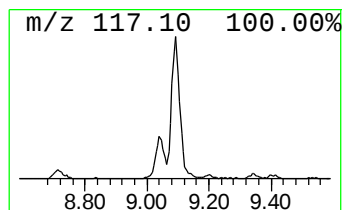
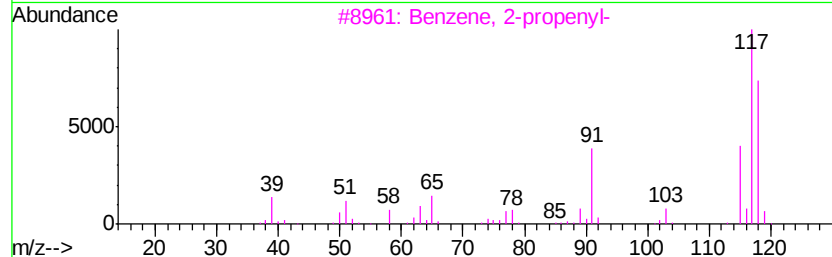
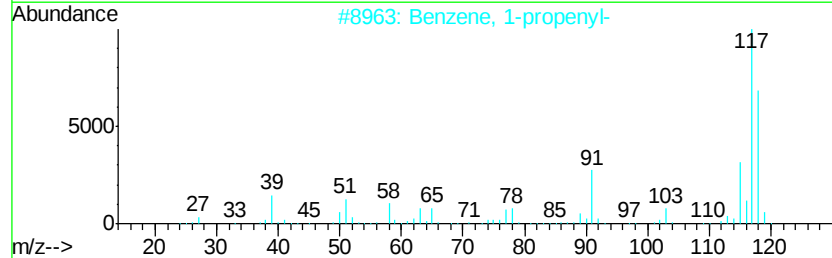
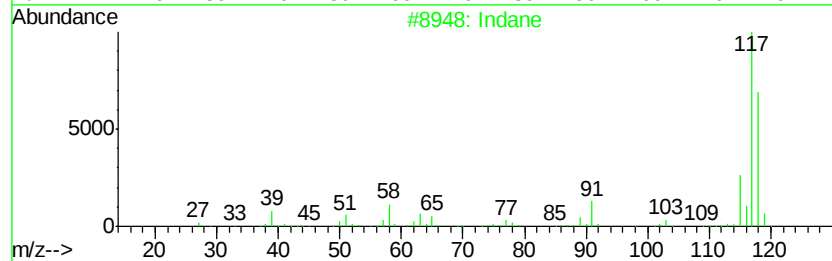
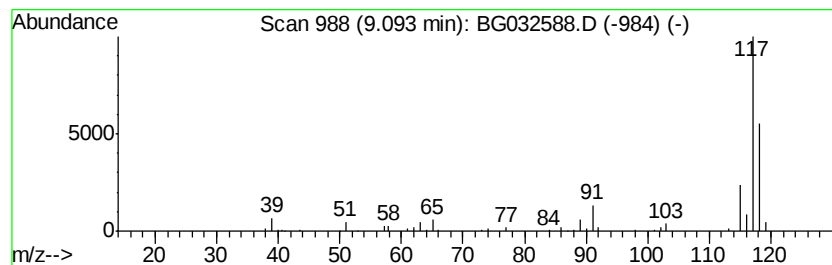
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 Indane Concentration Rank 7

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 9.09 | 34.52 ng | 207995 | 1,4-Dichlorobenzene-d4 | 8.71 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#         | Qual |
|------|----|----------------------------------|-----|---------|--------------|------|
| 1    | 5  | Indane                           | 118 | C9H10   | 000496-11-7  | 87   |
| 2    |    | Benzene, 1-propenyl-             | 118 | C9H10   | 000637-50-3  | 83   |
| 3    |    | Benzene, 2-propenyl-             | 118 | C9H10   | 000300-57-2  | 83   |
| 4    |    | Tetracyclo[3.3.1.0(2,8).0(4,6)]- | 118 | C9H10   | 1000191-13-7 | 80   |
| 5    |    | (Z)-1-Phenylpropene              | 118 | C9H10   | 000766-90-5  | 74   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
Data File : BG032588.D  
Acq On : 7 Feb 2018 7:26  
Operator : SJ/JU  
Sample : J1455-08  
Misc : MDL-W-8 ppm  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-03-OW

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

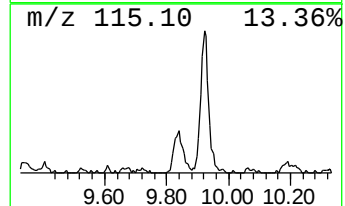
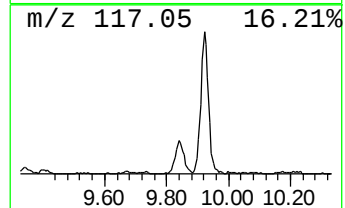
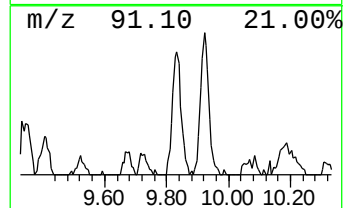
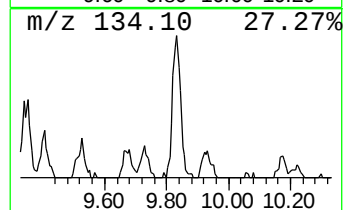
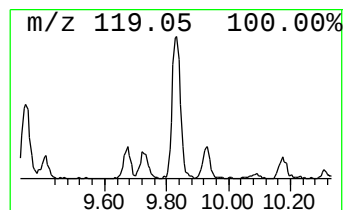
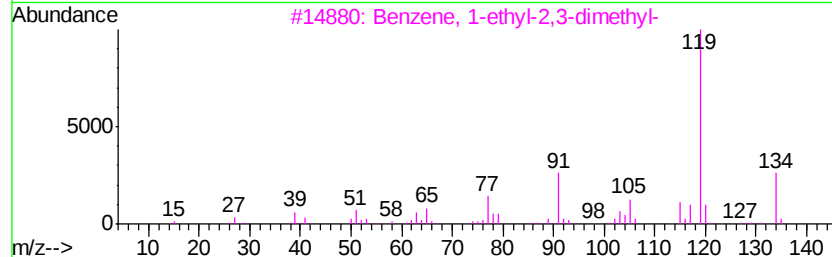
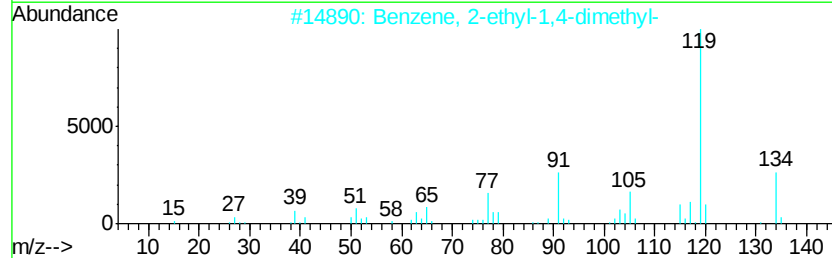
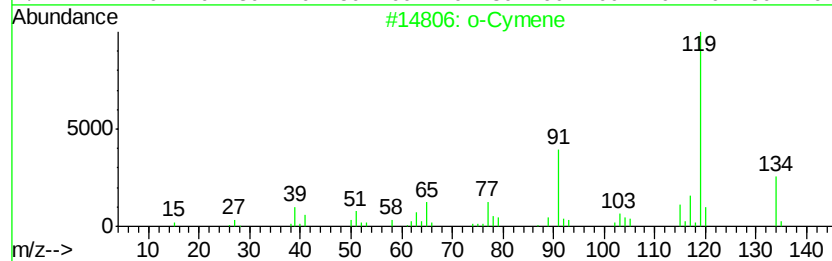
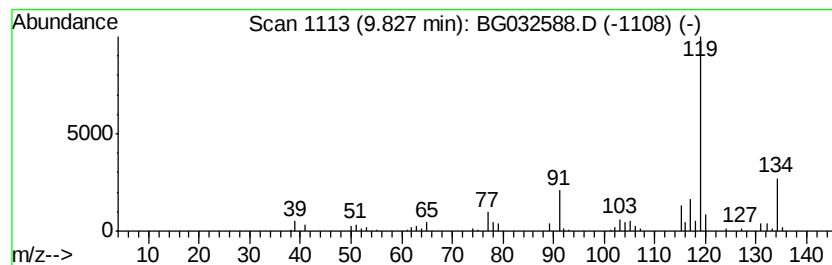
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 o-Cymene Concentration Rank 10

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 9.83 | 21.44 ng | 129228 | 1,4-Dichlorobenzene-d4 | 8.71 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 90   |
| 2    |    | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 87   |
| 3    |    | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 87   |
| 4    |    | p-Cymene                       | 134 | C10H14  | 000099-87-6 | 87   |
| 5    |    | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 87   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
Data File : BG032588.D  
Acq On : 7 Feb 2018 7:26  
Operator : SJ/JU  
Sample : J1455-08  
Misc : MDL-W-8 ppm  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-03-OW

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

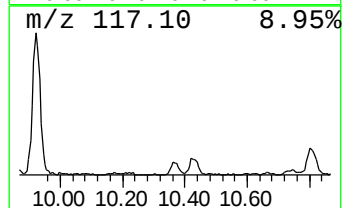
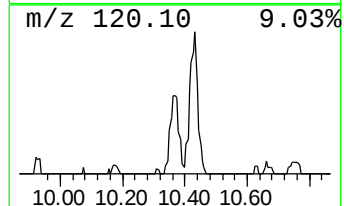
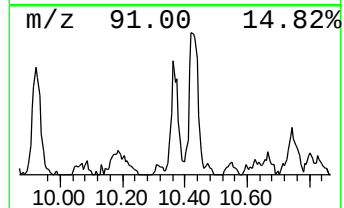
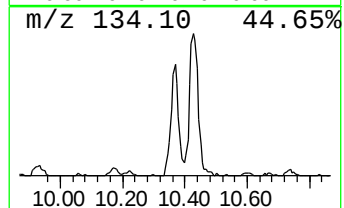
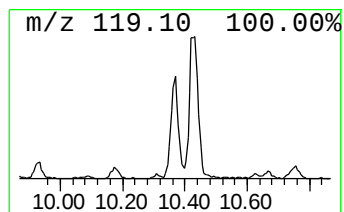
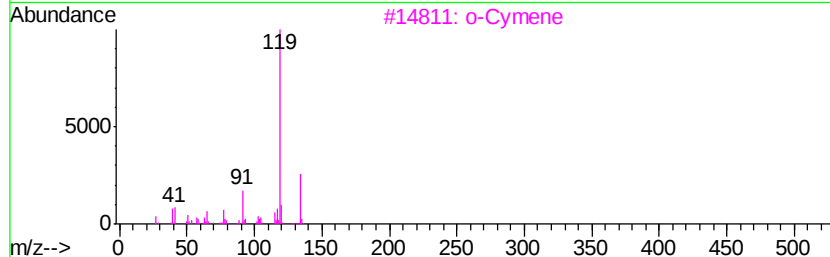
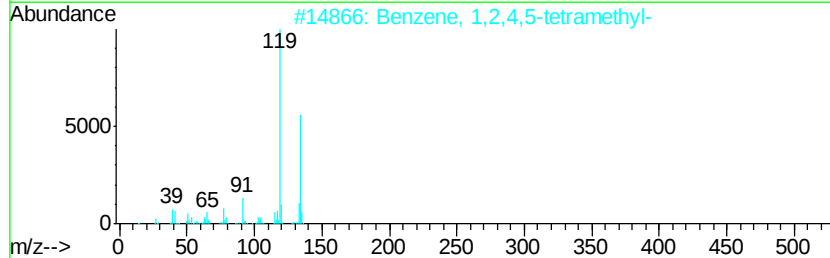
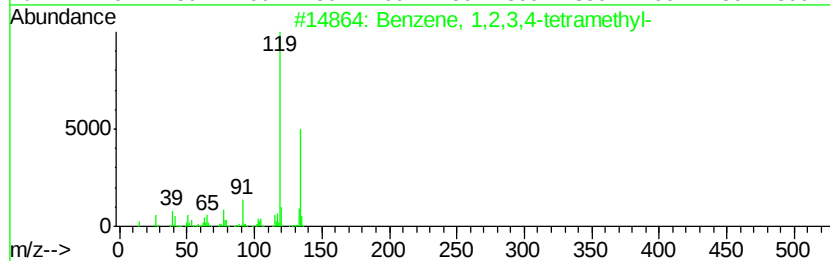
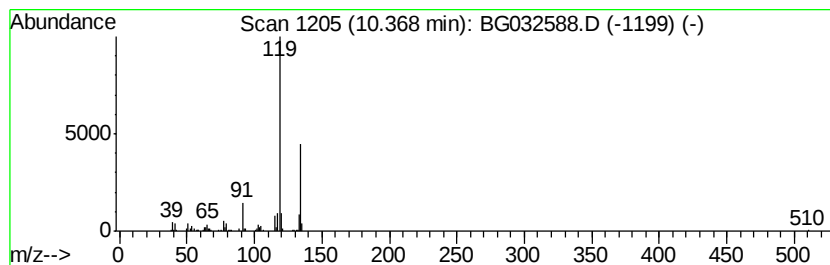
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 8

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 10.37 | 22.87 ng | 138276 | Naphthalene-d8   | 11.58 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 94   |
| 2    |    | Benzene, 1,2,4,5-tetramethyl-  | 134 | C10H14  | 000095-93-2 | 94   |
| 3    |    | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 93   |
| 4    |    | Benzene, 1,2,3,5-tetramethyl-  | 134 | C10H14  | 000527-53-7 | 91   |
| 5    |    | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 91   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
Data File : BG032588.D  
Acq On : 7 Feb 2018 7:26  
Operator : SJ/JU  
Sample : J1455-08  
Misc : MDL-W-8 ppm  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-03-OW

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

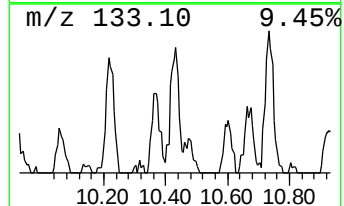
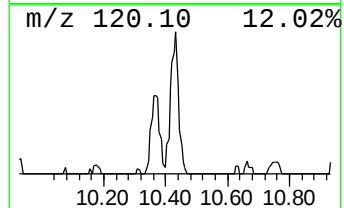
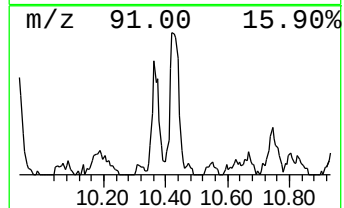
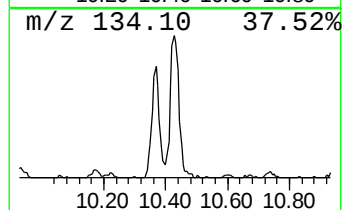
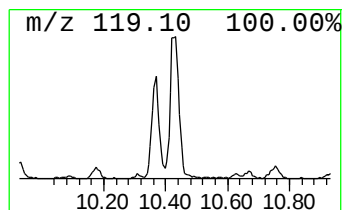
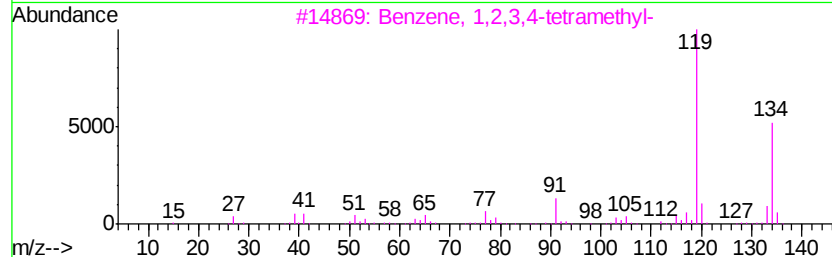
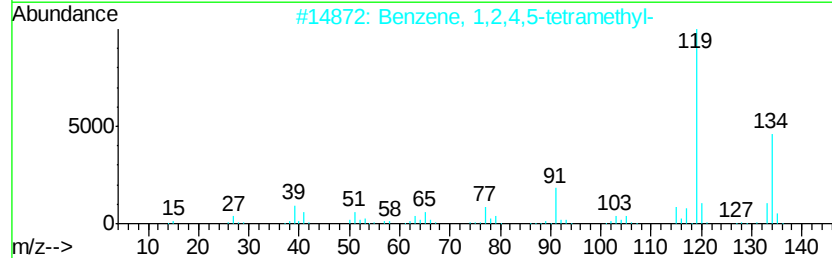
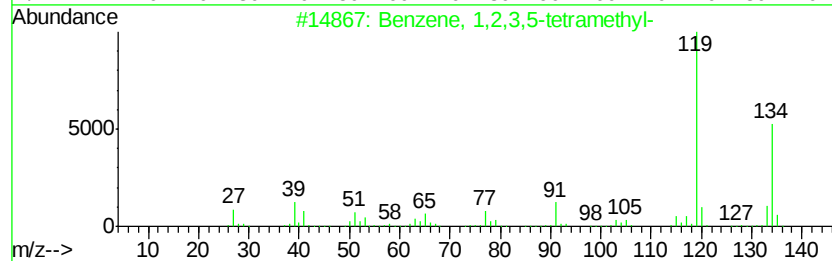
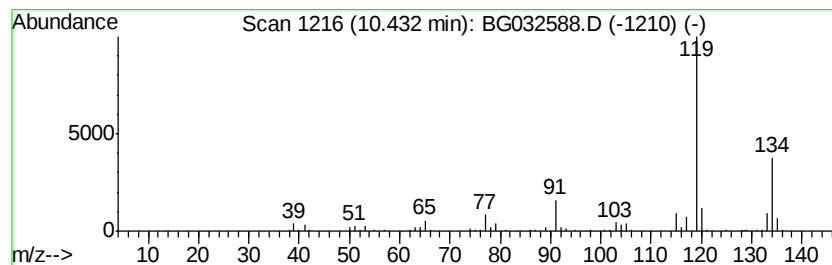
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 6

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 10.43 | 37.11 ng | 224313 | Naphthalene-d8   | 11.58 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3,5-tetramethyl-  | 134 | C10H14  | 000527-53-7 | 95   |
| 2    |    | Benzene, 1,2,4,5-tetramethyl-  | 134 | C10H14  | 000095-93-2 | 95   |
| 3    |    | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 94   |
| 4    |    | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 94   |
| 5    |    | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 94   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
Data File : BG032588.D  
Acq On : 7 Feb 2018 7:26  
Operator : SJ/JU  
Sample : J1455-08  
Misc : MDL-W-8 ppm  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-03-OW

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

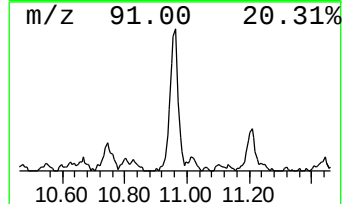
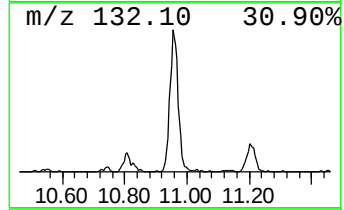
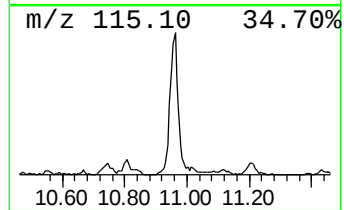
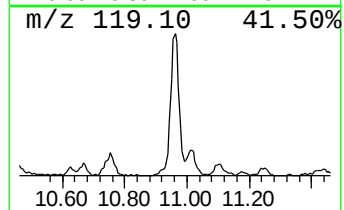
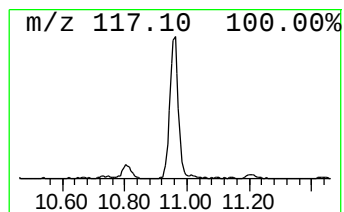
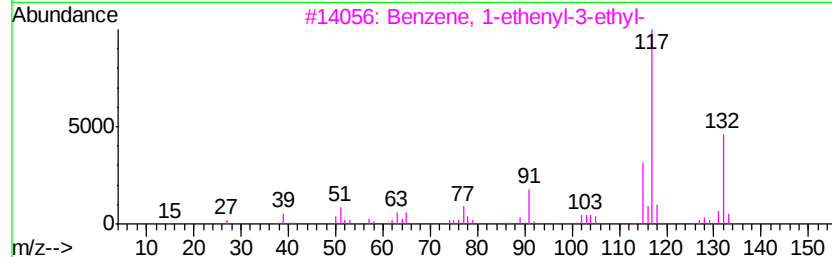
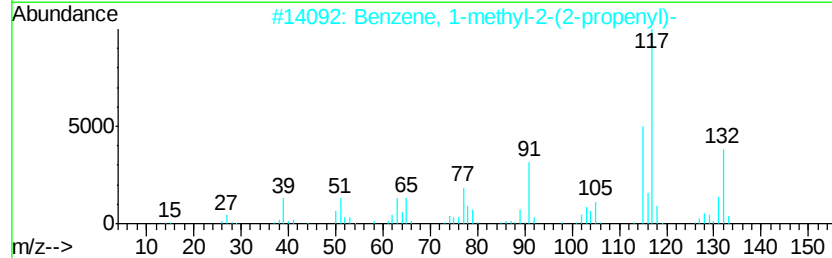
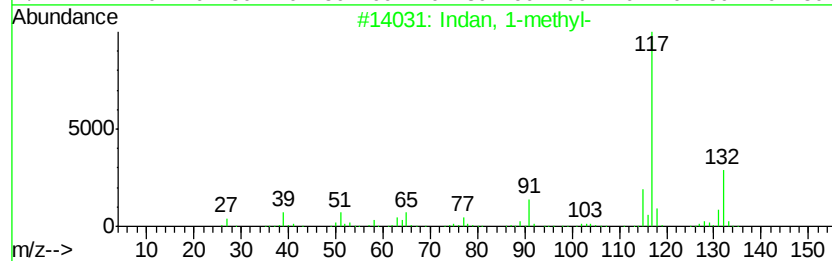
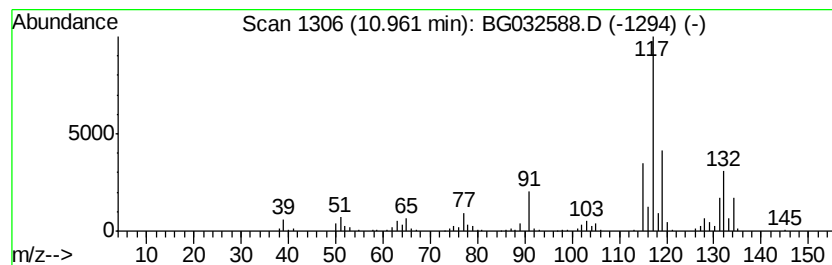
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Indan, 1-methyl- Concentration Rank 3

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 10.96 | 75.57 ng | 456852 | Naphthalene-d8   | 11.58 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Indan, 1-methyl-                    | 132 | C10H12  | 000767-58-8 | 76   |
| 2    |    | Benzene, 1-methyl-2-(2-propenyl)-   | 132 | C10H12  | 001587-04-8 | 76   |
| 3    |    | Benzene, 1-ethenyl-3-ethyl-         | 132 | C10H12  | 007525-62-4 | 70   |
| 4    |    | 1-Phenyl-1-butene                   | 132 | C10H12  | 000824-90-8 | 70   |
| 5    |    | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000768-00-3 | 68   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
Data File : BG032588.D  
Acq On : 7 Feb 2018 7:26  
Operator : SJ/JU  
Sample : J1455-08  
Misc : MDL-W-8 ppm  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-03-OW

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

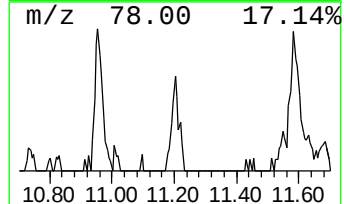
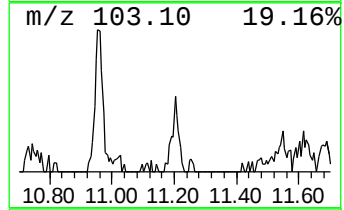
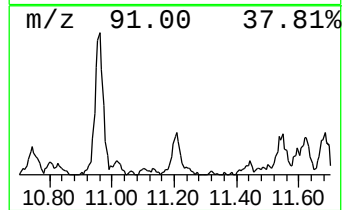
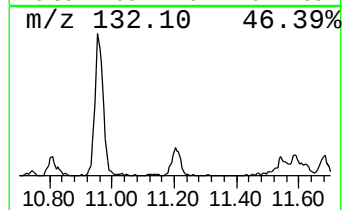
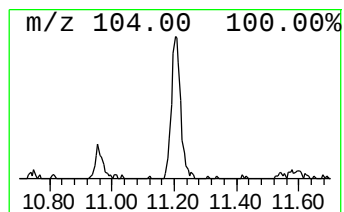
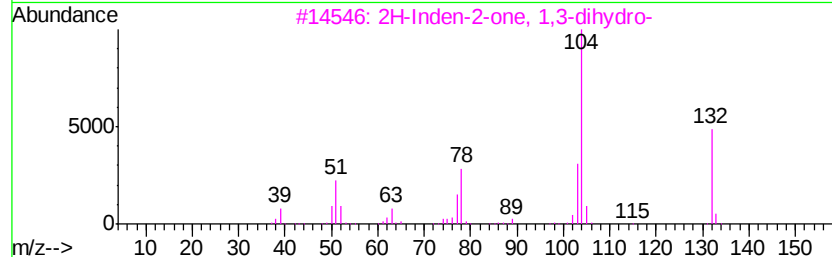
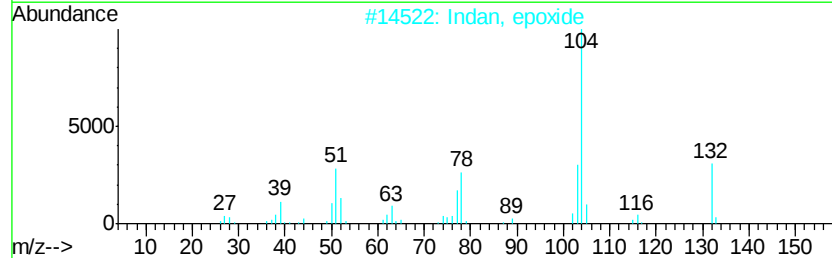
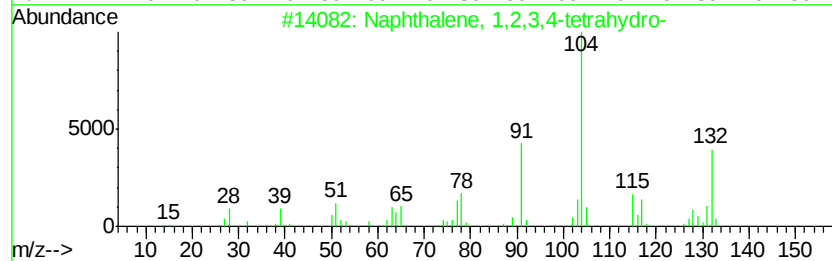
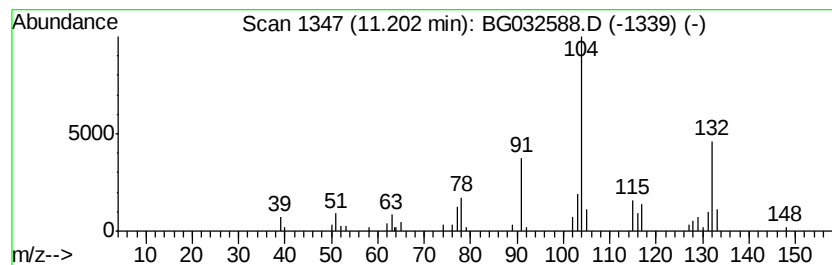
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 19

| R.T.  | EstConc  | Area  | Relative to ISTD | R.T.  |
|-------|----------|-------|------------------|-------|
| 11.20 | 12.00 ng | 72540 | Naphthalene-d8   | 11.58 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,2,3,4-tetrahydro-  | 132 | C10H12  | 000119-64-2 | 93   |
| 2    |    |   | Indan, epoxide                    | 132 | C9H8O   | 000768-22-9 | 58   |
| 3    |    |   | 2H-Inden-2-one, 1,3-dihydro-      | 132 | C9H8O   | 000615-13-4 | 52   |
| 4    |    |   | 1H-Inden-1-one, 2,3-dihydro-      | 132 | C9H8O   | 000083-33-0 | 49   |
| 5    |    |   | Isoquinoline, 1,2,3,4-tetrahydro- | 133 | C9H11N  | 000091-21-4 | 42   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
Data File : BG032588.D  
Acq On : 7 Feb 2018 7:26  
Operator : SJ/JU  
Sample : J1455-08  
Misc : MDL-W-8 ppm  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-03-OW

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

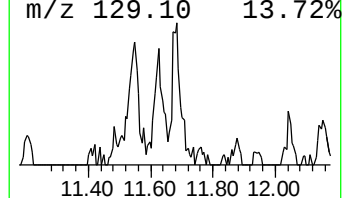
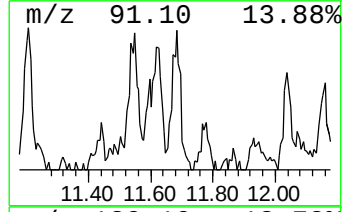
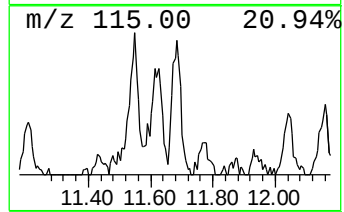
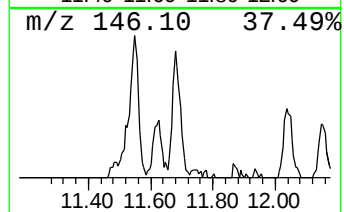
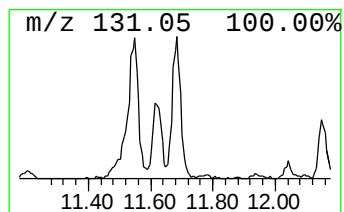
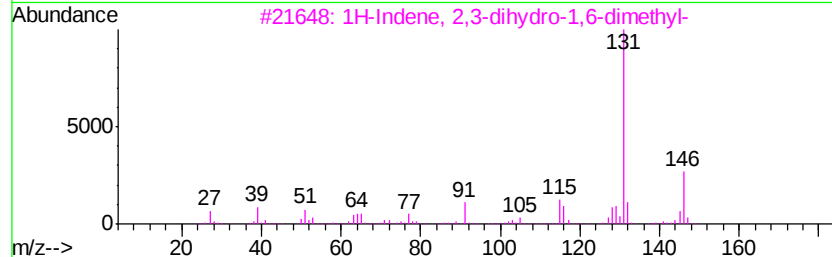
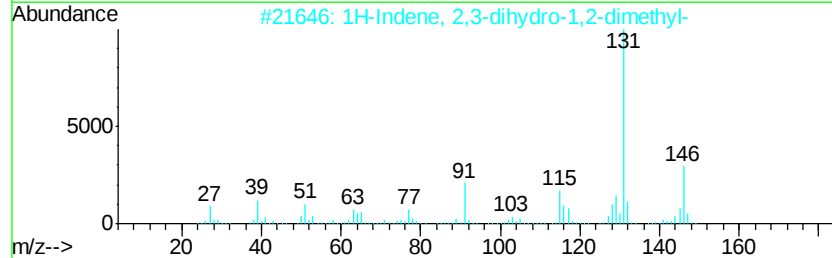
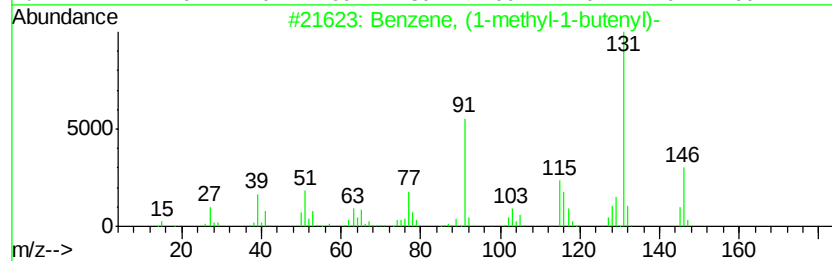
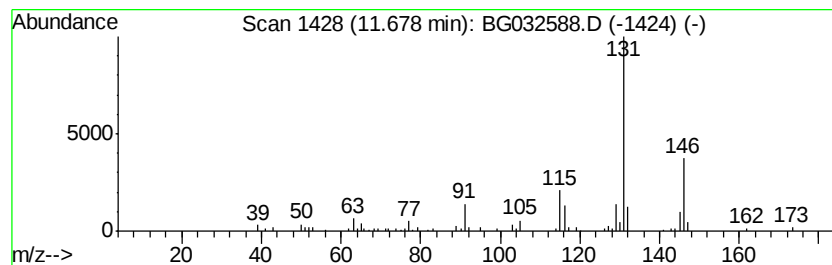
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Benzene, (1-methyl-1-butenyl)- Concentration Rank 11

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 11.68 | 21.44 ng | 129596 | Naphthalene-d8   | 11.58 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, (1-methyl-1-butenyl)-      | 146 | C11H14  | 053172-84-2 | 91   |
| 2    |    | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 | C11H14  | 017057-82-8 | 91   |
| 3    |    | 1H-Indene, 2,3-dihydro-1,6-dimet... | 146 | C11H14  | 017059-48-2 | 90   |
| 4    |    | Benzene, (1,2-dimethyl-1-propenyl)- | 146 | C11H14  | 000769-57-3 | 90   |
| 5    |    | Benzene, (2-methyl-1-butenyl)-      | 146 | C11H14  | 056253-64-6 | 90   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
Data File : BG032588.D  
Acq On : 7 Feb 2018 7:26  
Operator : SJ/JU  
Sample : J1455-08  
Misc : MDL-W-8 ppm  
ALS Vial : 22 Sample Multiplier: 1

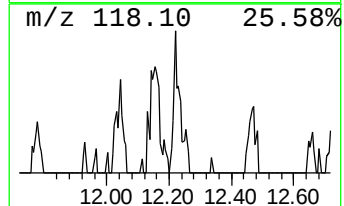
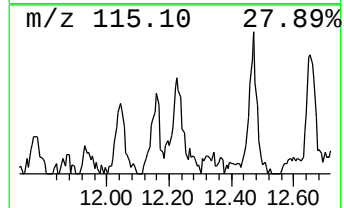
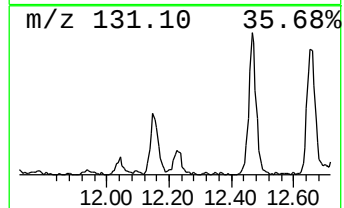
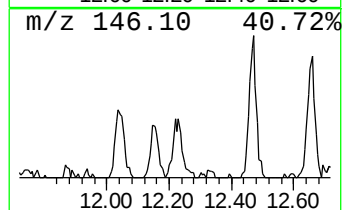
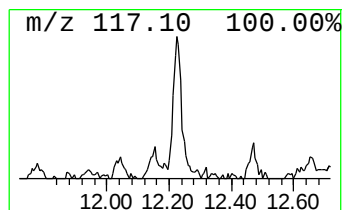
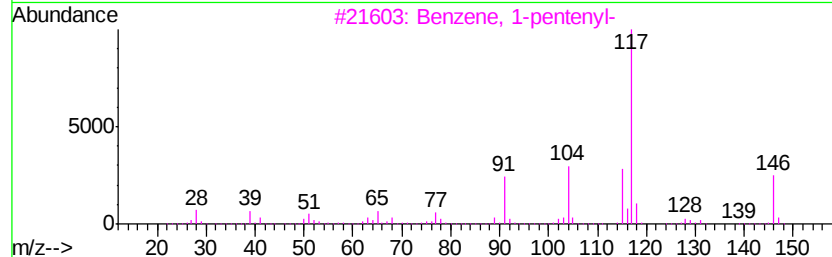
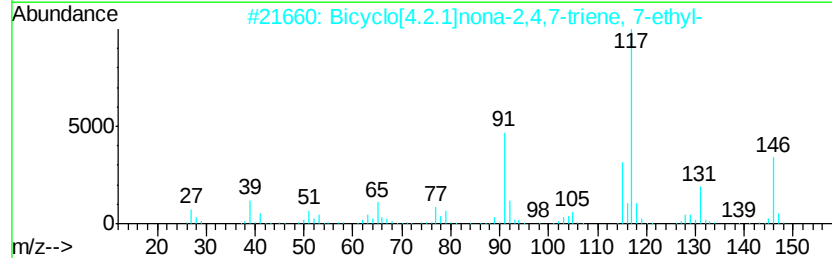
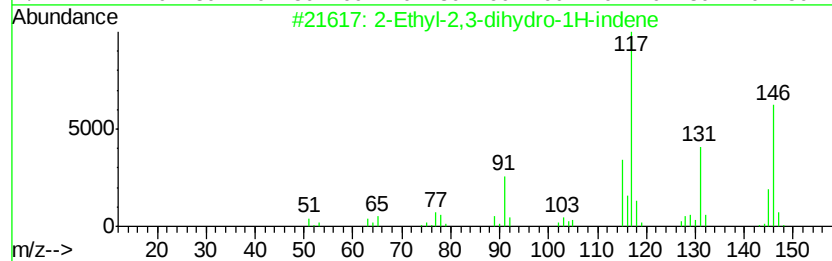
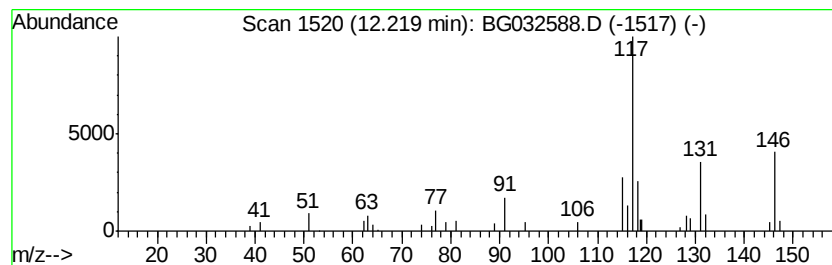
Instrument :  
BNA\_G  
ClientSampleId :  
LW-03-OW

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 12 2-Ethyl-2,3-dihydro-1H-indene Concentration Rank 17

| R.T.    | EstConc  | Area                                | Relative to ISTD | R.T.            |
|---------|----------|-------------------------------------|------------------|-----------------|
| 12.22   | 12.93 ng | 78174                               | Naphthalene-d8   | 11.58           |
| Hit# of | 5        | Tentative ID                        | MW MolForm       | CAS# Qual       |
| 1       |          | 2-Ethyl-2,3-dihydro-1H-indene       | 146 C11H14       | 056147-63-8 80  |
| 2       |          | Bicyclo[4.2.1]nona-2,4,7-triene,... | 146 C11H14       | 1000164-42-6 78 |
| 3       |          | Benzene, 1-pentenyl-                | 146 C11H14       | 000826-18-6 64  |
| 4       |          | 1H-Indene, 1-ethyl-2,3-dihydro-     | 146 C11H14       | 004830-99-3 58  |
| 5       |          | 1H-Imidazole, 4,5-dihydro-2-phenyl- | 146 C9H10N2      | 000936-49-2 50  |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
Data File : BG032588.D  
Acq On : 7 Feb 2018 7:26  
Operator : SJ/JU  
Sample : J1455-08  
Misc : MDL-W-8 ppm  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-03-OW

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

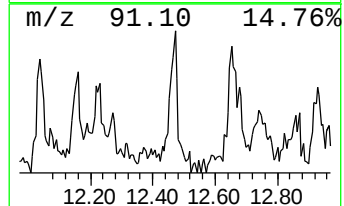
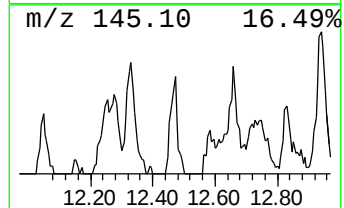
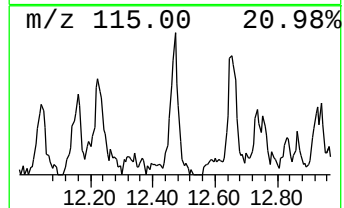
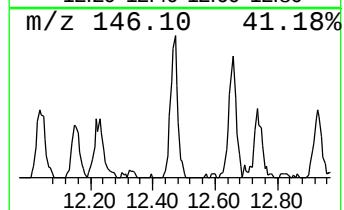
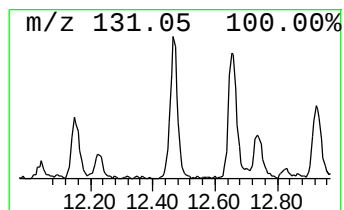
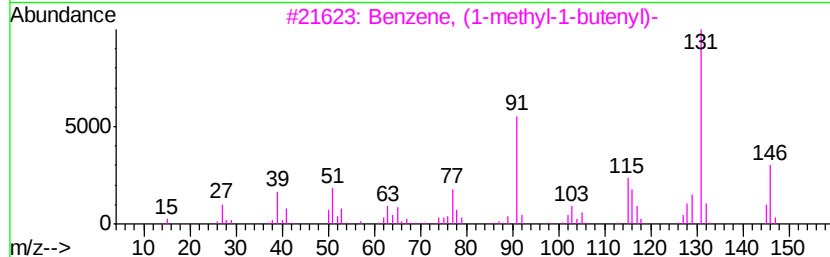
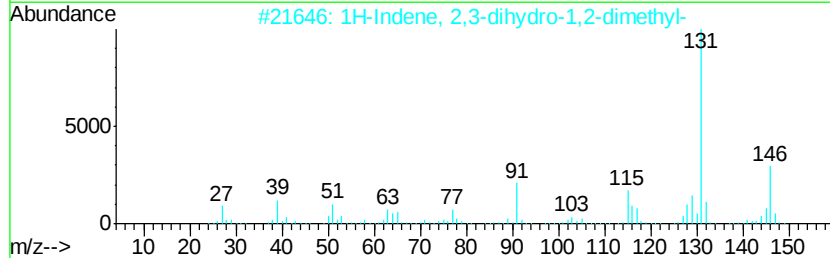
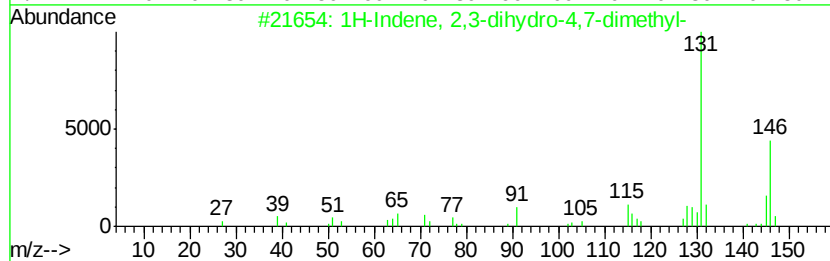
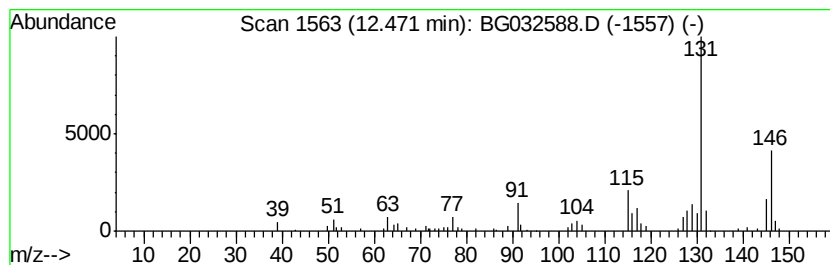
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 14

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 12.47 | 19.13 ng | 115672 | Naphthalene-d8   | 11.58 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 94   |
| 2    |    | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 | C11H14  | 017057-82-8 | 94   |
| 3    |    | Benzene, (1-methyl-1-butenyl)-      | 146 | C11H14  | 053172-84-2 | 93   |
| 4    |    | Benzene, (3-methyl-2-butenyl)-      | 146 | C11H14  | 004489-84-3 | 90   |
| 5    |    | Benzene, (1,2-dimethyl-1-propenyl)- | 146 | C11H14  | 000769-57-3 | 90   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
Data File : BG032588.D  
Acq On : 7 Feb 2018 7:26  
Operator : SJ/JU  
Sample : J1455-08  
Misc : MDL-W-8 ppm  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-03-OW

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

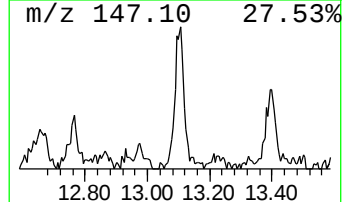
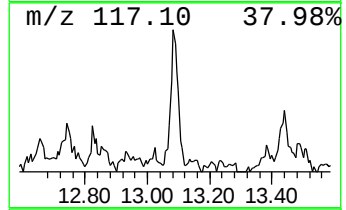
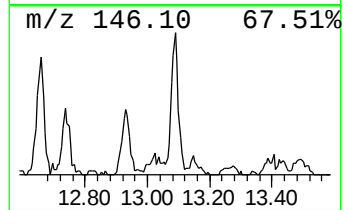
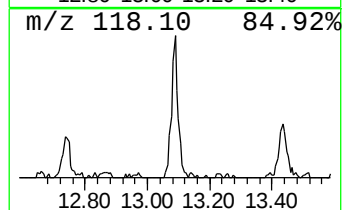
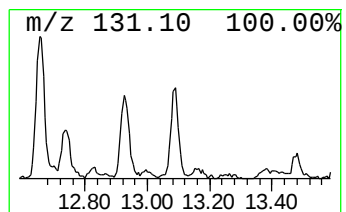
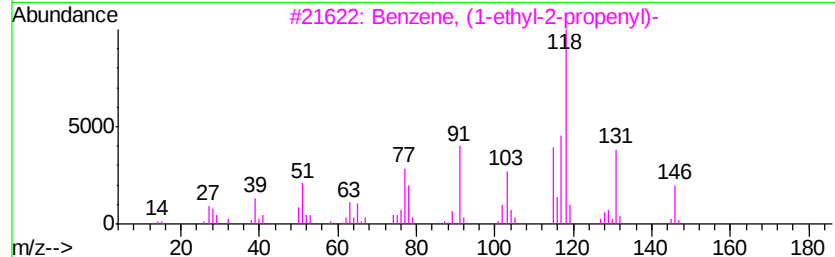
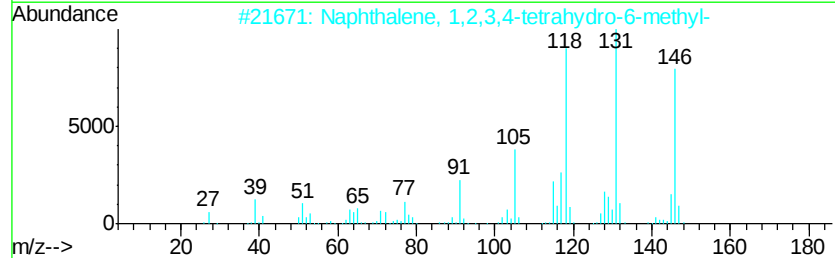
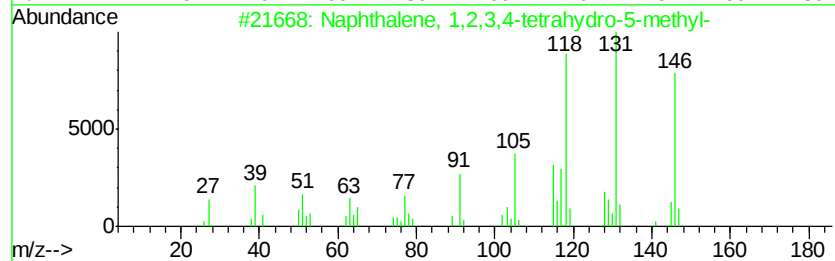
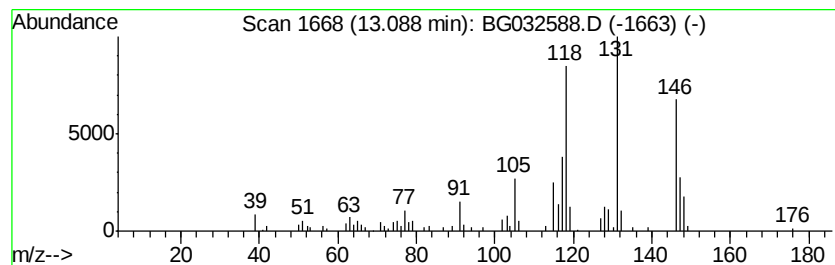
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Naphthalene, 1,2,3,4-tetra... Concentration Rank 9

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 13.09 | 21.83 ng | 131973 | Naphthalene-d8   | 11.58 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 002809-64-5  | 93   |
| 2    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 001680-51-9  | 91   |
| 3    |    | Benzene, (1-ethyl-2-propenyl)-      | 146 | C11H14  | 019947-22-9  | 89   |
| 4    |    | 2,3,4,5,6,7-Hexahydro-1H-cyclope... | 146 | C11H14  | 1000189-31-0 | 72   |
| 5    |    | Benzene, (1-methylenebutyl)-        | 146 | C11H14  | 005676-32-4  | 70   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
Data File : BG032588.D  
Acq On : 7 Feb 2018 7:26  
Operator : SJ/JU  
Sample : J1455-08  
Misc : MDL-W-8 ppm  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-03-OW

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

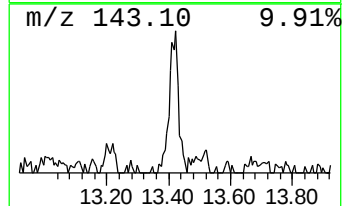
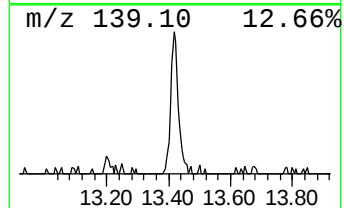
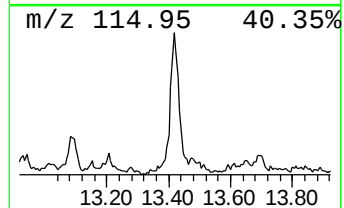
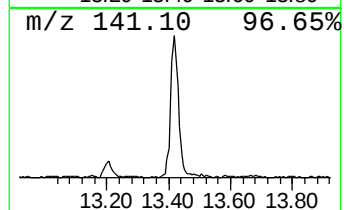
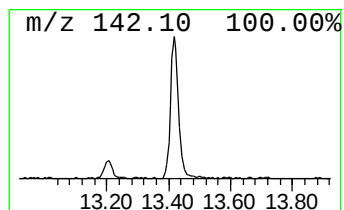
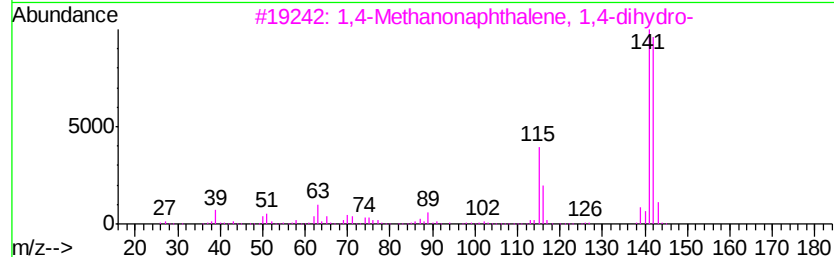
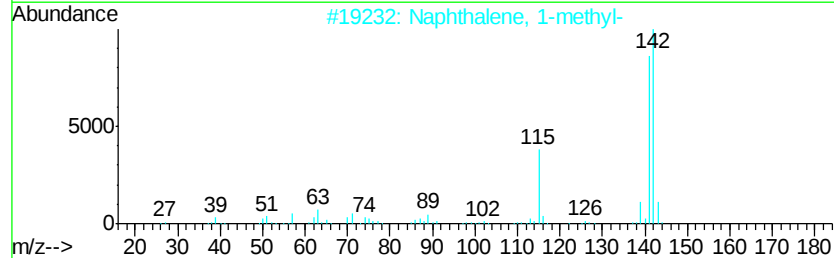
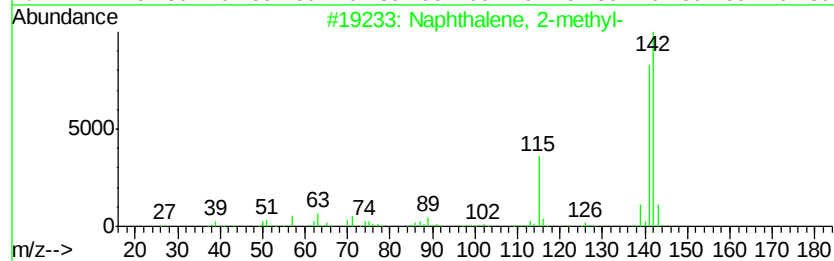
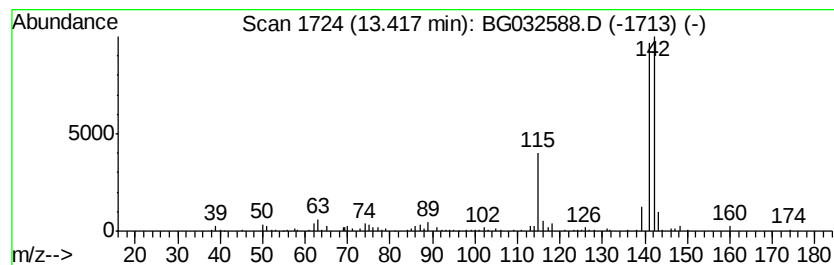
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Naphthalene, 1-methyl- Concentration Rank 4

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 13.42 | 49.00 ng | 296228 | Naphthalene-d8   | 11.58 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2-methyl-              | 142 | C11H10  | 000091-57-6 | 97   |
| 2    |    | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 97   |
| 3    |    | 1,4-Methanonaphthalene, 1,4-dihv... | 142 | C11H10  | 004453-90-1 | 91   |
| 4    |    | Benzocycloheptatriene               | 142 | C11H10  | 000264-09-5 | 91   |
| 5    |    | Bicyclo[4.4.1]undeca-1,3,5,7,9-p... | 142 | C11H10  | 002443-46-1 | 91   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
Data File : BG032588.D  
Acq On : 7 Feb 2018 7:26  
Operator : SJ/JU  
Sample : J1455-08  
Misc : MDL-W-8 ppm  
ALS Vial : 22 Sample Multiplier: 1

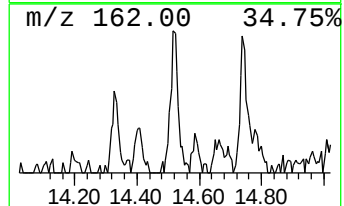
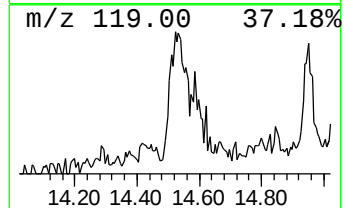
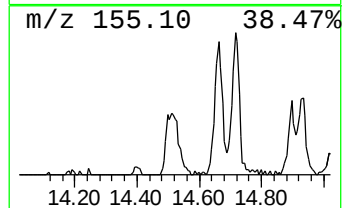
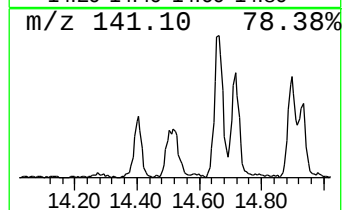
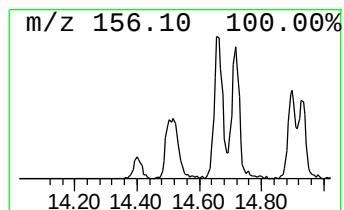
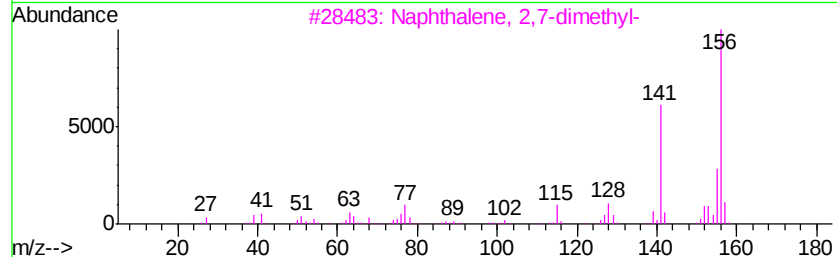
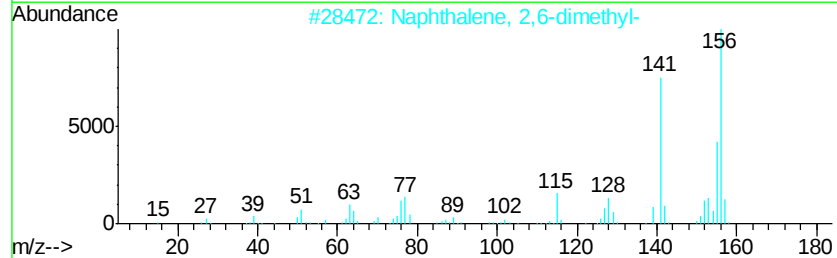
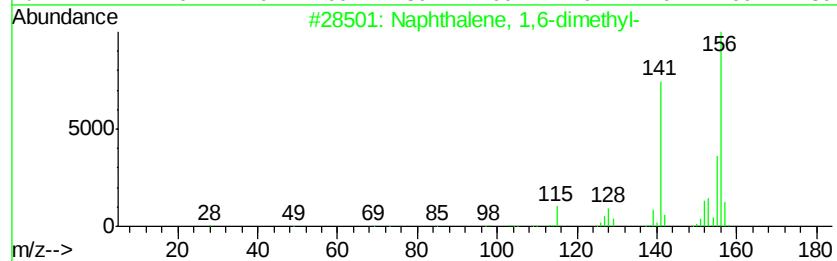
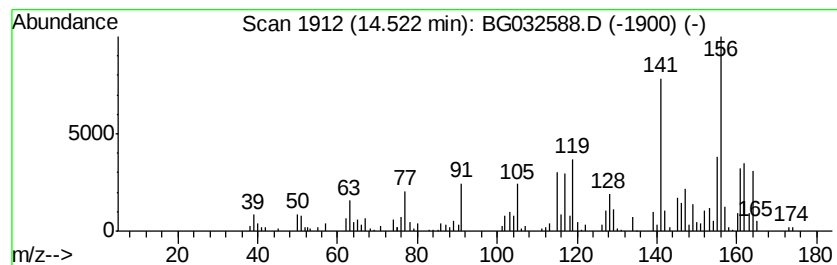
Instrument :  
BNA\_G  
ClientSampleId :  
LW-03-OW

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 18 Naphthalene, 1,6-dimethyl- Concentration Rank 16

| R.T.  | EstConc  | Area   | Relative to ISTD           |     | R.T.    |             |      |
|-------|----------|--------|----------------------------|-----|---------|-------------|------|
| 14.52 | 14.97 ng | 278898 | Acenaphthene-d10           |     | 15.34   |             |      |
| Hit#  | of       | 5      | Tentative ID               | MW  | MolForm | CAS#        | Qual |
| 1     |          |        | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 91   |
| 2     |          |        | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 91   |
| 3     |          |        | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 91   |
| 4     |          |        | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 83   |
| 5     |          |        | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 83   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
Data File : BG032588.D  
Acq On : 7 Feb 2018 7:26  
Operator : SJ/JU  
Sample : J1455-08  
Misc : MDL-W-8 ppm  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-03-OW

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

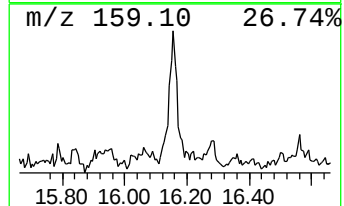
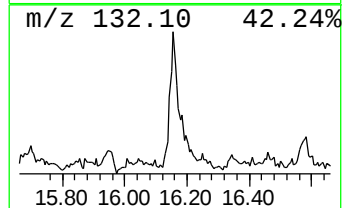
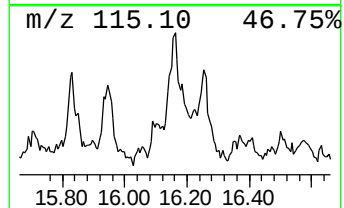
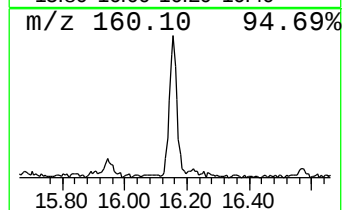
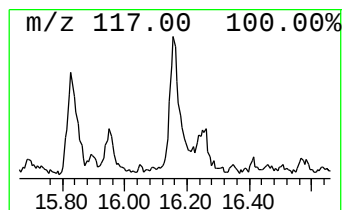
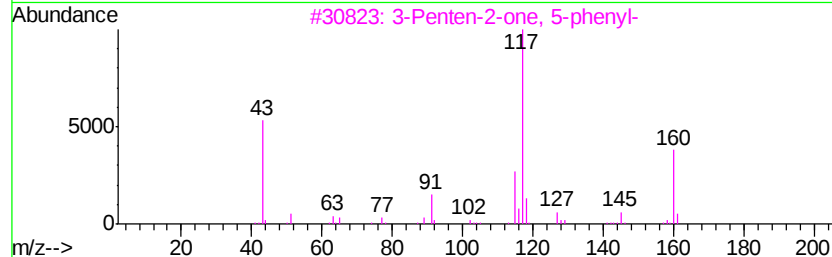
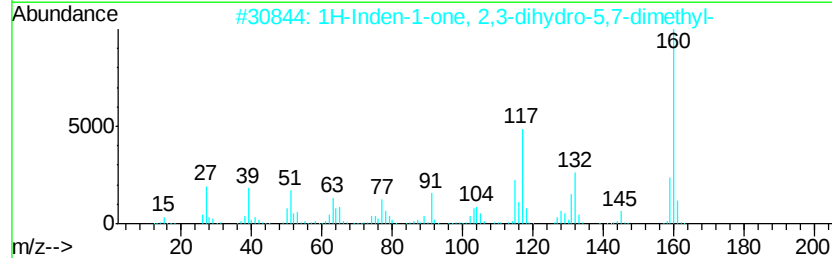
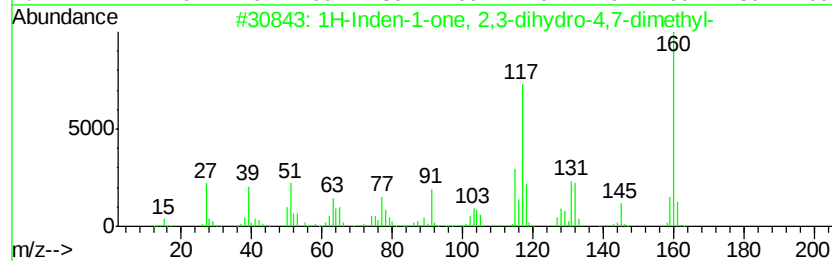
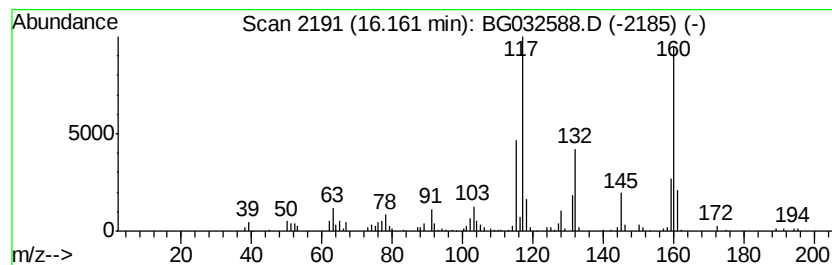
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 1H-Inden-1-one, 2,3-dihydro... Concentration Rank 20

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 16.16 | 11.08 ng | 206488 | Acenaphthene-d10 | 15.34 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1H-Inden-1-one, 2,3-dihydro-4,7-... | 160 | C11H12O | 005037-60-5 | 87   |
| 2    |    | 1H-Inden-1-one, 2,3-dihydro-5,7-... | 160 | C11H12O | 006682-69-5 | 86   |
| 3    |    | 3-Penten-2-one, 5-phenyl-           | 160 | C11H12O | 010521-97-8 | 58   |
| 4    |    | Benzene, 1-methyl-4-(2-propenyl)-   | 132 | C10H12  | 003333-13-9 | 45   |
| 5    |    | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000767-99-7 | 45   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
Data File : BG032588.D  
Acq On : 7 Feb 2018 7:26  
Operator : SJ/JU  
Sample : J1455-08  
Misc : MDL-W-8 ppm  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-03-OW

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

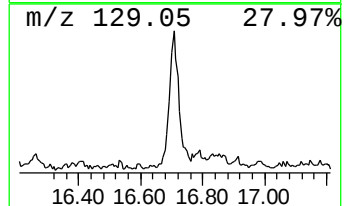
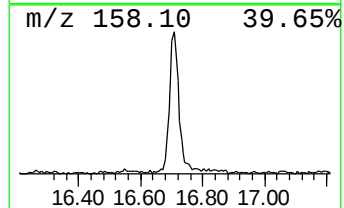
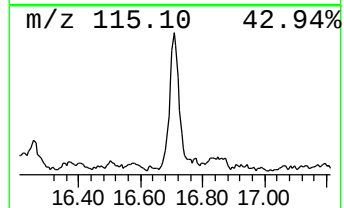
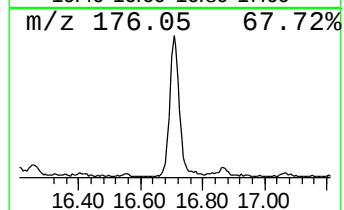
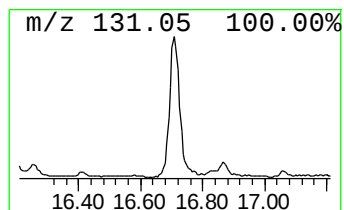
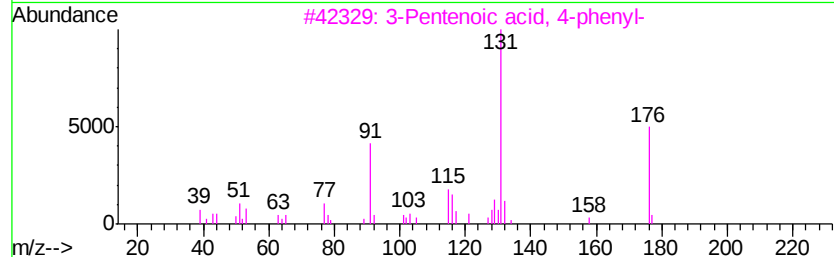
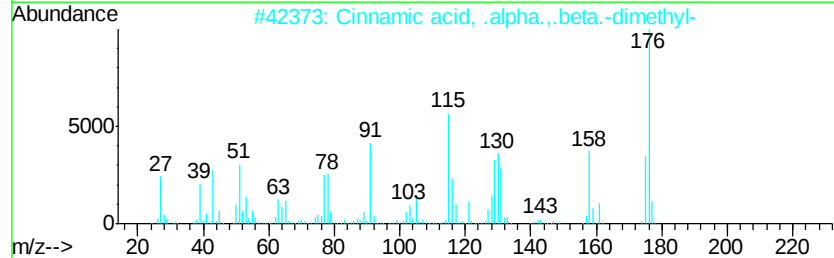
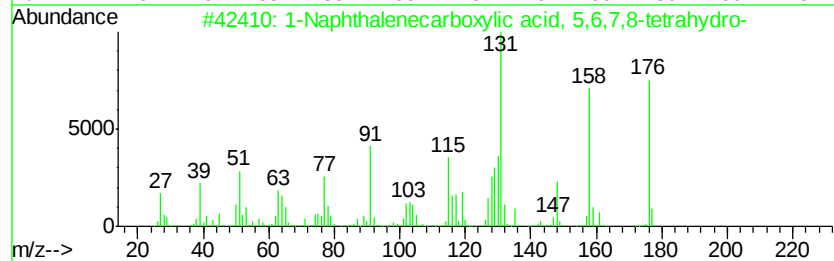
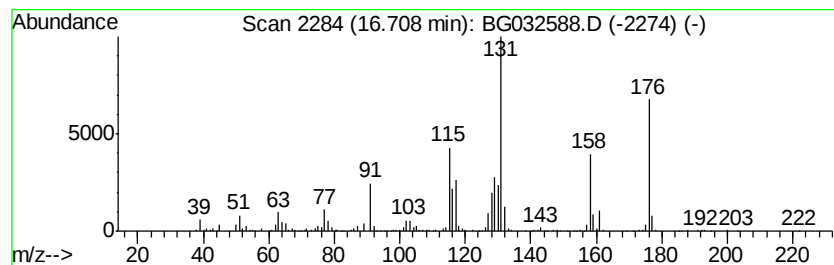
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 1-Naphthalenecarboxylic aci... Concentration Rank 5

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 16.71 | 43.95 ng | 818854 | Acenaphthene-d10 | 15.34 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 1-Naphthalenecarboxylic acid, 5,... | 176 | C11H12O2 | 004242-18-6  | 59   |
| 2    |    | Cinnamic acid, .alpha.,.beta.-di... | 176 | C11H12O2 | 1000151-56-4 | 58   |
| 3    |    | 3-Pentenoic acid, 4-phenyl-         | 176 | C11H12O2 | 053774-19-9  | 38   |
| 4    |    | Benzene, (3-chloro-1-methyl-1-pr... | 166 | C10H11Cl | 1000327-43-7 | 35   |
| 5    |    | Cyclopropane, 2-bromo-1-methyl-1... | 210 | C10H11Br | 055091-64-0  | 35   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG020618\  
 Data File : BG032588.D  
 Acq On : 7 Feb 2018 7:26  
 Operator : SJ/JU  
 Sample : J1455-08  
 Misc : MDL-W-8 ppm  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 LW-03-OW

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG012918.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| Butane, 2-methoxy... | 3.46  | 19.3    | ng    | 116340   | 1                     | 8.71  | 120521 | 20.0 |
| Mesitylene           | 7.88  | 12.9    | ng    | 77723    | 1                     | 8.71  | 120521 | 20.0 |
| Xenon                | 8.22  | 90.9    | ng    | 547968   | 1                     | 8.71  | 120521 | 20.0 |
| Indane               | 9.09  | 34.5    | ng    | 207995   | 1                     | 8.71  | 120521 | 20.0 |
| o-Cymene             | 9.83  | 21.4    | ng    | 129228   | 1                     | 8.71  | 120521 | 20.0 |
| Benzene, 1,2,3,4-... | 10.37 | 22.9    | ng    | 138276   | 2                     | 11.58 | 120903 | 20.0 |
| Benzene, 1,2,3,5-... | 10.43 | 37.1    | ng    | 224313   | 2                     | 11.58 | 120903 | 20.0 |
| Indan, 1-methyl-     | 10.96 | 75.6    | ng    | 456852   | 2                     | 11.58 | 120903 | 20.0 |
| Naphthalene, 1,2,... | 11.20 | 12.0    | ng    | 72540    | 2                     | 11.58 | 120903 | 20.0 |
| Benzene, (1-methy... | 11.68 | 21.4    | ng    | 129596   | 2                     | 11.58 | 120903 | 20.0 |
| 2-Ethyl-2,3-dihyd... | 12.22 | 12.9    | ng    | 78174    | 2                     | 11.58 | 120903 | 20.0 |
| 1H-Indene, 2,3-di... | 12.47 | 19.1    | ng    | 115672   | 2                     | 11.58 | 120903 | 20.0 |
| Naphthalene, 1,2,... | 13.09 | 21.8    | ng    | 131973   | 2                     | 11.58 | 120903 | 20.0 |
| Naphthalene, 1-me... | 13.42 | 49.0    | ng    | 296228   | 2                     | 11.58 | 120903 | 20.0 |
| Naphthalene, 1,6-... | 14.52 | 15.0    | ng    | 278898   | 3                     | 15.34 | 372623 | 20.0 |
| 1H-Inden-1-one, 2... | 16.16 | 11.1    | ng    | 206488   | 3                     | 15.34 | 372623 | 20.0 |
| 1-Naphthalenecarb... | 16.71 | 44.0    | ng    | 818854   | 3                     | 15.34 | 372623 | 20.0 |