

Data Path : Z:\HPCHEM1\BNA G\DATA\BG120515\  
 Data File : BG019952.D  
 Acq On : 5 Dec 2015 21:31  
 Operator : UM/NP  
 Sample : G4630-11  
 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MW-18

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-BG120215.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         | 5.189       | 394           | 400         | 405          | rBV      | 50659          | 77868         | 2.43%           | 0.325%        |
| 2         | 5.659       | 473           | 480         | 491          | rBV      | 299731         | 485564        | 15.18%          | 2.025%        |
| 3         | 6.593       | 631           | 639         | 647          | rBV      | 39181          | 75505         | 2.36%           | 0.315%        |
| 4         | 7.086       | 715           | 723         | 729          | rBV      | 52819          | 94918         | 2.97%           | 0.396%        |
| 5         | 7.257       | 745           | 752         | 759          | rBV      | 211685         | 367364        | 11.48%          | 1.532%        |
| 6         | 7.639       | 810           | 817         | 826          | rVB      | 551675         | 966045        | 30.20%          | 4.029%        |
| 7         | 8.015       | 877           | 881         | 886          | rVB      | 27186          | 40238         | 1.26%           | 0.168%        |
| 8         | 8.115       | 889           | 898         | 904          | rVB      | 127522         | 219176        | 6.85%           | 0.914%        |
| 9         | 8.438       | 942           | 953         | 959          | rBV      | 543969         | 967781        | 30.25%          | 4.036%        |
| 10        | 8.491       | 959           | 962         | 969          | rVB      | 91167          | 134913        | 4.22%           | 0.563%        |
| 11        | 8.608       | 975           | 982         | 987          | rBV2     | 50155          | 77076         | 2.41%           | 0.321%        |
| 12        | 8.761       | 1002          | 1008        | 1012         | rBV2     | 42619          | 79635         | 2.49%           | 0.332%        |
| 13        | 8.808       | 1012          | 1016        | 1022         | rVB4     | 38743          | 74478         | 2.33%           | 0.311%        |
| 14        | 9.225       | 1082          | 1087        | 1091         | rBV2     | 109958         | 180701        | 5.65%           | 0.754%        |
| 15        | 9.284       | 1091          | 1097        | 1111         | rVB2     | 475847         | 1056042       | 33.01%          | 4.404%        |
| 16        | 9.460       | 1119          | 1127        | 1137         | rBV9     | 29333          | 83540         | 2.61%           | 0.348%        |
| 17        | 9.578       | 1142          | 1147        | 1159         | rVB3     | 28694          | 64620         | 2.02%           | 0.269%        |
| 18        | 9.701       | 1159          | 1168        | 1172         | rBV7     | 21875          | 51910         | 1.62%           | 0.216%        |
| 19        | 9.748       | 1172          | 1176        | 1182         | rVB      | 94557          | 149952        | 4.69%           | 0.625%        |
| 20        | 9.848       | 1189          | 1193        | 1201         | rVB4     | 19284          | 37620         | 1.18%           | 0.157%        |
| 21        | 9.924       | 1201          | 1206        | 1213         | rBV6     | 15874          | 37557         | 1.17%           | 0.157%        |
| 22        | 10.012      | 1215          | 1221        | 1225         | rVV2     | 22738          | 34168         | 1.07%           | 0.142%        |
| 23        | 10.118      | 1232          | 1239        | 1245         | rBV4     | 53470          | 131641        | 4.12%           | 0.549%        |
| 24        | 10.324      | 1268          | 1274        | 1282         | rVV2     | 174888         | 320764        | 10.03%          | 1.338%        |
| 25        | 10.512      | 1296          | 1306        | 1311         | rBV10    | 27792          | 80901         | 2.53%           | 0.337%        |
| 26        | 10.624      | 1318          | 1325        | 1336         | rBV3     | 43915          | 112873        | 3.53%           | 0.471%        |
| 27        | 10.782      | 1347          | 1352        | 1359         | rBV9     | 23479          | 50481         | 1.58%           | 0.211%        |
| 28        | 10.929      | 1366          | 1377        | 1382         | rBV2     | 204668         | 474173        | 14.82%          | 1.977%        |
| 29        | 10.976      | 1383          | 1385        | 1391         | rVB4     | 48298          | 85349         | 2.67%           | 0.356%        |
| 30        | 11.041      | 1391          | 1396        | 1405         | rVB2     | 81363          | 148936        | 4.66%           | 0.621%        |
| 31        | 11.141      | 1408          | 1413        | 1421         | rBV      | 25433          | 48455         | 1.51%           | 0.202%        |
| 32        | 11.299      | 1435          | 1440        | 1445         | rBV5     | 30416          | 59522         | 1.86%           | 0.248%        |
| 33        | 11.399      | 1453          | 1457        | 1465         | rVB3     | 52357          | 103199        | 3.23%           | 0.430%        |
| 34        | 11.511      | 1466          | 1476        | 1481         | rBV3     | 52282          | 122260        | 3.82%           | 0.510%        |

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 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

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 BNA\_G  
 ClientSampleId :  
 MW-18

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-BG120215.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |       |         |         |        |        |
|----|--------|------|------|------|-------|---------|---------|--------|--------|
| 35 | 11.710 | 1504 | 1510 | 1513 | rBV6  | 19257   | 36426   | 1.14%  | 0.152% |
| 36 | 11.763 | 1514 | 1519 | 1526 | rVV10 | 31307   | 89295   | 2.79%  | 0.372% |
| 37 | 11.840 | 1526 | 1532 | 1538 | rVB4  | 53327   | 112159  | 3.51%  | 0.468% |
| 38 | 11.945 | 1546 | 1550 | 1553 | rBV2  | 32268   | 48702   | 1.52%  | 0.203% |
| 39 | 12.034 | 1560 | 1565 | 1572 | rVB2  | 49396   | 89235   | 2.79%  | 0.372% |
| 40 | 12.116 | 1574 | 1579 | 1585 | rBV2  | 48871   | 83181   | 2.60%  | 0.347% |
| 41 | 12.175 | 1585 | 1589 | 1593 | rBV2  | 33219   | 48928   | 1.53%  | 0.204% |
| 42 | 12.374 | 1619 | 1623 | 1627 | rBV3  | 28291   | 46408   | 1.45%  | 0.194% |
| 43 | 12.468 | 1634 | 1639 | 1644 | rVB4  | 92123   | 150955  | 4.72%  | 0.630% |
| 44 | 12.527 | 1644 | 1649 | 1664 | rVB10 | 68958   | 233241  | 7.29%  | 0.973% |
| 45 | 12.727 | 1678 | 1683 | 1687 | rBV2  | 75140   | 103117  | 3.22%  | 0.430% |
| 46 | 12.797 | 1688 | 1695 | 1700 | rVV2  | 128372  | 269204  | 8.42%  | 1.123% |
| 47 | 12.868 | 1704 | 1707 | 1711 | rVB5  | 34713   | 53067   | 1.66%  | 0.221% |
| 48 | 13.091 | 1741 | 1745 | 1750 | rVB4  | 34138   | 55958   | 1.75%  | 0.233% |
| 49 | 13.138 | 1750 | 1753 | 1757 | rVB6  | 23993   | 33790   | 1.06%  | 0.141% |
| 50 | 13.203 | 1760 | 1764 | 1766 | rVV5  | 31346   | 53361   | 1.67%  | 0.223% |
| 51 | 13.232 | 1766 | 1769 | 1773 | rVV4  | 60329   | 95969   | 3.00%  | 0.400% |
| 52 | 13.279 | 1773 | 1777 | 1783 | rVB3  | 94322   | 151583  | 4.74%  | 0.632% |
| 53 | 13.373 | 1786 | 1793 | 1797 | rVV   | 1333596 | 2072070 | 64.77% | 8.641% |
| 54 | 13.420 | 1797 | 1801 | 1805 | rVB   | 121599  | 178039  | 5.57%  | 0.742% |
| 55 | 13.696 | 1842 | 1848 | 1852 | rBV7  | 38593   | 77974   | 2.44%  | 0.325% |
| 56 | 13.737 | 1852 | 1855 | 1858 | rBV5  | 26802   | 40794   | 1.28%  | 0.170% |
| 57 | 13.861 | 1873 | 1876 | 1879 | rBV5  | 23331   | 35206   | 1.10%  | 0.147% |
| 58 | 13.908 | 1880 | 1884 | 1889 | rVV4  | 75560   | 145991  | 4.56%  | 0.609% |
| 59 | 13.961 | 1889 | 1893 | 1898 | rVV5  | 77819   | 133952  | 4.19%  | 0.559% |
| 60 | 14.008 | 1898 | 1901 | 1902 | rVV2  | 37712   | 46179   | 1.44%  | 0.193% |
| 61 | 14.067 | 1903 | 1911 | 1916 | rVV5  | 140884  | 391709  | 12.24% | 1.634% |
| 62 | 14.119 | 1916 | 1920 | 1929 | rVB4  | 93366   | 187709  | 5.87%  | 0.783% |
| 63 | 14.219 | 1934 | 1937 | 1940 | rVB4  | 44388   | 53215   | 1.66%  | 0.222% |
| 64 | 14.296 | 1948 | 1950 | 1954 | rBV2  | 29410   | 36550   | 1.14%  | 0.152% |
| 65 | 14.372 | 1959 | 1963 | 1970 | rVB4  | 129345  | 235418  | 7.36%  | 0.982% |
| 66 | 14.466 | 1970 | 1979 | 1986 | rBV5  | 76140   | 226604  | 7.08%  | 0.945% |
| 67 | 14.554 | 1989 | 1994 | 2004 | rVB3  | 123552  | 238666  | 7.46%  | 0.995% |
| 68 | 14.742 | 2019 | 2026 | 2033 | rVB2  | 312144  | 588381  | 18.39% | 2.454% |
| 69 | 14.824 | 2036 | 2040 | 2044 | rBV5  | 84809   | 128202  | 4.01%  | 0.535% |
| 70 | 14.901 | 2049 | 2053 | 2066 | rVB5  | 111672  | 303113  | 9.48%  | 1.264% |
| 71 | 15.065 | 2077 | 2081 | 2086 | rBV7  | 40793   | 91299   | 2.85%  | 0.381% |

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Integration Parameters: rteint.p  
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 Filtering: 5  
 Min Area: 3 % of largest Peak  
 Max Peaks: 100  
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 Peak separation: 5

Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG120215.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|     |        |      |      |      |      |         |         |         |         |
|-----|--------|------|------|------|------|---------|---------|---------|---------|
| 72  | 15.130 | 2088 | 2092 | 2101 | rVB5 | 90336   | 209273  | 6.54%   | 0.873%  |
| 73  | 15.218 | 2101 | 2107 | 2110 | rBV5 | 104561  | 201412  | 6.30%   | 0.840%  |
| 74  | 15.383 | 2127 | 2135 | 2141 | rBV5 | 174518  | 418610  | 13.09%  | 1.746%  |
| 75  | 15.559 | 2161 | 2165 | 2168 | rVV4 | 54897   | 71920   | 2.25%   | 0.300%  |
| 76  | 15.600 | 2168 | 2172 | 2178 | rVB6 | 72535   | 109926  | 3.44%   | 0.458%  |
| 77  | 15.782 | 2198 | 2203 | 2208 | rBV5 | 100507  | 175917  | 5.50%   | 0.734%  |
| 78  | 15.841 | 2209 | 2213 | 2218 | rVB6 | 51752   | 91805   | 2.87%   | 0.383%  |
| 79  | 15.994 | 2235 | 2239 | 2245 | rBV7 | 49692   | 102815  | 3.21%   | 0.429%  |
| 80  | 16.229 | 2273 | 2279 | 2284 | rBV  | 1127442 | 1705639 | 53.32%  | 7.113%  |
| 81  | 16.464 | 2316 | 2319 | 2323 | rBV6 | 64011   | 96895   | 3.03%   | 0.404%  |
| 82  | 16.658 | 2348 | 2352 | 2361 | rVB5 | 86940   | 180629  | 5.65%   | 0.753%  |
| 83  | 16.834 | 2378 | 2382 | 2386 | rVB2 | 81348   | 102023  | 3.19%   | 0.425%  |
| 84  | 17.269 | 2453 | 2456 | 2460 | rBV4 | 44907   | 80981   | 2.53%   | 0.338%  |
| 85  | 17.357 | 2467 | 2471 | 2476 | rBV2 | 65502   | 105667  | 3.30%   | 0.441%  |
| 86  | 17.486 | 2486 | 2493 | 2501 | rBV  | 425364  | 667070  | 20.85%  | 2.782%  |
| 87  | 17.756 | 2536 | 2539 | 2548 | rVB6 | 72879   | 118962  | 3.72%   | 0.496%  |
| 88  | 17.956 | 2570 | 2573 | 2578 | rVB3 | 46213   | 58093   | 1.82%   | 0.242%  |
| 89  | 18.367 | 2639 | 2643 | 2646 | rBV  | 104543  | 134256  | 4.20%   | 0.560%  |
| 90  | 19.631 | 2854 | 2858 | 2864 | rBV  | 93666   | 119839  | 3.75%   | 0.500%  |
| 91  | 20.089 | 2930 | 2936 | 2941 | rBV  | 2503468 | 3198983 | 100.00% | 13.341% |
| 92  | 20.770 | 3048 | 3052 | 3058 | rVB  | 53939   | 75896   | 2.37%   | 0.317%  |
| 93  | 20.982 | 3084 | 3088 | 3093 | rBV  | 32362   | 49895   | 1.56%   | 0.208%  |
| 94  | 21.035 | 3093 | 3097 | 3100 | rVB  | 55396   | 74770   | 2.34%   | 0.312%  |
| 95  | 21.099 | 3105 | 3108 | 3113 | rVB3 | 30759   | 40504   | 1.27%   | 0.169%  |
| 96  | 21.387 | 3154 | 3157 | 3167 | rVB4 | 35660   | 49947   | 1.56%   | 0.208%  |
| 97  | 21.540 | 3179 | 3183 | 3188 | rVB  | 46943   | 60309   | 1.89%   | 0.252%  |
| 98  | 21.758 | 3213 | 3220 | 3228 | rVB  | 480651  | 826785  | 25.85%  | 3.448%  |
| 99  | 22.092 | 3273 | 3277 | 3284 | rVB2 | 21233   | 37046   | 1.16%   | 0.154%  |
| 100 | 25.030 | 3768 | 3777 | 3791 | rVB  | 264847  | 752074  | 23.51%  | 3.136%  |

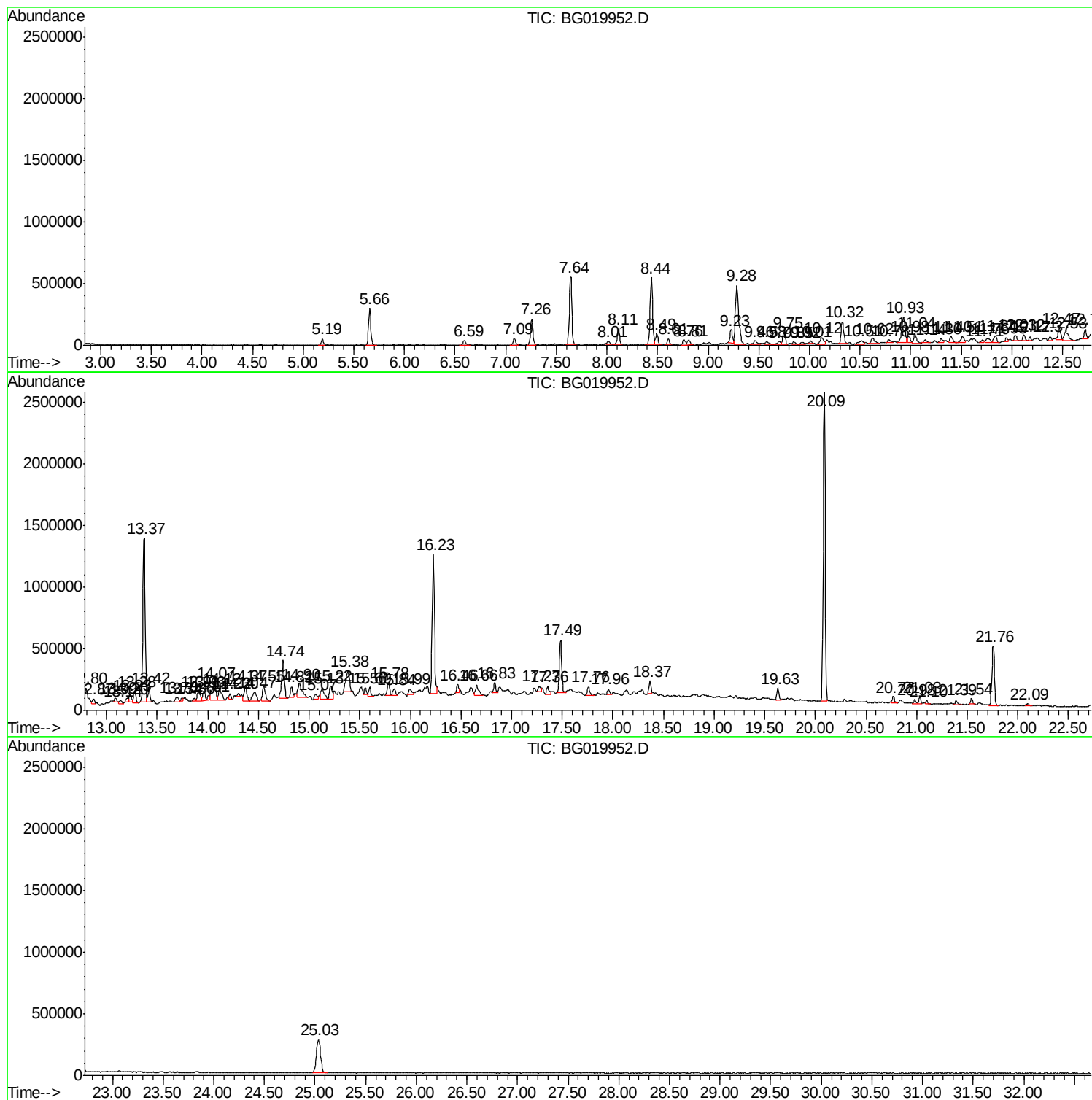
Sum of corrected areas: 23978816

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ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-18

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120515\  
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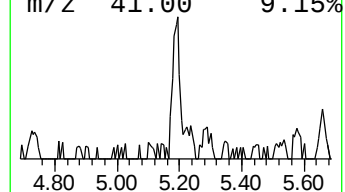
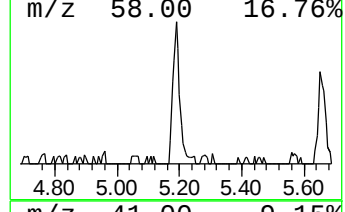
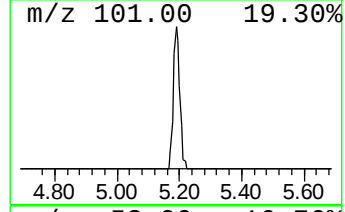
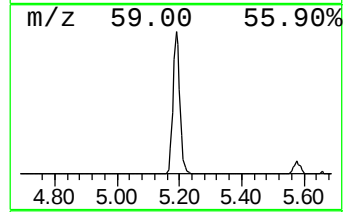
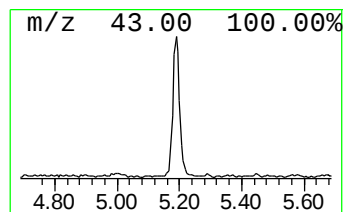
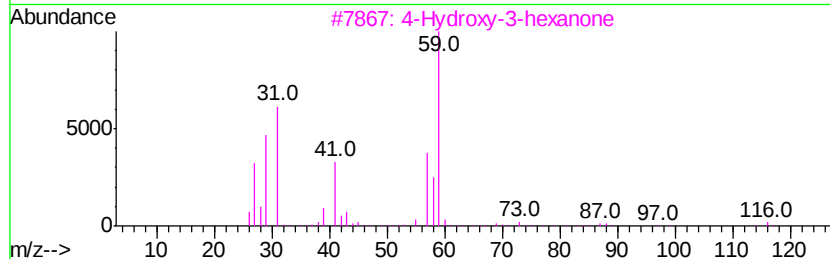
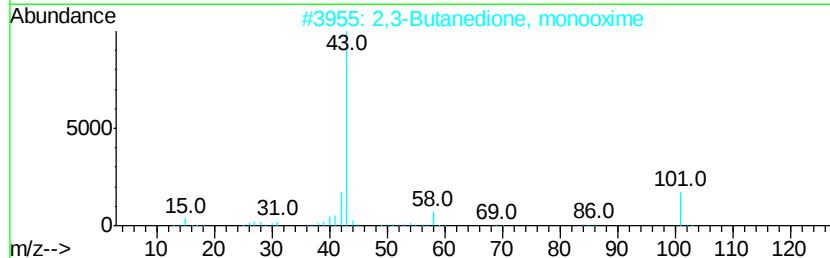
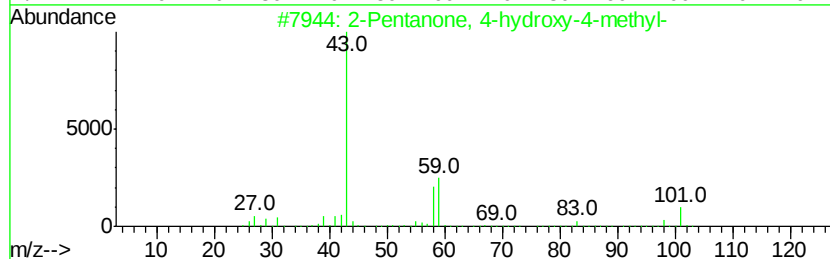
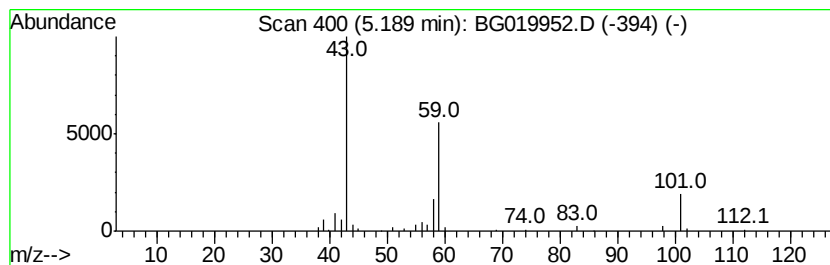
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 18

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 5.19 | 7.11 ng | 77868 | 1,4-Dichlorobenzene-d4 | 8.11 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-     | 116 | C6H12O2  | 000123-42-2 | 56   |
| 2    |    |   | 2,3-Butanedione, monooxime           | 101 | C4H7NO2  | 000057-71-6 | 25   |
| 3    |    |   | 4-Hydroxy-3-hexanone                 | 116 | C6H12O2  | 004984-85-4 | 23   |
| 4    |    |   | Acetic acid, cyano-, 1,1-dimethyl... | 141 | C7H11NO2 | 001116-98-9 | 17   |
| 5    |    |   | Butyl aldoxime, 3-methyl-, syn-      | 101 | C5H11NO  | 005780-40-5 | 17   |



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ALS Vial : 41 Sample Multiplier: 1

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

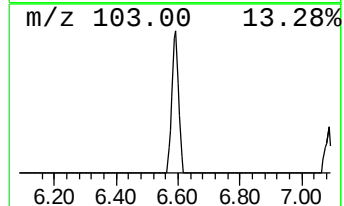
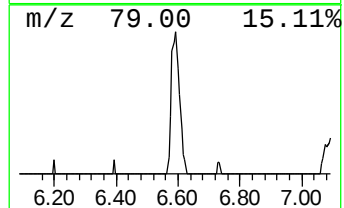
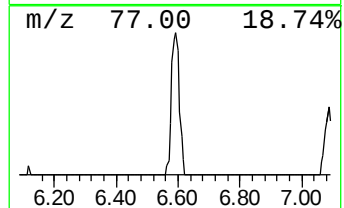
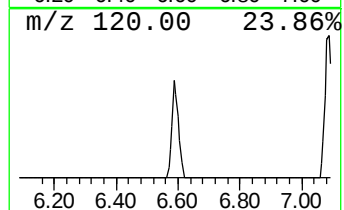
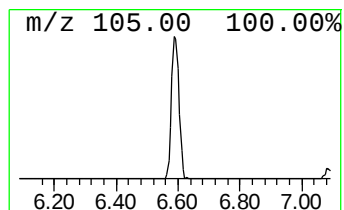
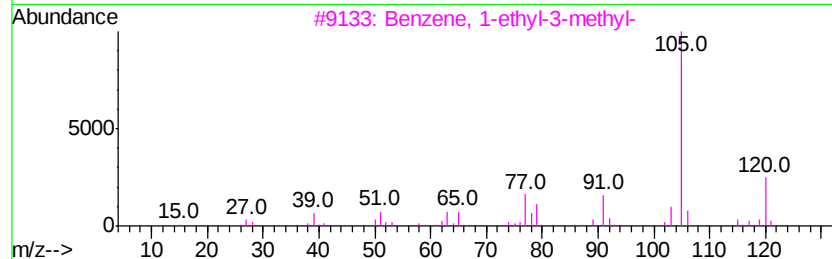
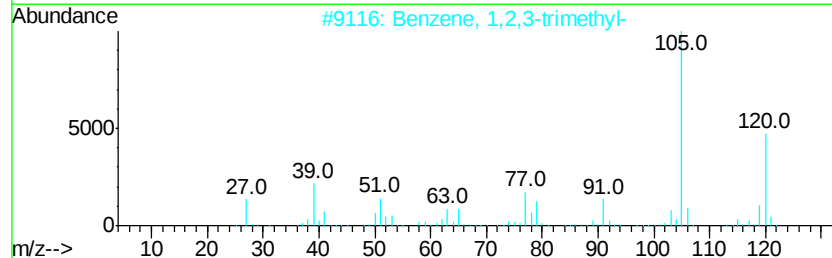
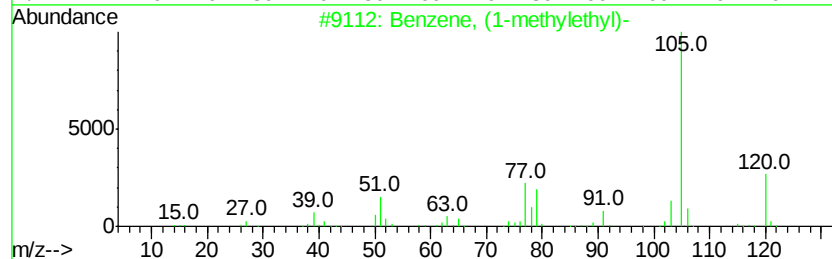
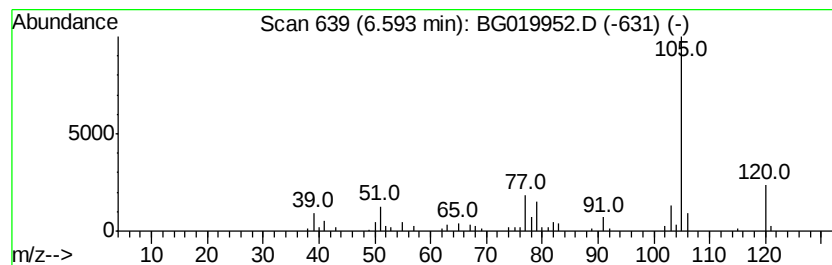
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 Benzene, (1-methylethyl)- Concentration Rank 20

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 6.59 | 6.89 ng | 75505 | 1,4-Dichlorobenzene-d4 | 8.11 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, (1-methylethyl)-  | 120 | C9H12   | 000098-82-8 | 97   |
| 2    |    | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |
| 3    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 83   |
| 4    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 80   |
| 5    |    | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 72   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120515\  
Data File : BG019952.D  
Acq On : 5 Dec 2015 21:31  
Operator : UM/NP  
Sample : G4630-11  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-18

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

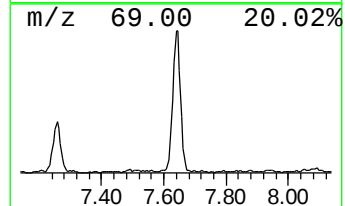
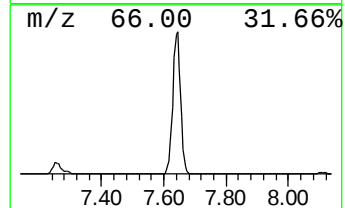
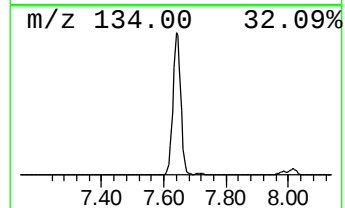
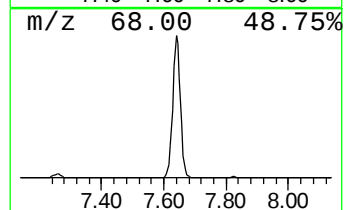
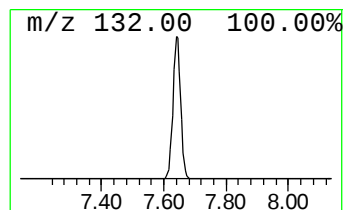
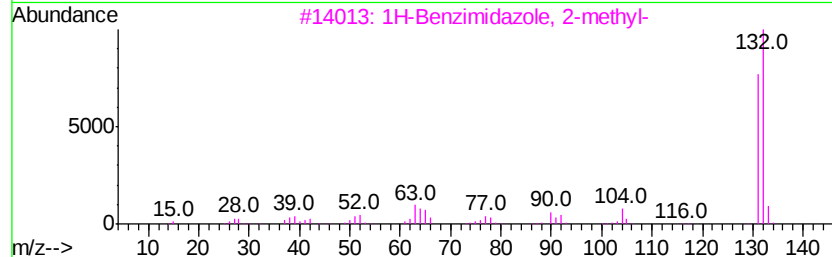
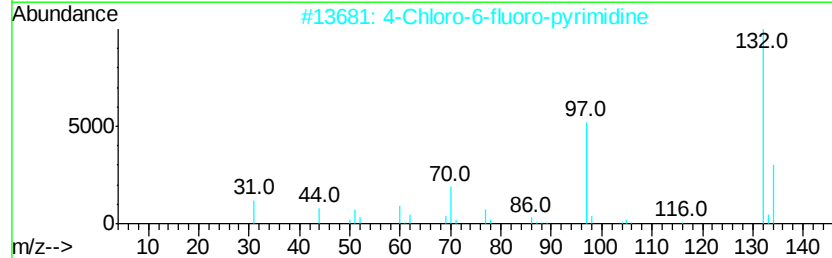
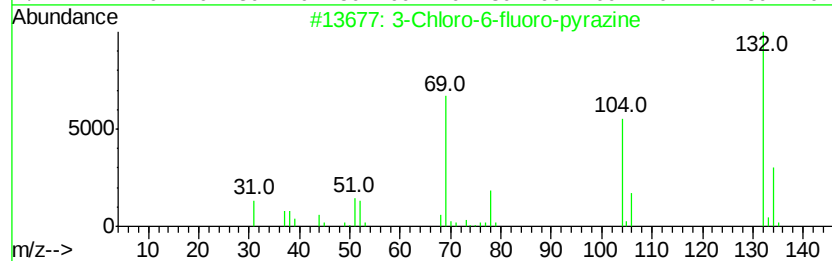
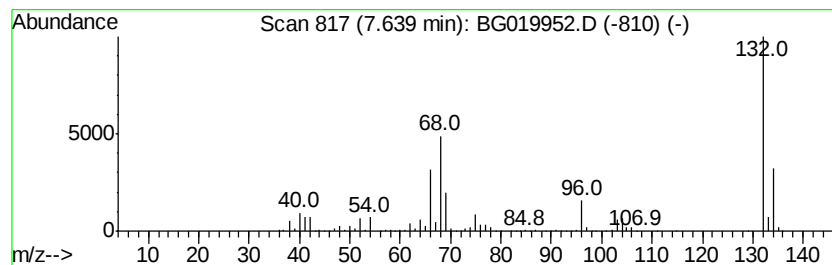
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 unknown7.64 Concentration Rank 2

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 7.64 | 88.15 ng | 966045 | 1,4-Dichlorobenzene-d4 | 8.11 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | 4-Chloro-6-fluoro-pyrimidine     | 132 | C4H2ClFN2 | 051422-01-6  | 27   |
| 3    |    | 1H-Benzimidazole, 2-methyl-      | 132 | C8H8N2    | 000615-15-6  | 22   |
| 4    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 22   |
| 5    |    | 1,3-Cyclopentanedione, 2-chloro- | 132 | C5H5ClO2  | 014203-19-1  | 17   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120515\  
Data File : BG019952.D  
Acq On : 5 Dec 2015 21:31  
Operator : UM/NP  
Sample : G4630-11  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-18

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

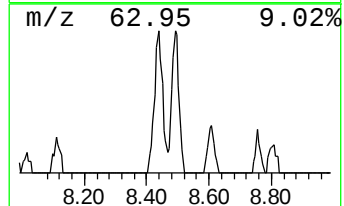
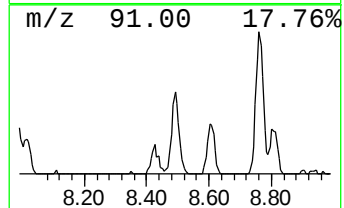
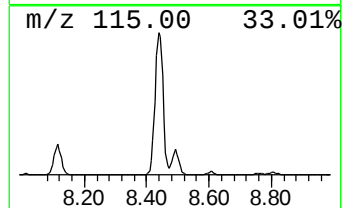
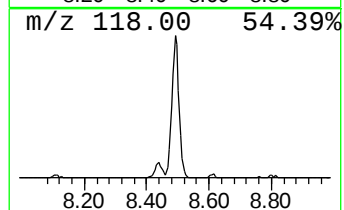
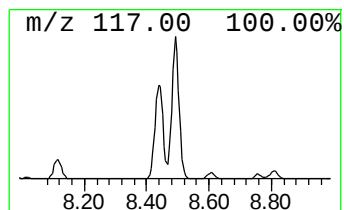
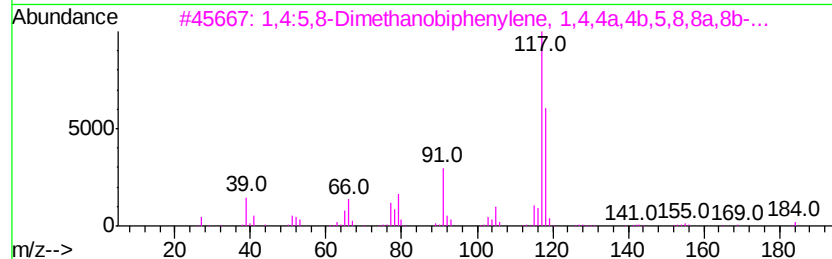
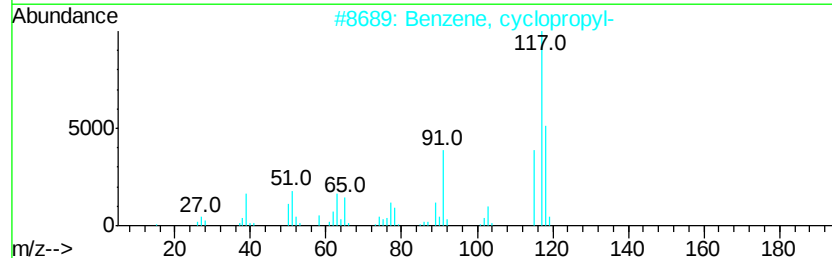
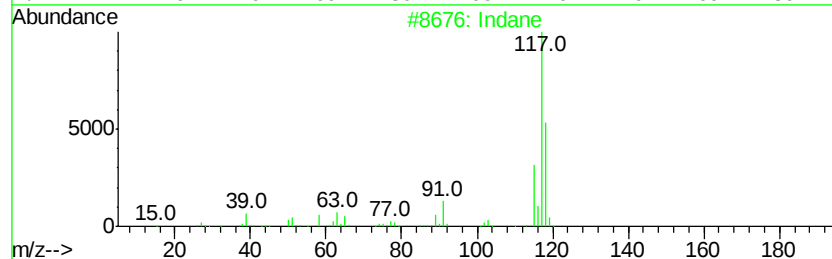
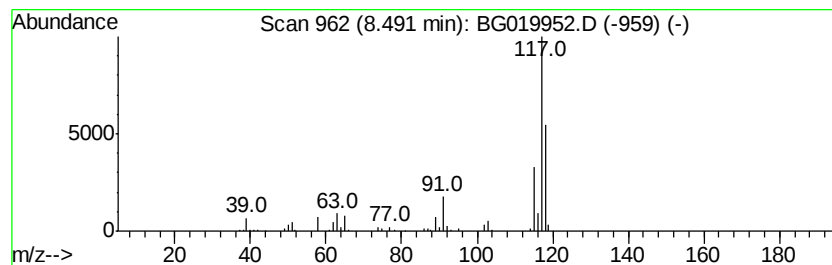
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 Indane Concentration Rank 7

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 8.49 | 12.31 ng | 134913 | 1,4-Dichlorobenzene-d4 | 8.11 |

| Hit# of | 5                                   | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|-------------------------------------|--------------|-----|---------|-------------|------|
| 1       | Indane                              |              | 118 | C9H10   | 000496-11-7 | 91   |
| 2       | Benzene, cyclopropyl-               |              | 118 | C9H10   | 000873-49-4 | 87   |
| 3       | 1,4:5,8-Dimethanobiphenylene, 1,... |              | 184 | C14H16  | 001624-13-1 | 72   |
| 4       | Benzene, 2-propenyl-                |              | 118 | C9H10   | 000300-57-2 | 64   |
| 5       | Deltacyclene                        |              | 118 | C9H10   | 007785-10-6 | 64   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120515\  
Data File : BG019952.D  
Acq On : 5 Dec 2015 21:31  
Operator : UM/NP  
Sample : G4630-11  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-18

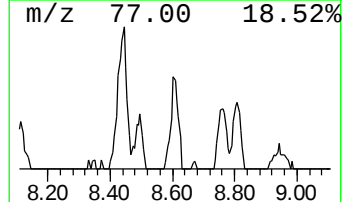
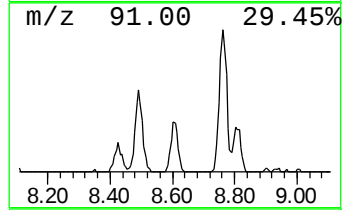
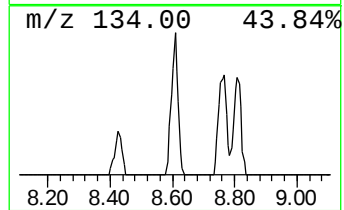
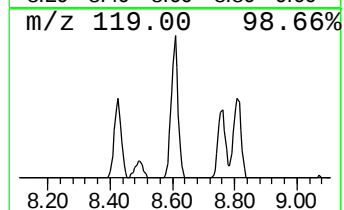
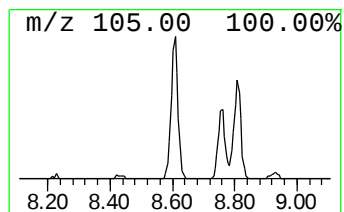
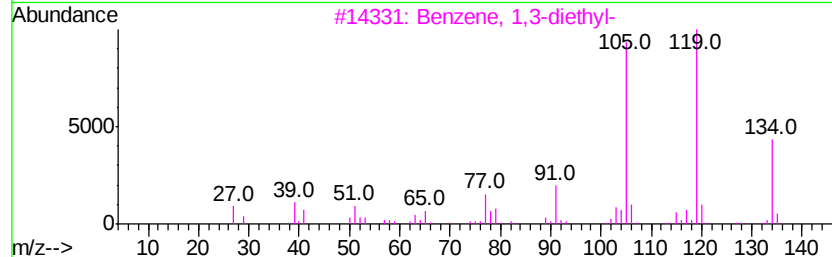
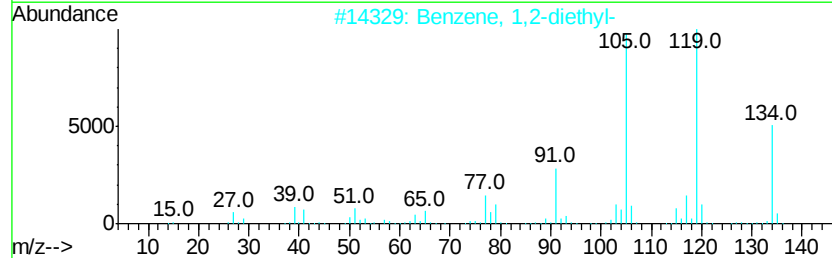
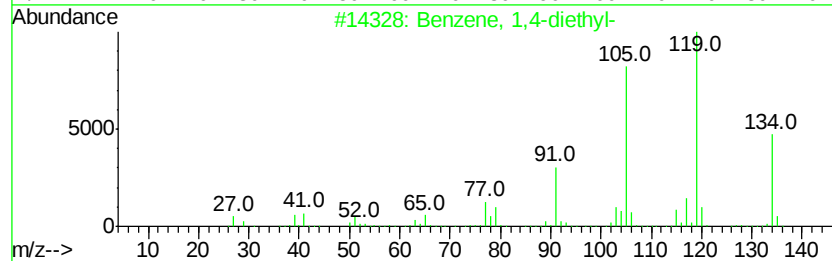
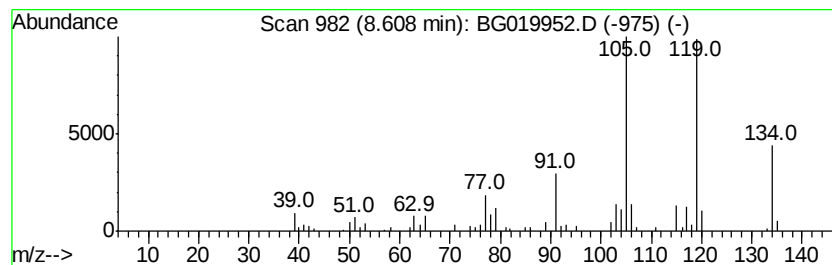
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 Benzene, 1,4-diethyl- Concentration Rank 19

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 8.61 | 7.03 ng | 77076 | 1,4-Dichlorobenzene-d4 | 8.11 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,4-diethyl-          | 134 | C10H14  | 000105-05-5 | 97   |
| 2    |    |   | Benzene, 1,2-diethyl-          | 134 | C10H14  | 000135-01-3 | 96   |
| 3    |    |   | Benzene, 1,3-diethyl-          | 134 | C10H14  | 000141-93-5 | 94   |
| 4    |    |   | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 91   |
| 5    |    |   | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 83   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120515\  
Data File : BG019952.D  
Acq On : 5 Dec 2015 21:31  
Operator : UM/NP  
Sample : G4630-11  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-18

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

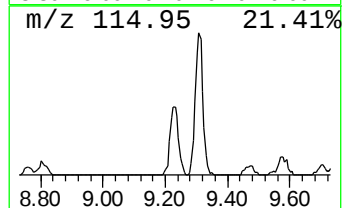
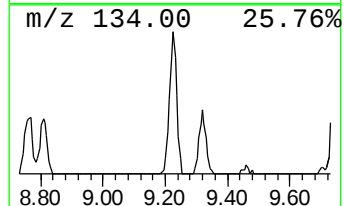
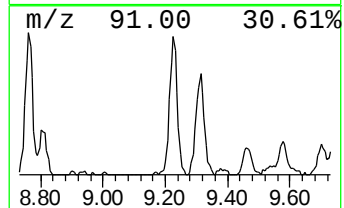
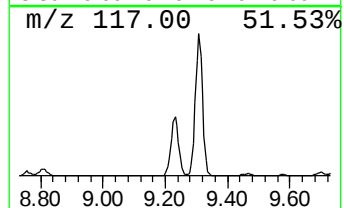
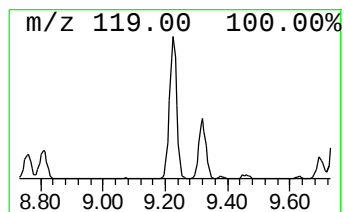
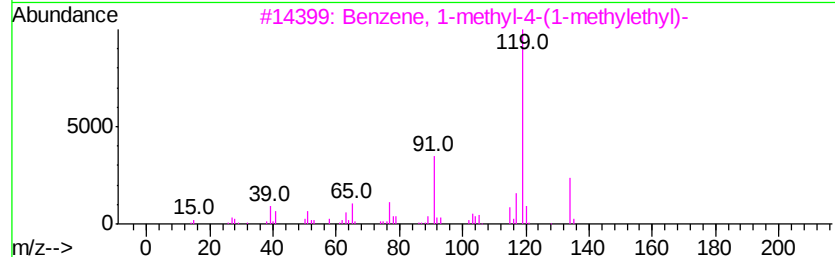
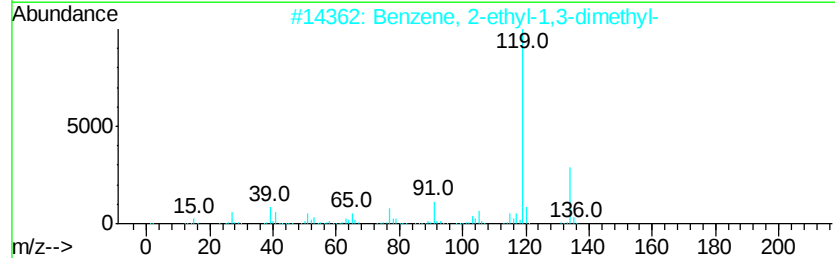
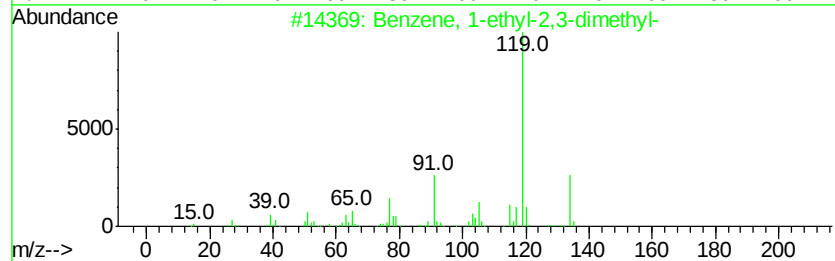
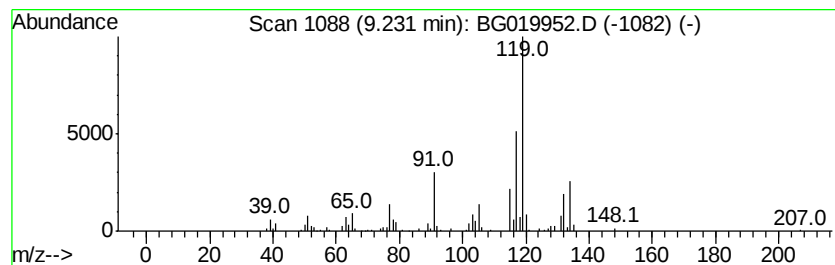
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 3

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 9.23 | 16.49 ng | 180701 | 1,4-Dichlorobenzene-d4 | 8.11 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-2,3-dimethyl-       | 134 | C10H14  | 000933-98-2 | 92   |
| 2    |    | Benzene, 2-ethyl-1,3-dimethyl-       | 134 | C10H14  | 002870-04-4 | 70   |
| 3    |    | Benzene, 1-methyl-4-(1-methylethyl)- | 134 | C10H14  | 000099-87-6 | 70   |
| 4    |    | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 70   |
| 5    |    | Benzene, 1-ethyl-2,4-dimethyl-       | 134 | C10H14  | 000874-41-9 | 70   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120515\  
Data File : BG019952.D  
Acq On : 5 Dec 2015 21:31  
Operator : UM/NP  
Sample : G4630-11  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampled :  
MW-18

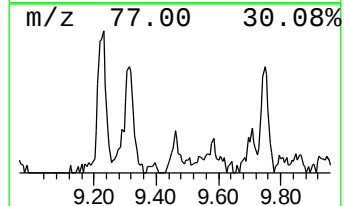
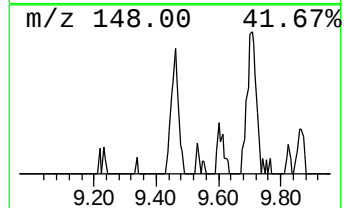
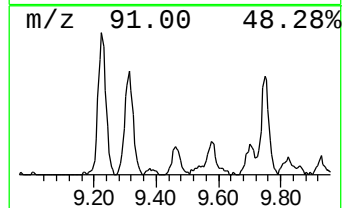
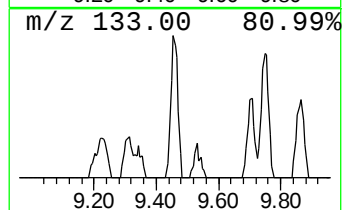
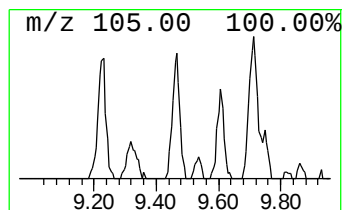
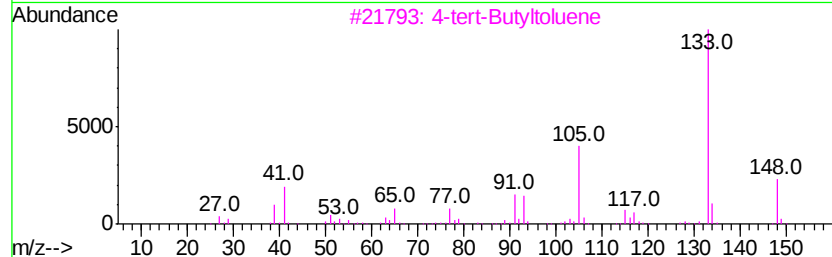
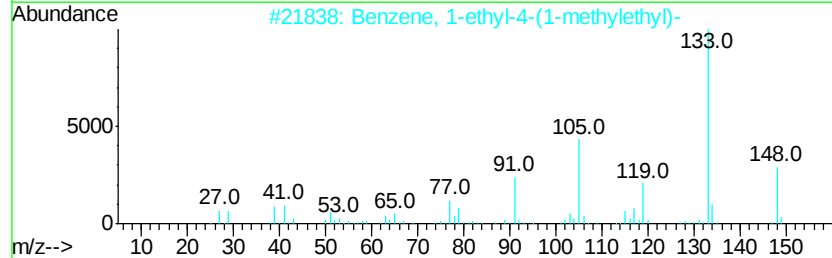
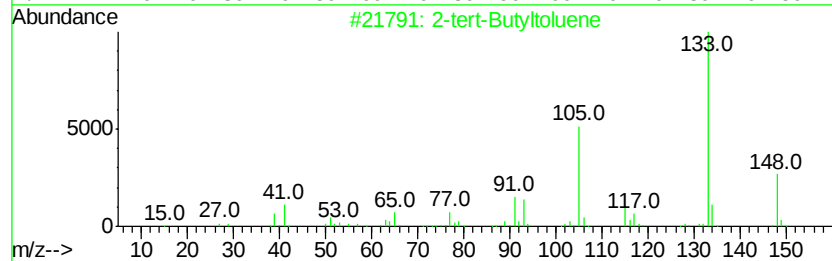
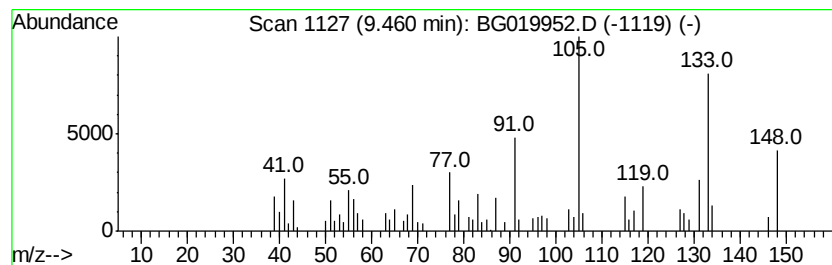
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 2-tert-Butyltoluene Concentration Rank 15

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 9.46 | 7.62 ng | 83540 | 1,4-Dichlorobenzene-d4 | 8.11 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | 2-tert-Butyltoluene                 | 148 | C11H16    | 001074-92-6 | 58   |
| 2    |    |   | Benzene, 1-ethyl-4-(1-methylethyl)- | 148 | C11H16    | 004218-48-8 | 50   |
| 3    |    |   | 4-tert-Butyltoluene                 | 148 | C11H16    | 000098-51-1 | 47   |
| 4    |    |   | Ethanone, 1-(4-ethylphenyl)-        | 148 | C10H12O   | 000937-30-4 | 43   |
| 5    |    |   | 2-Nitrophenyl-.beta.-phenylpropi... | 271 | C15H13NO4 | 040123-51-1 | 38   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120515\  
Data File : BG019952.D  
Acq On : 5 Dec 2015 21:31  
Operator : UM/NP  
Sample : G4630-11  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

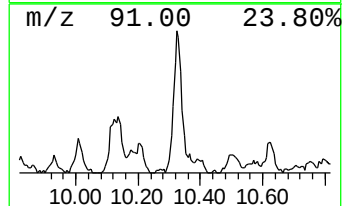
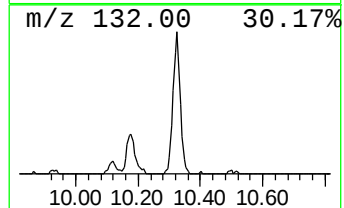
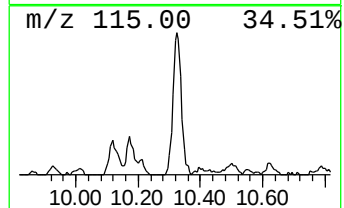
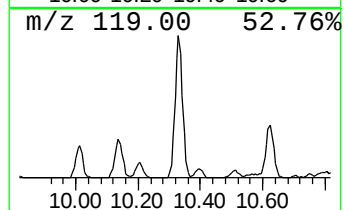
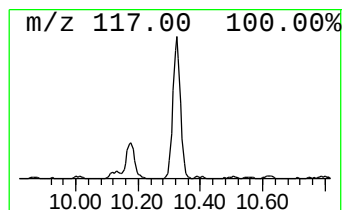
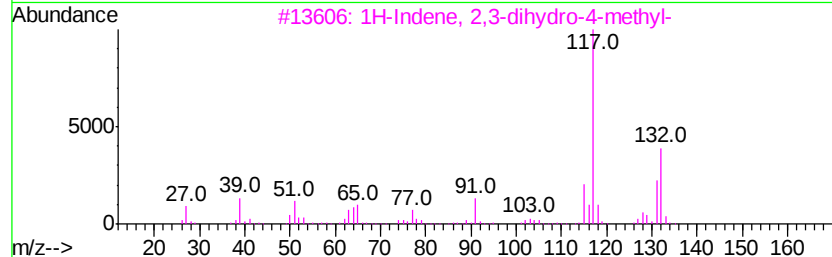
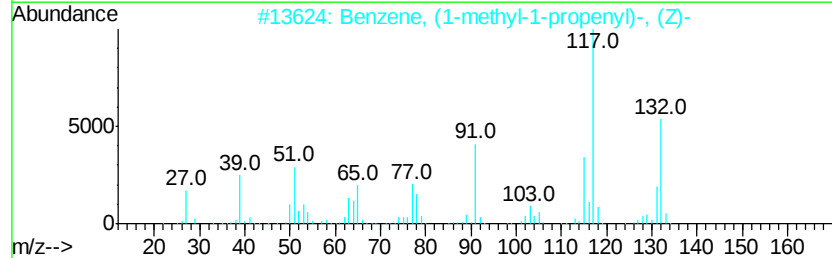
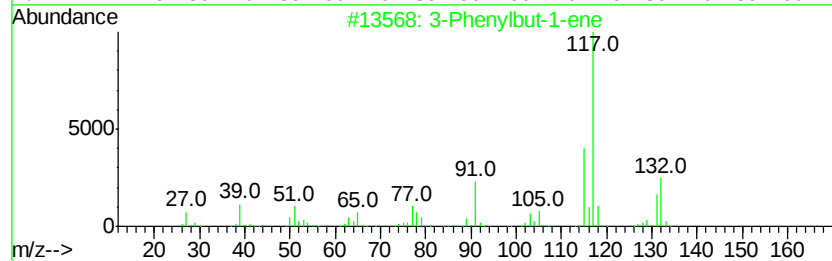
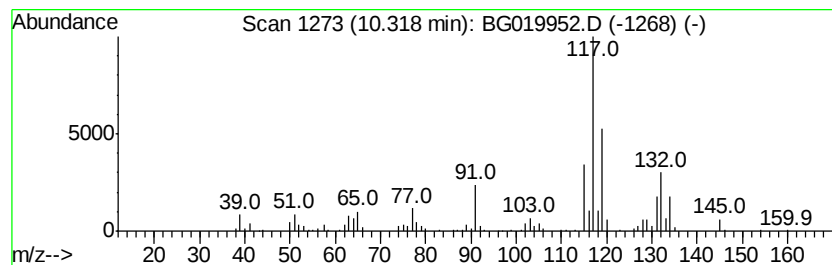
Instrument :  
BNA\_G  
ClientSampleId :  
MW-18

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 11 3-Phenylbut-1-ene Concentration Rank 5

| R.T.    | EstConc  | Area                                 | Relative to ISTD | R.T.           |
|---------|----------|--------------------------------------|------------------|----------------|
| 10.32   | 13.53 ng | 320764                               | Naphthalene-d8   | 10.93          |
| Hit# of | 5        | Tentative ID                         | MW MolForm       | CAS# Qual      |
| 1       |          | 3-Phenylbut-1-ene                    | 132 C10H12       | 000934-10-1 70 |
| 2       |          | Benzene, (1-methyl-1-propenyl)-, ... | 132 C10H12       | 000767-99-7 64 |
| 3       |          | 1H-Indene, 2,3-dihydro-4-methyl-     | 132 C10H12       | 000824-22-6 64 |
| 4       |          | Benzene, 2-ethenyl-1,3-dimethyl-     | 132 C10H12       | 002039-90-9 62 |
| 5       |          | 1-Phenyl-1-butene                    | 132 C10H12       | 000824-90-8 62 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120515\  
Data File : BG019952.D  
Acq On : 5 Dec 2015 21:31  
Operator : UM/NP  
Sample : G4630-11  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-18

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

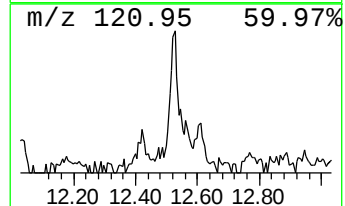
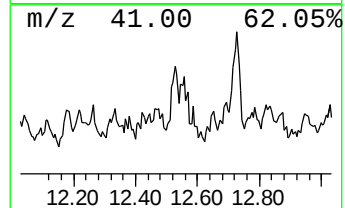
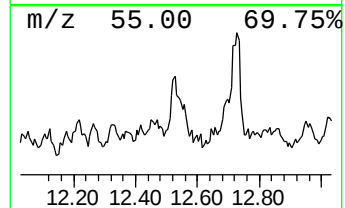
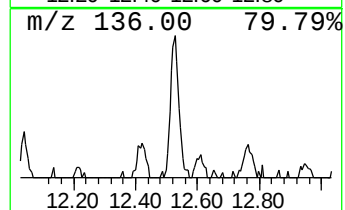
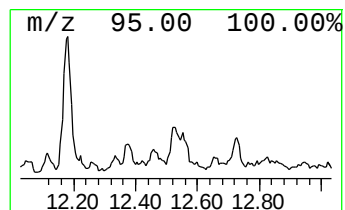
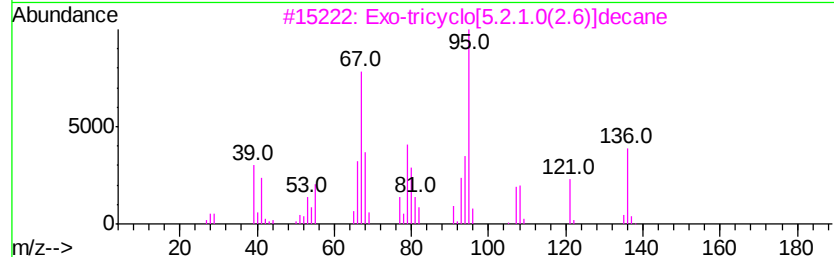
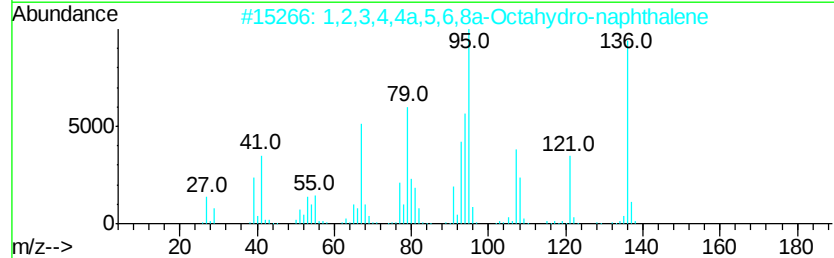
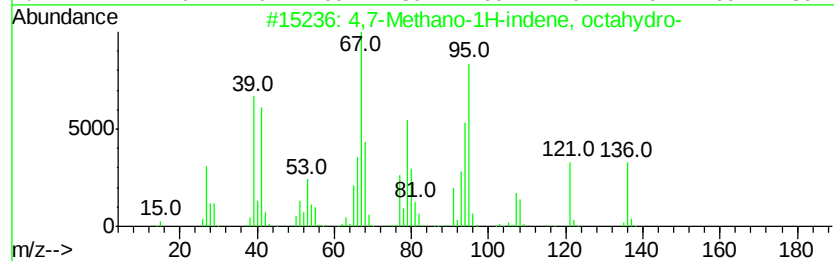
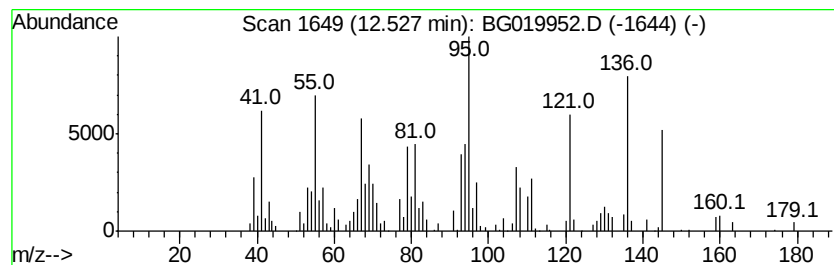
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 4,7-Methano-1H-indene, octa... Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.53 | 9.84 ng | 233241 | Naphthalene-d8   | 10.93 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | 4,7-Methano-1H-indene, octahydro-   | 136 | C10H16  | 006004-38-2  | 95   |
| 2    |    | 1,2,3,4,4a,5,6,8a-Octahydro-naph... | 136 | C10H16  | 031244-58-3  | 86   |
| 3    |    | Exo-tricyclo[5.2.1.0(2.6)]decane    | 136 | C10H16  | 1000215-28-7 | 64   |
| 4    |    | Endo-tricyclo[5.2.1.0(2.6)]decane   | 136 | C10H16  | 1000215-28-8 | 64   |
| 5    |    | Tricyclo[5.2.1.0(4,8)]decane        | 136 | C10H16  | 042836-61-3  | 53   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120515\  
Data File : BG019952.D  
Acq On : 5 Dec 2015 21:31  
Operator : UM/NP  
Sample : G4630-11  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-18

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

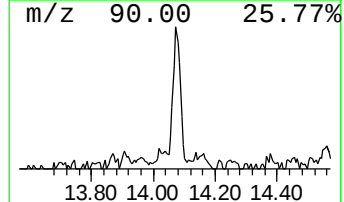
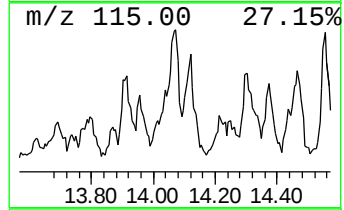
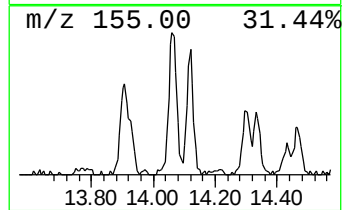
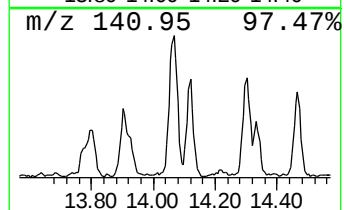
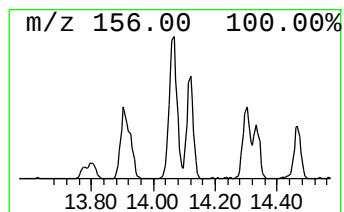
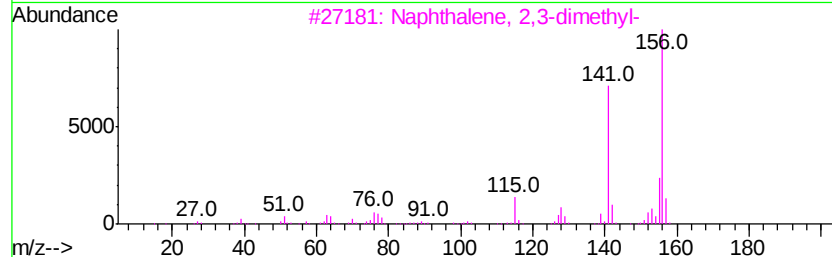
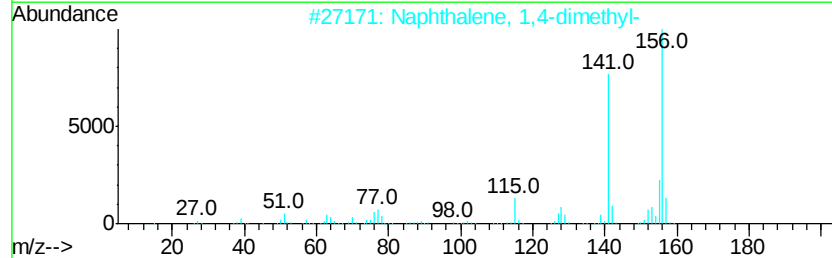
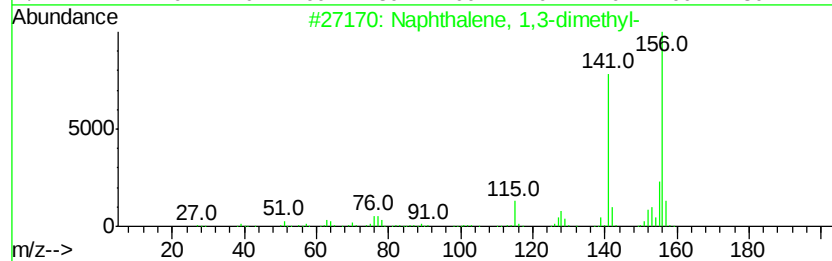
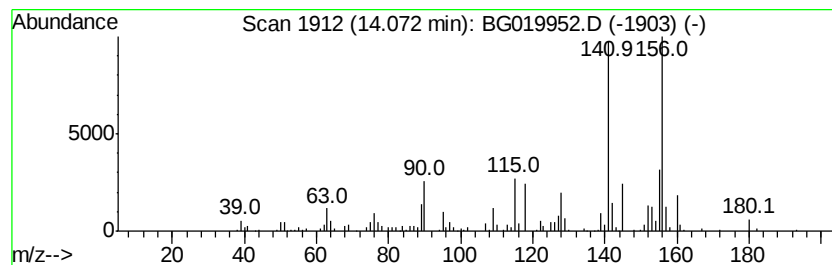
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Naphthalene, 1,3-dimethyl- Concentration Rank 6

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 14.07 | 13.31 ng | 391709 | Acenaphthene-d10 | 14.74 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 97   |
| 2    |    | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 96   |
| 3    |    | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 96   |
| 4    |    | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 95   |
| 5    |    | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 95   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120515\  
Data File : BG019952.D  
Acq On : 5 Dec 2015 21:31  
Operator : UM/NP  
Sample : G4630-11  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-18

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

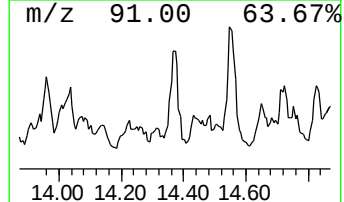
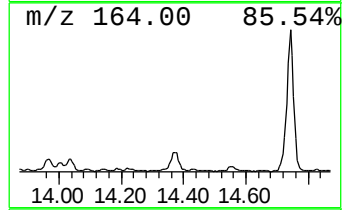
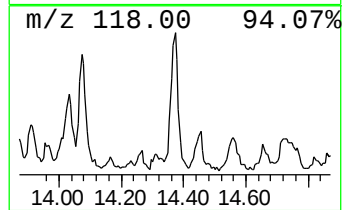
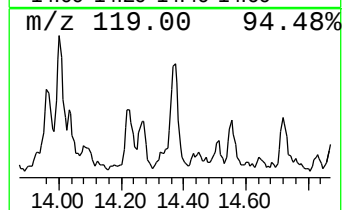
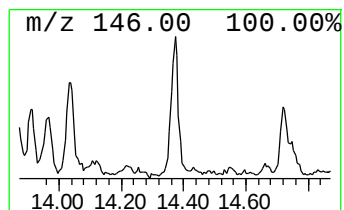
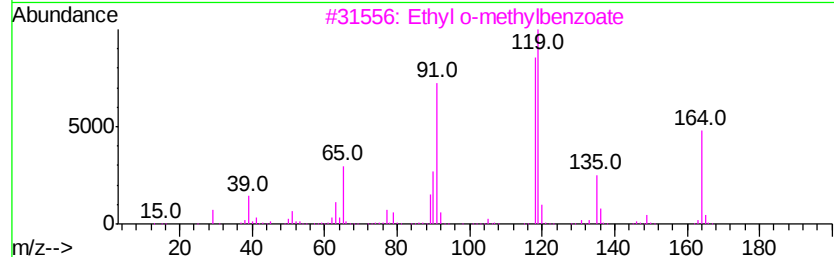
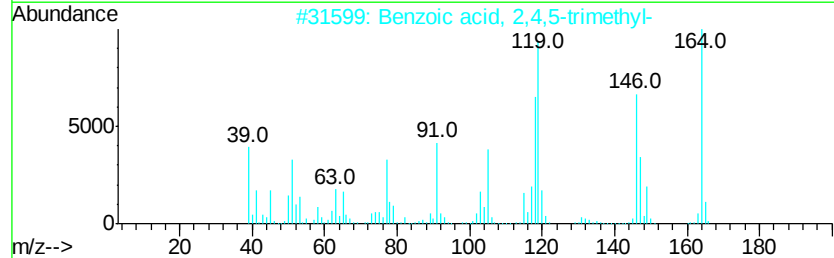
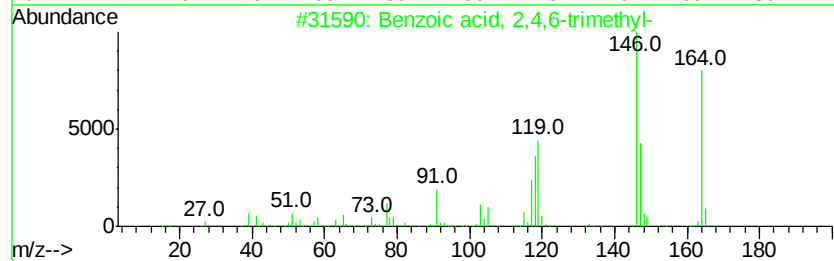
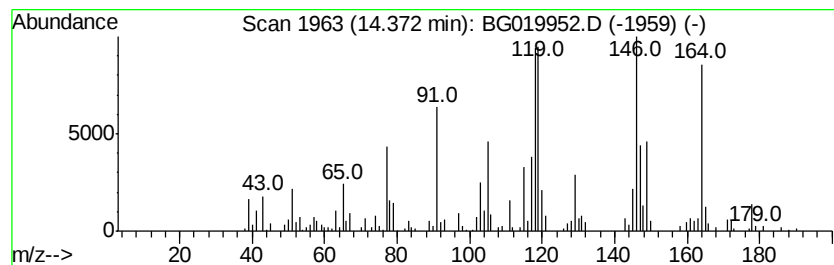
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 Benzoic acid, 2,4,6-trimethyl- Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.37 | 8.00 ng | 235418 | Acenaphthene-d10 | 14.74 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Benzoic acid, 2,4,6-trimethyl-      | 164 | C10H12O2 | 000480-63-7 | 70   |
| 2    |    | Benzoic acid, 2,4,5-trimethyl-      | 164 | C10H12O2 | 000528-90-5 | 62   |
| 3    |    | Ethyl o-methylbenzoate              | 164 | C10H12O2 | 000087-24-1 | 45   |
| 4    |    | 2-Propenoic acid, 3-(3-hydroxyph... | 164 | C9H8O3   | 000588-30-7 | 43   |
| 5    |    | Naphthalene, 1,2,3,4-tetrahydro...  | 146 | C11H14   | 001680-51-9 | 38   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120515\  
Data File : BG019952.D  
Acq On : 5 Dec 2015 21:31  
Operator : UM/NP  
Sample : G4630-11  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-18

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

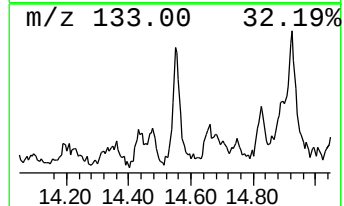
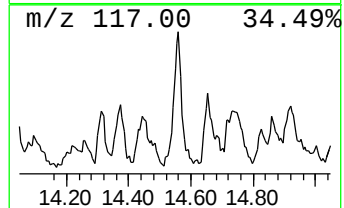
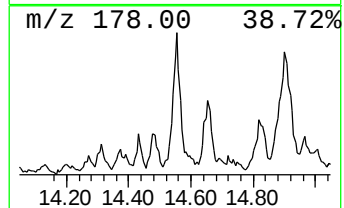
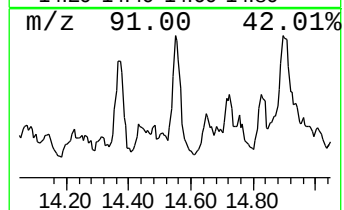
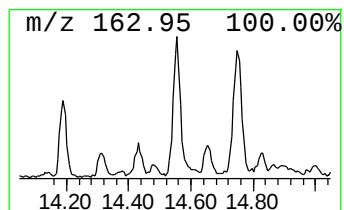
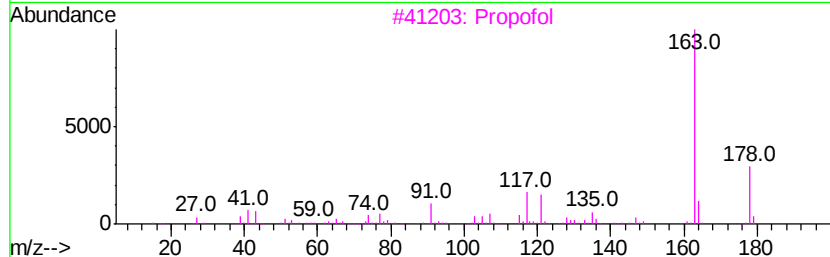
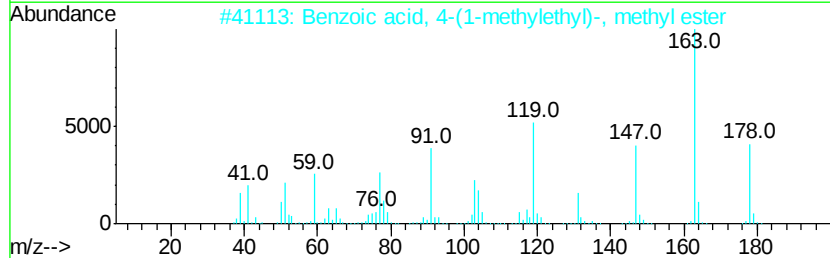
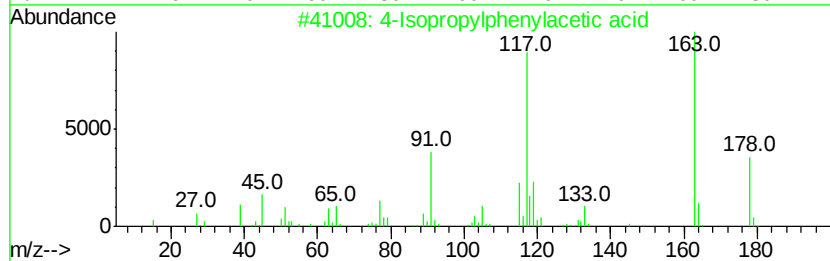
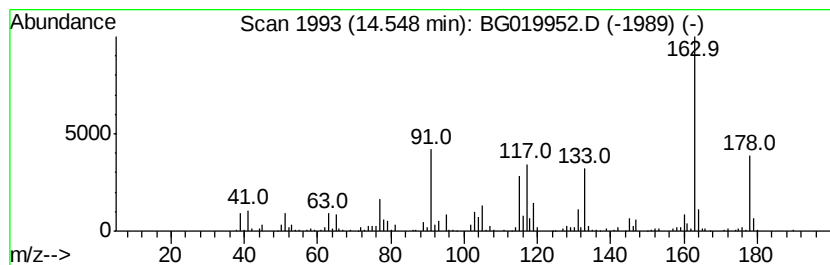
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 4-Isopropylphenylacetic acid Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.55 | 8.11 ng | 238666 | Acenaphthene-d10 | 14.74 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 4-Isopropylphenylacetic acid        | 178 | C11H14O2 | 004476-28-2 | 64   |
| 2    |    | Benzoic acid, 4-(1-methylethyl)-... | 178 | C11H14O2 | 020185-55-1 | 53   |
| 3    |    | Propofol                            | 178 | C12H18O  | 002078-54-8 | 52   |
| 4    |    | Benzoic acid, p-tert-butyl-         | 178 | C11H14O2 | 000098-73-7 | 46   |
| 5    |    | p-Allylnitrobenzene                 | 163 | C9H9NO2  | 053483-17-3 | 46   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120515\  
Data File : BG019952.D  
Acq On : 5 Dec 2015 21:31  
Operator : UM/NP  
Sample : G4630-11  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-18

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

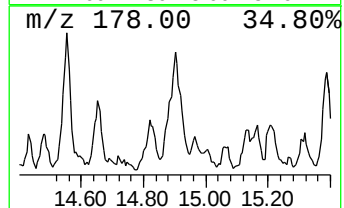
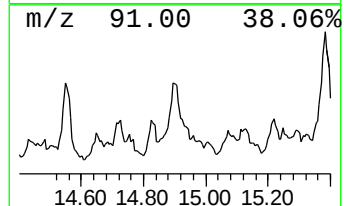
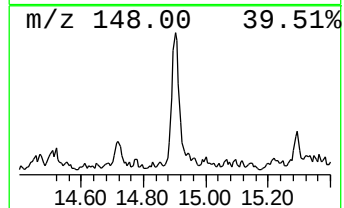
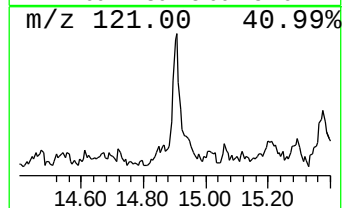
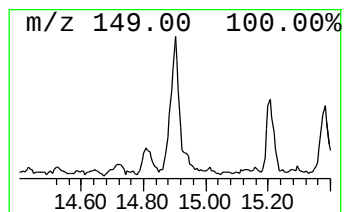
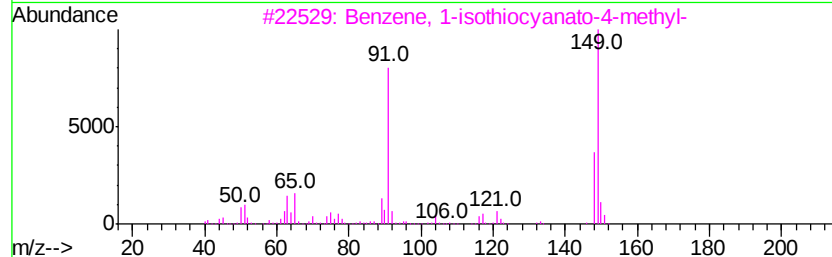
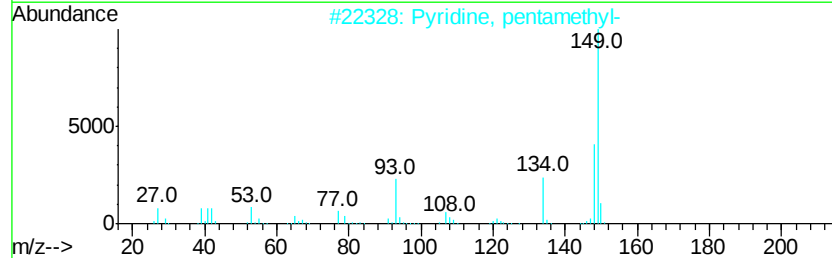
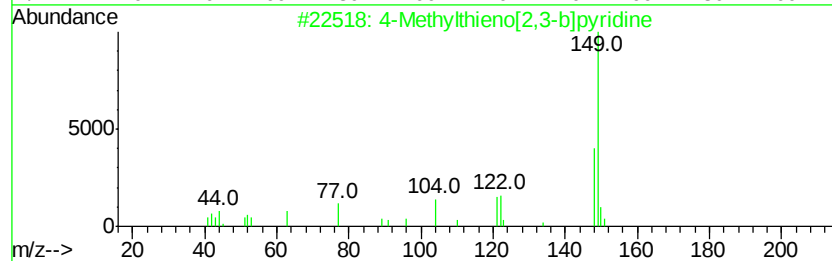
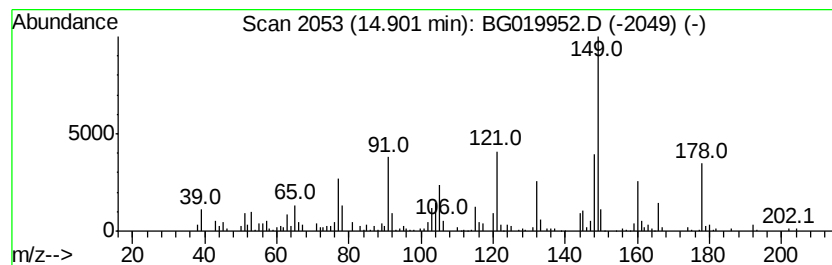
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 unknown14.90 Concentration Rank 9

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 14.90 | 10.30 ng | 303113 | Acenaphthene-d10 | 14.74 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 4-Methylthieno[2,3-b]pyridine       | 149 | C8H7NS       | 013362-81-7 | 47      |      |      |
| 2    | Pyridine, pentamethyl-              | 149 | C10H15N      | 003748-83-2 | 46      |      |      |
| 3    | Benzene, 1-isothiocyanato-4-methyl- | 149 | C8H7NS       | 000622-59-3 | 46      |      |      |
| 4    | 3-(3-Pyridyl)propenoic acid         | 149 | C8H7NO2      | 001126-74-5 | 43      |      |      |
| 5    | 6-Methylthieno[2,3-b]pyridine       | 149 | C8H7NS       | 001759-30-4 | 38      |      |      |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120515\  
Data File : BG019952.D  
Acq On : 5 Dec 2015 21:31  
Operator : UM/NP  
Sample : G4630-11  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-18

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

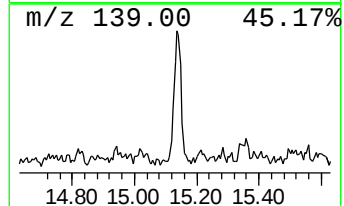
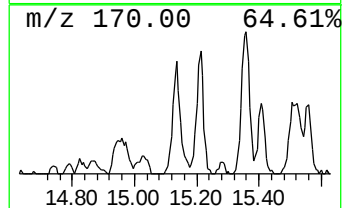
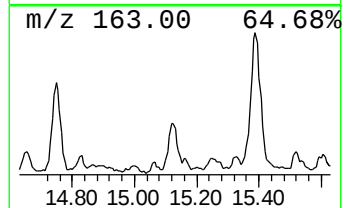
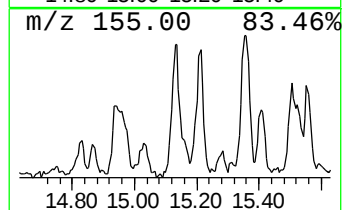
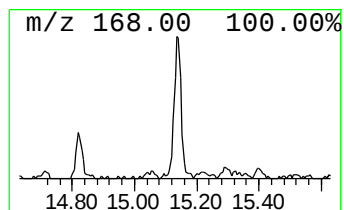
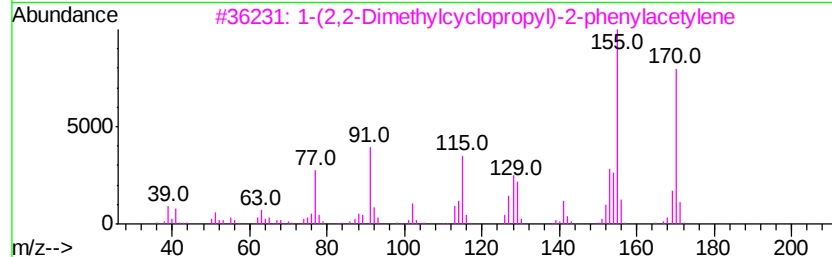
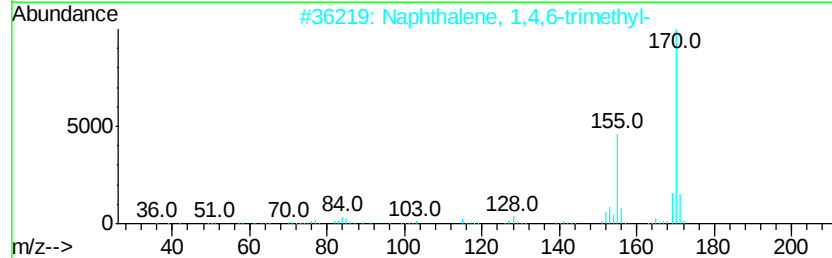
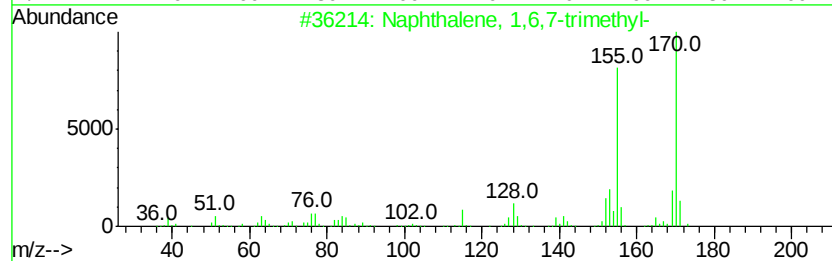
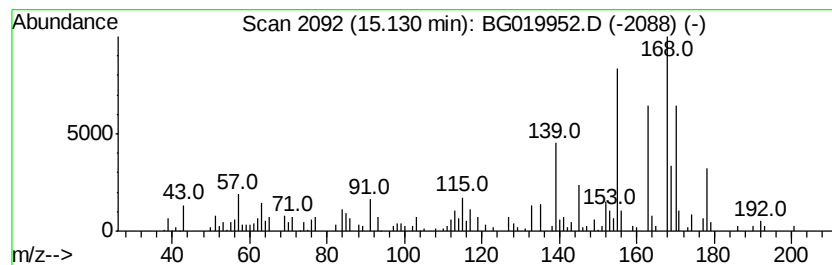
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 unknown15.13 Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.13 | 7.11 ng | 209273 | Acenaphthene-d10 | 14.74 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Naphthalene, 1,6,7-trimethyl-       | 170 | C13H14  | 002245-38-7  | 42   |
| 2    |    | Naphthalene, 1,4,6-trimethyl-       | 170 | C13H14  | 002131-42-2  | 38   |
| 3    |    | 1-(2,2-Dimethylcyclopropyl)-2-ph... | 170 | C13H14  | 1000294-91-1 | 38   |
| 4    |    | 3-(2-Methyl-propenyl)-1H-indene     | 170 | C13H14  | 1000187-78-5 | 38   |
| 5    |    | Azulene, 4,6,8-trimethyl-           | 170 | C13H14  | 000941-81-1  | 38   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG120515\  
Data File : BG019952.D  
Acq On : 5 Dec 2015 21:31  
Operator : UM/NP  
Sample : G4630-11  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

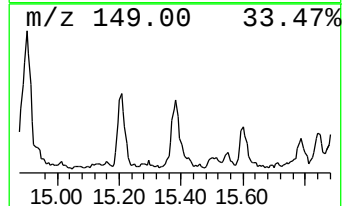
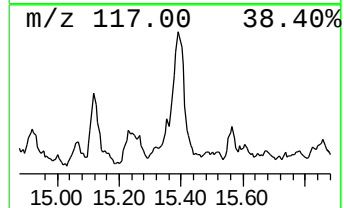
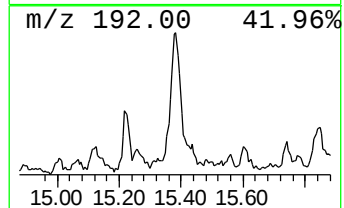
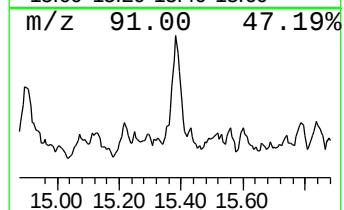
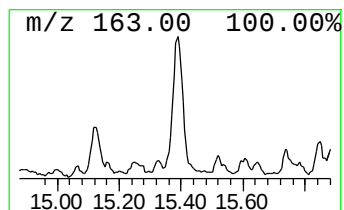
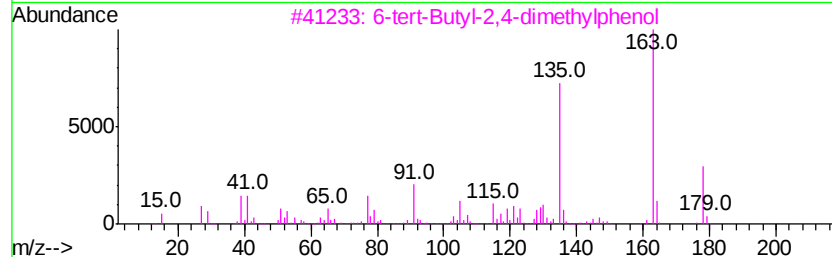
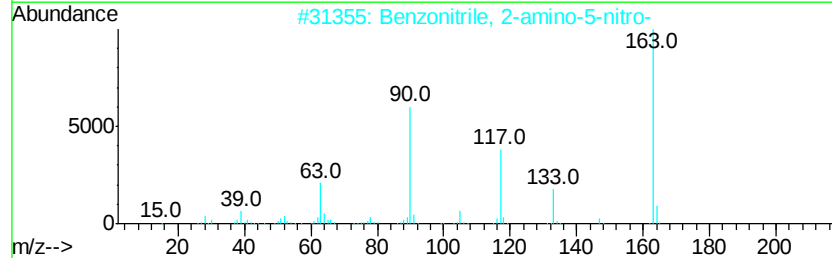
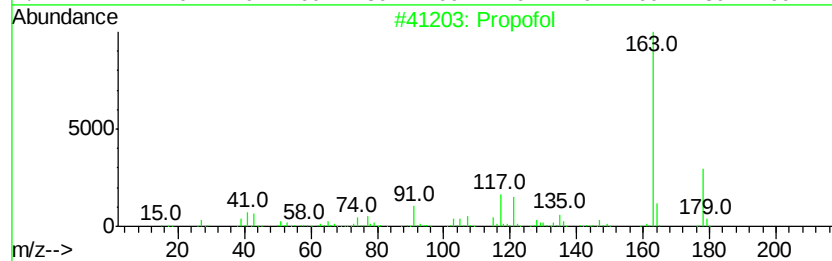
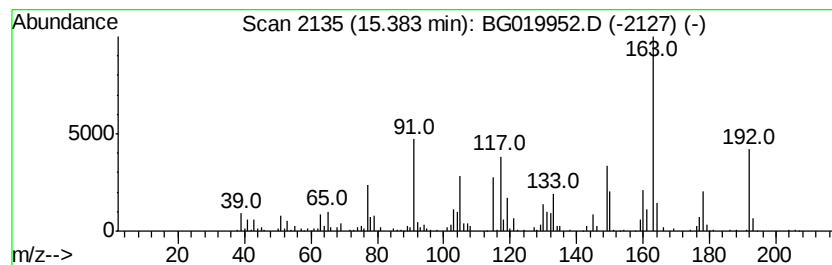
Instrument :  
BNA\_G  
ClientSampleId :  
MW-18

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 20 unknown15.38 Concentration Rank 4

| R.T.    | EstConc                         | Area         | Relative to ISTD | R.T.           |
|---------|---------------------------------|--------------|------------------|----------------|
| 15.38   | 14.23 ng                        | 418610       | Acenaphthene-d10 | 14.74          |
| Hit# of | 5                               | Tentative ID | MW MolForm       | CAS# Qual      |
| 1       | Propofol                        |              | 178 C12H18O      | 002078-54-8 38 |
| 2       | Benzonitrile, 2-amino-5-nitro-  |              | 163 C7H5N3O2     | 017420-30-3 38 |
| 3       | 6-tert-Butyl-2,4-dimethylphenol |              | 178 C12H18O      | 001879-09-0 30 |
| 4       | Benzonitrile, 2,4-dimethoxy-    |              | 163 C9H9NO2      | 004107-65-7 30 |
| 5       | Benzoic acid, p-tert-butyl-     |              | 178 C11H14O2     | 000098-73-7 27 |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG120515\  
 Data File : BG019952.D  
 Acq On : 5 Dec 2015 21:31  
 Operator : UM/NP  
 Sample : G4630-11  
 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MW-18

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG120215.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT 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 |
| 2-Pentanone, 4-hy... | 5.19  | 7.1     | ng    | 77868    | 1                     | 8.11  | 219176 | 20.0 |
| Benzene, (1-methy... | 6.59  | 6.9     | ng    | 75505    | 1                     | 8.11  | 219176 | 20.0 |
| unknown7.64          | 7.64  | 88.2    | ng    | 966045   | 1                     | 8.11  | 219176 | 20.0 |
| Indane               | 8.49  | 12.3    | ng    | 134913   | 1                     | 8.11  | 219176 | 20.0 |
| Benzene, 1,4-diet... | 8.61  | 7.0     | ng    | 77076    | 1                     | 8.11  | 219176 | 20.0 |
| Benzene, 1-ethyl-... | 9.23  | 16.5    | ng    | 180701   | 1                     | 8.11  | 219176 | 20.0 |
| 2-tert-Butyltoluene  | 9.46  | 7.6     | ng    | 83540    | 1                     | 8.11  | 219176 | 20.0 |
| 3-Phenylbut-1-ene    | 10.32 | 13.5    | ng    | 320764   | 2                     | 10.93 | 474173 | 20.0 |
| 4,7-Methano-1H-in... | 12.53 | 9.8     | ng    | 233241   | 2                     | 10.93 | 474173 | 20.0 |
| Naphthalene, 1,3-... | 14.07 | 13.3    | ng    | 391709   | 3                     | 14.74 | 588381 | 20.0 |
| Benzoic acid, 2,4... | 14.37 | 8.0     | ng    | 235418   | 3                     | 14.74 | 588381 | 20.0 |
| 4-Isopropylphenyl... | 14.55 | 8.1     | ng    | 238666   | 3                     | 14.74 | 588381 | 20.0 |
| unknown14.90         | 14.90 | 10.3    | ng    | 303113   | 3                     | 14.74 | 588381 | 20.0 |
| unknown15.13         | 15.13 | 7.1     | ng    | 209273   | 3                     | 14.74 | 588381 | 20.0 |
| unknown15.38         | 15.38 | 14.2    | ng    | 418610   | 3                     | 14.74 | 588381 | 20.0 |