

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
 Data File : BG048872.D  
 Acq On : 23 Apr 2021 20:53  
 Operator : CG/JU  
 Sample : M2107-19  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SB-34-9.5-10.0

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG041421.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG048872.D\data.ms

| 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         | 4.180       | 138           | 143         | 155          | rVB3     | 6954           | 14156         | 1.73%           | 0.141%        |
| 2         | 5.608       | 379           | 386         | 395          | rBV      | 336392         | 539401        | 66.05%          | 5.377%        |
| 3         | 6.924       | 603           | 610         | 617          | rVB5     | 4093           | 9591          | 1.17%           | 0.096%        |
| 4         | 7.223       | 649           | 661         | 671          | rBV      | 351229         | 626565        | 76.73%          | 6.246%        |
| 5         | 7.541       | 710           | 715         | 722          | rVB6     | 9728           | 19055         | 2.33%           | 0.190%        |
| 6         | 8.017       | 791           | 796         | 798          | rBV3     | 12735          | 17966         | 2.20%           | 0.179%        |
| 7         | 8.058       | 798           | 803         | 808          | rVV2     | 84520          | 138416        | 16.95%          | 1.380%        |
| 8         | 8.105       | 810           | 811         | 817          | rVB4     | 6572           | 8824          | 1.08%           | 0.088%        |
| 9         | 8.187       | 817           | 825         | 829          | rBV2     | 32673          | 54795         | 6.71%           | 0.546%        |
| 10        | 8.275       | 837           | 840         | 846          | rBV7     | 7864           | 14869         | 1.82%           | 0.148%        |
| 11        | 8.340       | 846           | 851         | 856          | rVB4     | 14397          | 21345         | 2.61%           | 0.213%        |
| 12        | 8.833       | 930           | 935         | 940          | rVB6     | 11419          | 21481         | 2.63%           | 0.214%        |
| 13        | 8.892       | 940           | 945         | 951          | rBV5     | 10979          | 18796         | 2.30%           | 0.187%        |
| 14        | 9.133       | 981           | 986         | 987          | rBV3     | 6019           | 8611          | 1.05%           | 0.086%        |
| 15        | 9.168       | 987           | 992         | 996          | rVV6     | 12883          | 26792         | 3.28%           | 0.267%        |
| 16        | 9.227       | 996           | 1002        | 1014         | rVB      | 228982         | 415114        | 50.83%          | 4.138%        |
| 17        | 9.380       | 1022          | 1028        | 1029         | rBV3     | 6400           | 9682          | 1.19%           | 0.097%        |
| 18        | 9.415       | 1030          | 1034        | 1042         | rVB5     | 8835           | 17915         | 2.19%           | 0.179%        |
| 19        | 9.697       | 1079          | 1082        | 1086         | rBV4     | 6548           | 9362          | 1.15%           | 0.093%        |
| 20        | 9.791       | 1094          | 1098        | 1101         | rBV2     | 7646           | 10055         | 1.23%           | 0.100%        |
| 21        | 9.991       | 1129          | 1132        | 1137         | rVB      | 11255          | 15271         | 1.87%           | 0.152%        |
| 22        | 10.049      | 1137          | 1142        | 1147         | rBV5     | 10728          | 19781         | 2.42%           | 0.197%        |
| 23        | 10.666      | 1238          | 1247        | 1249         | rBV3     | 4290           | 8738          | 1.07%           | 0.087%        |
| 24        | 10.884      | 1278          | 1284        | 1289         | rVV      | 114146         | 205533        | 25.17%          | 2.049%        |
| 25        | 10.937      | 1289          | 1293        | 1298         | rVB2     | 21795          | 37506         | 4.59%           | 0.374%        |
| 26        | 11.025      | 1304          | 1308        | 1312         | rVB2     | 11122          | 20202         | 2.47%           | 0.201%        |
| 27        | 11.542      | 1390          | 1396        | 1398         | rBV5     | 5634           | 10347         | 1.27%           | 0.103%        |
| 28        | 11.577      | 1398          | 1402        | 1407         | rVB6     | 12719          | 20728         | 2.54%           | 0.207%        |
| 29        | 11.700      | 1422          | 1423        | 1429         | rVB5     | 5845           | 9516          | 1.17%           | 0.095%        |
| 30        | 11.894      | 1447          | 1456        | 1460         | rBV2     | 50929          | 90736         | 11.11%          | 0.905%        |
| 31        | 11.930      | 1460          | 1462        | 1466         | rVV4     | 12540          | 17484         | 2.14%           | 0.174%        |
| 32        | 11.977      | 1466          | 1470        | 1476         | rVB7     | 9475           | 18913         | 2.32%           | 0.189%        |
| 33        | 12.212      | 1505          | 1510        | 1516         | rBV8     | 6210           | 12614         | 1.54%           | 0.126%        |
| 34        | 12.270      | 1516          | 1520        | 1526         | rBV6     | 8497           | 20661         | 2.53%           | 0.206%        |
| 35        | 12.364      | 1533          | 1536        | 1540         | rBV5     | 7537           | 13369         | 1.64%           | 0.133%        |
| 36        | 12.499      | 1553          | 1559        | 1564         | rVB7     | 12820          | 26939         | 3.30%           | 0.269%        |

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 Sample : M2107-19  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SB-34-9.5-10.0

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG041421.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 37 | 12.729 | 1594 | 1598 | 1600 | rBV5 | 5516   | 9427   | 1.15%  | 0.094% |
| 38 | 12.917 | 1626 | 1630 | 1633 | rBV5 | 12783  | 16968  | 2.08%  | 0.169% |
| 39 | 12.981 | 1637 | 1641 | 1650 | rVB3 | 30980  | 59126  | 7.24%  | 0.589% |
| 40 | 13.111 | 1661 | 1663 | 1667 | rVB5 | 10183  | 9843   | 1.21%  | 0.098% |
| 41 | 13.234 | 1679 | 1684 | 1691 | rVB2 | 103868 | 151425 | 18.54% | 1.510% |
| 42 | 13.334 | 1693 | 1701 | 1712 | rVB  | 426359 | 668469 | 81.86% | 6.664% |
| 43 | 13.498 | 1721 | 1729 | 1731 | rBV6 | 24580  | 46431  | 5.69%  | 0.463% |
| 44 | 13.592 | 1741 | 1745 | 1749 | rBV2 | 48547  | 79643  | 9.75%  | 0.794% |
| 45 | 13.763 | 1772 | 1774 | 1780 | rVB7 | 17716  | 23284  | 2.85%  | 0.232% |
| 46 | 13.986 | 1807 | 1812 | 1816 | rBV7 | 24490  | 33439  | 4.09%  | 0.333% |
| 47 | 14.039 | 1816 | 1821 | 1825 | rVV6 | 28314  | 50135  | 6.14%  | 0.500% |
| 48 | 14.086 | 1826 | 1829 | 1833 | rVV4 | 24065  | 31069  | 3.80%  | 0.310% |
| 49 | 14.133 | 1833 | 1837 | 1840 | rBV4 | 46560  | 66485  | 8.14%  | 0.663% |
| 50 | 14.180 | 1840 | 1845 | 1851 | rVB3 | 170299 | 235168 | 28.80% | 2.344% |
| 51 | 14.227 | 1851 | 1853 | 1855 | rBV3 | 9514   | 12077  | 1.48%  | 0.120% |
| 52 | 14.397 | 1880 | 1882 | 1885 | rVB4 | 9803   | 8618   | 1.06%  | 0.086% |
| 53 | 14.444 | 1886 | 1890 | 1891 | rBV4 | 19711  | 22092  | 2.71%  | 0.220% |
| 54 | 14.538 | 1901 | 1906 | 1911 | rVB5 | 52574  | 84097  | 10.30% | 0.838% |
| 55 | 14.679 | 1924 | 1930 | 1932 | rBV5 | 36229  | 71193  | 8.72%  | 0.710% |
| 56 | 14.715 | 1932 | 1936 | 1942 | rVV  | 165219 | 275071 | 33.68% | 2.742% |
| 57 | 14.762 | 1942 | 1944 | 1947 | rVV3 | 11009  | 15071  | 1.85%  | 0.150% |
| 58 | 14.797 | 1947 | 1950 | 1955 | rVB7 | 18610  | 26141  | 3.20%  | 0.261% |
| 59 | 14.926 | 1967 | 1972 | 1980 | rVB8 | 31568  | 71834  | 8.80%  | 0.716% |
| 60 | 15.002 | 1982 | 1985 | 1986 | rBV3 | 14595  | 16539  | 2.03%  | 0.165% |
| 61 | 15.102 | 1997 | 2002 | 2008 | rBV4 | 43195  | 74339  | 9.10%  | 0.741% |
| 62 | 15.185 | 2013 | 2016 | 2019 | rBV2 | 24566  | 34233  | 4.19%  | 0.341% |
| 63 | 15.331 | 2036 | 2041 | 2046 | rBV7 | 39204  | 75629  | 9.26%  | 0.754% |
| 64 | 15.378 | 2047 | 2049 | 2054 | rVB5 | 21186  | 26353  | 3.23%  | 0.263% |
| 65 | 15.431 | 2054 | 2058 | 2062 | rVB7 | 23839  | 32328  | 3.96%  | 0.322% |
| 66 | 15.502 | 2064 | 2070 | 2073 | rBV7 | 41783  | 69474  | 8.51%  | 0.693% |
| 67 | 15.602 | 2083 | 2087 | 2091 | rBV7 | 15748  | 30342  | 3.72%  | 0.302% |
| 68 | 15.760 | 2110 | 2114 | 2117 | rBV5 | 18719  | 32676  | 4.00%  | 0.326% |
| 69 | 15.819 | 2121 | 2124 | 2128 | rVB4 | 42994  | 56941  | 6.97%  | 0.568% |
| 70 | 15.948 | 2138 | 2146 | 2158 | rBV  | 106282 | 210281 | 25.75% | 2.096% |
| 71 | 16.213 | 2186 | 2191 | 2198 | rVB  | 389240 | 544781 | 66.71% | 5.431% |
| 72 | 16.436 | 2221 | 2229 | 2235 | rBV  | 250650 | 389749 | 47.73% | 3.885% |
| 73 | 16.530 | 2241 | 2245 | 2249 | rBV6 | 13268  | 29431  | 3.60%  | 0.293% |
| 74 | 16.565 | 2249 | 2251 | 2253 | rBV3 | 12634  | 13585  | 1.66%  | 0.135% |
| 75 | 17.118 | 2343 | 2345 | 2347 | rBV3 | 9769   | 9191   | 1.13%  | 0.092% |
| 76 | 17.259 | 2364 | 2369 | 2373 | rBV2 | 75069  | 113479 | 13.90% | 1.131% |

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 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SB-34-9.5-10.0

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG041421.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|     |        |      |      |      |      |        |        |         |        |
|-----|--------|------|------|------|------|--------|--------|---------|--------|
| 77  | 17.476 | 2400 | 2406 | 2411 | rBV  | 198692 | 313984 | 38.45%  | 3.130% |
| 78  | 17.870 | 2470 | 2473 | 2476 | rBV5 | 22783  | 30977  | 3.79%   | 0.309% |
| 79  | 18.592 | 2594 | 2596 | 2601 | rVB6 | 25904  | 23821  | 2.92%   | 0.237% |
| 80  | 18.757 | 2619 | 2624 | 2629 | rVB5 | 78536  | 130291 | 15.96%  | 1.299% |
| 81  | 18.945 | 2652 | 2656 | 2663 | rVB8 | 62708  | 96909  | 11.87%  | 0.966% |
| 82  | 19.086 | 2677 | 2680 | 2687 | rVB6 | 46763  | 81266  | 9.95%   | 0.810% |
| 83  | 19.215 | 2697 | 2702 | 2708 | rBV8 | 29671  | 73807  | 9.04%   | 0.736% |
| 84  | 19.268 | 2708 | 2711 | 2716 | rVB6 | 34714  | 44854  | 5.49%   | 0.447% |
| 85  | 19.844 | 2806 | 2809 | 2813 | rBV2 | 39459  | 48358  | 5.92%   | 0.482% |
| 86  | 20.085 | 2846 | 2850 | 2855 | rVB  | 675289 | 816610 | 100.00% | 8.141% |
| 87  | 20.302 | 2883 | 2887 | 2892 | rVB  | 71286  | 95653  | 11.71%  | 0.954% |
| 88  | 20.378 | 2896 | 2900 | 2904 | rVB3 | 63433  | 75422  | 9.24%   | 0.752% |
| 89  | 20.872 | 2981 | 2984 | 2991 | rVB  | 65185  | 85034  | 10.41%  | 0.848% |
| 90  | 21.389 | 3068 | 3072 | 3078 | rVB  | 129662 | 184124 | 22.55%  | 1.835% |
| 91  | 21.577 | 3100 | 3104 | 3111 | rVB3 | 76760  | 110567 | 13.54%  | 1.102% |
| 92  | 21.771 | 3132 | 3137 | 3143 | rVB2 | 189980 | 323619 | 39.63%  | 3.226% |
| 93  | 21.947 | 3164 | 3167 | 3174 | rVB3 | 58101  | 84127  | 10.30%  | 0.839% |
| 94  | 22.570 | 3269 | 3273 | 3280 | rBV3 | 61816  | 127098 | 15.56%  | 1.267% |
| 95  | 23.287 | 3391 | 3395 | 3400 | rVB6 | 34140  | 57274  | 7.01%   | 0.571% |
| 96  | 24.121 | 3531 | 3537 | 3543 | rVB2 | 110856 | 225329 | 27.59%  | 2.246% |
| 97  | 25.108 | 3695 | 3705 | 3712 | rBV3 | 133837 | 379127 | 46.43%  | 3.779% |
| 98  | 26.277 | 3897 | 3904 | 3913 | rVB3 | 68669  | 211829 | 25.94%  | 2.112% |
| 99  | 28.011 | 4197 | 4199 | 4202 | rBV4 | 10952  | 15022  | 1.84%   | 0.150% |
| 100 | 31.847 | 4848 | 4852 | 4853 | rBV4 | 14156  | 18626  | 2.28%   | 0.186% |

Sum of corrected areas: 10031314



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Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

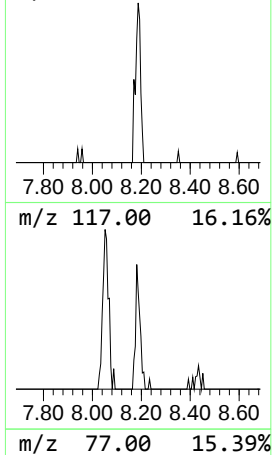
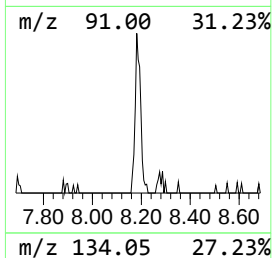
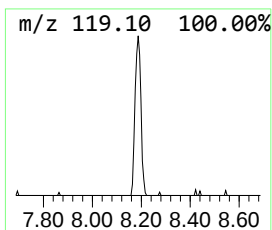
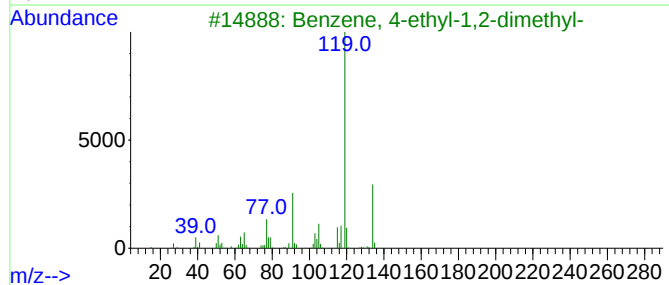
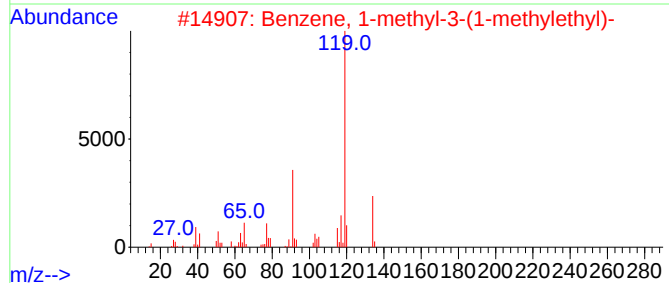
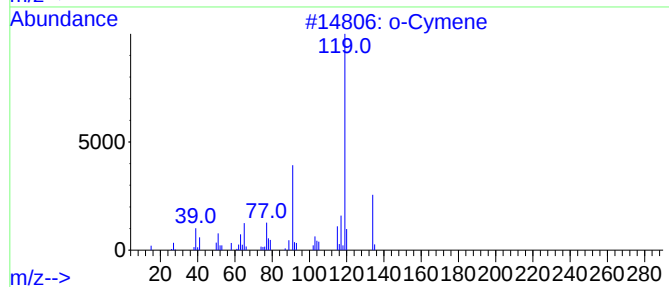
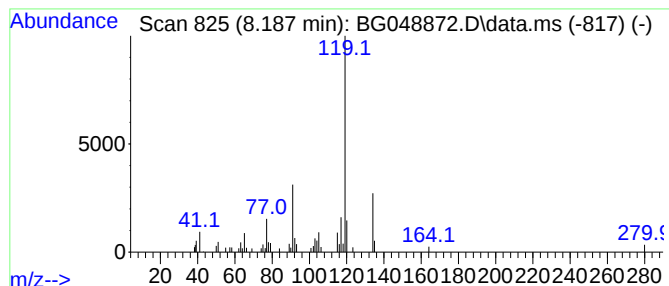
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG041421.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 o-Cymene Concentration Rank 9

| R.T.  | EstConc | Area  | Relative to ISTD       | R.T.  |
|-------|---------|-------|------------------------|-------|
| 8.187 | 7.92 ng | 54795 | 1,4-Dichlorobenzene-d4 | 8.058 |

| Hit# | of 5 | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|------|--------------------------------------|-----|---------|-------------|------|
| 1    |      | o-Cymene                             | 134 | C10H14  | 000527-84-4 | 93   |
| 2    |      | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 93   |
| 3    |      | Benzene, 4-ethyl-1,2-dimethyl-       | 134 | C10H14  | 000934-80-5 | 93   |
| 4    |      | Benzene, 1,2,4,5-tetramethyl-        | 134 | C10H14  | 000095-93-2 | 93   |
| 5    |      | Benzene, 2-ethyl-1,3-dimethyl-       | 134 | C10H14  | 002870-04-4 | 93   |



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ALS Vial : 17 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
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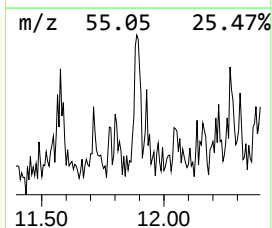
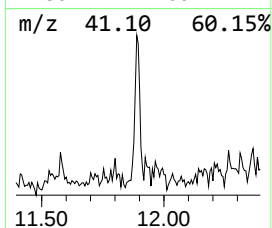
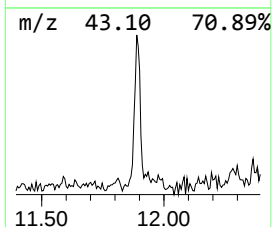
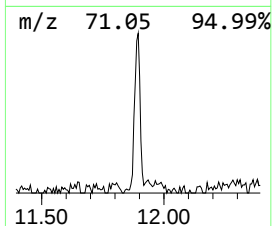
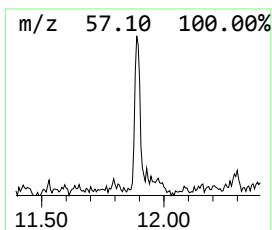
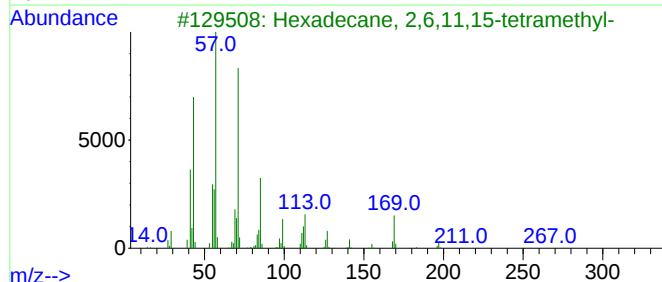
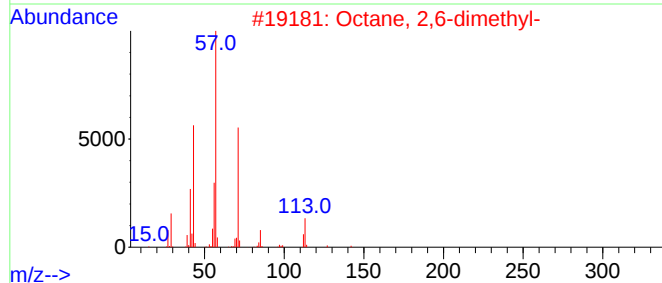
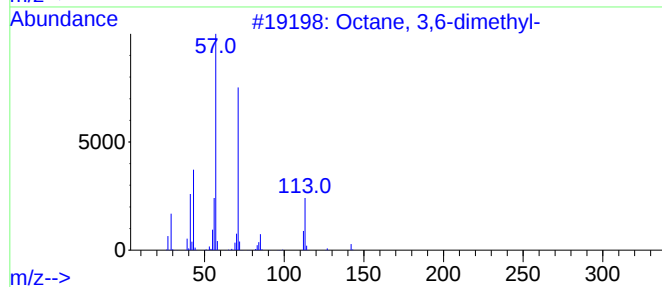
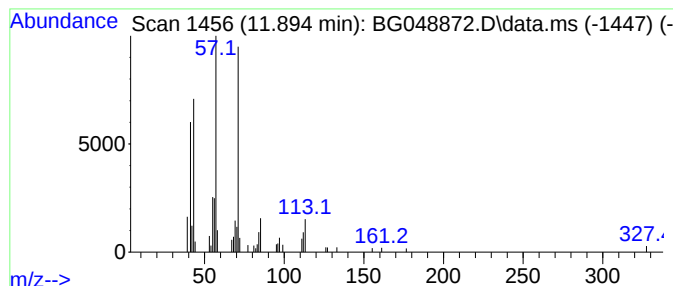
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG041421.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 Octane, 3,6-dimethyl- Concentration Rank 7

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 11.894 | 8.83 ng | 90736 | Naphthalene-d8   | 10.884 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Octane, 3,6-dimethyl-               | 142 | C10H22    | 015869-94-0  | 78   |
| 2    |    |   | Octane, 2,6-dimethyl-               | 142 | C10H22    | 002051-30-1  | 78   |
| 3    |    |   | Hexadecane, 2,6,11,15-tetramethyl-  | 282 | C20H42    | 000504-44-9  | 78   |
| 4    |    |   | Dodecane, 2,6,10-trimethyl-         | 212 | C15H32    | 003891-98-3  | 72   |
| 5    |    |   | Sulfurous acid, 2-ethylhexyl non... | 320 | C17H36O3S | 1000309-19-2 | 72   |



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TIC Integration Parameters: LSCINT.P

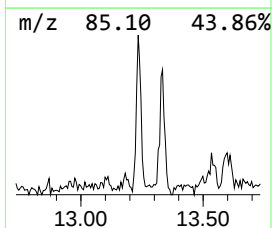
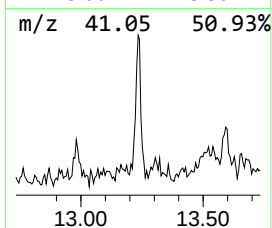
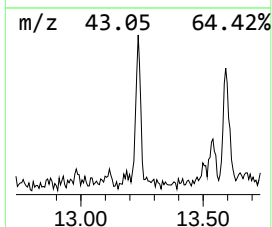
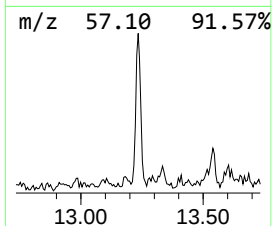
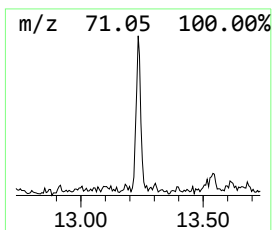
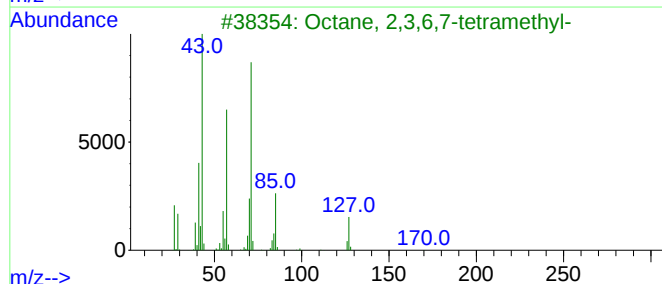
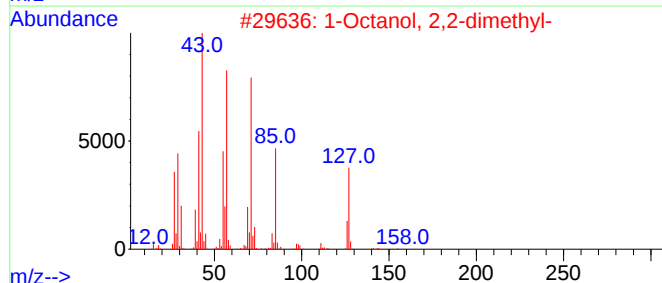
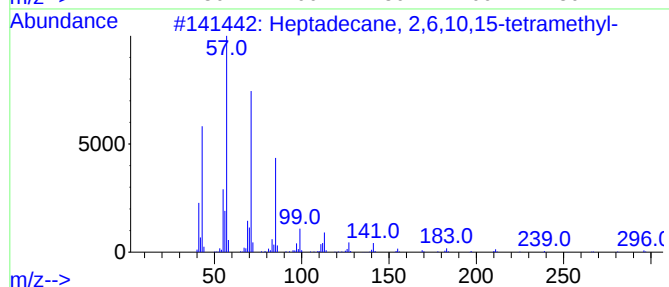
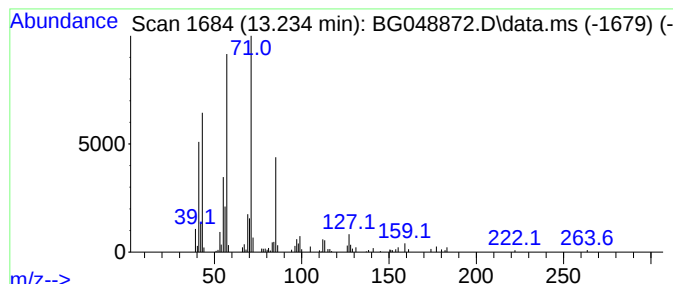
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Peak Number 3 Heptadecane, 2,6,10,15-tetr... Concentration Rank 6

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 13.234 | 11.01 ng | 151425 | Acenaphthene-d10 | 14.720 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|------|-------------------------------------|-----|-----------|--------------|------|
| 1    |      | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44    | 054833-48-6  | 86   |
| 2    |      | 1-Octanol, 2,2-dimethyl-            | 158 | C10H22O   | 002370-14-1  | 72   |
| 3    |      | Octane, 2,3,6,7-tetramethyl-        | 170 | C12H26    | 052670-34-5  | 59   |
| 4    |      | Sulfurous acid, hexyl pentyl ester  | 236 | C11H24O3S | 1000309-14-1 | 53   |
| 5    |      | Sulfurous acid, 2-ethylhexyl tri... | 376 | C21H44O3S | 1000309-19-6 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
Data File : BG048872.D  
Acq On : 23 Apr 2021 20:53  
Operator : CG/JU  
Sample : M2107-19  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

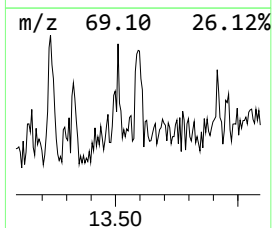
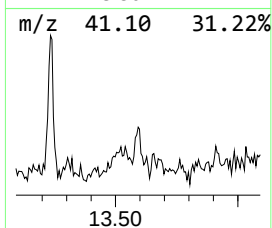
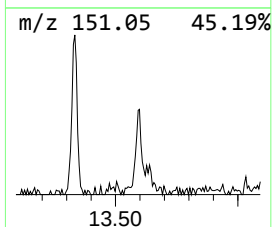
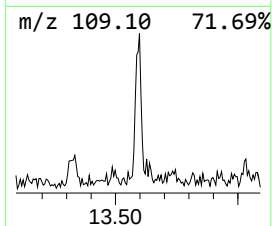
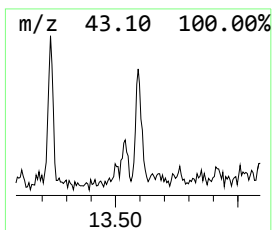
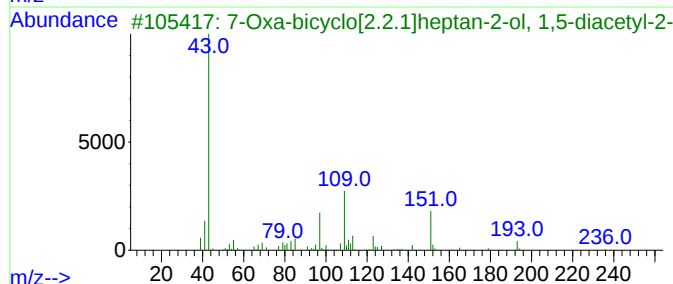
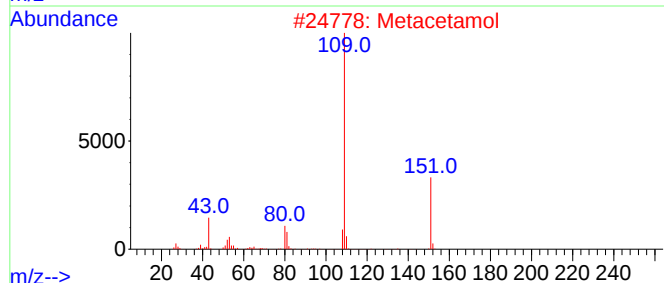
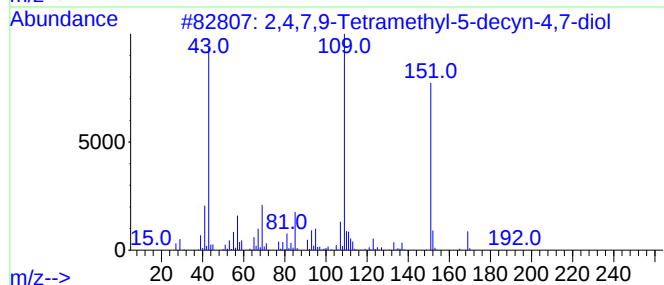
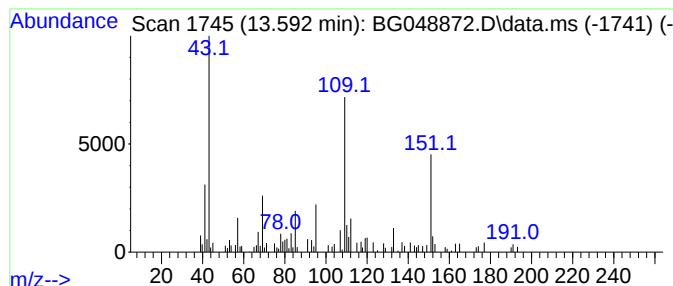
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 4 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 16

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 13.592 | 5.79 ng | 79643 | Acenaphthene-d10 | 14.720 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2,4,7,9-Tetramethyl-5-decyn-4,7-... | 226 | C14H26O2     | 000126-86-3 | 50      |      |      |
| 2    | Metacetamol                         | 151 | C8H9NO2      | 000621-42-1 | 43      |      |      |
| 3    | 7-Oxa-bicyclo[2.2.1]heptan-2-ol,... | 254 | C14H22O4     | 329362-13-2 | 38      |      |      |
| 4    | 3-Pyridinemethanol, acetate         | 151 | C8H9NO2      | 010072-09-0 | 38      |      |      |
| 5    | Camphorsulfonic acid                | 232 | C10H16O4S    | 003144-16-9 | 37      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
Data File : BG048872.D  
Acq On : 23 Apr 2021 20:53  
Operator : CG/JU  
Sample : M2107-19  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG041421.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

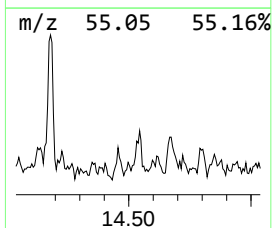
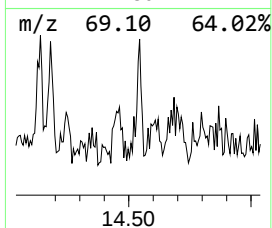
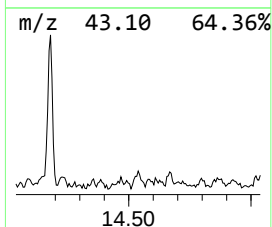
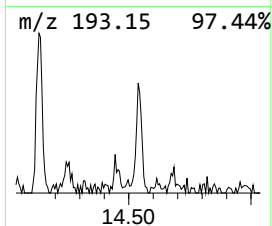
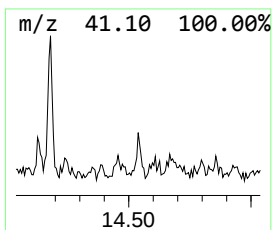
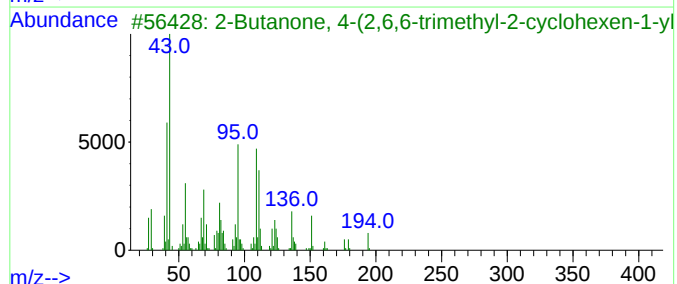
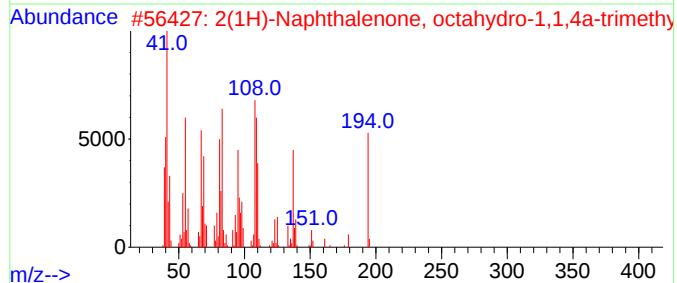
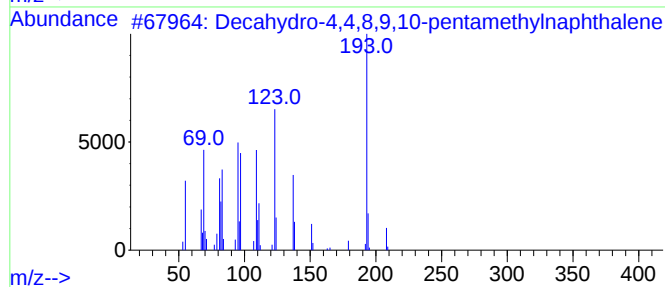
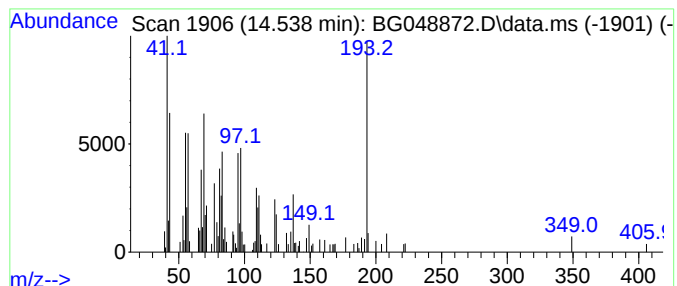
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Peak Number 5 Decahydro-4,4,8,9,10-pentam... Concentration Rank 14

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 14.538 | 6.11 ng | 84097 | Acenaphthene-d10 | 14.720 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | Decahydro-4,4,8,9,10-pentamethyl... | 208 | C15H28    | 080655-44-3 | 62   |
| 2    |    |   | 2(1H)-Naphthalenone, octahydro-1... | 194 | C13H22O   | 000775-54-2 | 25   |
| 3    |    |   | 2-Butanone, 4-(2,6,6-trimethyl-2... | 194 | C13H22O   | 039721-65-8 | 18   |
| 4    |    |   | 4-Pyrimidinamine, 2-methoxy-6-(t... | 193 | C6H6F3N3O | 054824-10-1 | 14   |
| 5    |    |   | [1,1'-Biphenyl]-4-acetonitrile      | 193 | C14H11N   | 031603-77-7 | 14   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
Data File : BG048872.D  
Acq On : 23 Apr 2021 20:53  
Operator : CG/JU  
Sample : M2107-19  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG041421.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

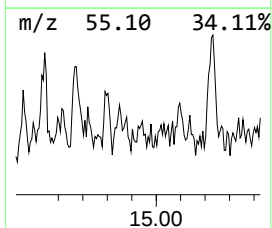
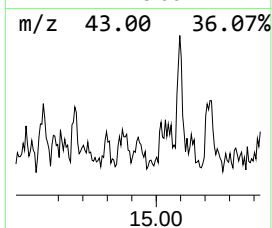
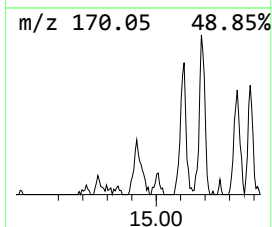
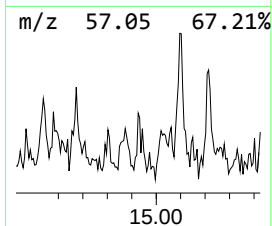
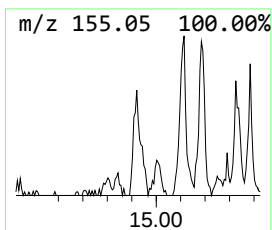
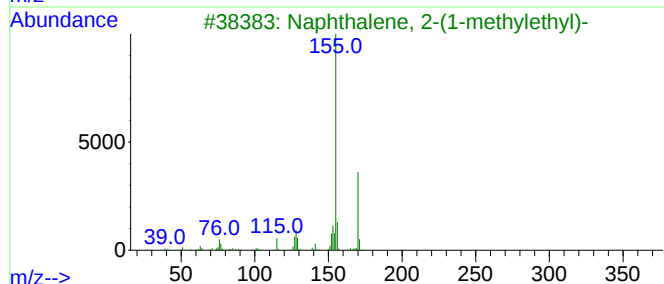
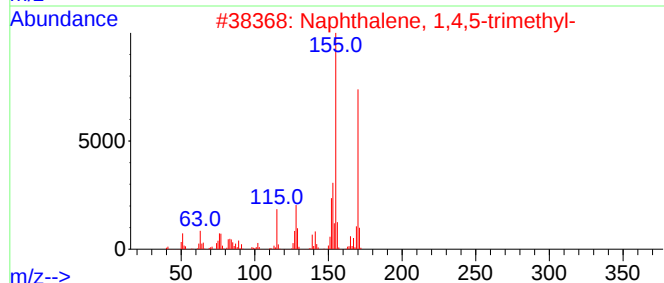
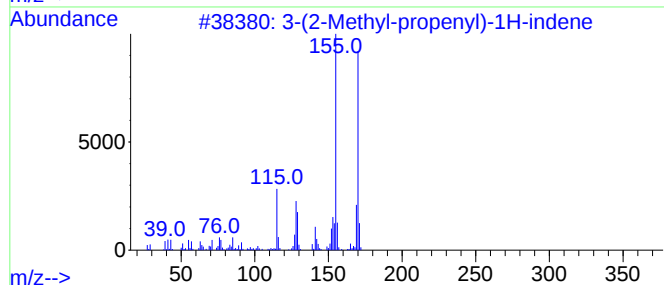
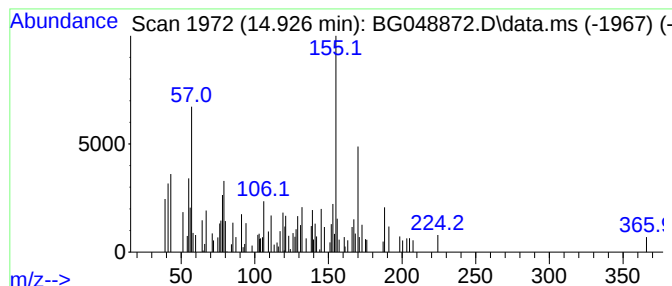
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 3-(2-Methyl-propenyl)-1H-in... Concentration Rank 20

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 14.926 | 5.22 ng | 71834 | Acenaphthene-d10 | 14.720 |

| Hit# | of                              | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|---------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 3-(2-Methyl-propenyl)-1H-indene | 170 | C13H14       | 1000187-78-5 | 70      |      |      |
| 2    | Naphthalene, 1,4,5-trimethyl-   | 170 | C13H14       | 002131-41-1  | 60      |      |      |
| 3    | Naphthalene, 2-(1-methylethyl)- | 170 | C13H14       | 002027-17-0  | 55      |      |      |
| 4    | 2-Naphthyl methyl ketone        | 170 | C12H10O      | 000093-08-3  | 43      |      |      |
| 5    | Ethanone, 1-(1-Naphthalenyl)-   | 170 | C12H10O      | 000941-98-0  | 43      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
Data File : BG048872.D  
Acq On : 23 Apr 2021 20:53  
Operator : CG/JU  
Sample : M2107-19  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG041421.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

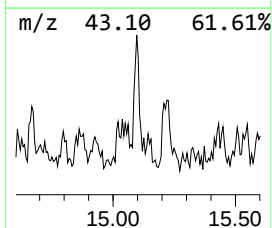
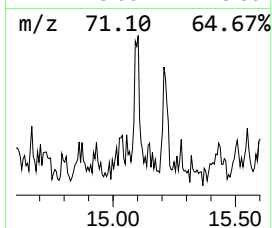
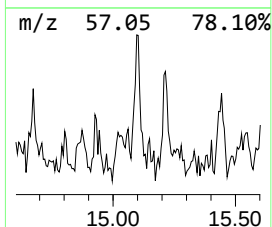
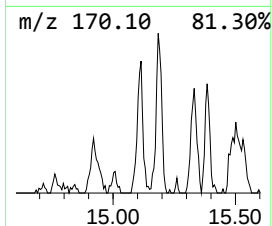
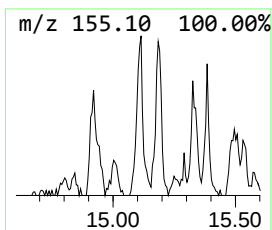
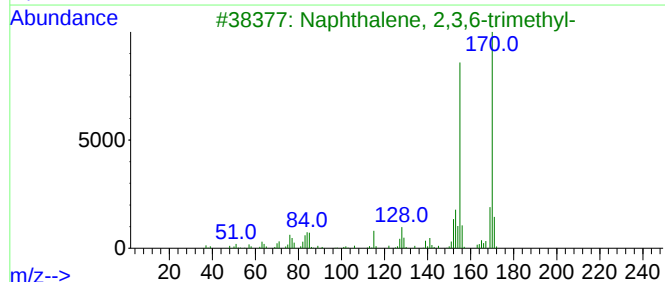
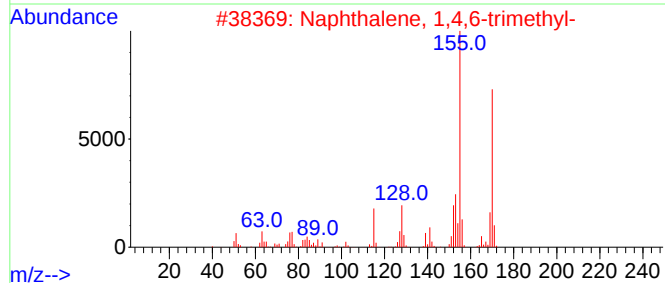
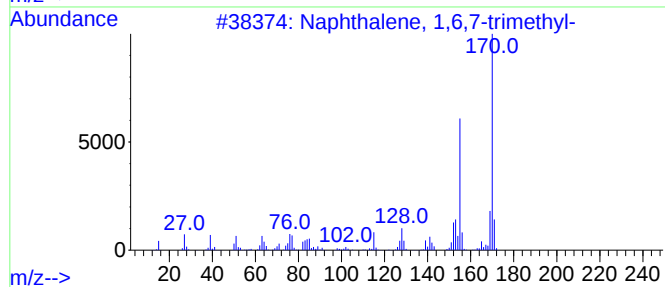
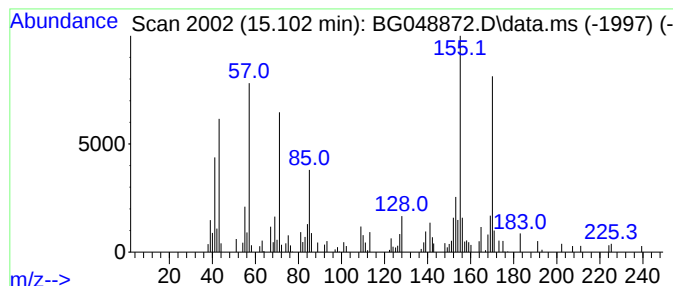
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Naphthalene, 1,6,7-trimethyl- Concentration Rank 18

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 15.102 | 5.41 ng | 74339 | Acenaphthene-d10 | 14.720 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 95   |
| 2    |    |   | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 94   |
| 3    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 93   |
| 4    |    |   | 4,6,8-Trimethylazulene        | 170 | C13H14  | 000941-81-1 | 76   |
| 5    |    |   | Naphthalene, 1,4,5-trimethyl- | 170 | C13H14  | 002131-41-1 | 76   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
Data File : BG048872.D  
Acq On : 23 Apr 2021 20:53  
Operator : CG/JU  
Sample : M2107-19  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

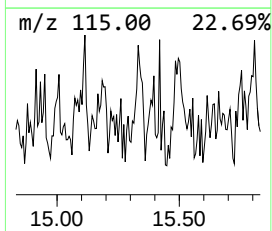
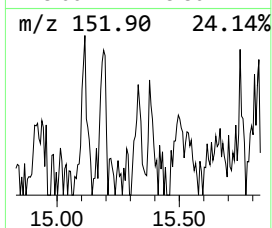
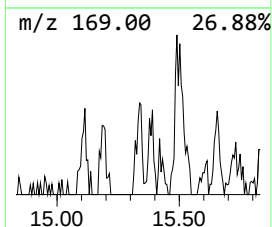
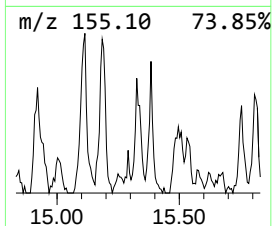
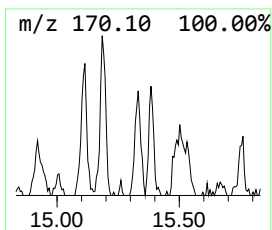
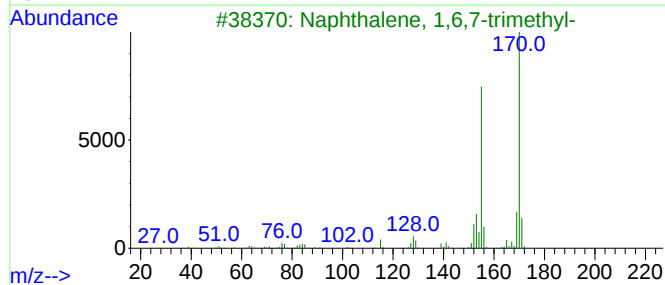
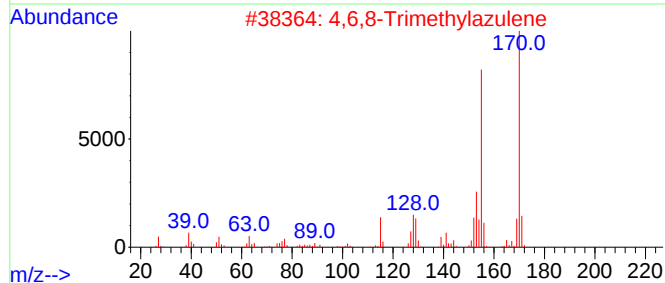
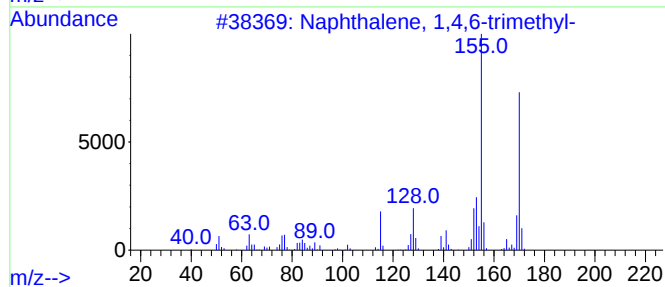
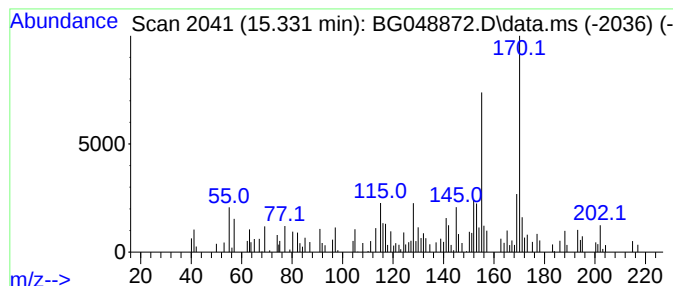
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 Naphthalene, 1,4,6-trimethyl- Concentration Rank 17

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 15.332 | 5.50 ng | 75629 | Acenaphthene-d10 | 14.720 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm | CAS#         | Qual |
|------|----|---|---------------------------------|-----|---------|--------------|------|
| 1    |    |   | Naphthalene, 1,4,6-trimethyl-   | 170 | C13H14  | 002131-42-2  | 90   |
| 2    |    |   | 4,6,8-Trimethylazulene          | 170 | C13H14  | 000941-81-1  | 81   |
| 3    |    |   | Naphthalene, 1,6,7-trimethyl-   | 170 | C13H14  | 002245-38-7  | 81   |
| 4    |    |   | Naphthalene, 2,3,6-trimethyl-   | 170 | C13H14  | 000829-26-5  | 81   |
| 5    |    |   | 3-(2-Methyl-propenyl)-1H-indene | 170 | C13H14  | 1000187-78-5 | 76   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
Data File : BG048872.D  
Acq On : 23 Apr 2021 20:53  
Operator : CG/JU  
Sample : M2107-19  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

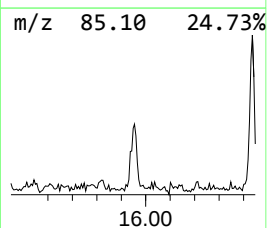
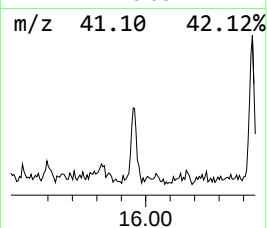
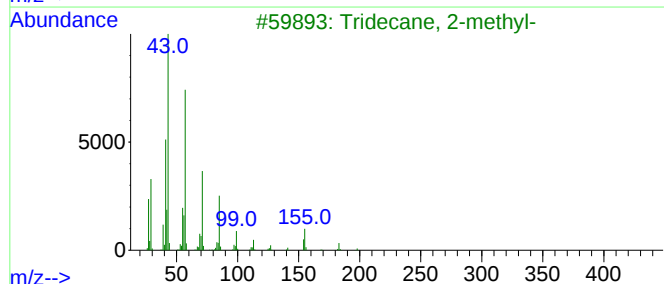
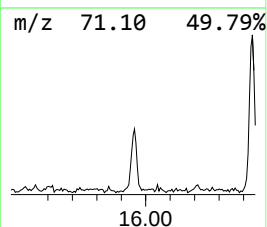
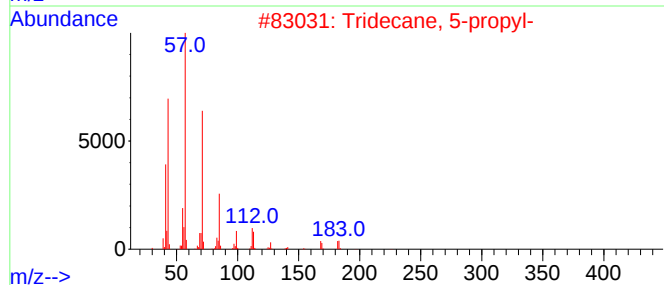
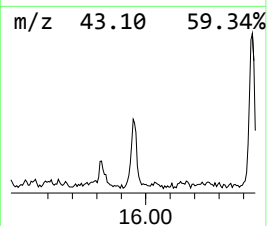
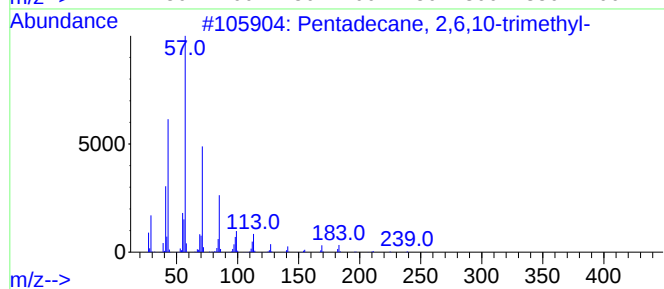
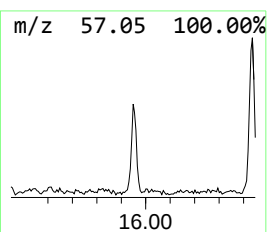
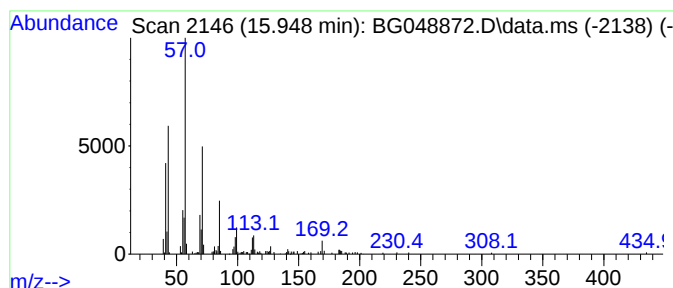
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 Pentadecane, 2,6,10-trimethyl- Concentration Rank 2

| R.T.      | EstConc                        | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|--------------------------------|--------|------------------|---------|-------------|------|
| 15.948    | 15.29 ng                       | 210281 | Acenaphthene-d10 |         | 14.720      |      |
| Hit# of 5 | Tentative ID                   |        | MW               | MolForm | CAS#        | Qual |
| 1         | Pentadecane, 2,6,10-trimethyl- |        | 254              | C18H38  | 003892-00-0 | 87   |
| 2         | Tridecane, 5-propyl-           |        | 226              | C16H34  | 055045-11-9 | 86   |
| 3         | Tridecane, 2-methyl-           |        | 198              | C14H30  | 001560-96-9 | 83   |
| 4         | Tridecane, 6-methyl-           |        | 198              | C14H30  | 013287-21-3 | 81   |
| 5         | Heptacosane                    |        | 380              | C27H56  | 000593-49-7 | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
Data File : BG048872.D  
Acq On : 23 Apr 2021 20:53  
Operator : CG/JU  
Sample : M2107-19  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

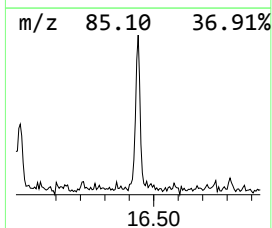
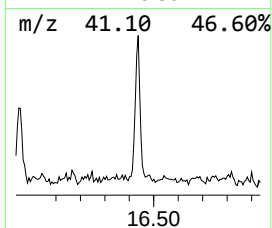
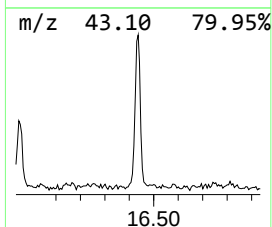
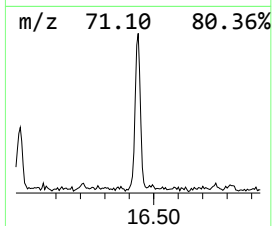
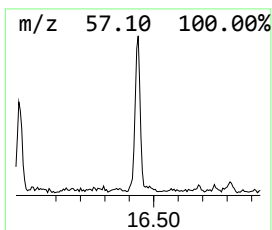
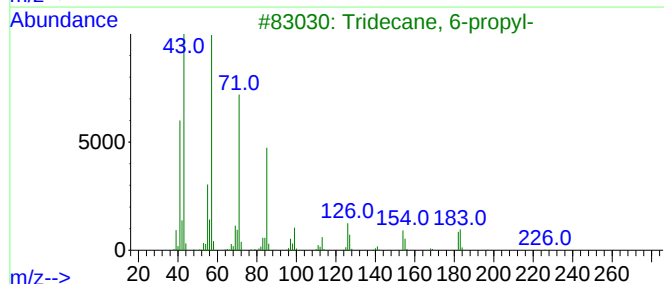
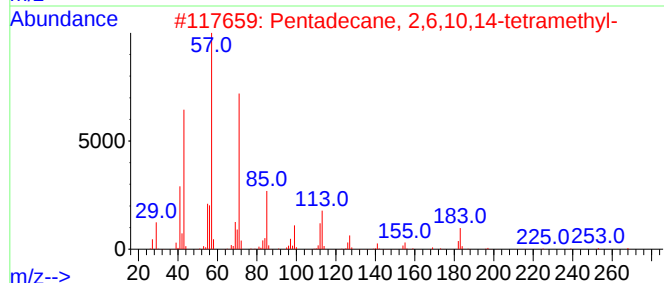
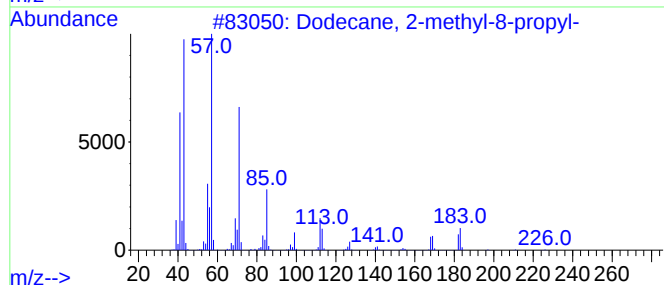
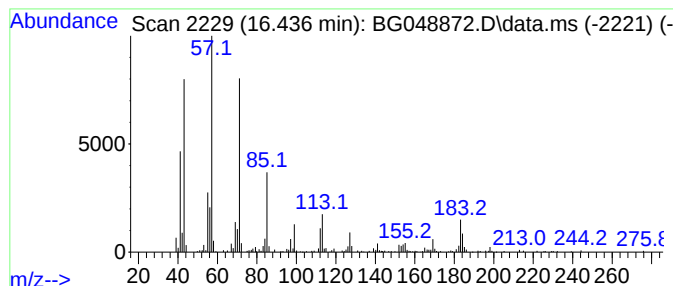
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 Dodecane, 2-methyl-8-propyl- Concentration Rank 1

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 16.436 | 24.83 ng | 389749 | Phenanthrene-d10 | 17.476 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Dodecane, 2-methyl-8-propyl-        | 226 | C16H34  | 055045-07-3 | 93   |
| 2    |    |   | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40  | 001921-70-6 | 91   |
| 3    |    |   | Tridecane, 6-propyl-                | 226 | C16H34  | 055045-10-8 | 90   |
| 4    |    |   | Hexadecane, 2,6,10,14-tetramethyl-  | 282 | C20H42  | 000638-36-8 | 90   |
| 5    |    |   | Hexadecane, 2,6,11,15-tetramethyl-  | 282 | C20H42  | 000504-44-9 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
Data File : BG048872.D  
Acq On : 23 Apr 2021 20:53  
Operator : CG/JU  
Sample : M2107-19  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG041421.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

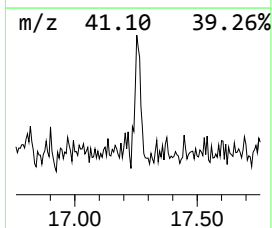
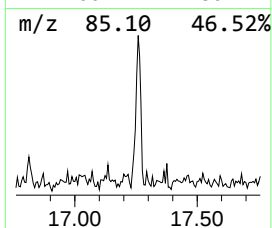
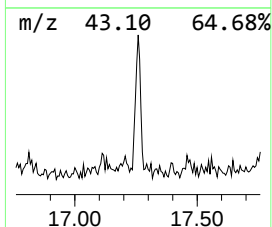
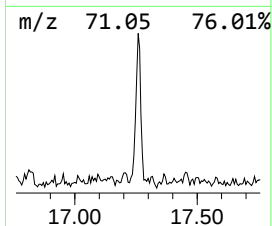
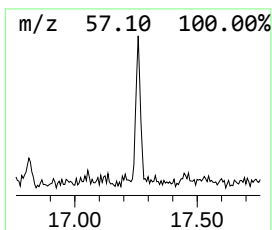
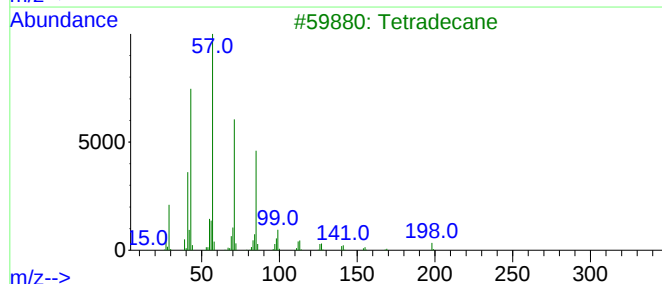
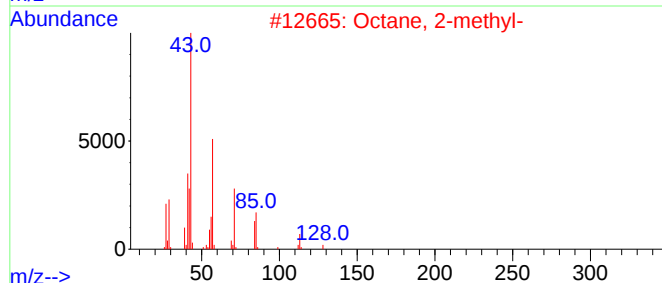
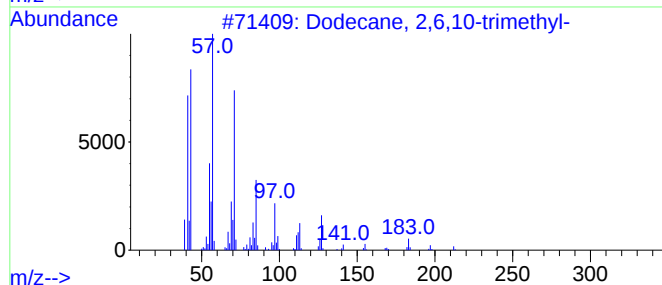
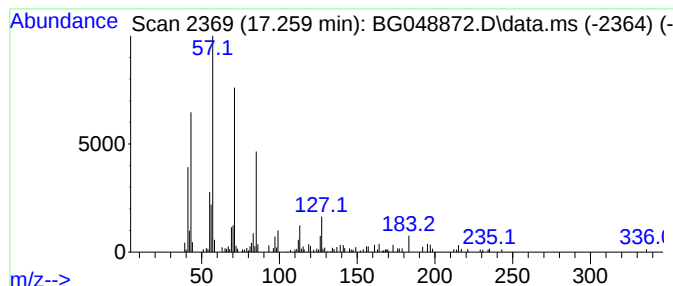
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Dodecane, 2,6,10-trimethyl- Concentration Rank 11

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 17.259 | 7.23 ng | 113479 | Phenanthrene-d10 | 17.476 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Dodecane, 2,6,10-trimethyl-         | 212 | C15H32    | 003891-98-3  | 90   |
| 2    |    |   | Octane, 2-methyl-                   | 128 | C9H20     | 003221-61-2  | 83   |
| 3    |    |   | Tetradecane                         | 198 | C14H30    | 000629-59-4  | 80   |
| 4    |    |   | Dodecane, 2,6,11-trimethyl-         | 212 | C15H32    | 031295-56-4  | 74   |
| 5    |    |   | Sulfurous acid, 2-ethylhexyl hex... | 278 | C14H30O3S | 1000309-20-2 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
Data File : BG048872.D  
Acq On : 23 Apr 2021 20:53  
Operator : CG/JU  
Sample : M2107-19  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

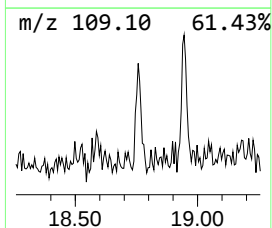
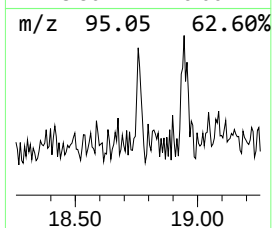
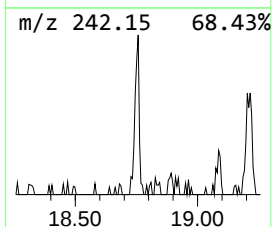
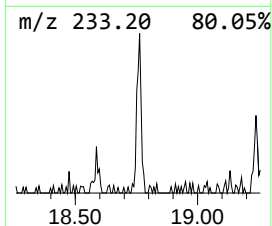
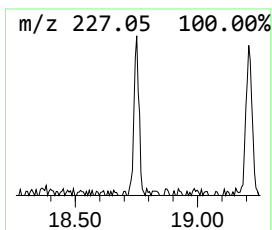
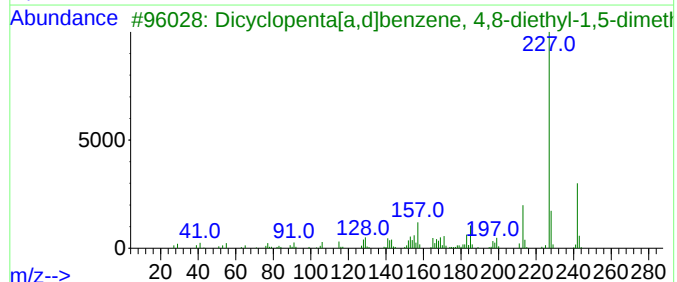
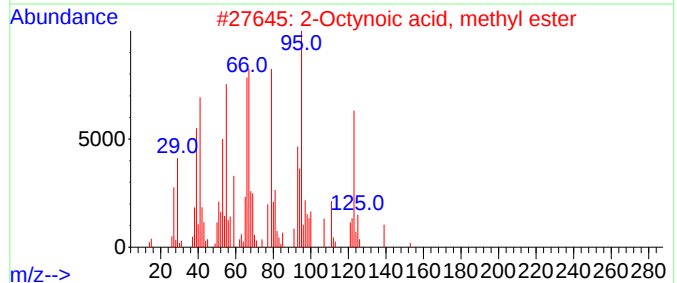
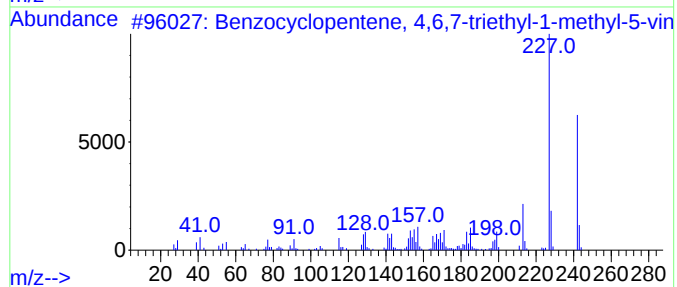
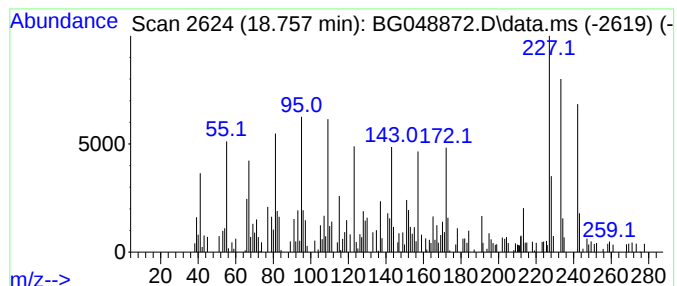
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 12 unknown18.757 Concentration Rank 8

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 18.757 | 8.30 ng | 130291 | Phenanthrene-d10 | 17.476 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Benzocyclopentene, 4,6,7-triethy... | 242 | C18H26     | 1000156-41-5 | 25   |
| 2    |    |   | 2-Octynoic acid, methyl ester       | 154 | C9H14O2    | 000111-12-6  | 14   |
| 3    |    |   | Dicyclopenta[a,d]benzene, 4,8-di... | 242 | C18H26     | 1000156-41-6 | 14   |
| 4    |    |   | Cis-2-methyl-4-n-butylthiane        | 172 | C10H20S    | 076097-67-1  | 11   |
| 5    |    |   | 4,5-Dimethyl-4-nitro-2-phenyl-2,... | 233 | C11H11N3O3 | 1000300-94-0 | 11   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
Data File : BG048872.D  
Acq On : 23 Apr 2021 20:53  
Operator : CG/JU  
Sample : M2107-19  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

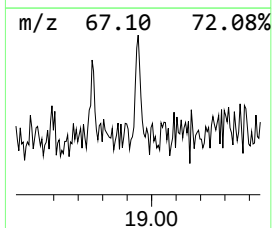
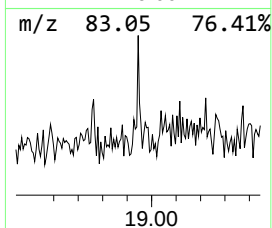
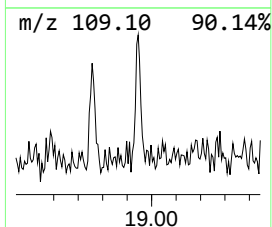
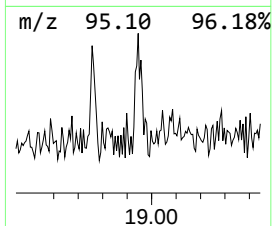
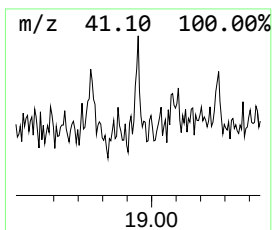
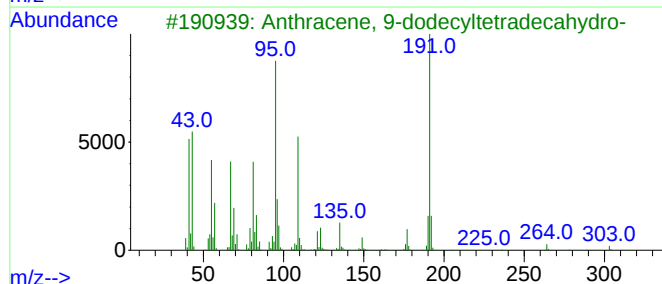
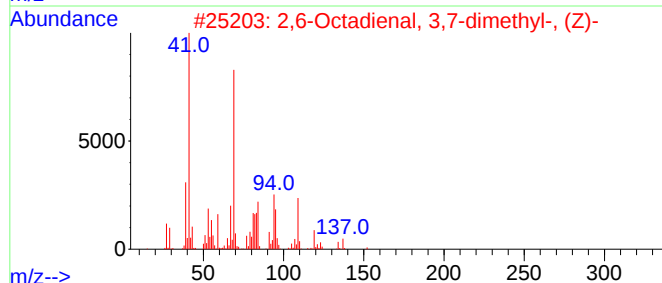
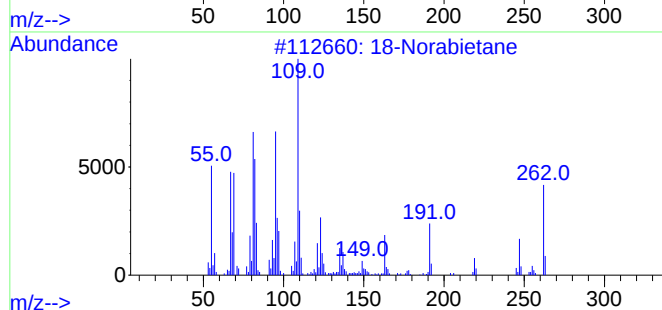
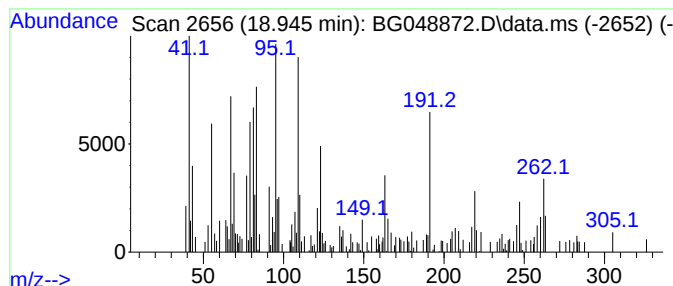
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 13 18-Norabietane Concentration Rank 13

| R.T.      | EstConc                             | Area  | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|-------|------------------|--------------|------|
| 18.945    | 6.17 ng                             | 96909 | Phenanthrene-d10 | 17.476       |      |
| Hit# of 5 | Tentative ID                        | MW    | MolForm          | CAS#         | Qual |
| 1         | 18-Norabietane                      | 262   | C19H34           | 1000293-16-6 | 50   |
| 2         | 2,6-Octadienal, 3,7-dimethyl-, (Z)- | 152   | C10H16O          | 000106-26-3  | 47   |
| 3         | Anthracene, 9-dodecyltetradeca...   | 360   | C26H48           | 055401-75-7  | 35   |
| 4         | 8-Thiabicyclo[4.3.0]nonane S-oxide  | 158   | C8H14OS          | 1000215-10-0 | 27   |
| 5         | Divinylbis(cyclopropyl)silane       | 164   | C10H16Si         | 1000109-95-2 | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
Data File : BG048872.D  
Acq On : 23 Apr 2021 20:53  
Operator : CG/JU  
Sample : M2107-19  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

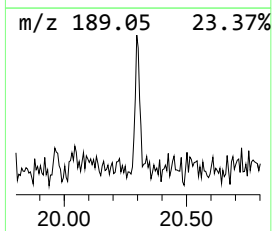
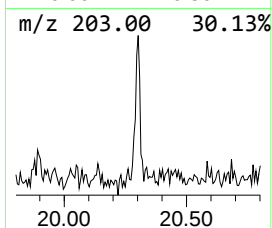
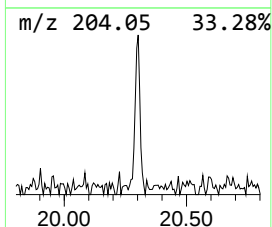
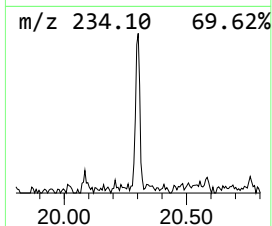
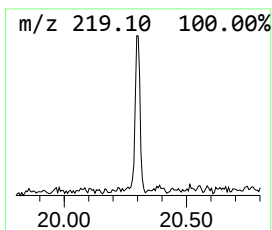
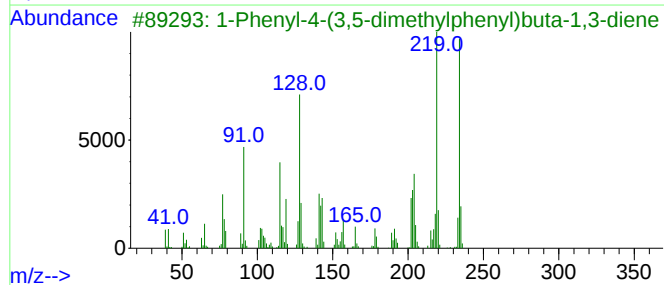
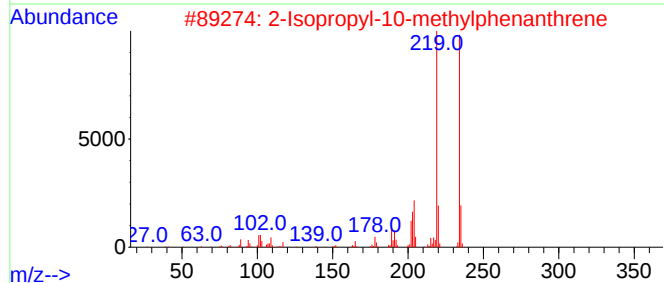
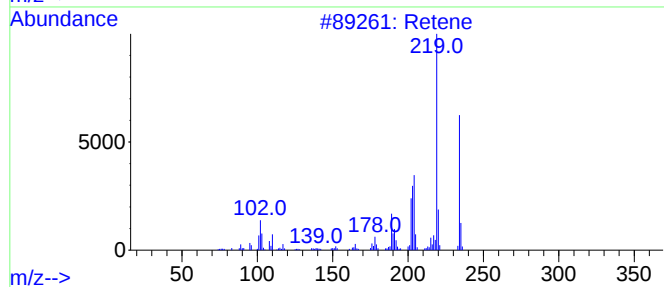
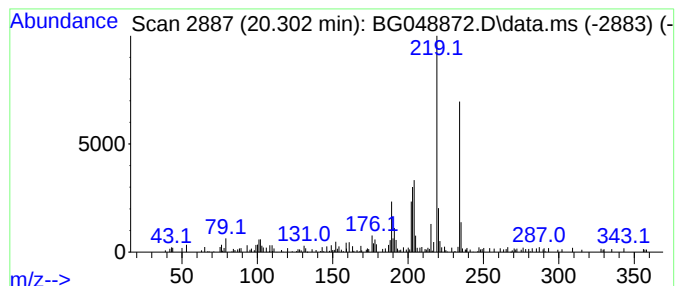
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 14 Retene Concentration Rank 15

| R.T.      | EstConc                             | Area  | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|-------|------------------|--------------|------|
| 20.302    | 5.91 ng                             | 95653 | Chrysene-d12     | 21.771       |      |
| Hit# of 5 | Tentative ID                        | MW    | MolForm          | CAS#         | Qual |
| 1         | Retene                              | 234   | C18H18           | 000483-65-8  | 95   |
| 2         | 2-Isopropyl-10-methylphenanthrene   | 234   | C18H18           | 066552-97-4  | 93   |
| 3         | 1-Phenyl-4-(3,5-dimethylphenyl)b... | 234   | C18H18           | 1000202-15-0 | 83   |
| 4         | Phenanthrene, 3,4,5,6-tetramethyl-  | 234   | C18H18           | 007343-06-8  | 64   |
| 5         | Anthracene, 2-(1,1-dimethylethyl)-  | 234   | C18H18           | 018801-00-8  | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
Data File : BG048872.D  
Acq On : 23 Apr 2021 20:53  
Operator : CG/JU  
Sample : M2107-19  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

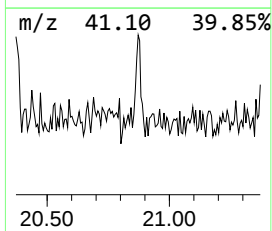
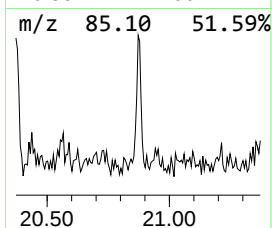
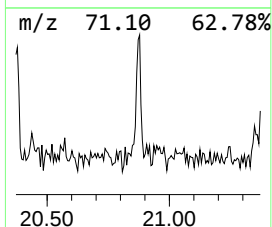
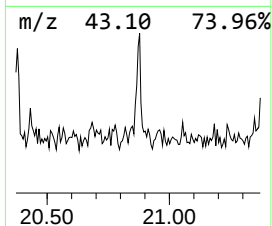
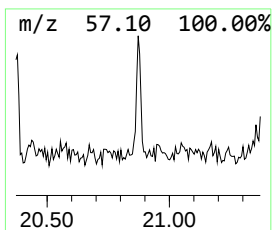
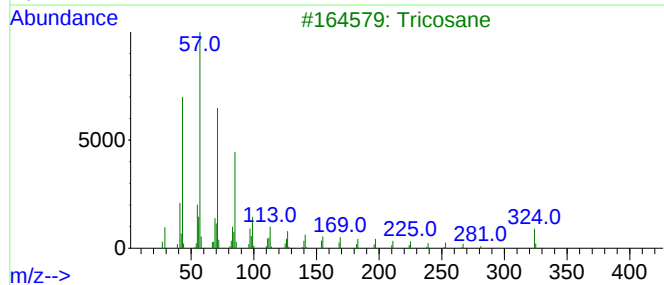
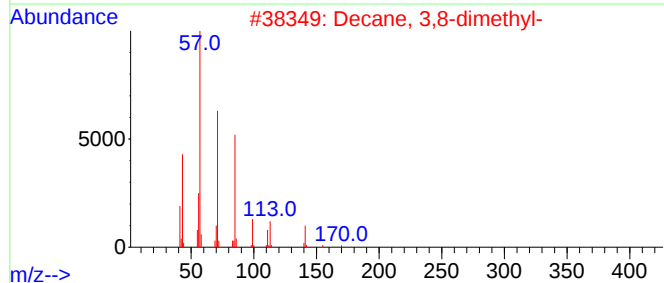
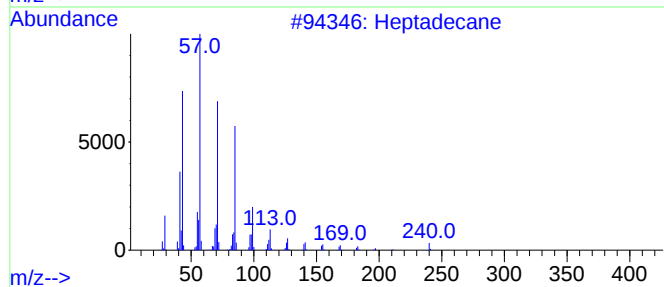
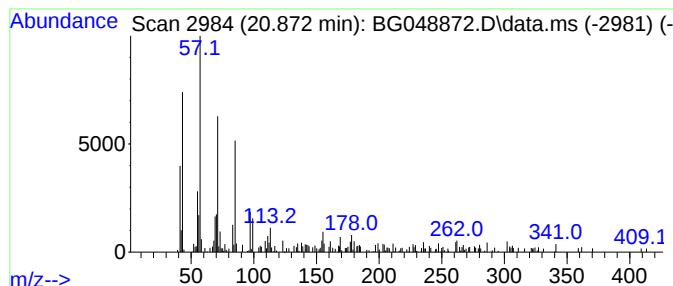
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 15 Heptadecane Concentration Rank 19

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 20.872 | 5.26 ng | 85034 | Chrysene-d12     | 21.771 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane           | 240 | C17H36  | 000629-78-7 | 86   |
| 2    |    |   | Decane, 3,8-dimethyl- | 170 | C12H26  | 017312-55-9 | 86   |
| 3    |    |   | Tricosane             | 324 | C23H48  | 000638-67-5 | 62   |
| 4    |    |   | Undecane              | 156 | C11H24  | 001120-21-4 | 62   |
| 5    |    |   | Tridecane             | 184 | C13H28  | 000629-50-5 | 62   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
Data File : BG048872.D  
Acq On : 23 Apr 2021 20:53  
Operator : CG/JU  
Sample : M2107-19  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG041421.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

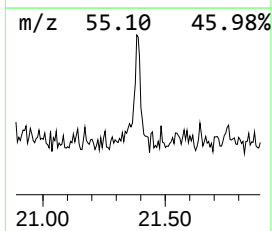
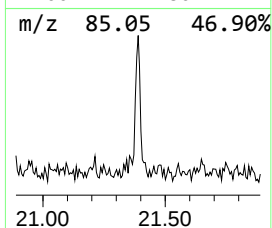
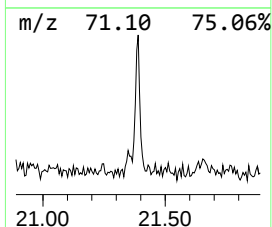
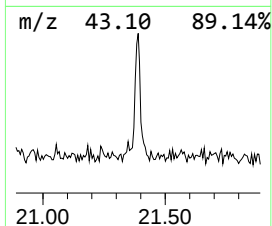
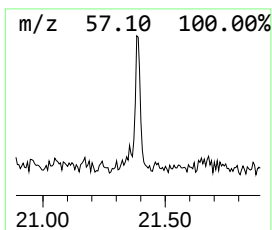
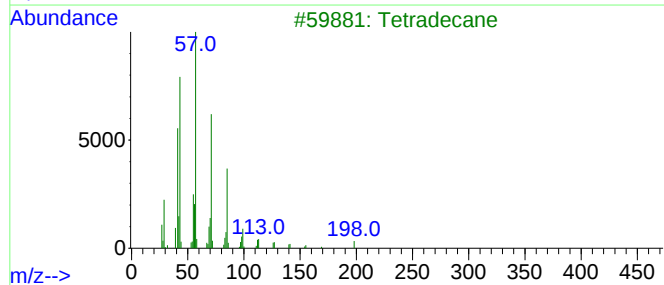
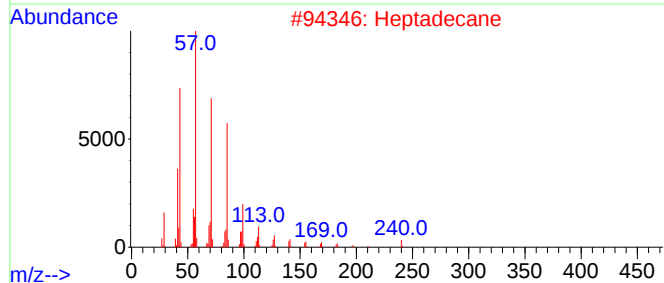
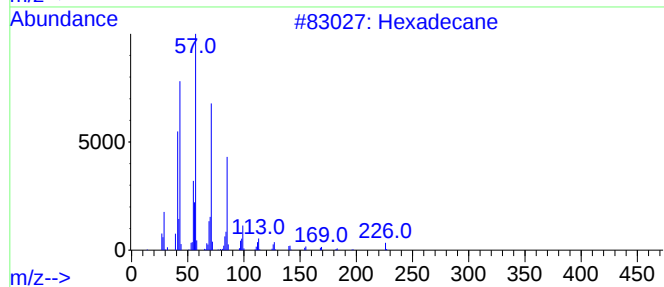
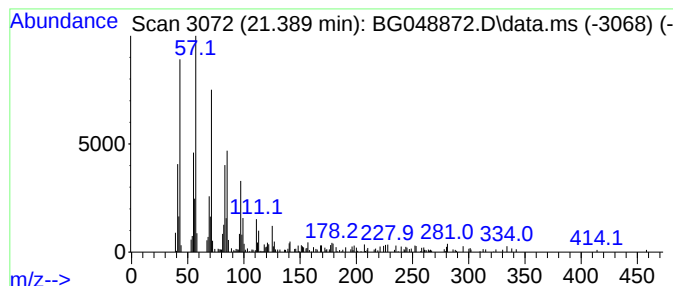
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Hexadecane Concentration Rank 4

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 21.389 | 11.38 ng | 184124 | Chrysene-d12     | 21.771 |

| Hit# | of 5 | Tentative ID      | MW  | MolForm | CAS#         | Qual |
|------|------|-------------------|-----|---------|--------------|------|
| 1    |      | Hexadecane        | 226 | C16H34  | 000544-76-3  | 95   |
| 2    |      | Heptadecane       | 240 | C17H36  | 000629-78-7  | 95   |
| 3    |      | Tetradecane       | 198 | C14H30  | 000629-59-4  | 90   |
| 4    |      | E-15-Heptadecenal | 252 | C17H32O | 1000130-97-9 | 86   |
| 5    |      | 4-Methyldocosane  | 324 | C23H48  | 025117-30-0  | 81   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
Data File : BG048872.D  
Acq On : 23 Apr 2021 20:53  
Operator : CG/JU  
Sample : M2107-19  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG041421.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

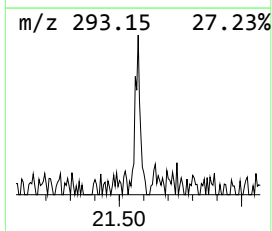
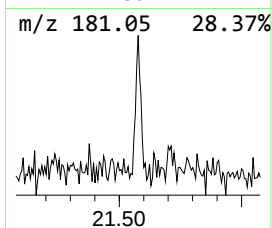
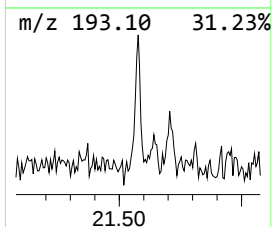
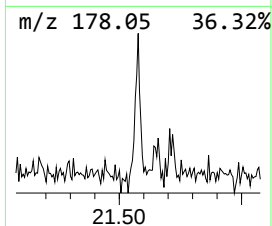
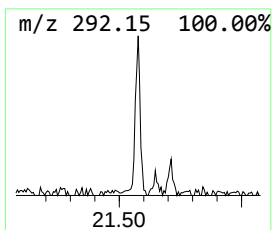
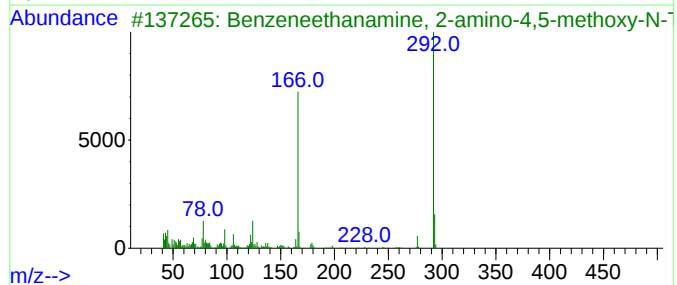
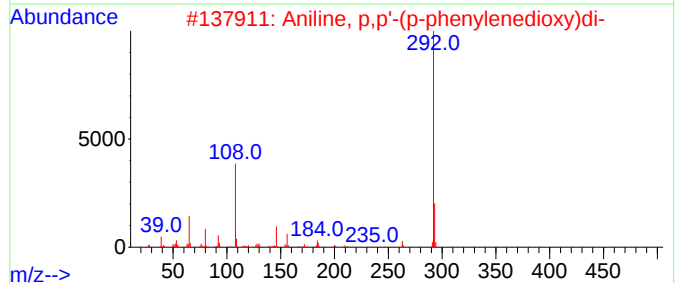
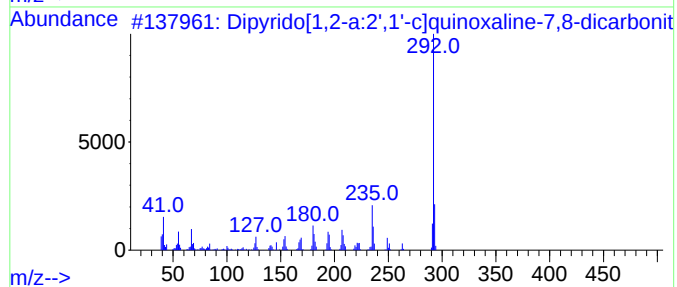
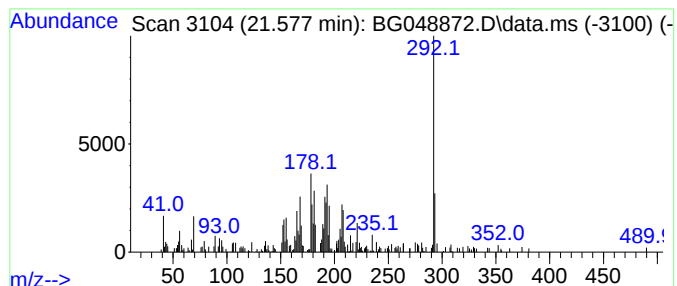
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 unknown21.577 Concentration Rank 12

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 21.577 | 6.83 ng | 110567 | Chrysene-d12     | 21.771 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|--------------|--------------|------|
| 1    |    |   | Dipyrido[1,2-a:2',1'-c]quinoxali... | 292 | C18H20N4     | 1000303-96-3 | 27   |
| 2    |    |   | Aniline, p,p'-(p-phenylenedioxy)di- | 292 | C18H16N2O2   | 003491-12-1  | 22   |
| 3    |    |   | Benzeneethanamine, 2-amino-4,5-m... | 292 | C12H15F3N2O3 | 1000125-99-1 | 22   |
| 4    |    |   | 4-(p-(Dimethylamino)benzylidene)... | 292 | C18H16N2O2   | 065156-10-7  | 22   |
| 5    |    |   | 4,6-Di(4-methoxyphenyl)pyrimidine   | 292 | C18H16N2O2   | 183670-49-7  | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
Data File : BG048872.D  
Acq On : 23 Apr 2021 20:53  
Operator : CG/JU  
Sample : M2107-19  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG041421.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

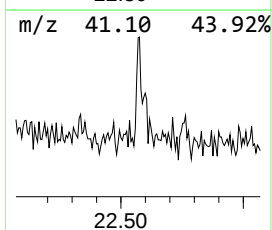
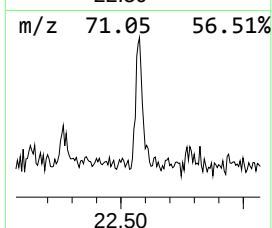
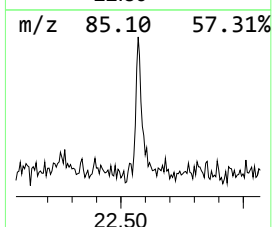
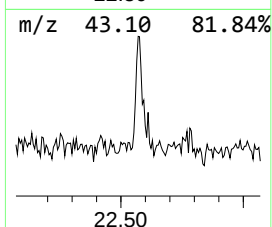
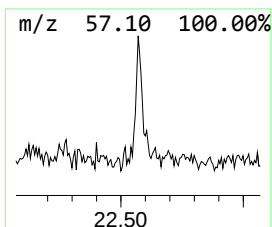
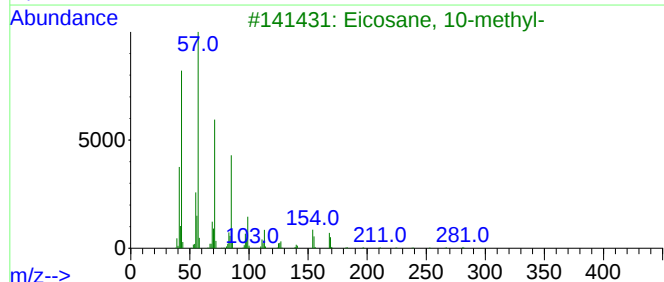
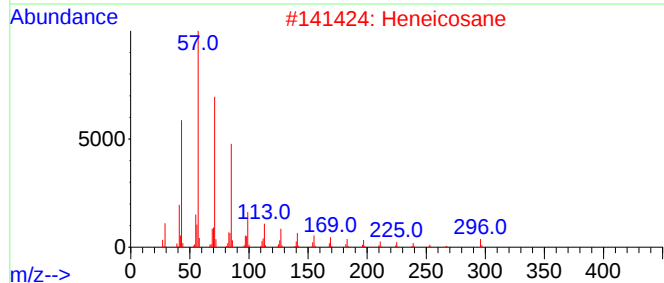
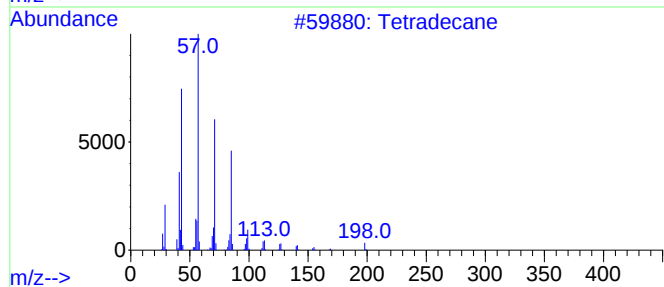
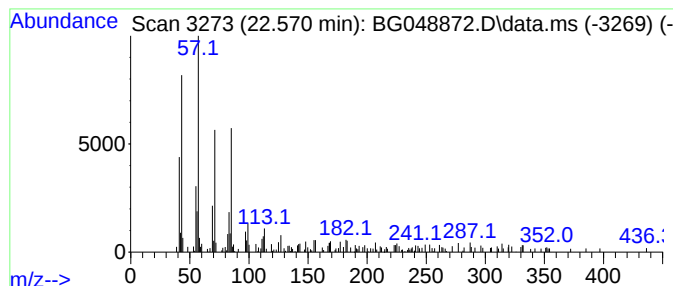
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Tetradecane Concentration Rank 10

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 22.570 | 7.85 ng | 127098 | Chrysene-d12     | 21.771 |

| Hit# | of 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|------|----------------------|-----|---------|-------------|------|
| 1    |      | Tetradecane          | 198 | C14H30  | 000629-59-4 | 86   |
| 2    |      | Heneicosane          | 296 | C21H44  | 000629-94-7 | 70   |
| 3    |      | Eicosane, 10-methyl- | 296 | C21H44  | 054833-23-7 | 58   |
| 4    |      | Tetracosane          | 338 | C24H50  | 000646-31-1 | 55   |
| 5    |      | Tricosane, 2-methyl- | 338 | C24H50  | 001928-30-9 | 55   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
Data File : BG048872.D  
Acq On : 23 Apr 2021 20:53  
Operator : CG/JU  
Sample : M2107-19  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG041421.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

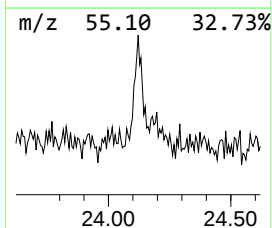
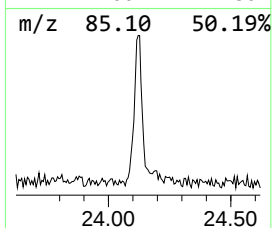
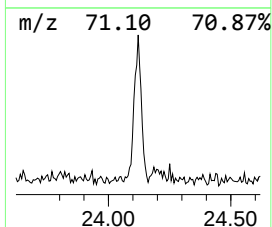
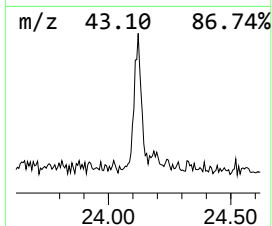
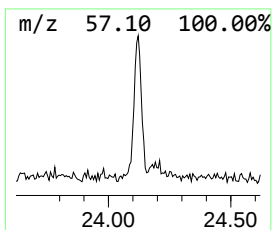
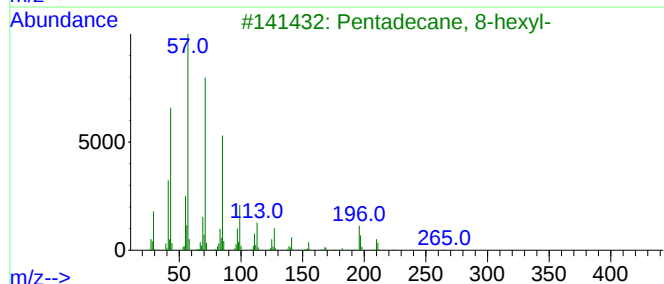
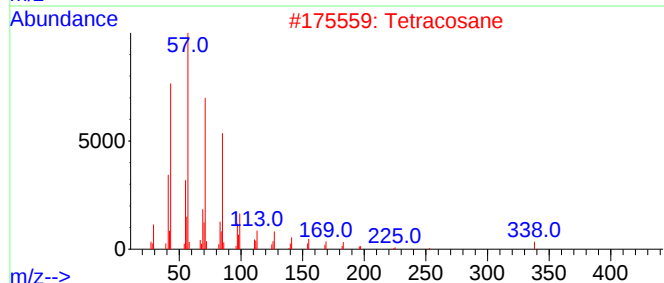
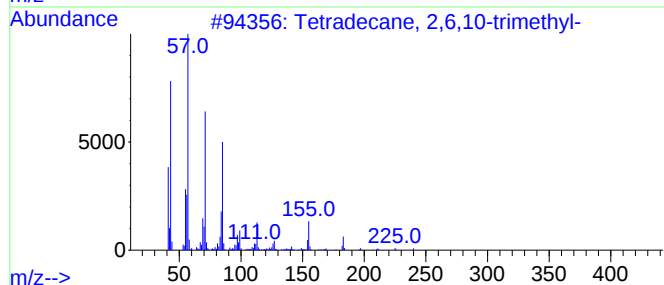
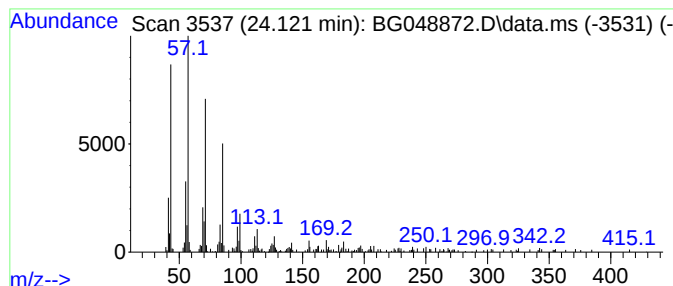
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Tetradecane, 2,6,10-trimethyl- Concentration Rank 3

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 24.121 | 11.89 ng | 225329 | Perylene-d12     | 25.102 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#         | Qual |
|------|----|---|--------------------------------|-----|---------|--------------|------|
| 1    |    |   | Tetradecane, 2,6,10-trimethyl- | 240 | C17H36  | 014905-56-7  | 93   |
| 2    |    |   | Tetracosane                    | 338 | C24H50  | 000646-31-1  | 90   |
| 3    |    |   | Pentadecane, 8-hexyl-          | 296 | C21H44  | 013475-75-7  | 87   |
| 4    |    |   | 2-methyloctacosane             | 408 | C29H60  | 1000376-72-8 | 87   |
| 5    |    |   | Nonadecane                     | 268 | C19H40  | 000629-92-5  | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
Data File : BG048872.D  
Acq On : 23 Apr 2021 20:53  
Operator : CG/JU  
Sample : M2107-19  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SB-34-9.5-10.0

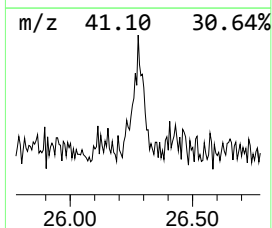
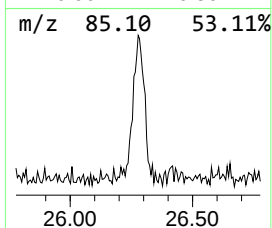
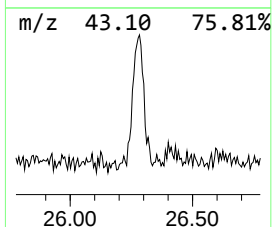
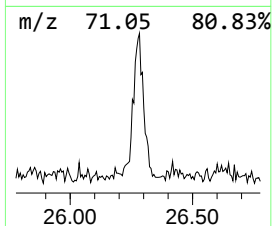
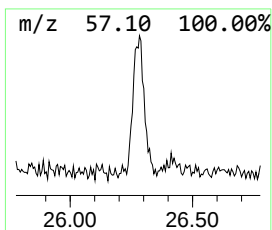
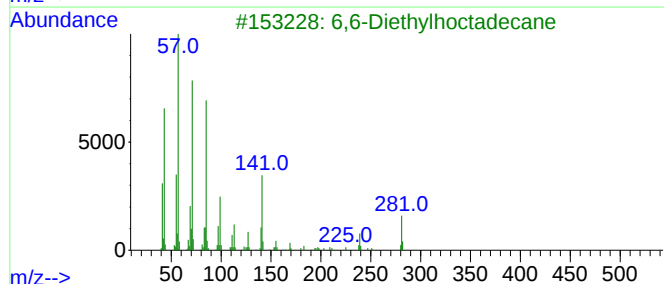
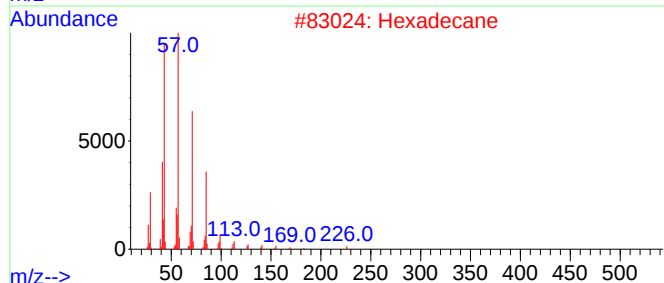
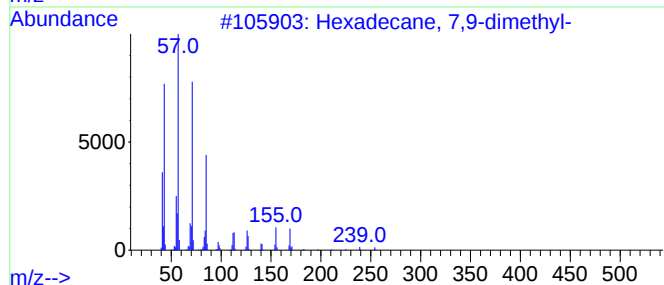
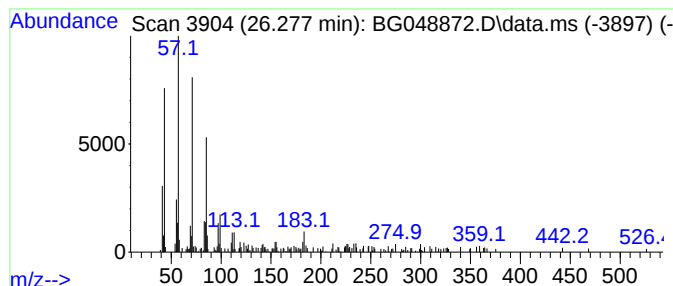
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 20 Hexadecane, 7,9-dimethyl- Concentration Rank 5

| R.T.      | EstConc                   | Area   | Relative to ISTD |         | R.T.         |      |
|-----------|---------------------------|--------|------------------|---------|--------------|------|
| 26.278    | 11.17 ng                  | 211829 | Perylene-d12     |         | 25.102       |      |
| Hit# of 5 | Tentative ID              |        | MW               | MolForm | CAS#         | Qual |
| 1         | Hexadecane, 7,9-dimethyl- |        | 254              | C18H38  | 021164-95-4  | 81   |
| 2         | Hexadecane                |        | 226              | C16H34  | 000544-76-3  | 74   |
| 3         | 6,6-Diethylhooctadecane   |        | 310              | C22H46  | 1000360-41-8 | 72   |
| 4         | Hexacosane                |        | 366              | C26H54  | 000630-01-3  | 72   |
| 5         | Dodecane, 1-iodo-         |        | 296              | C12H25I | 004292-19-7  | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042321\  
 Data File : BG048872.D  
 Acq On : 23 Apr 2021 20:53  
 Operator : CG/JU  
 Sample : M2107-19  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SB-34-9.5-10.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG041421.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 |
| o-Cymene           | 8.187  | 7.9     | ng    | 54795    | 1                     | 8.058  | 138416 | 20.0 |
| Octane, 3,6-dim... | 11.894 | 8.8     | ng    | 90736    | 2                     | 10.884 | 205533 | 20.0 |
| Heptadecane, 2,... | 13.234 | 11.0    | ng    | 151425   | 3                     | 14.720 | 275071 | 20.0 |
| 2,4,7,9-Tetrame... | 13.592 | 5.8     | ng    | 79643    | 3                     | 14.720 | 275071 | 20.0 |
| Decahydro-4,4,8... | 14.538 | 6.1     | ng    | 84097    | 3                     | 14.720 | 275071 | 20.0 |
| 3-(2-Methyl-pro... | 14.926 | 5.2     | ng    | 71834    | 3                     | 14.720 | 275071 | 20.0 |
| Naphthalene, 1,... | 15.102 | 5.4     | ng    | 74339    | 3                     | 14.720 | 275071 | 20.0 |
| Naphthalene, 1,... | 15.332 | 5.5     | ng    | 75629    | 3                     | 14.720 | 275071 | 20.0 |
| Pentadecane, 2,... | 15.948 | 15.3    | ng    | 210281   | 3                     | 14.720 | 275071 | 20.0 |
| Dodecane, 2-met... | 16.436 | 24.8    | ng    | 389749   | 4                     | 17.476 | 313984 | 20.0 |
| Dodecane, 2,6,1... | 17.259 | 7.2     | ng    | 113479   | 4                     | 17.476 | 313984 | 20.0 |
| unknown18.757      | 18.757 | 8.3     | ng    | 130291   | 4                     | 17.476 | 313984 | 20.0 |
| 18-Norabietane     | 18.945 | 6.2     | ng    | 96909    | 4                     | 17.476 | 313984 | 20.0 |
| Retene             | 20.302 | 5.9     | ng    | 95653    | 5                     | 21.771 | 323619 | 20.0 |
| Heptadecane        | 20.872 | 5.3     | ng    | 85034    | 5                     | 21.771 | 323619 | 20.0 |
| Hexadecane         | 21.389 | 11.4    | ng    | 184124   | 5                     | 21.771 | 323619 | 20.0 |
| unknown21.577      | 21.577 | 6.8     | ng    | 110567   | 5                     | 21.771 | 323619 | 20.0 |
| Tetradecane        | 22.570 | 7.8     | ng    | 127098   | 5                     | 21.771 | 323619 | 20.0 |
| Tetradecane, 2,... | 24.121 | 11.9    | ng    | 225329   | 6                     | 25.102 | 379127 | 20.0 |
| Hexadecane, 7,9... | 26.278 | 11.2    | ng    | 211829   | 6                     | 25.102 | 379127 | 20.0 |