

Data Path : Z:\HPCHEM1\BNA G\DATA\BG030116\  
 Data File : BG021197.D  
 Acq On : 1 Mar 2016 23:43  
 Operator : UM/SJ  
 Sample : H1565-08  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 CEB-5(2.5-5)20160223

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-BG022916.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 3.061       | 32            | 38          | 57           | rVV      | 1934928        | 2750207       | 100.00%         | 16.554%       |
| 2         | 3.202       | 57            | 62          | 71           | rVB2     | 18991          | 37193         | 1.35%           | 0.224%        |
| 3         | 5.241       | 401           | 409         | 421          | rBV      | 406997         | 624495        | 22.71%          | 3.759%        |
| 4         | 5.705       | 482           | 488         | 494          | rVB      | 67357          | 109593        | 3.98%           | 0.660%        |
| 5         | 5.775       | 494           | 500         | 509          | rBV      | 54090          | 118494        | 4.31%           | 0.713%        |
| 6         | 6.146       | 555           | 563         | 570          | rVB2     | 52782          | 90675         | 3.30%           | 0.546%        |
| 7         | 6.627       | 635           | 645         | 652          | rBV2     | 18474          | 34892         | 1.27%           | 0.210%        |
| 8         | 7.127       | 718           | 730         | 735          | rBV3     | 55572          | 103603        | 3.77%           | 0.624%        |
| 9         | 7.227       | 742           | 747         | 752          | rBV      | 106310         | 167199        | 6.08%           | 1.006%        |
| 10        | 7.297       | 752           | 759         | 765          | rVV2     | 300625         | 536154        | 19.50%          | 3.227%        |
| 11        | 7.362       | 765           | 770         | 779          | rVB      | 74069          | 122449        | 4.45%           | 0.737%        |
| 12        | 7.532       | 793           | 799         | 806          | rVB      | 44993          | 71586         | 2.60%           | 0.431%        |
| 13        | 7.679       | 816           | 824         | 832          | rVB      | 152443         | 270460        | 9.83%           | 1.628%        |
| 14        | 7.791       | 832           | 843         | 854          | rVB      | 235493         | 401526        | 14.60%          | 2.417%        |
| 15        | 8.120       | 891           | 899         | 901          | rBV      | 35558          | 62430         | 2.27%           | 0.376%        |
| 16        | 8.149       | 901           | 904         | 909          | rVB      | 43926          | 66154         | 2.41%           | 0.398%        |
| 17        | 8.261       | 917           | 923         | 930          | rVB2     | 75000          | 124667        | 4.53%           | 0.750%        |
| 18        | 8.372       | 937           | 942         | 949          | rBV      | 33115          | 57787         | 2.10%           | 0.348%        |
| 19        | 8.472       | 952           | 959         | 964          | rVV      | 241224         | 420012        | 15.27%          | 2.528%        |
| 20        | 8.525       | 964           | 968         | 977          | rVB      | 24252          | 45057         | 1.64%           | 0.271%        |
| 21        | 8.631       | 979           | 986         | 990          | rBV3     | 23405          | 46396         | 1.69%           | 0.279%        |
| 22        | 8.684       | 990           | 995         | 1006         | rVB2     | 69312          | 150641        | 5.48%           | 0.907%        |
| 23        | 8.784       | 1006          | 1012        | 1025         | rBV4     | 88785          | 202641        | 7.37%           | 1.220%        |
| 24        | 9.095       | 1060          | 1065        | 1070         | rVV      | 33056          | 49388         | 1.80%           | 0.297%        |
| 25        | 9.148       | 1070          | 1074        | 1081         | rVB      | 32739          | 52009         | 1.89%           | 0.313%        |
| 26        | 9.254       | 1086          | 1092        | 1096         | rBV2     | 76520          | 125131        | 4.55%           | 0.753%        |
| 27        | 9.312       | 1096          | 1102        | 1109         | rBV      | 227664         | 437875        | 15.92%          | 2.636%        |
| 28        | 9.589       | 1142          | 1149        | 1152         | rBV2     | 16864          | 33012         | 1.20%           | 0.199%        |
| 29        | 9.771       | 1175          | 1180        | 1185         | rBV      | 40083          | 66822         | 2.43%           | 0.402%        |
| 30        | 9.835       | 1185          | 1191        | 1199         | rVB      | 61656          | 99893         | 3.63%           | 0.601%        |
| 31        | 10.147      | 1238          | 1244        | 1248         | rBV2     | 17947          | 28889         | 1.05%           | 0.174%        |
| 32        | 10.206      | 1248          | 1254        | 1262         | rVB6     | 35186          | 77834         | 2.83%           | 0.469%        |
| 33        | 10.294      | 1262          | 1269        | 1273         | rBV2     | 16939          | 37298         | 1.36%           | 0.225%        |
| 34        | 10.352      | 1273          | 1279        | 1285         | rVV2     | 69986          | 136708        | 4.97%           | 0.823%        |

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 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 CEB-5(2.5-5)20160223

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-BG022916.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 35 | 10.411 | 1285 | 1289 | 1294 | rVB2 | 22993  | 35449  | 1.29%  | 0.213% |
| 36 | 10.523 | 1301 | 1308 | 1315 | rVB5 | 13541  | 32855  | 1.19%  | 0.198% |
| 37 | 10.911 | 1369 | 1374 | 1376 | rBV  | 29461  | 49437  | 1.80%  | 0.298% |
| 38 | 10.958 | 1377 | 1382 | 1386 | rVV  | 78402  | 160629 | 5.84%  | 0.967% |
| 39 | 11.005 | 1386 | 1390 | 1396 | rVB  | 77009  | 146620 | 5.33%  | 0.883% |
| 40 | 11.862 | 1529 | 1536 | 1541 | rBV2 | 23068  | 45762  | 1.66%  | 0.275% |
| 41 | 12.344 | 1610 | 1618 | 1624 | rBV3 | 31268  | 57414  | 2.09%  | 0.346% |
| 42 | 12.597 | 1656 | 1661 | 1669 | rVB  | 65490  | 111774 | 4.06%  | 0.673% |
| 43 | 12.814 | 1691 | 1698 | 1708 | rVB2 | 44910  | 77767  | 2.83%  | 0.468% |
| 44 | 13.384 | 1788 | 1795 | 1801 | rVB  | 498114 | 749710 | 27.26% | 4.513% |
| 45 | 13.566 | 1816 | 1826 | 1834 | rBV2 | 18340  | 33157  | 1.21%  | 0.200% |
| 46 | 14.077 | 1904 | 1913 | 1918 | rBV  | 24247  | 43756  | 1.59%  | 0.263% |
| 47 | 14.195 | 1928 | 1933 | 1941 | rVB  | 73694  | 112781 | 4.10%  | 0.679% |
| 48 | 14.753 | 2016 | 2028 | 2034 | rVV2 | 114103 | 193229 | 7.03%  | 1.163% |
| 49 | 14.818 | 2034 | 2039 | 2047 | rVB  | 27090  | 46409  | 1.69%  | 0.279% |
| 50 | 15.147 | 2089 | 2095 | 2100 | rBV2 | 23604  | 39046  | 1.42%  | 0.235% |
| 51 | 15.576 | 2164 | 2168 | 2173 | rVB  | 28362  | 34911  | 1.27%  | 0.210% |
| 52 | 15.799 | 2199 | 2206 | 2214 | rBV  | 33322  | 57840  | 2.10%  | 0.348% |
| 53 | 16.234 | 2275 | 2280 | 2289 | rVB  | 66626  | 101308 | 3.68%  | 0.610% |
| 54 | 16.433 | 2310 | 2314 | 2317 | rBV2 | 25035  | 35198  | 1.28%  | 0.212% |
| 55 | 16.786 | 2364 | 2374 | 2379 | rBV7 | 13499  | 35228  | 1.28%  | 0.212% |
| 56 | 17.227 | 2443 | 2449 | 2455 | rBV4 | 21853  | 40395  | 1.47%  | 0.243% |
| 57 | 17.309 | 2460 | 2463 | 2471 | rVB2 | 20213  | 28957  | 1.05%  | 0.174% |
| 58 | 17.491 | 2488 | 2494 | 2497 | rBV  | 135497 | 209096 | 7.60%  | 1.259% |
| 59 | 17.532 | 2497 | 2501 | 2508 | rVV  | 365636 | 548357 | 19.94% | 3.301% |
| 60 | 17.620 | 2508 | 2516 | 2522 | rVV  | 52541  | 87469  | 3.18%  | 0.526% |
| 61 | 17.885 | 2551 | 2561 | 2570 | rBV2 | 26309  | 51070  | 1.86%  | 0.307% |
| 62 | 18.361 | 2635 | 2642 | 2646 | rBV2 | 87780  | 143217 | 5.21%  | 0.862% |
| 63 | 18.408 | 2646 | 2650 | 2657 | rVB  | 79613  | 123700 | 4.50%  | 0.745% |
| 64 | 18.490 | 2657 | 2664 | 2669 | rBV  | 24283  | 40965  | 1.49%  | 0.247% |
| 65 | 18.555 | 2669 | 2675 | 2679 | rVV2 | 72531  | 155146 | 5.64%  | 0.934% |
| 66 | 18.590 | 2679 | 2681 | 2688 | rVB3 | 48064  | 65905  | 2.40%  | 0.397% |
| 67 | 18.848 | 2720 | 2725 | 2729 | rBV  | 44960  | 66399  | 2.41%  | 0.400% |
| 68 | 18.895 | 2729 | 2733 | 2738 | rVB  | 45023  | 67837  | 2.47%  | 0.408% |
| 69 | 19.265 | 2791 | 2796 | 2801 | rBV  | 45128  | 78764  | 2.86%  | 0.474% |
| 70 | 19.406 | 2816 | 2820 | 2828 | rBV4 | 32548  | 65884  | 2.40%  | 0.397% |
| 71 | 19.530 | 2835 | 2841 | 2847 | rVB  | 428371 | 610267 | 22.19% | 3.673% |

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Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
CEB-5(2.5-5)20160223

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:\HPCHEM1\BNA\_G\METHODS\8270-BG022916.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 72 | 19.618 | 2853 | 2856 | 2862 | rVB7 | 22762  | 33036  | 1.20%  | 0.199% |
| 73 | 19.771 | 2878 | 2882 | 2886 | rBV  | 21456  | 30361  | 1.10%  | 0.183% |
| 74 | 19.859 | 2891 | 2897 | 2898 | rBV  | 67236  | 96208  | 3.50%  | 0.579% |
| 75 | 19.894 | 2898 | 2903 | 2910 | rVV  | 415230 | 628354 | 22.85% | 3.782% |
| 76 | 19.959 | 2910 | 2914 | 2920 | rVB3 | 28946  | 48041  | 1.75%  | 0.289% |
| 77 | 20.082 | 2929 | 2935 | 2940 | rBV  | 772332 | 967940 | 35.20% | 5.826% |
| 78 | 20.206 | 2951 | 2956 | 2957 | rBV4 | 33118  | 47324  | 1.72%  | 0.285% |
| 79 | 20.376 | 2981 | 2985 | 2989 | rVB2 | 67771  | 105178 | 3.82%  | 0.633% |
| 80 | 20.476 | 2998 | 3002 | 3005 | rBV  | 36273  | 49918  | 1.82%  | 0.300% |
| 81 | 20.535 | 3006 | 3012 | 3016 | rVV2 | 43127  | 85716  | 3.12%  | 0.516% |
| 82 | 20.670 | 3032 | 3035 | 3040 | rBV  | 38558  | 57289  | 2.08%  | 0.345% |
| 83 | 20.717 | 3040 | 3043 | 3046 | rVV  | 37032  | 49113  | 1.79%  | 0.296% |
| 84 | 20.752 | 3046 | 3049 | 3055 | rVB2 | 39195  | 56578  | 2.06%  | 0.341% |
| 85 | 21.134 | 3112 | 3114 | 3122 | rVB  | 39821  | 57584  | 2.09%  | 0.347% |
| 86 | 21.369 | 3151 | 3154 | 3159 | rVB2 | 56024  | 79217  | 2.88%  | 0.477% |
| 87 | 21.422 | 3160 | 3163 | 3168 | rVB2 | 28324  | 39676  | 1.44%  | 0.239% |
| 88 | 21.622 | 3193 | 3197 | 3203 | rVB  | 134195 | 196425 | 7.14%  | 1.182% |
| 89 | 21.739 | 3211 | 3217 | 3224 | rBV2 | 208159 | 569913 | 20.72% | 3.430% |
| 90 | 21.804 | 3224 | 3228 | 3235 | rVB  | 170796 | 281827 | 10.25% | 1.696% |
| 91 | 23.989 | 3592 | 3600 | 3608 | rBV  | 97100  | 342937 | 12.47% | 2.064% |
| 92 | 24.871 | 3743 | 3750 | 3765 | rVB  | 85478  | 257253 | 9.35%  | 1.548% |
| 93 | 25.035 | 3770 | 3778 | 3786 | rVV  | 70529  | 190595 | 6.93%  | 1.147% |

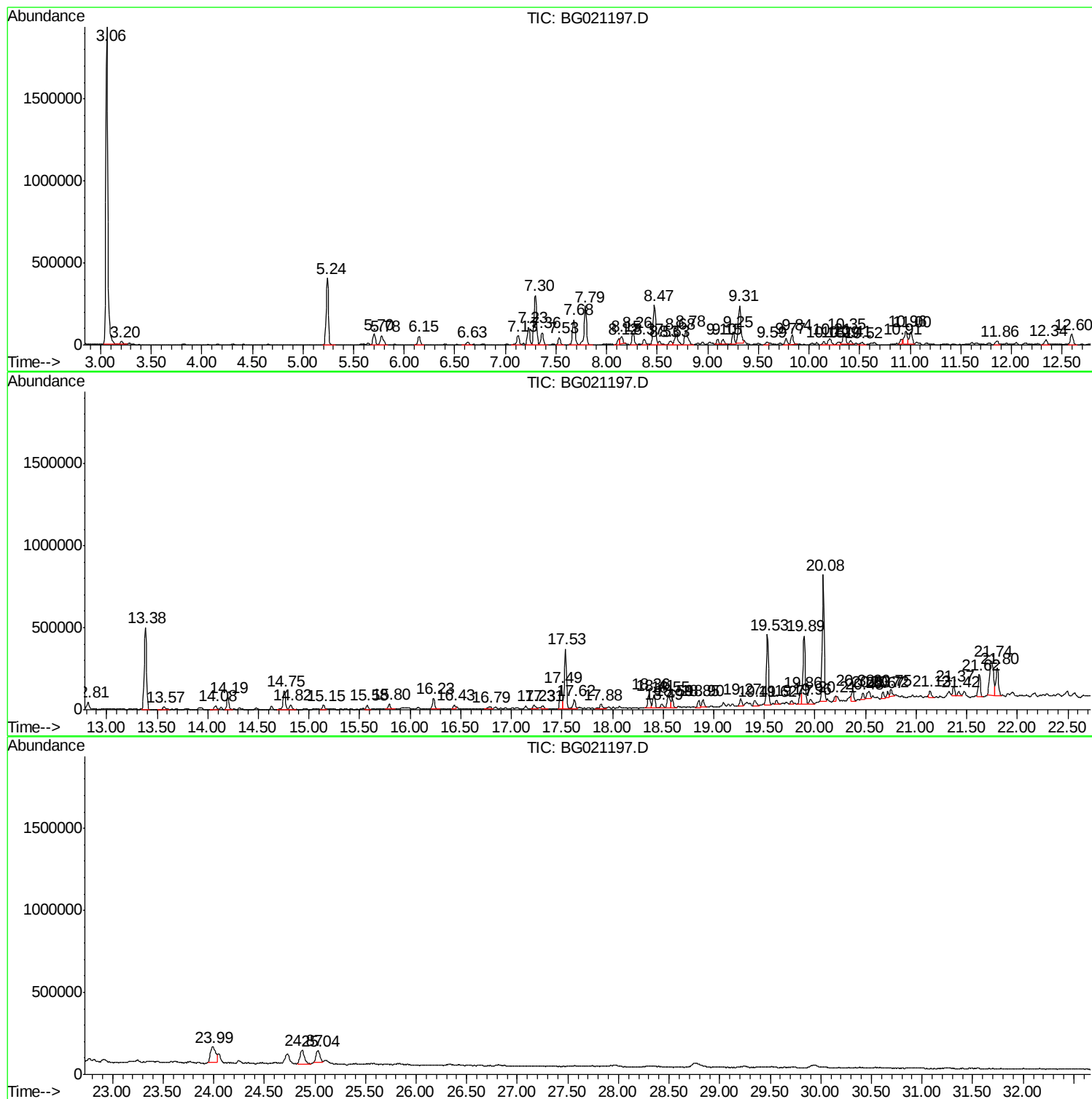
Sum of corrected areas: 16613361

Data Path : Z:\HPCHEM1\BNA G\DATA\BG030116\  
Data File : BG021197.D  
Acq On : 1 Mar 2016 23:43  
Operator : UM/SJ  
Sample : H1565-08  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
CEB-5(2.5-5)20160223

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.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\BG030116\  
Data File : BG021197.D  
Acq On : 1 Mar 2016 23:43  
Operator : UM/SJ  
Sample : H1565-08  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
CEB-5(2.5-5)20160223

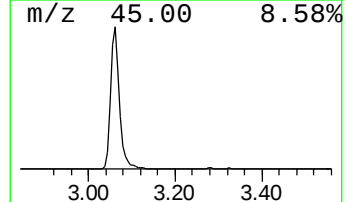
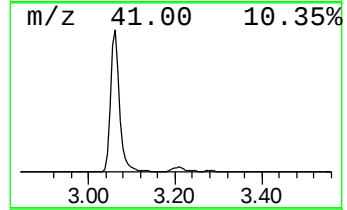
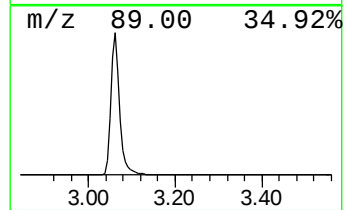
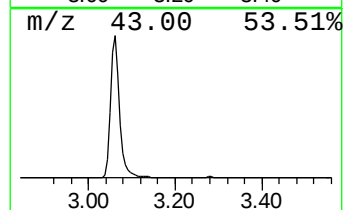
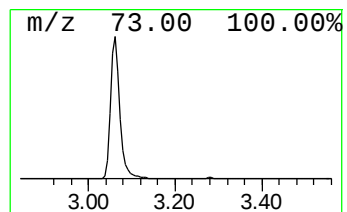
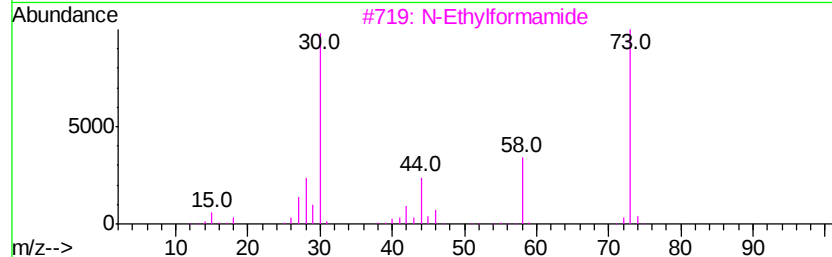
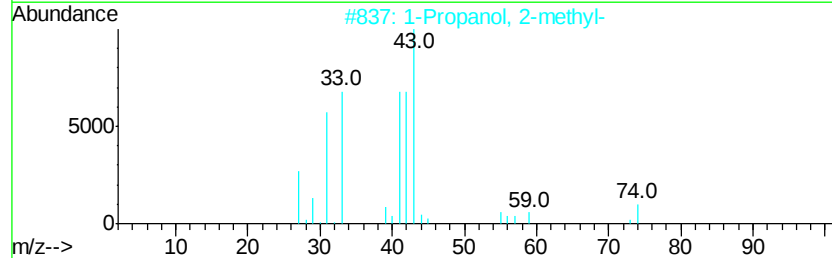
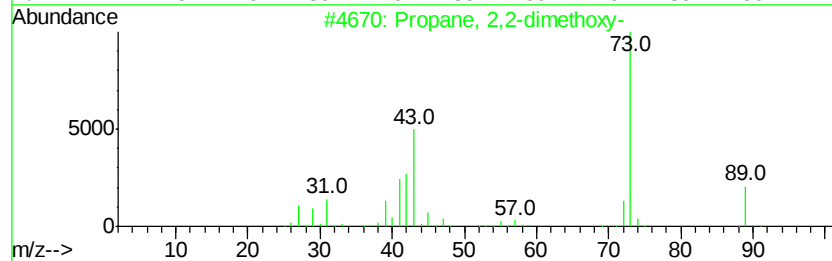
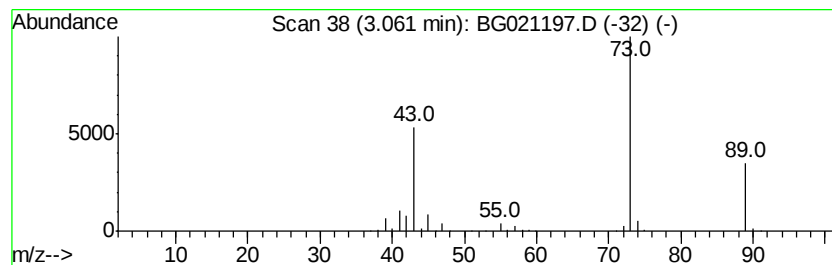
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 1

| R.T. | EstConc   | Area    | Relative to ISTD       | R.T. |
|------|-----------|---------|------------------------|------|
| 3.06 | 831.46 ng | 2750210 | 1,4-Dichlorobenzene-d4 | 8.15 |

| Hit# | of | Tentative ID                        | MW  | MolForm       | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------------|--------------|------|
| 1    | 5  | Propane, 2,2-dimethoxy-             | 104 | C5H12O2       | 000077-76-9  | 56   |
| 2    |    | 1-Propanol, 2-methyl-               | 74  | C4H10O        | 000078-83-1  | 9    |
| 3    |    | N-Ethylformamide                    | 73  | C3H7NO        | 000627-45-2  | 9    |
| 4    |    | 1,3-Diaminoquanidine                | 89  | CH7N5         | 004364-78-7  | 9    |
| 5    |    | 1,3-Dioxolane, 2-(3-bromo-5,5,5-... | 352 | C10H16BrCl3O2 | 1000115-31-4 | 9    |



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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

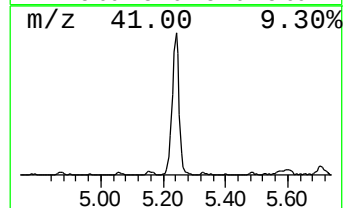
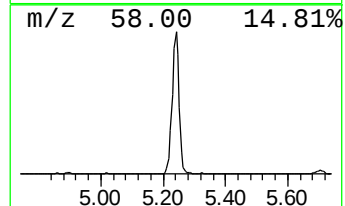
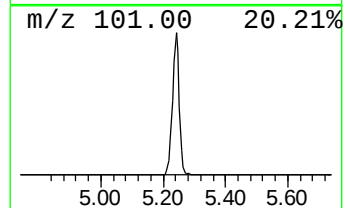
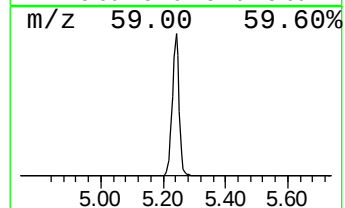
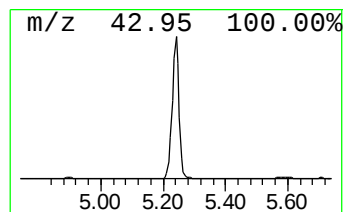
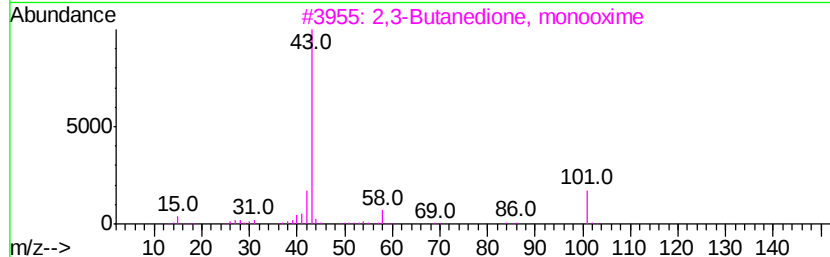
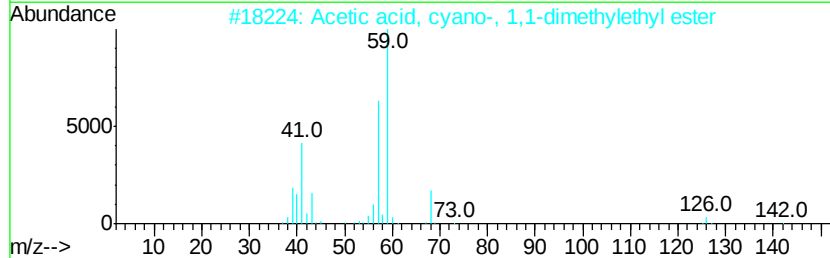
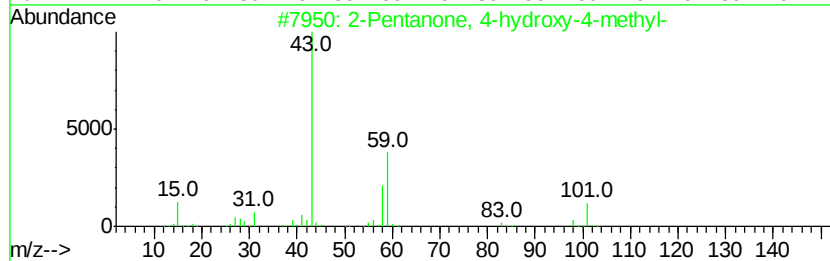
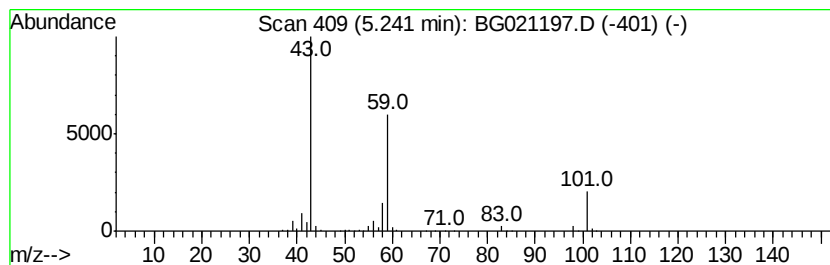
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc   | Area   | Relative to ISTD       | R.T. |
|------|-----------|--------|------------------------|------|
| 5.24 | 188.80 ng | 624495 | 1,4-Dichlorobenzene-d4 | 8.15 |

| Hit# | of | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-     | 116 | C6H12O2  | 000123-42-2 | 56   |
| 2    |    | Acetic acid, cyano-, 1,1-dimethyl... | 141 | C7H11NO2 | 001116-98-9 | 25   |
| 3    |    | 2,3-Butanedione, monooxime           | 101 | C4H7NO2  | 000057-71-6 | 12   |
| 4    |    | Butane, 1-ethoxy-                    | 102 | C6H14O   | 000628-81-9 | 9    |
| 5    |    | Morpholine, 4-methyl-                | 101 | C5H11NO  | 000109-02-4 | 9    |



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ClientSampleId :  
CEB-5(2.5-5)20160223

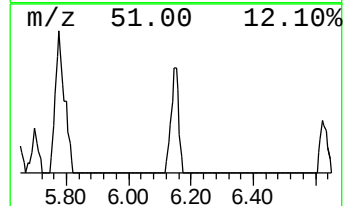
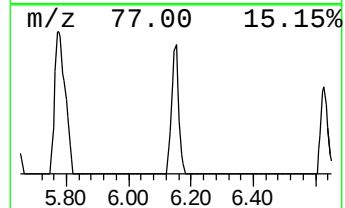
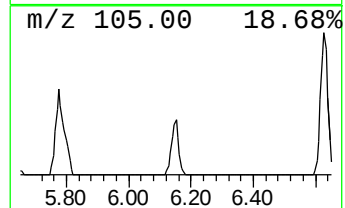
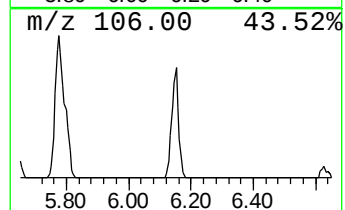
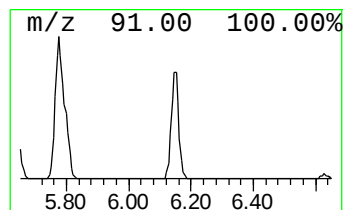
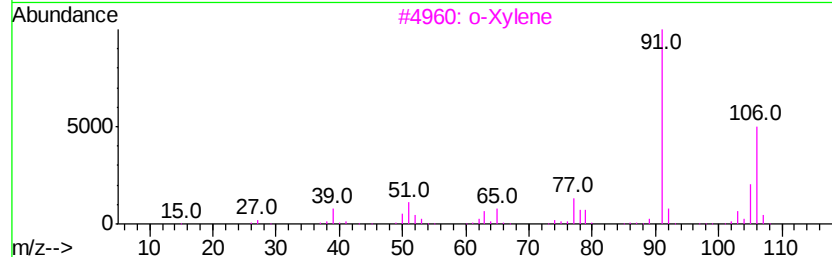
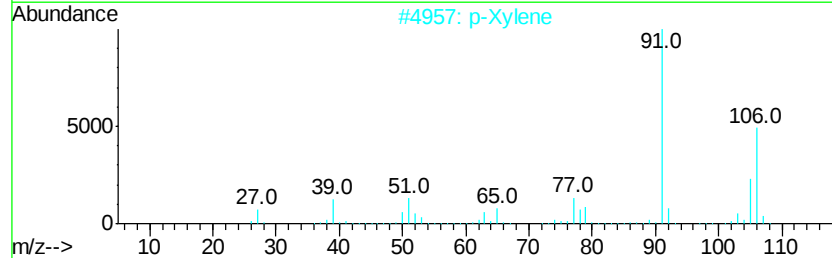
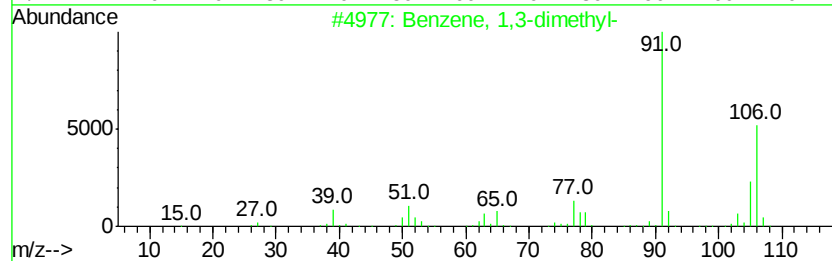
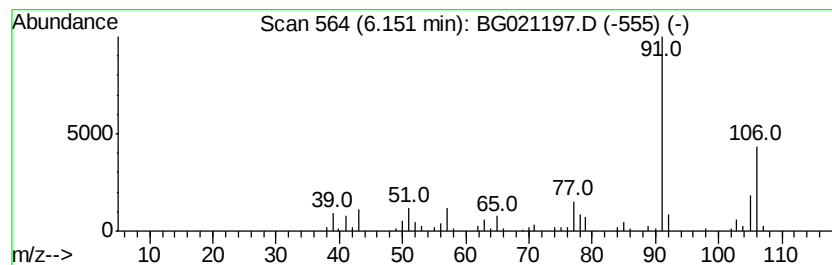
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 4 Benzene, 1,3-dimethyl- Concentration Rank 14

| R.T. | EstConc  | Area  | Relative to ISTD       | R.T. |
|------|----------|-------|------------------------|------|
| 6.15 | 27.41 ng | 90675 | 1,4-Dichlorobenzene-d4 | 8.15 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,3-dimethyl-              | 106 | C8H10   | 000108-38-3 | 95   |
| 2    |    | p-Xylene                            | 106 | C8H10   | 000106-42-3 | 95   |
| 3    |    | o-Xylene                            | 106 | C8H10   | 000095-47-6 | 95   |
| 4    |    | Ethylbenzene                        | 106 | C8H10   | 000100-41-4 | 87   |
| 5    |    | 1,3-Cyclopentadiene, 5-(1-methyl... | 106 | C8H10   | 002175-91-9 | 74   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG030116\  
Data File : BG021197.D  
Acq On : 1 Mar 2016 23:43  
Operator : UM/SJ  
Sample : H1565-08  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
CEB-5(2.5-5)20160223

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

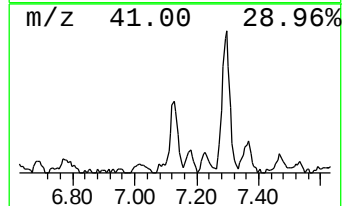
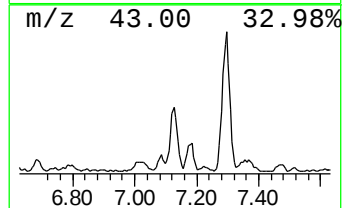
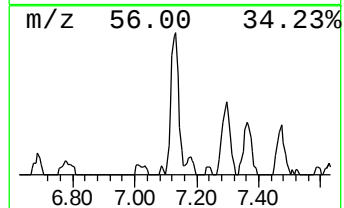
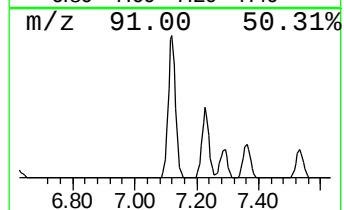
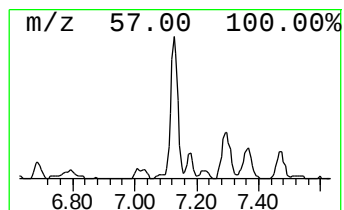
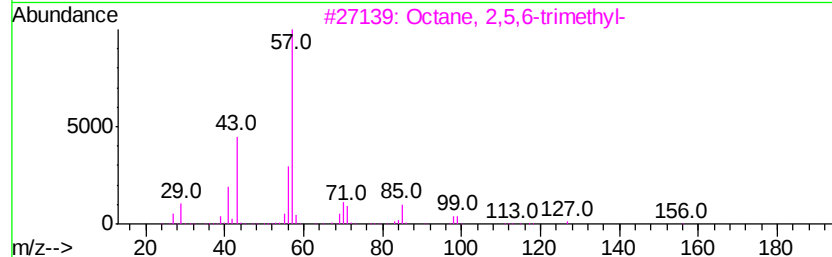
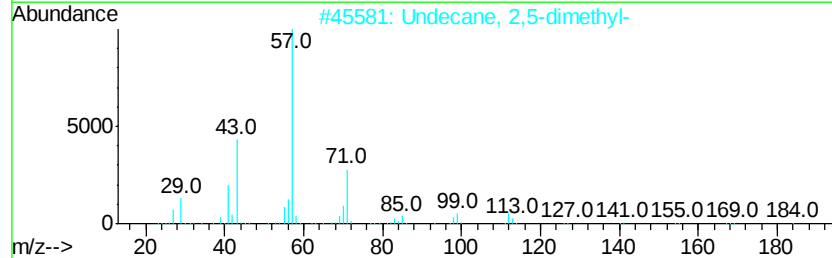
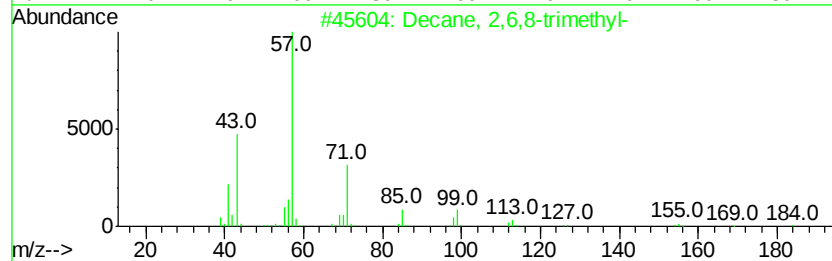
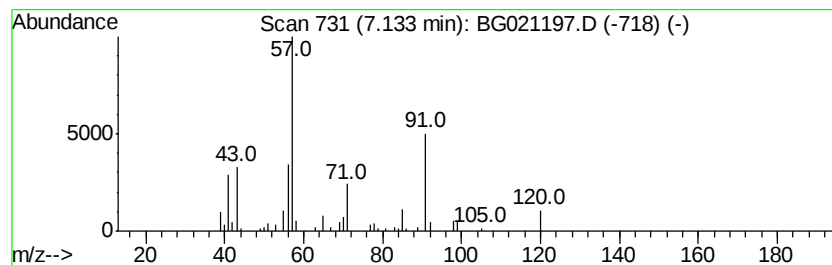
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 unknown7.13 Concentration Rank 13

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 7.13 | 31.32 ng | 103603 | 1,4-Dichlorobenzene-d4 | 8.15 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Decane, 2,6,8-trimethyl- | 184 | C13H28  | 062108-26-3 | 47   |
| 2    |    | Undecane, 2,5-dimethyl-  | 184 | C13H28  | 017301-22-3 | 43   |
| 3    |    | Octane, 2,5,6-trimethyl- | 156 | C11H24  | 062016-14-2 | 43   |
| 4    |    | Benzene, propyl-         | 120 | C9H12   | 000103-65-1 | 35   |
| 5    |    | Octane, 3-methyl-        | 128 | C9H20   | 002216-33-3 | 35   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG030116\  
Data File : BG021197.D  
Acq On : 1 Mar 2016 23:43  
Operator : UM/SJ  
Sample : H1565-08  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
CEB-5(2.5-5)20160223

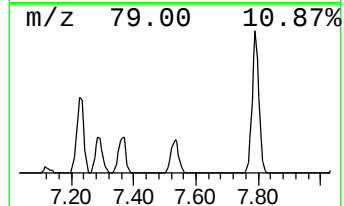
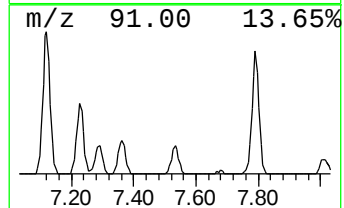
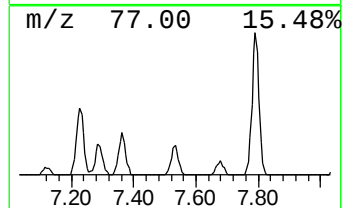
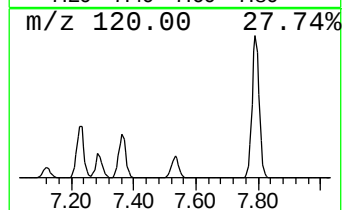
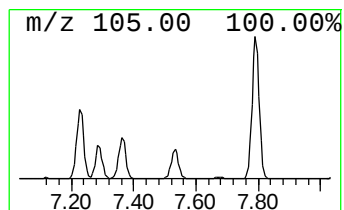
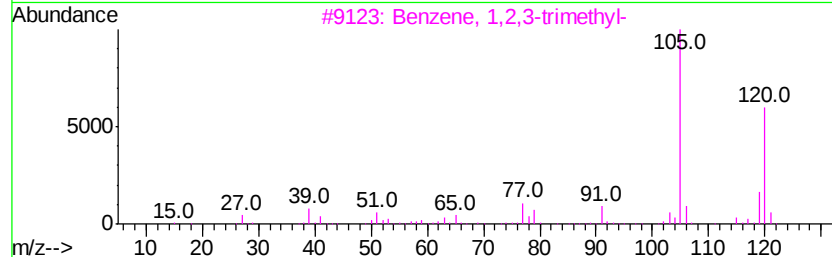
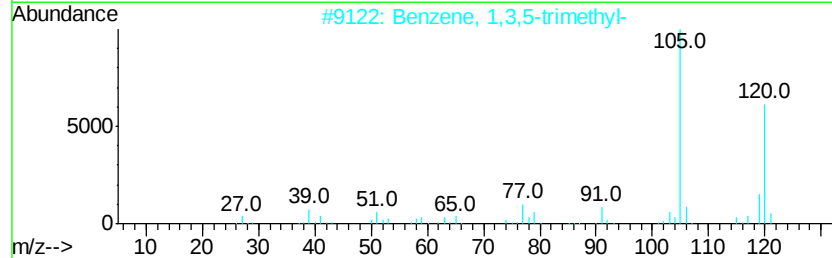
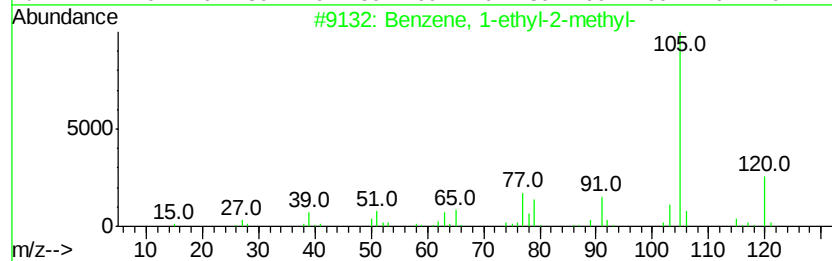
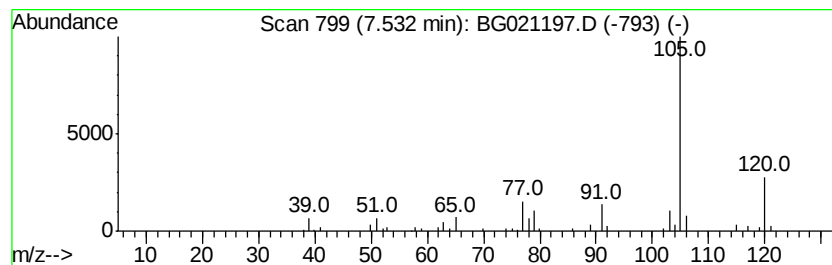
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 Benzene, 1-ethyl-2-methyl- Concentration Rank 15

| R.T. | EstConc  | Area  | Relative to ISTD       | R.T. |
|------|----------|-------|------------------------|------|
| 7.53 | 21.64 ng | 71586 | 1,4-Dichlorobenzene-d4 | 8.15 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 93   |
| 2    |    | Benzene, 1,3,5-trimethyl-  | 120 | C9H12   | 000108-67-8 | 91   |
| 3    |    | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |
| 4    |    | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 91   |
| 5    |    | Benzene, (1-methylethyl)-  | 120 | C9H12   | 000098-82-8 | 90   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG030116\  
Data File : BG021197.D  
Acq On : 1 Mar 2016 23:43  
Operator : UM/SJ  
Sample : H1565-08  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
CEB-5(2.5-5)20160223

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

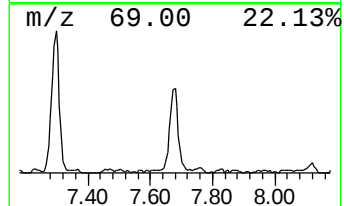
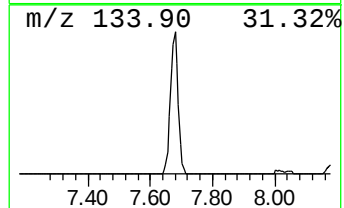
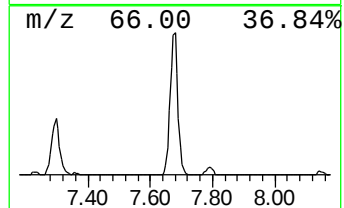
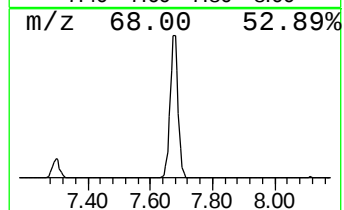
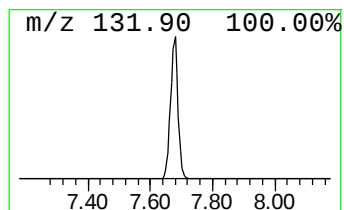
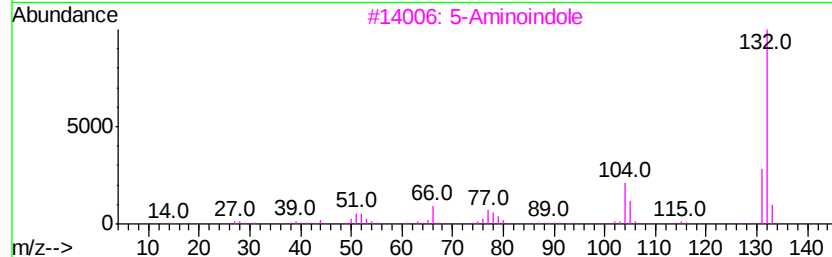
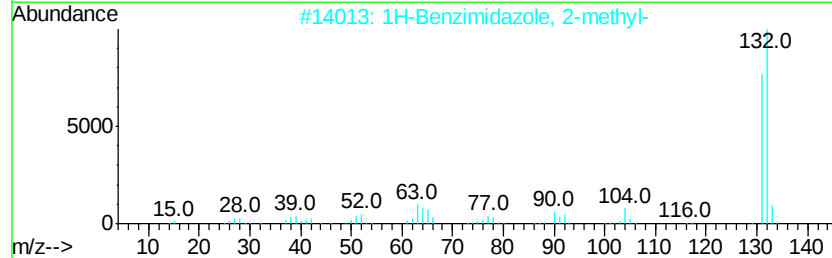
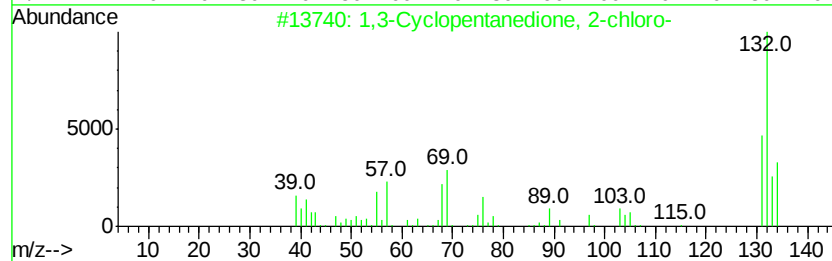
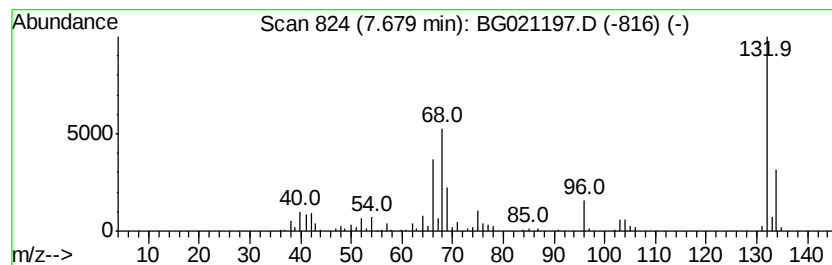
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TIC Integration Parameters: LSCINT.P

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Peak Number 9 unknown7.68 Concentration Rank 5

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 7.68 | 81.77 ng | 270460 | 1,4-Dichlorobenzene-d4 | 8.15 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | 1,3-Cyclopentanedione, 2-chloro-    | 132 | C5H5ClO2  | 014203-19-1 | 35   |
| 2    |    | 1H-Benzimidazole, 2-methyl-         | 132 | C8H8N2    | 000615-15-6 | 22   |
| 3    |    | 5-Aminoindole                       | 132 | C8H8N2    | 005192-03-0 | 22   |
| 4    |    | 5-Fluoro-2-chloropyrimidine         | 132 | C4H2ClFN2 | 062802-42-0 | 16   |
| 5    |    | 2H-Cyclopenta[d]pyridazine, 2-me... | 132 | C8H8N2    | 022291-85-6 | 12   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG030116\  
Data File : BG021197.D  
Acq On : 1 Mar 2016 23:43  
Operator : UM/SJ  
Sample : H1565-08  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
CEB-5(2.5-5)20160223

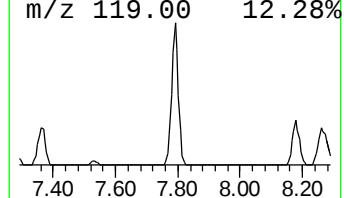
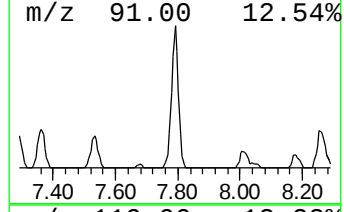
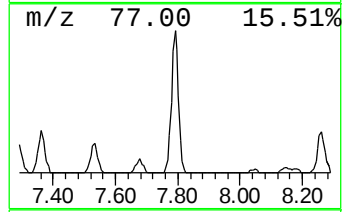
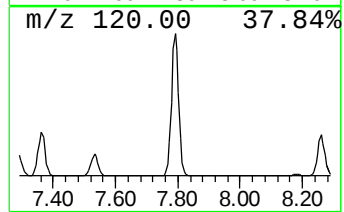
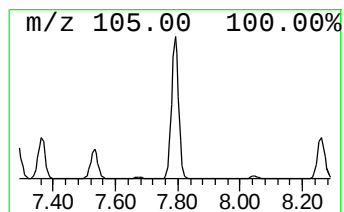
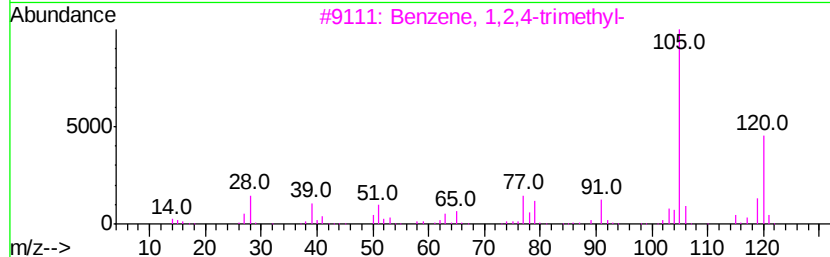
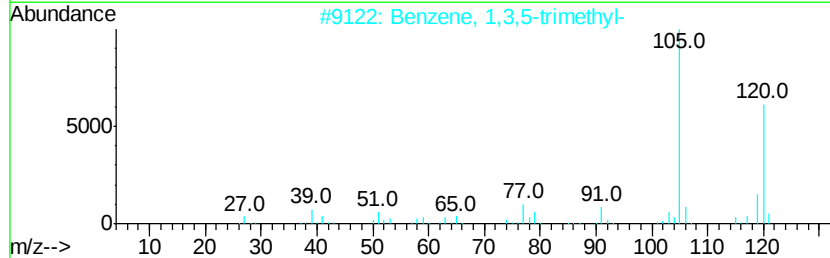
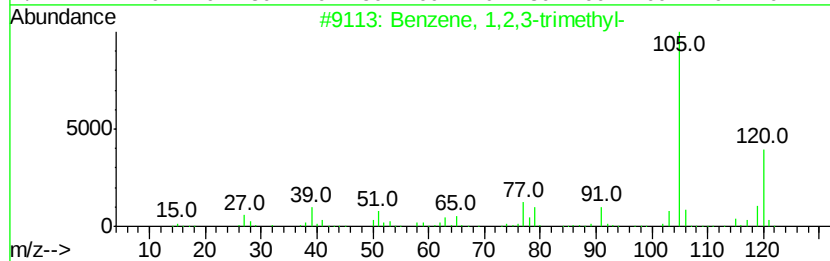
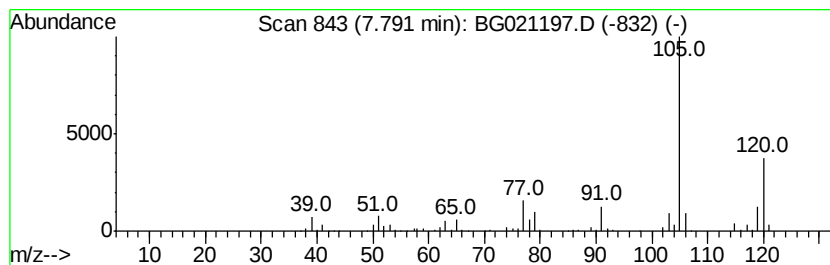
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 Benzene, 1,2,3-trimethyl- Concentration Rank 4

| R.T. | EstConc   | Area   | Relative to ISTD       | R.T. |
|------|-----------|--------|------------------------|------|
| 7.79 | 121.39 ng | 401526 | 1,4-Dichlorobenzene-d4 | 8.15 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 95   |
| 2    |    | Benzene, 1,3,5-trimethyl-  | 120 | C9H12   | 000108-67-8 | 94   |
| 3    |    | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 94   |
| 4    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |
| 5    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 90   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG030116\  
Data File : BG021197.D  
Acq On : 1 Mar 2016 23:43  
Operator : UM/SJ  
Sample : H1565-08  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
CEB-5(2.5-5)20160223

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

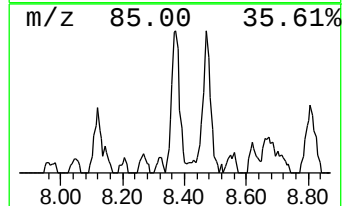
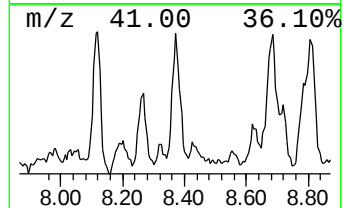
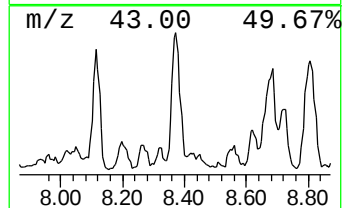
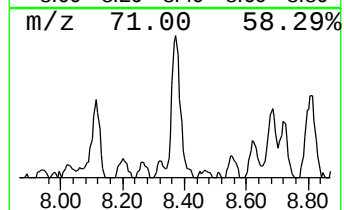
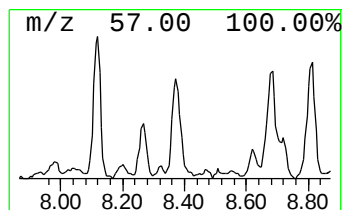
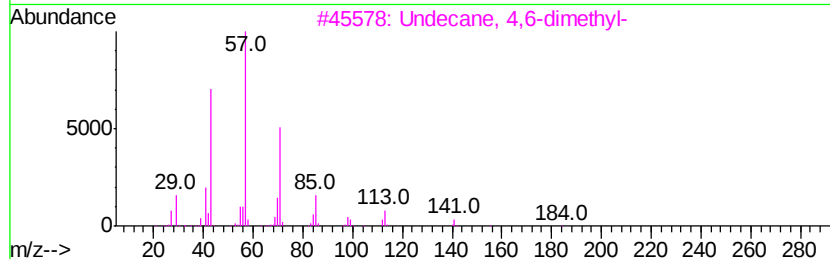
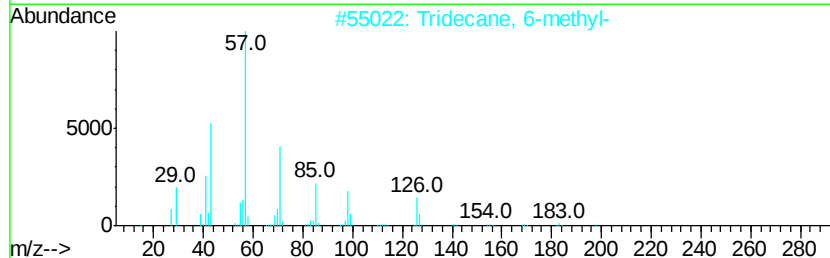
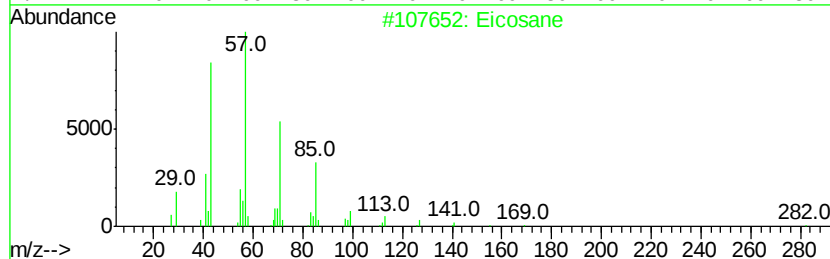
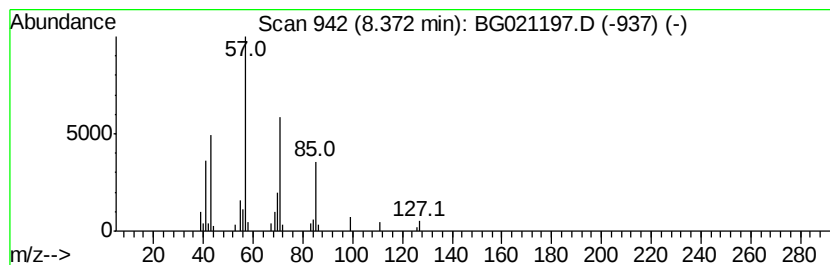
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Eicosane Concentration Rank 16

| R.T. | EstConc  | Area  | Relative to ISTD       | R.T. |
|------|----------|-------|------------------------|------|
| 8.37 | 17.47 ng | 57787 | 1,4-Dichlorobenzene-d4 | 8.15 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Eicosane                | 282 | C20H42  | 000112-95-8 | 72   |
| 2    |    | Tridecane, 6-methyl-    | 198 | C14H30  | 013287-21-3 | 72   |
| 3    |    | Undecane, 4,6-dimethyl- | 184 | C13H28  | 017312-82-2 | 72   |
| 4    |    | Hexadecane              | 226 | C16H34  | 000544-76-3 | 64   |
| 5    |    | Tetradecane             | 198 | C14H30  | 000629-59-4 | 59   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG030116\  
Data File : BG021197.D  
Acq On : 1 Mar 2016 23:43  
Operator : UM/SJ  
Sample : H1565-08  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
CEB-5(2.5-5)20160223

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

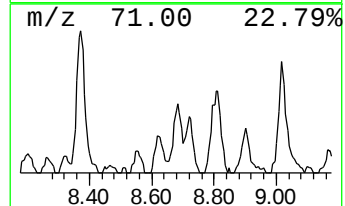
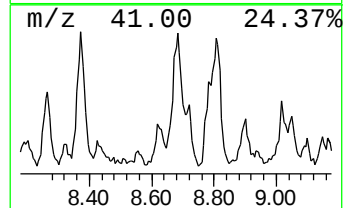
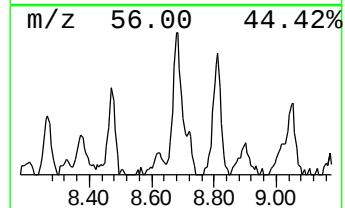
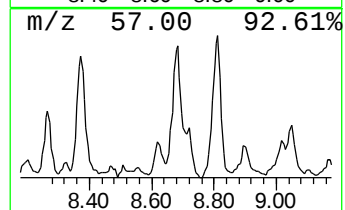
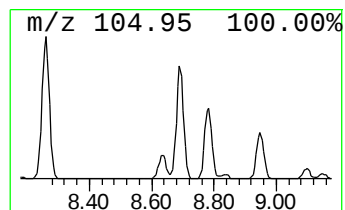
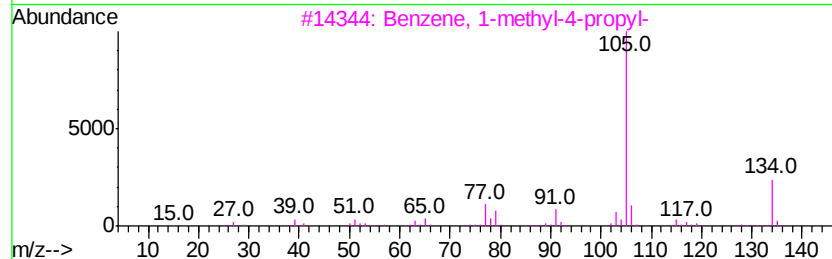
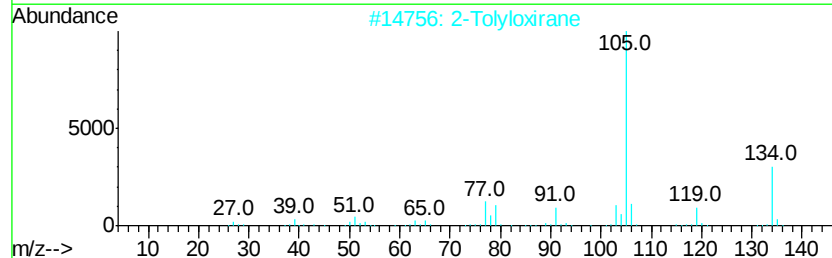
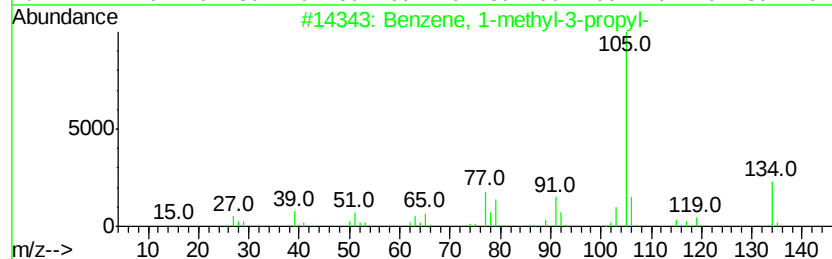
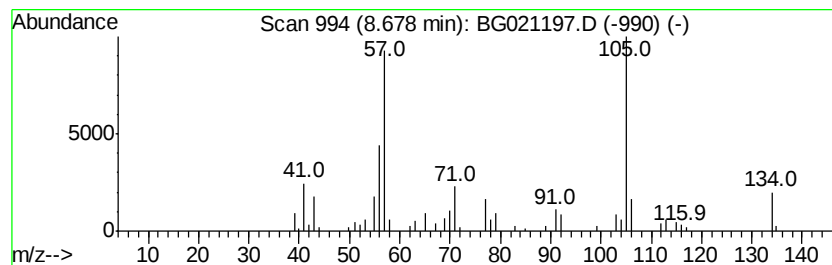
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Benzene, 1-methyl-3-propyl- Concentration Rank 8

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 8.68 | 45.54 ng | 150641 | 1,4-Dichlorobenzene-d4 | 8.15 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-3-propyl- | 134 | C10H14  | 001074-43-7 | 93   |
| 2    |    | 2-Tolyloxirane              | 134 | C9H10O  | 002783-26-8 | 70   |
| 3    |    | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 70   |
| 4    |    | Benzene, (1-methylpropyl)-  | 134 | C10H14  | 000135-98-8 | 70   |
| 5    |    | Benzene, 1-methyl-2-propyl- | 134 | C10H14  | 001074-17-5 | 53   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG030116\  
Data File : BG021197.D  
Acq On : 1 Mar 2016 23:43  
Operator : UM/SJ  
Sample : H1565-08  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
CEB-5(2.5-5)20160223

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

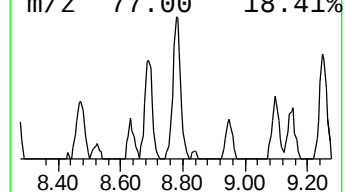
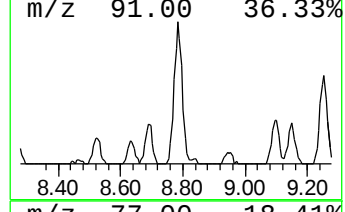
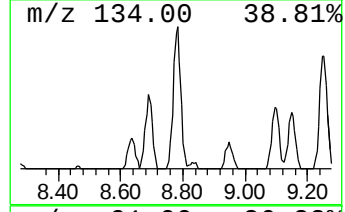
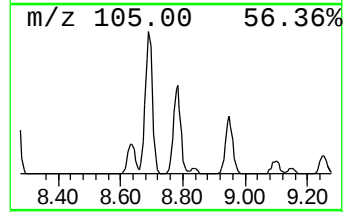
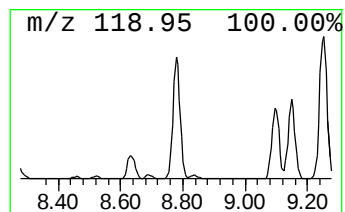
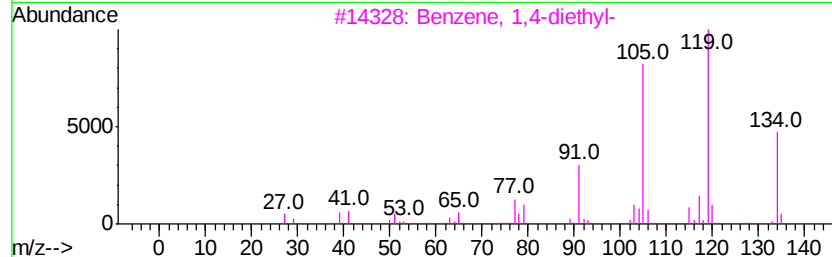
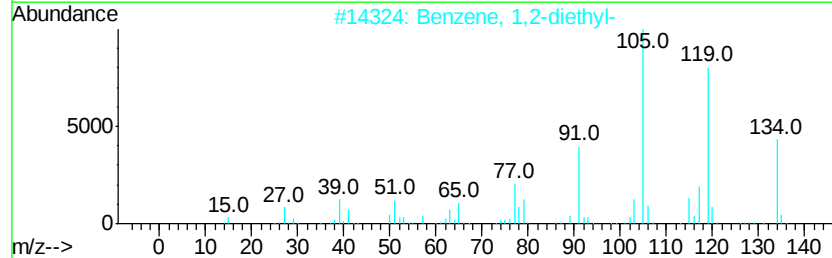
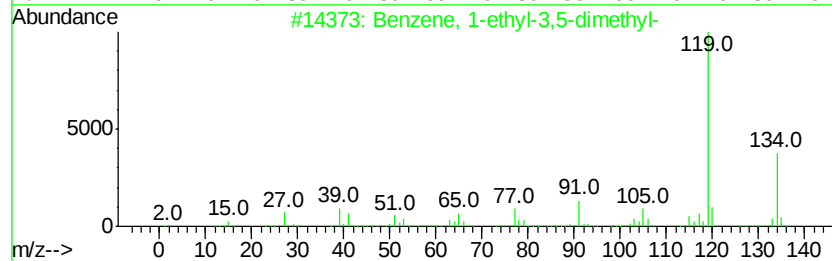
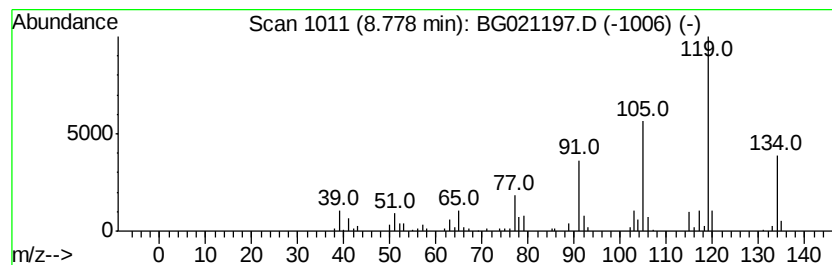
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 6

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 8.78 | 61.26 ng | 202641 | 1,4-Dichlorobenzene-d4 | 8.15 |

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



Data Path : Z:\HPCHEM1\BNA G\DATA\BG030116\  
Data File : BG021197.D  
Acq On : 1 Mar 2016 23:43  
Operator : UM/SJ  
Sample : H1565-08  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
CEB-5(2.5-5)20160223

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

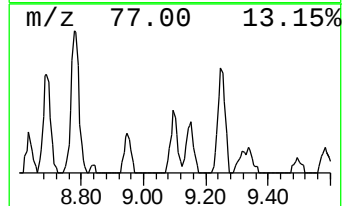
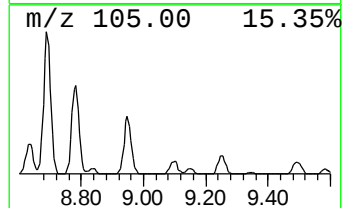
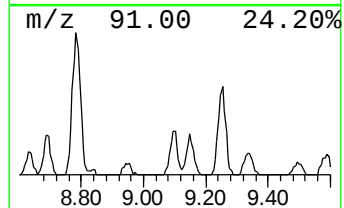
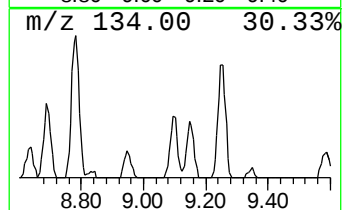
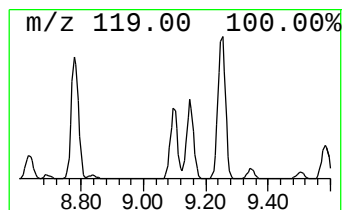
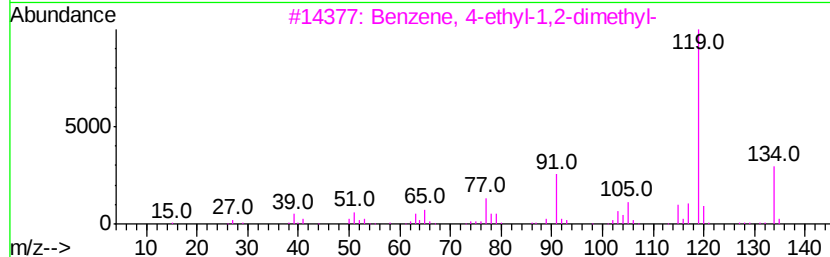
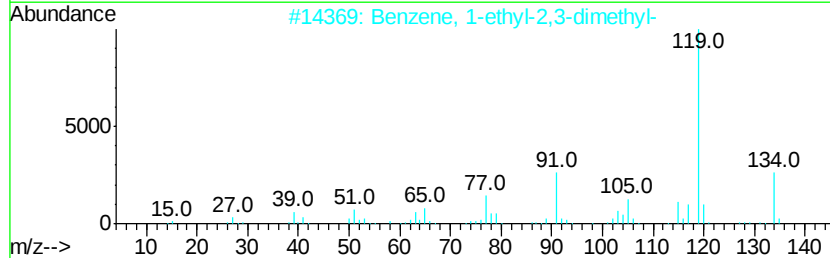
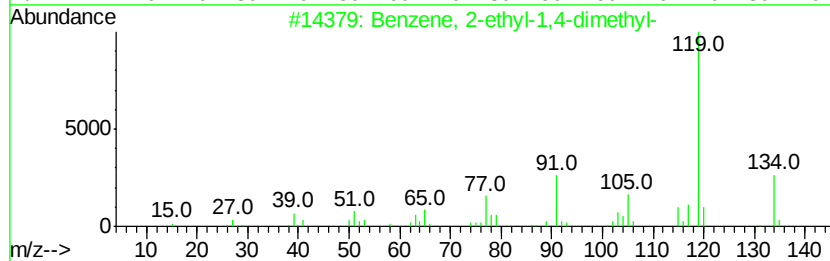
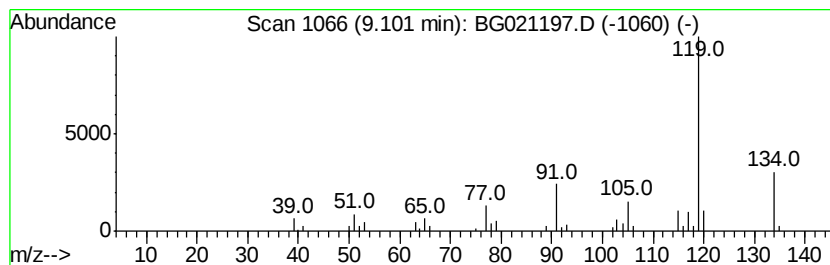
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 19

| R.T. | EstConc  | Area  | Relative to ISTD       | R.T. |
|------|----------|-------|------------------------|------|
| 9.10 | 14.93 ng | 49388 | 1,4-Dichlorobenzene-d4 | 8.15 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 96   |
| 2    |    | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 96   |
| 3    |    | Benzene, 4-ethyl-1,2-dimethyl-      | 134 | C10H14  | 000934-80-5 | 96   |
| 4    |    | Benzene, 2-ethyl-1,3-dimethyl-      | 134 | C10H14  | 002870-04-4 | 95   |
| 5    |    | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 95   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG030116\  
Data File : BG021197.D  
Acq On : 1 Mar 2016 23:43  
Operator : UM/SJ  
Sample : H1565-08  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
CEB-5(2.5-5)20160223

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

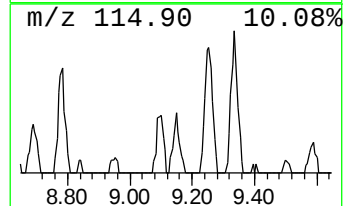
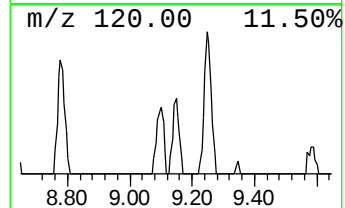
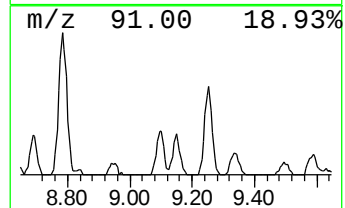
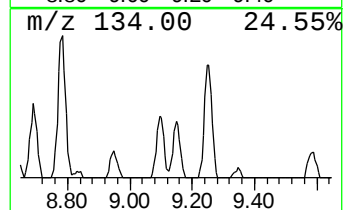
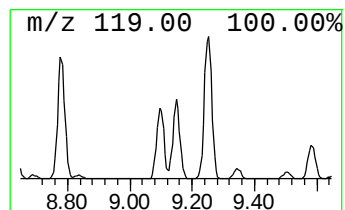
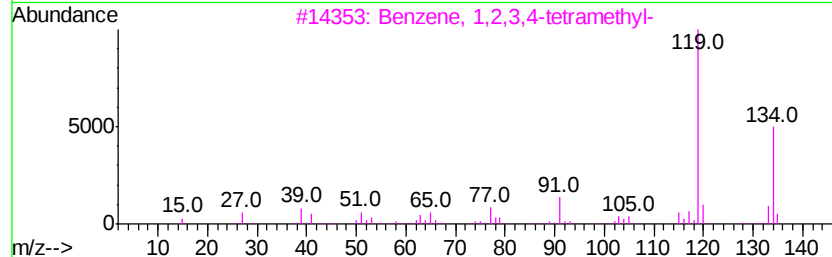
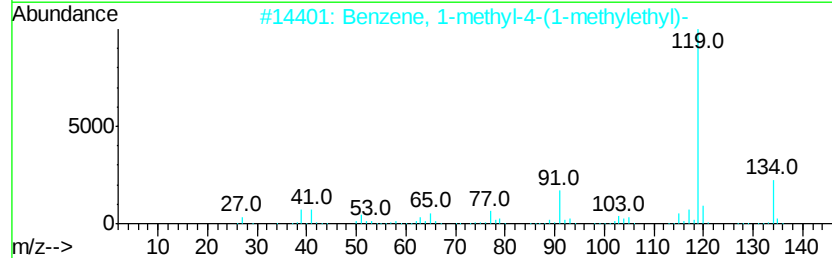
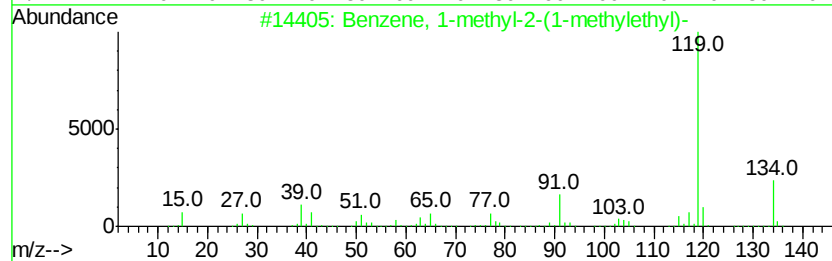
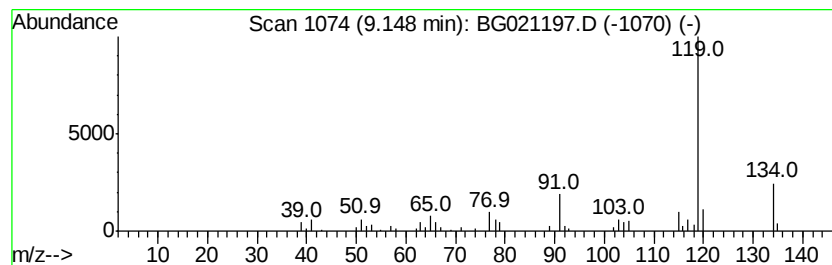
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Benzene, 1-methyl-2-(1-meth... Concentration Rank 18

| R.T. | EstConc  | Area  | Relative to ISTD       | R.T. |
|------|----------|-------|------------------------|------|
| 9.15 | 15.72 ng | 52009 | 1,4-Dichlorobenzene-d4 | 8.15 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 94   |
| 2    |    | Benzene, 1-methyl-4-(1-methyleth... | 134 | C10H14  | 000099-87-6 | 94   |
| 3    |    | Benzene, 1,2,3,4-tetramethyl-       | 134 | C10H14  | 000488-23-3 | 93   |
| 4    |    | Benzene, 1,2,4,5-tetramethyl-       | 134 | C10H14  | 000095-93-2 | 91   |
| 5    |    | Benzene, 4-ethyl-1,2-dimethyl-      | 134 | C10H14  | 000934-80-5 | 91   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG030116\  
Data File : BG021197.D  
Acq On : 1 Mar 2016 23:43  
Operator : UM/SJ  
Sample : H1565-08  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
CEB-5(2.5-5)20160223

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

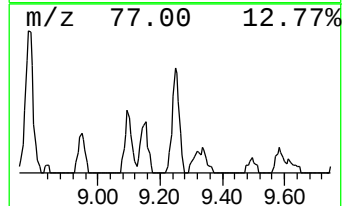
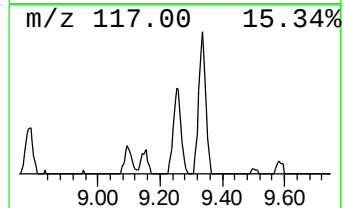
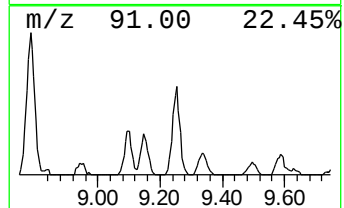
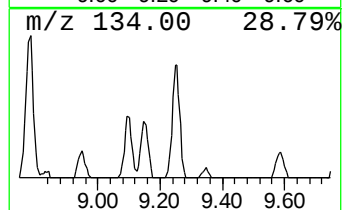
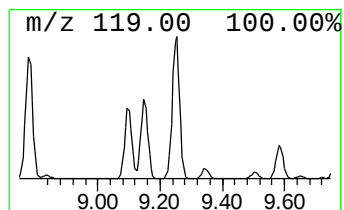
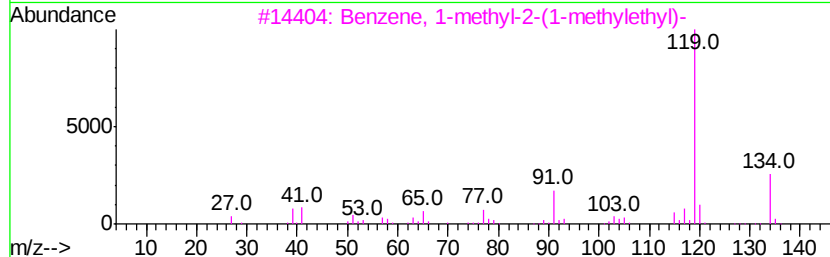
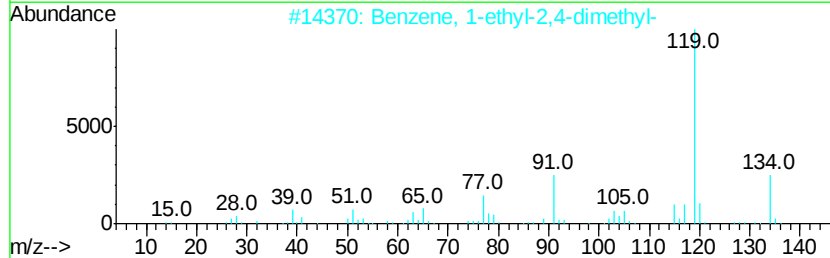
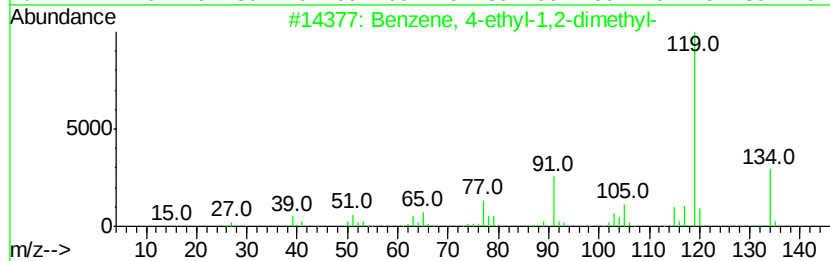
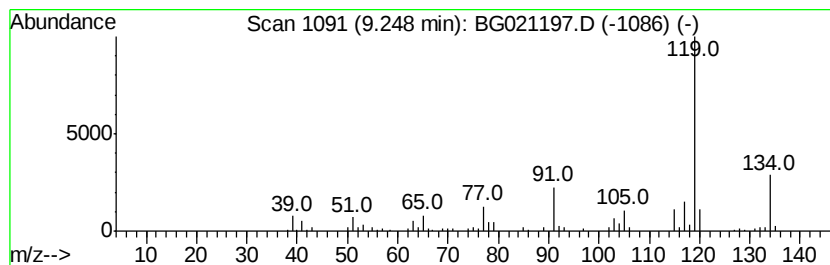
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 9

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 9.25 | 37.83 ng | 125131 | 1,4-Dichlorobenzene-d4 | 8.15 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 4-ethyl-1,2-dimethyl-       | 134 | C10H14  | 000934-80-5 | 95   |
| 2    |    | Benzene, 1-ethyl-2,4-dimethyl-       | 134 | C10H14  | 000874-41-9 | 95   |
| 3    |    | Benzene, 1-methyl-2-(1-methylethyl)- | 134 | C10H14  | 000527-84-4 | 95   |
| 4    |    | Benzene, 1-ethyl-2,3-dimethyl-       | 134 | C10H14  | 000933-98-2 | 94   |
| 5    |    | Benzene, 2-ethyl-1,3-dimethyl-       | 134 | C10H14  | 002870-04-4 | 94   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG030116\  
Data File : BG021197.D  
Acq On : 1 Mar 2016 23:43  
Operator : UM/SJ  
Sample : H1565-08  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
CEB-5(2.5-5)20160223

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

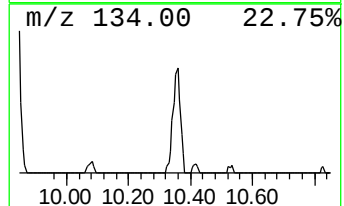
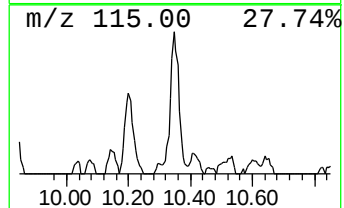
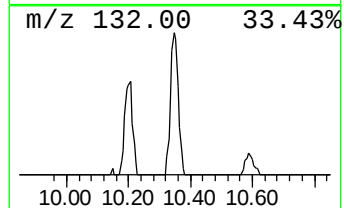
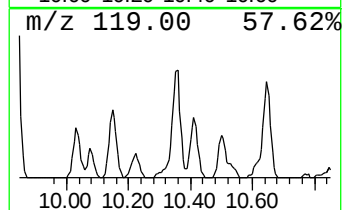
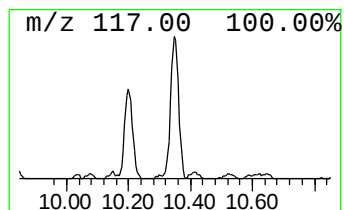
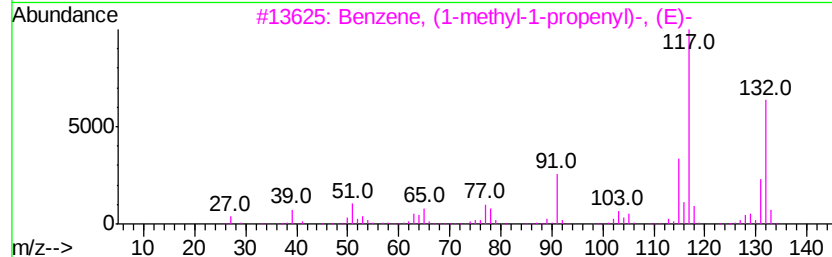
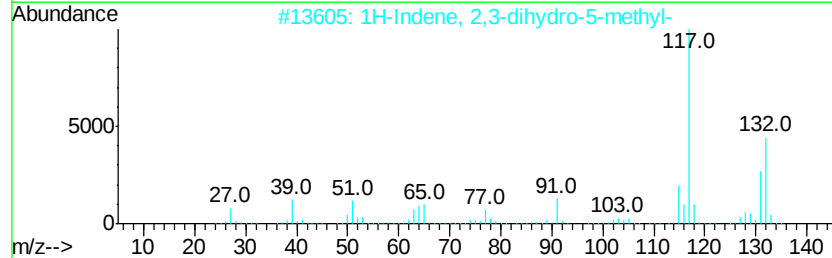
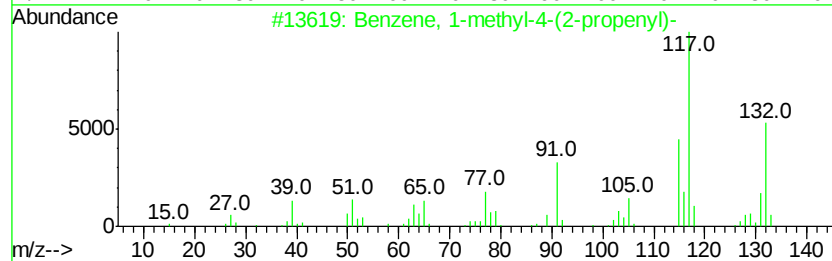
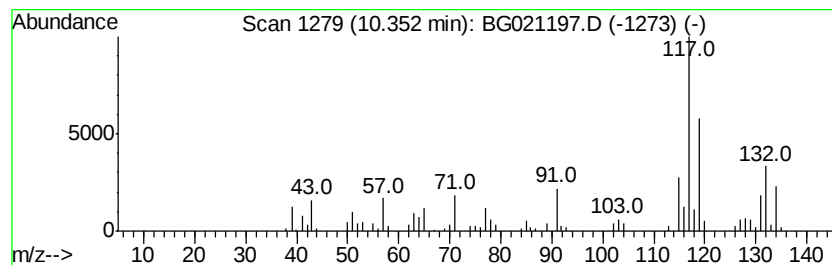
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 Benzene, 1-methyl-4-(2-prop... Concentration Rank 17

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 10.35 | 17.02 ng | 136708 | Naphthalene-d8   | 10.96 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-4-(2-propenyl)-   | 132 | C10H12  | 003333-13-9 | 87   |
| 2    |    | 1H-Indene, 2,3-dihydro-5-methyl-    | 132 | C10H12  | 000874-35-1 | 86   |
| 3    |    | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000768-00-3 | 70   |
| 4    |    | Benzene, 1-methyl-2-(2-propenyl)-   | 132 | C10H12  | 001587-04-8 | 70   |
| 5    |    | Benzene, 1-ethenyl-3-ethyl-         | 132 | C10H12  | 007525-62-4 | 70   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG030116\  
Data File : BG021197.D  
Acq On : 1 Mar 2016 23:43  
Operator : UM/SJ  
Sample : H1565-08  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

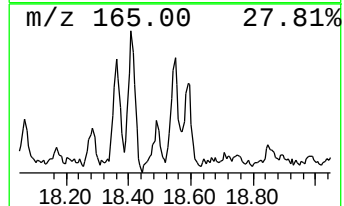
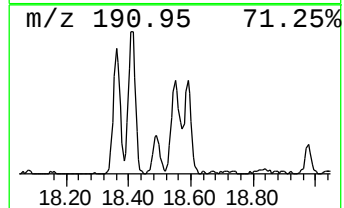
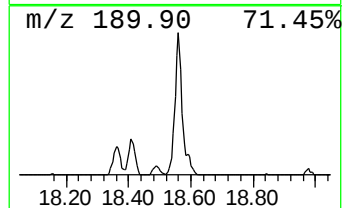
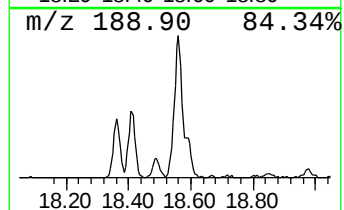
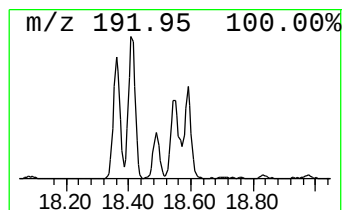
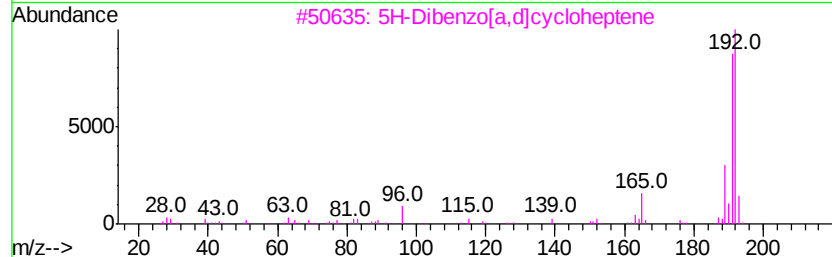
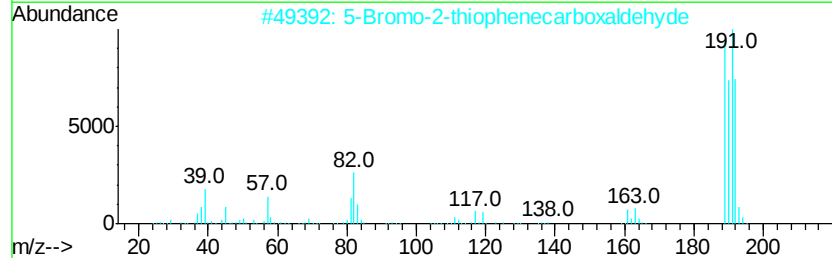
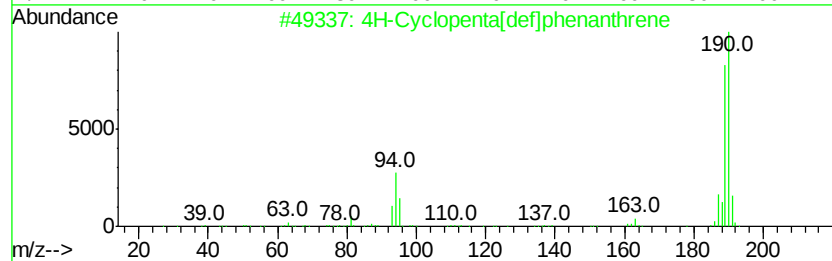
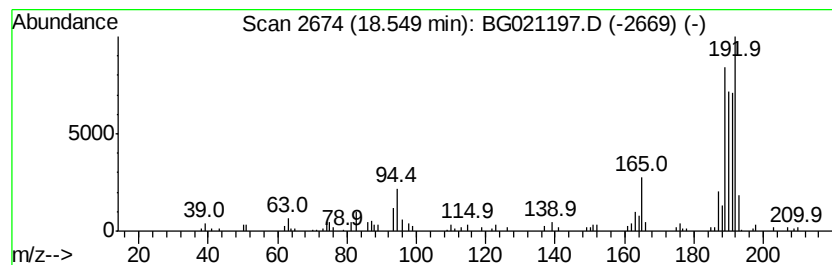
Instrument :  
BNA\_G  
ClientSampleId :  
CEB-5(2.5-5)20160223

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 20 4H-Cyclopenta[def]phenanthrene Concentration Rank 20

| R.T.    | EstConc  | Area                               | Relative to ISTD | R.T.     |             |      |
|---------|----------|------------------------------------|------------------|----------|-------------|------|
| 18.55   | 14.84 ng | 155146                             | Phenanthrene-d10 | 17.49    |             |      |
| Hit# of | 5        | Tentative ID                       | MW               | MolForm  | CAS#        | Qual |
| 1       |          | 4H-Cyclopenta[def]phenanthrene     | 190              | C15H10   | 000203-64-5 | 58   |
| 2       |          | 5-Bromo-2-thiophenecarboxaldehyde  | 190              | C5H3BrOS | 004701-17-1 | 52   |
| 3       |          | 5H-Dibenzo[a,d]cycloheptene        | 192              | C15H12   | 000256-81-5 | 51   |
| 4       |          | 6H-Cyclobuta[ik]phenanthrene       | 190              | C15H10   | 083469-43-6 | 50   |
| 5       |          | Benzene, 1,1'-(1,2-propadienyli... | 192              | C15H12   | 014251-57-1 | 38   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG030116\  
 Data File : BG021197.D  
 Acq On : 1 Mar 2016 23:43  
 Operator : UM/SJ  
 Sample : H1565-08  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 CEB-5(2.5-5)20160223

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG022916.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 |
| Propane, 2,2-dime... | 3.06  | 831.5   | ng    | 2750210  | 1                     | 8.15  | 66154  | 20.0 |
| 2-Pentanone, 4-hy... | 5.24  | 188.8   | ng    | 624495   | 1                     | 8.15  | 66154  | 20.0 |
| Benzene, 1,3-dime... | 6.15  | 27.4    | ng    | 90675    | 1                     | 8.15  | 66154  | 20.0 |
| unknown7.13          | 7.13  | 31.3    | ng    | 103603   | 1                     | 8.15  | 66154  | 20.0 |
| Benzene, 1-ethyl-... | 7.53  | 21.6    | ng    | 71586    | 1                     | 8.15  | 66154  | 20.0 |
| unknown7.68          | 7.68  | 81.8    | ng    | 270460   | 1                     | 8.15  | 66154  | 20.0 |
| Benzene, 1,2,3-tr... | 7.79  | 121.4   | ng    | 401526   | 1                     | 8.15  | 66154  | 20.0 |
| Eicosane             | 8.37  | 17.5    | ng    | 57787    | 1                     | 8.15  | 66154  | 20.0 |
| Benzene, 1-methyl... | 8.68  | 45.5    | ng    | 150641   | 1                     | 8.15  | 66154  | 20.0 |
| Benzene, 1-ethyl-... | 8.78  | 61.3    | ng    | 202641   | 1                     | 8.15  | 66154  | 20.0 |
| Benzene, 2-ethyl-... | 9.10  | 14.9    | ng    | 49388    | 1                     | 8.15  | 66154  | 20.0 |
| Benzene, 1-methyl... | 9.15  | 15.7    | ng    | 52009    | 1                     | 8.15  | 66154  | 20.0 |
| Benzene, 4-ethyl-... | 9.25  | 37.8    | ng    | 125131   | 1                     | 8.15  | 66154  | 20.0 |
| Benzene, 1-methyl... | 10.35 | 17.0    | ng    | 136708   | 2                     | 10.96 | 160629 | 20.0 |
| 4H-Cyclopenta[def... | 18.55 | 14.8    | ng    | 155146   | 4                     | 17.49 | 209096 | 20.0 |