

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061716\  
 Data File : BG022437.D  
 Acq On : 18 Jun 2016 23:58  
 Operator : UM/SJ  
 Sample : H3471-25  
 Misc :  
 ALS Vial : 46 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 STA-1160-(0-4)

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-BG060616.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         | 2.873       | 3             | 6           | 23           | rVB      | 851326         | 1298780       | 100.00%         | 9.797%        |
| 2         | 4.694       | 308           | 316         | 325          | rBV2     | 26501          | 51406         | 3.96%           | 0.388%        |
| 3         | 5.105       | 380           | 386         | 395          | rVB      | 22856          | 44109         | 3.40%           | 0.333%        |
| 4         | 5.505       | 449           | 454         | 461          | rVB2     | 8569           | 13632         | 1.05%           | 0.103%        |
| 5         | 5.605       | 464           | 471         | 489          | rBV      | 285005         | 633430        | 48.77%          | 4.778%        |
| 6         | 6.016       | 535           | 541         | 551          | rVB2     | 25555          | 49097         | 3.78%           | 0.370%        |
| 7         | 7.003       | 701           | 709         | 715          | rBV      | 21991          | 42000         | 3.23%           | 0.317%        |
| 8         | 7.109       | 721           | 727         | 732          | rBV      | 54296          | 97500         | 7.51%           | 0.735%        |
| 9         | 7.167       | 732           | 737         | 742          | rVV      | 41164          | 79964         | 6.16%           | 0.603%        |
| 10        | 7.232       | 742           | 748         | 761          | rVV      | 332692         | 680746        | 52.41%          | 5.135%        |
| 11        | 7.326       | 761           | 764         | 772          | rVB9     | 5823           | 14227         | 1.10%           | 0.107%        |
| 12        | 7.414       | 774           | 779         | 786          | rBV      | 25563          | 46707         | 3.60%           | 0.352%        |
| 13        | 7.579       | 800           | 807         | 814          | rBV      | 374353         | 709974        | 54.66%          | 5.355%        |
| 14        | 7.638       | 814           | 817         | 819          | rVV      | 38394          | 56344         | 4.34%           | 0.425%        |
| 15        | 7.679       | 819           | 824         | 831          | rVB      | 97554          | 180184        | 13.87%          | 1.359%        |
| 16        | 7.937       | 858           | 868         | 873          | rBV9     | 6682           | 20643         | 1.59%           | 0.156%        |
| 17        | 7.996       | 873           | 878         | 880          | rVV3     | 12082          | 22259         | 1.71%           | 0.168%        |
| 18        | 8.043       | 880           | 886         | 898          | rVV      | 110413         | 233348        | 17.97%          | 1.760%        |
| 19        | 8.149       | 898           | 904         | 912          | rVV2     | 49336          | 101589        | 7.82%           | 0.766%        |
| 20        | 8.366       | 934           | 941         | 953          | rVB      | 288234         | 567656        | 43.71%          | 4.282%        |
| 21        | 8.525       | 964           | 968         | 971          | rBV2     | 10600          | 18078         | 1.39%           | 0.136%        |
| 22        | 8.583       | 972           | 978         | 988          | rVB      | 51780          | 104547        | 8.05%           | 0.789%        |
| 23        | 8.677       | 988           | 994         | 1000         | rBV3     | 53642          | 109058        | 8.40%           | 0.823%        |
| 24        | 8.842       | 1017          | 1022        | 1033         | rVV2     | 18094          | 45886         | 3.53%           | 0.346%        |
| 25        | 8.995       | 1042          | 1048        | 1051         | rBV3     | 10058          | 16754         | 1.29%           | 0.126%        |
| 26        | 9.042       | 1051          | 1056        | 1062         | rVB2     | 15159          | 26133         | 2.01%           | 0.197%        |
| 27        | 9.142       | 1062          | 1073        | 1079         | rBV6     | 41182          | 98349         | 7.57%           | 0.742%        |
| 28        | 9.218       | 1079          | 1086        | 1103         | rVB      | 206882         | 501721        | 38.63%          | 3.785%        |
| 29        | 9.383       | 1105          | 1114        | 1123         | rBV      | 22471          | 65167         | 5.02%           | 0.492%        |
| 30        | 9.477       | 1124          | 1130        | 1137         | rBV3     | 11466          | 27323         | 2.10%           | 0.206%        |
| 31        | 9.670       | 1159          | 1163        | 1168         | rVB2     | 10877          | 16863         | 1.30%           | 0.127%        |
| 32        | 9.729       | 1168          | 1173        | 1185         | rVB2     | 19649          | 46076         | 3.55%           | 0.348%        |
| 33        | 10.041      | 1219          | 1226        | 1234         | rBV7     | 12613          | 35776         | 2.75%           | 0.270%        |
| 34        | 10.193      | 1244          | 1252        | 1256         | rBV6     | 9049           | 21051         | 1.62%           | 0.159%        |

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Integration Parameters: rteint.p  
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 Start Thrs: 0.2  
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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-BG060616.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |         |        |        |
|----|--------|------|------|------|------|--------|---------|--------|--------|
| 35 | 10.252 | 1256 | 1262 | 1269 | rVV2 | 32172  | 70789   | 5.45%  | 0.534% |
| 36 | 10.305 | 1269 | 1271 | 1276 | rVB3 | 8916   | 13223   | 1.02%  | 0.100% |
| 37 | 10.417 | 1283 | 1290 | 1292 | rBV4 | 7906   | 15568   | 1.20%  | 0.117% |
| 38 | 10.552 | 1306 | 1313 | 1320 | rVB2 | 14412  | 31389   | 2.42%  | 0.237% |
| 39 | 10.793 | 1346 | 1354 | 1355 | rBV6 | 10578  | 14397   | 1.11%  | 0.109% |
| 40 | 10.863 | 1359 | 1366 | 1371 | rVV  | 179036 | 360183  | 27.73% | 2.717% |
| 41 | 10.910 | 1371 | 1374 | 1382 | rVB  | 34762  | 63992   | 4.93%  | 0.483% |
| 42 | 11.856 | 1526 | 1535 | 1538 | rBV6 | 5787   | 15201   | 1.17%  | 0.115% |
| 43 | 12.244 | 1597 | 1601 | 1611 | rVB6 | 9721   | 19796   | 1.52%  | 0.149% |
| 44 | 12.514 | 1642 | 1647 | 1654 | rBV  | 14533  | 22406   | 1.73%  | 0.169% |
| 45 | 12.667 | 1666 | 1673 | 1679 | rBV2 | 12616  | 25402   | 1.96%  | 0.192% |
| 46 | 12.732 | 1679 | 1684 | 1690 | rVB2 | 10486  | 18633   | 1.43%  | 0.141% |
| 47 | 13.301 | 1773 | 1781 | 1795 | rBV  | 655113 | 1108419 | 85.34% | 8.361% |
| 48 | 13.484 | 1807 | 1812 | 1824 | rVB7 | 7563   | 18124   | 1.40%  | 0.137% |
| 49 | 14.001 | 1895 | 1900 | 1905 | rBV4 | 6971   | 13911   | 1.07%  | 0.105% |
| 50 | 14.130 | 1915 | 1922 | 1929 | rBV  | 71016  | 119122  | 9.17%  | 0.899% |
| 51 | 14.412 | 1964 | 1970 | 1977 | rBV  | 15649  | 33998   | 2.62%  | 0.256% |
| 52 | 14.547 | 1988 | 1993 | 1999 | rBV2 | 13159  | 18224   | 1.40%  | 0.137% |
| 53 | 14.682 | 2009 | 2016 | 2024 | rVB2 | 276131 | 475663  | 36.62% | 3.588% |
| 54 | 14.994 | 2064 | 2069 | 2071 | rBV3 | 9174   | 13631   | 1.05%  | 0.103% |
| 55 | 15.070 | 2079 | 2082 | 2092 | rVB6 | 7003   | 14020   | 1.08%  | 0.106% |
| 56 | 15.452 | 2141 | 2147 | 2151 | rBV7 | 7213   | 13109   | 1.01%  | 0.099% |
| 57 | 15.499 | 2151 | 2155 | 2160 | rVB  | 18921  | 26174   | 2.02%  | 0.197% |
| 58 | 15.893 | 2214 | 2222 | 2227 | rBV6 | 6339   | 14028   | 1.08%  | 0.106% |
| 59 | 16.175 | 2264 | 2270 | 2281 | rBV  | 354888 | 599455  | 46.16% | 4.522% |
| 60 | 16.357 | 2296 | 2301 | 2304 | rBV3 | 21160  | 39254   | 3.02%  | 0.296% |
| 61 | 16.498 | 2319 | 2325 | 2330 | rBV  | 11519  | 19957   | 1.54%  | 0.151% |
| 62 | 16.556 | 2330 | 2335 | 2341 | rVV3 | 8741   | 20560   | 1.58%  | 0.155% |
| 63 | 16.615 | 2341 | 2345 | 2350 | rVV3 | 9787   | 15883   | 1.22%  | 0.120% |
| 64 | 16.821 | 2374 | 2380 | 2384 | rBV6 | 10072  | 17722   | 1.36%  | 0.134% |
| 65 | 16.880 | 2387 | 2390 | 2395 | rVB  | 10711  | 15283   | 1.18%  | 0.115% |
| 66 | 17.085 | 2421 | 2425 | 2430 | rBV5 | 8959   | 13783   | 1.06%  | 0.104% |
| 67 | 17.144 | 2432 | 2435 | 2441 | rBV4 | 15860  | 26437   | 2.04%  | 0.199% |
| 68 | 17.426 | 2476 | 2483 | 2488 | rBV2 | 285065 | 484926  | 37.34% | 3.658% |
| 69 | 17.467 | 2488 | 2490 | 2497 | rVB  | 46693  | 65819   | 5.07%  | 0.496% |
| 70 | 17.561 | 2502 | 2506 | 2512 | rVV2 | 12226  | 20032   | 1.54%  | 0.151% |
| 71 | 17.884 | 2558 | 2561 | 2567 | rVB5 | 10891  | 15575   | 1.20%  | 0.117% |

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 STA-1160-(0-4)

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-BG060616.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |         |        |        |
|----|--------|------|------|------|------|--------|---------|--------|--------|
| 72 | 18.296 | 2626 | 2631 | 2635 | rBV2 | 15520  | 29374   | 2.26%  | 0.222% |
| 73 | 18.348 | 2637 | 2640 | 2647 | rVB3 | 20062  | 37884   | 2.92%  | 0.286% |
| 74 | 18.490 | 2659 | 2664 | 2669 | rBV4 | 18382  | 38835   | 2.99%  | 0.293% |
| 75 | 18.531 | 2669 | 2671 | 2675 | rVB2 | 11859  | 14108   | 1.09%  | 0.106% |
| 76 | 18.578 | 2675 | 2679 | 2682 | rBV6 | 8374   | 14814   | 1.14%  | 0.112% |
| 77 | 18.789 | 2711 | 2715 | 2721 | rBV4 | 11403  | 20371   | 1.57%  | 0.154% |
| 78 | 18.848 | 2721 | 2725 | 2731 | rVB5 | 12627  | 20525   | 1.58%  | 0.155% |
| 79 | 19.200 | 2781 | 2785 | 2791 | rBV3 | 25272  | 40955   | 3.15%  | 0.309% |
| 80 | 19.353 | 2805 | 2811 | 2814 | rBV6 | 15225  | 30298   | 2.33%  | 0.229% |
| 81 | 19.471 | 2825 | 2831 | 2837 | rBV  | 159798 | 244163  | 18.80% | 1.842% |
| 82 | 19.700 | 2865 | 2870 | 2875 | rBV2 | 36029  | 46584   | 3.59%  | 0.351% |
| 83 | 19.835 | 2885 | 2893 | 2900 | rVB  | 150898 | 259366  | 19.97% | 1.956% |
| 84 | 20.023 | 2919 | 2925 | 2931 | rBV  | 736853 | 1033929 | 79.61% | 7.799% |
| 85 | 20.734 | 3043 | 3046 | 3051 | rVB4 | 29672  | 40681   | 3.13%  | 0.307% |
| 86 | 21.086 | 3103 | 3106 | 3114 | rVB  | 27320  | 42817   | 3.30%  | 0.323% |
| 87 | 21.298 | 3140 | 3142 | 3149 | rVB5 | 43749  | 70485   | 5.43%  | 0.532% |
| 88 | 21.692 | 3201 | 3209 | 3214 | rBV3 | 272451 | 615663  | 47.40% | 4.644% |
| 89 | 23.901 | 3579 | 3585 | 3602 | rVB  | 85809  | 349444  | 26.91% | 2.636% |
| 90 | 24.935 | 3753 | 3761 | 3776 | rBV2 | 111782 | 336207  | 25.89% | 2.536% |

Sum of corrected areas: 13256968



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ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
STA-1160-(0-4)

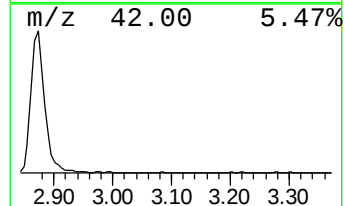
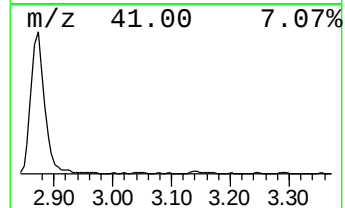
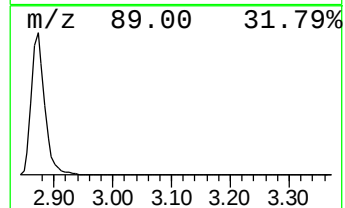
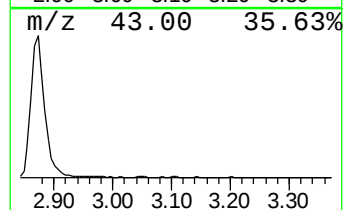
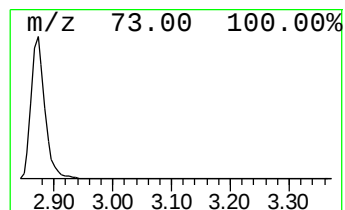
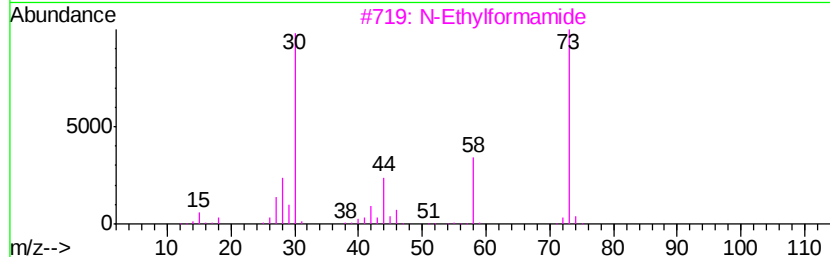
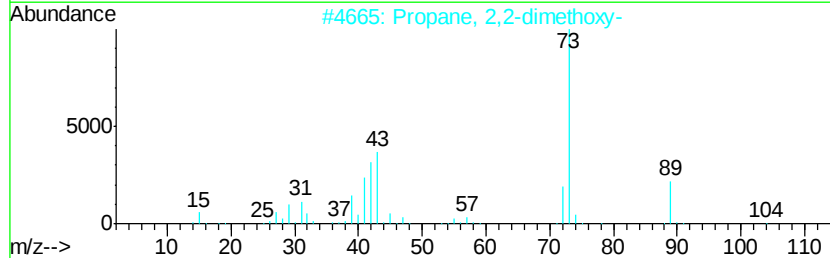
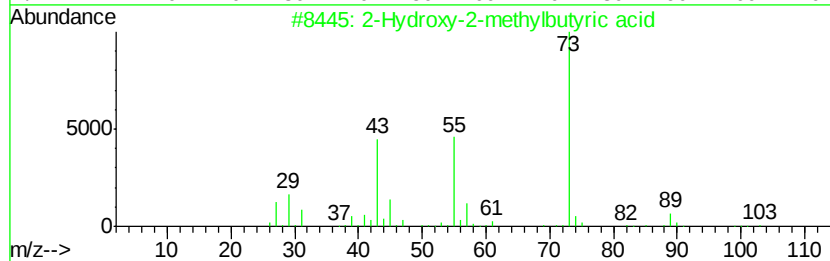
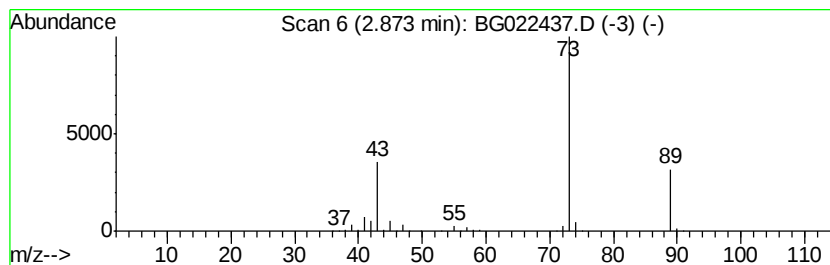
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 unknown2.87 Concentration Rank 1

| R.T. | EstConc   | Area    | Relative to ISTD       | R.T. |
|------|-----------|---------|------------------------|------|
| 2.87 | 111.32 ng | 1298780 | 1,4-Dichlorobenzene-d4 | 8.04 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Hydroxy-2-methylbutyric acid | 118 | C5H10O3 | 003739-30-8 | 28   |
| 2    |    | Propane, 2,2-dimethoxy-        | 104 | C5H12O2 | 000077-76-9 | 9    |
| 3    |    | N-Ethylformamide               | 73  | C3H7NO  | 000627-45-2 | 7    |
| 4    |    | Methane, isothiocyanato-       | 73  | C2H3NS  | 000556-61-6 | 7    |
| 5    |    | 1,3,5-Trioxane                 | 90  | C3H6O3  | 000110-88-3 | 4    |



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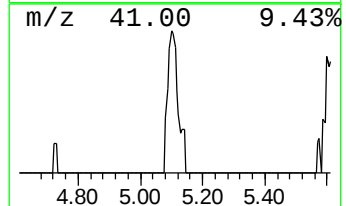
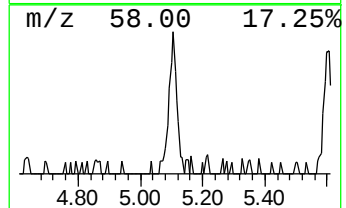
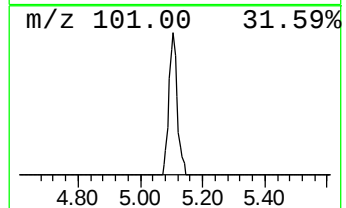
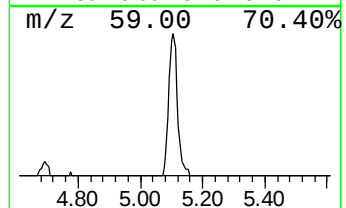
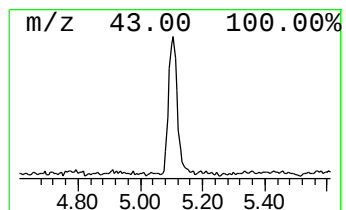
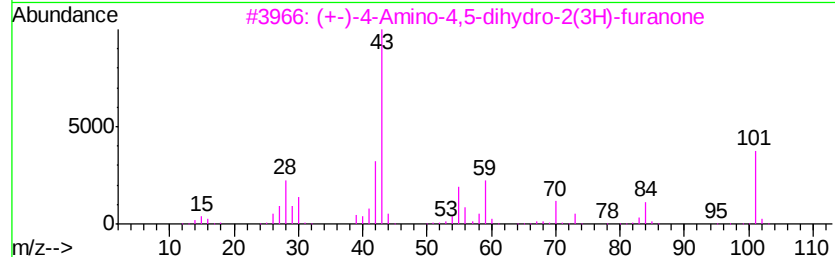
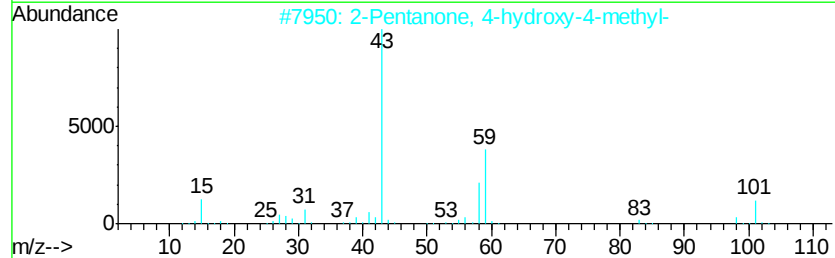
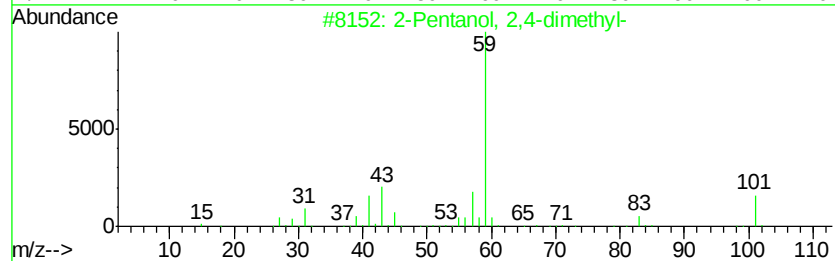
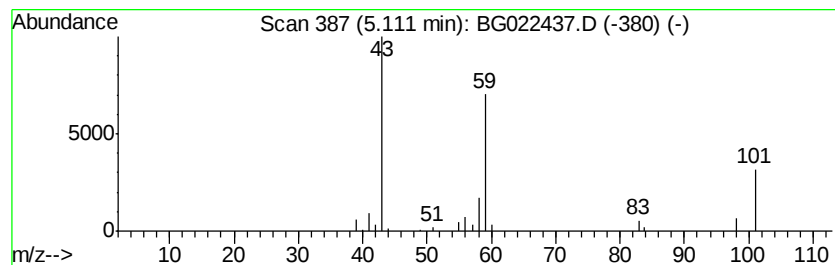
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 2-Pentanol, 2,4-dimethyl- Concentration Rank 18

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 5.11 | 3.78 ng | 44109 | 1,4-Dichlorobenzene-d4 | 8.04 |

| Hit# | of | 5 | Tentative ID                           | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 2-Pentanol, 2,4-dimethyl-              | 116 | C7H16O  | 000625-06-9 | 50   |
| 2    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-       | 116 | C6H12O2 | 000123-42-2 | 45   |
| 3    |    |   | (+)-4-Amino-4,5-dihydro-2(3H)-furanone | 101 | C4H7NO2 | 016504-58-8 | 43   |
| 4    |    |   | Propane, 1-ethoxy-2-methyl-            | 102 | C6H14O  | 000627-02-1 | 25   |
| 5    |    |   | Butane, 1-ethoxy-                      | 102 | C6H14O  | 000628-81-9 | 17   |



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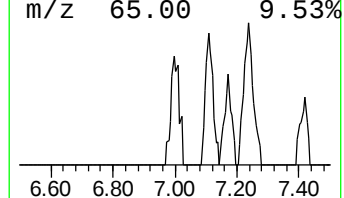
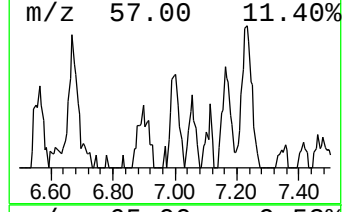
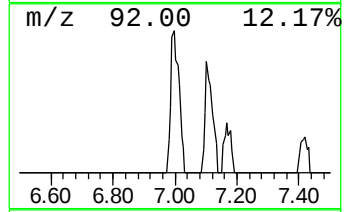
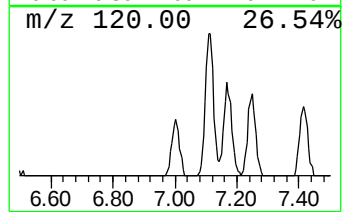
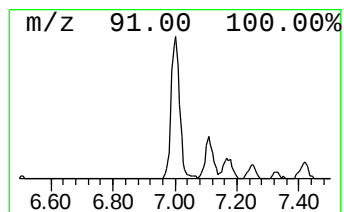
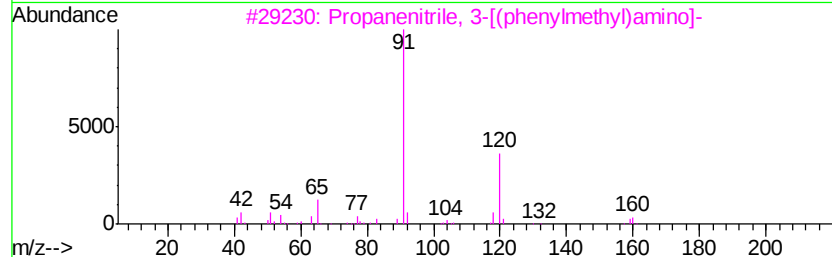
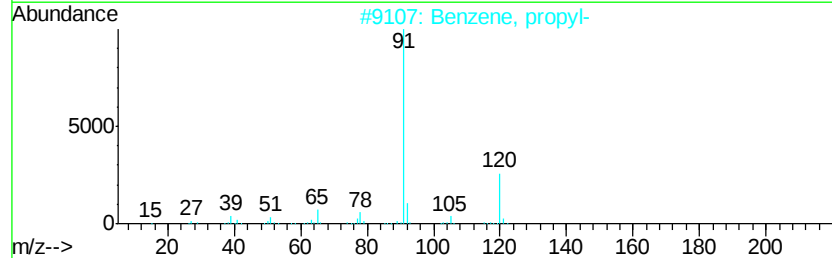
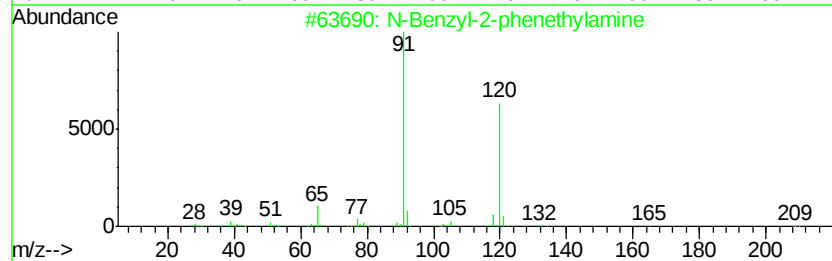
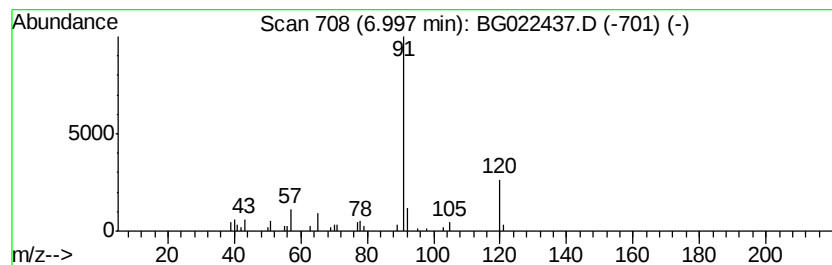
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 5 N-Benzyl-2-phenethylamine Concentration Rank 19

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 7.00 | 3.60 ng | 42000 | 1,4-Dichlorobenzene-d4 | 8.04 |

| Hit# | of | Tentative ID                             | MW  | MolForm   | CAS#        | Qual |
|------|----|------------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | N-Benzyl-2-phenethylamine                | 211 | C15H17N   | 003647-71-0 | 59   |
| 2    |    | Benzene, propyl-                         | 120 | C9H12     | 000103-65-1 | 49   |
| 3    |    | Propanenitrile, 3-[(phenylmethyl)amino]- | 160 | C10H12N2  | 000706-03-6 | 42   |
| 4    |    | 2,4-Difluorobenzene, 1-benzoyloxy-       | 220 | C13H10F2O | 152434-86-1 | 38   |
| 5    |    | 1,3,5-Cycloheptatriene, 7-ethyl-         | 120 | C9H12     | 017634-51-4 | 27   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061716\  
Data File : BG022437.D  
Acq On : 18 Jun 2016 23:58  
Operator : UM/SJ  
Sample : H3471-25  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
STA-1160-(0-4)

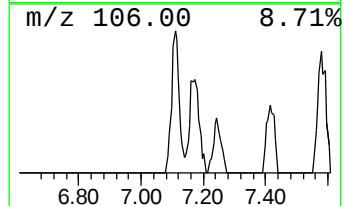
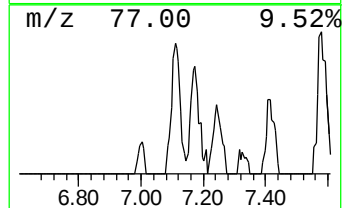
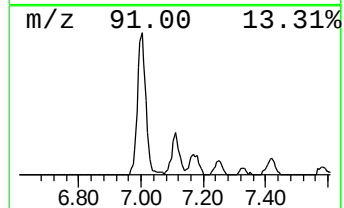
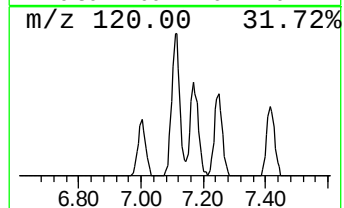
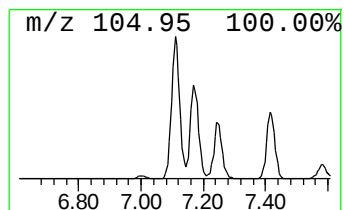
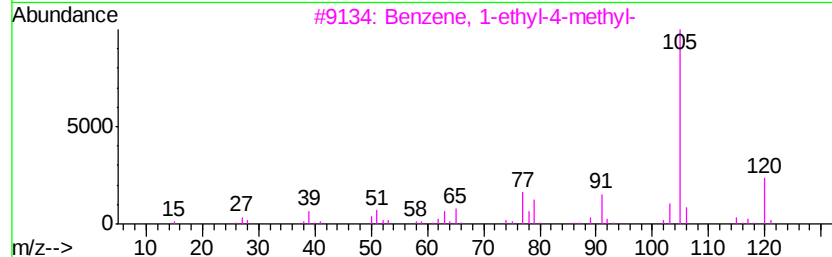
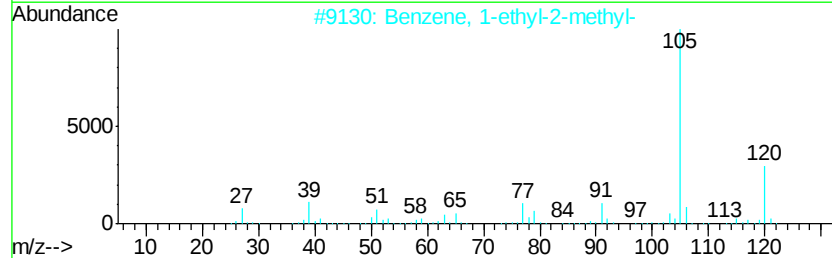
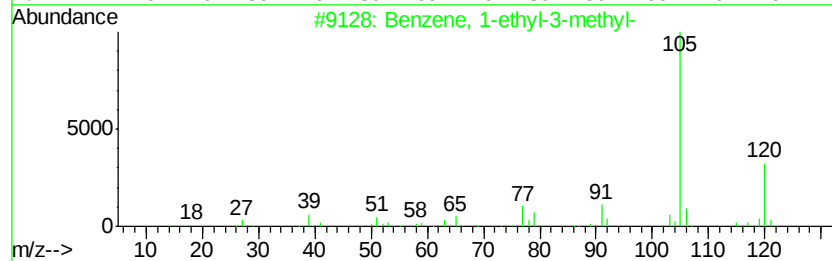
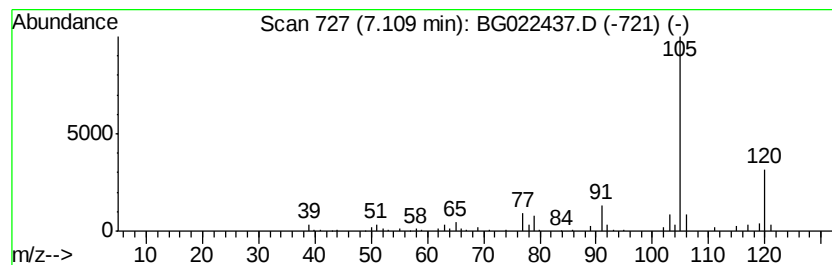
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 Benzene, 1-ethyl-3-methyl- Concentration Rank 9

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 7.11 | 8.36 ng | 97500 | 1,4-Dichlorobenzene-d4 | 8.04 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 97   |
| 2    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 95   |
| 3    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 93   |
| 4    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 90   |
| 5    |    |   | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 90   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061716\  
Data File : BG022437.D  
Acq On : 18 Jun 2016 23:58  
Operator : UM/SJ  
Sample : H3471-25  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
STA-1160-(0-4)

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

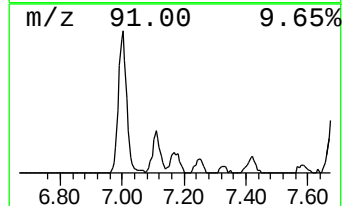
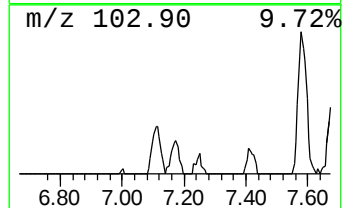
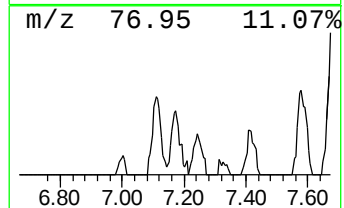
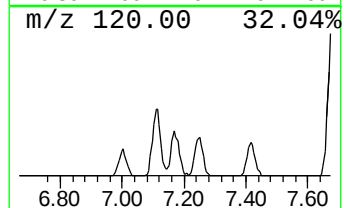
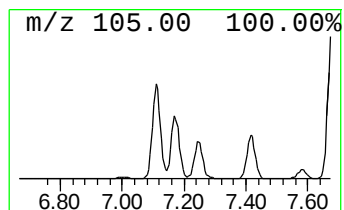
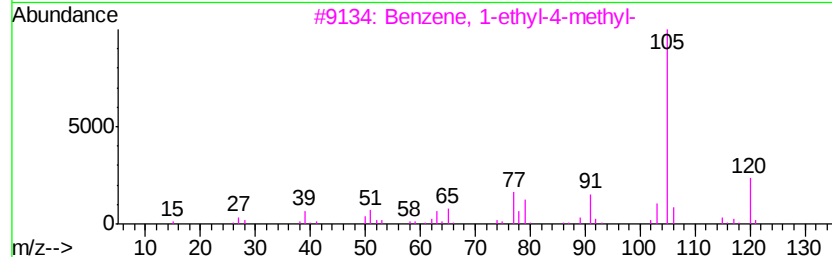
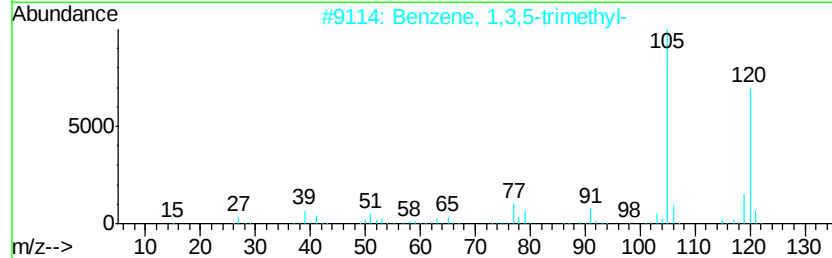
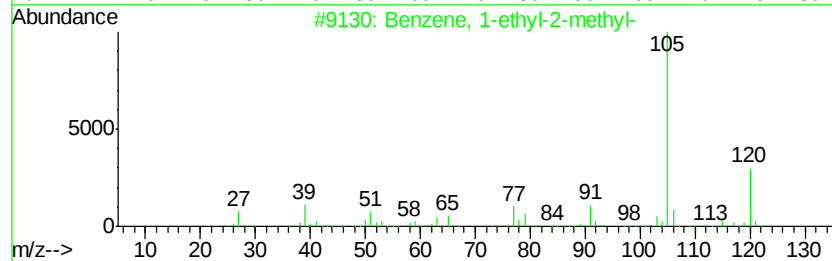
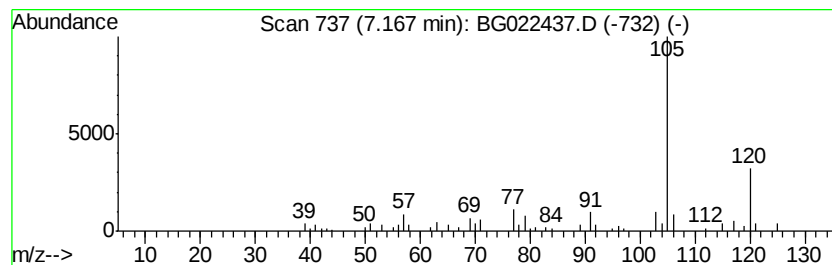
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 7.17 | 6.85 ng | 79964 | 1,4-Dichlorobenzene-d4 | 8.04 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 90   |
| 2    |    |   | Benzene, 1,3,5-trimethyl-  | 120 | C9H12   | 000108-67-8 | 90   |
| 3    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 90   |
| 4    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 87   |
| 5    |    |   | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 86   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061716\  
Data File : BG022437.D  
Acq On : 18 Jun 2016 23:58  
Operator : UM/SJ  
Sample : H3471-25  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
STA-1160-(0-4)

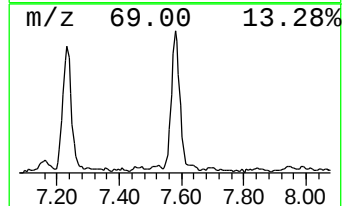
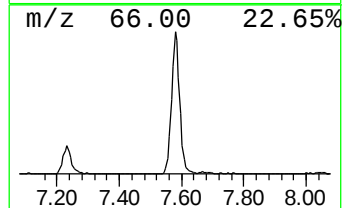
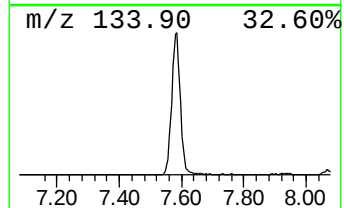
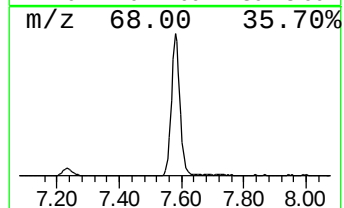
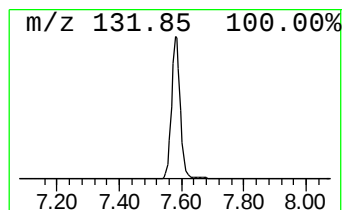
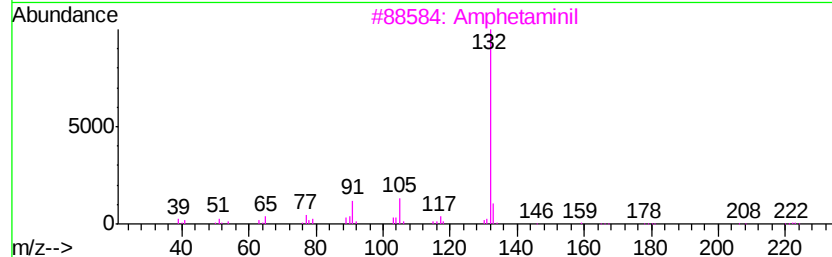
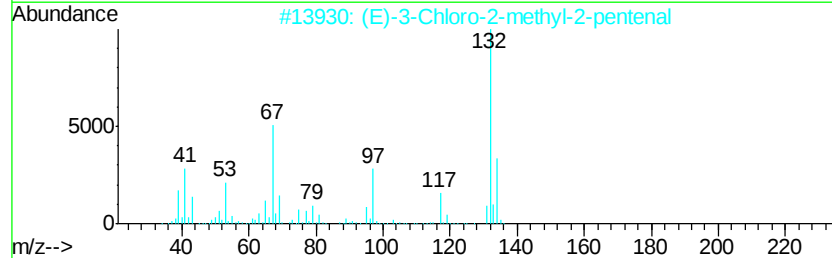
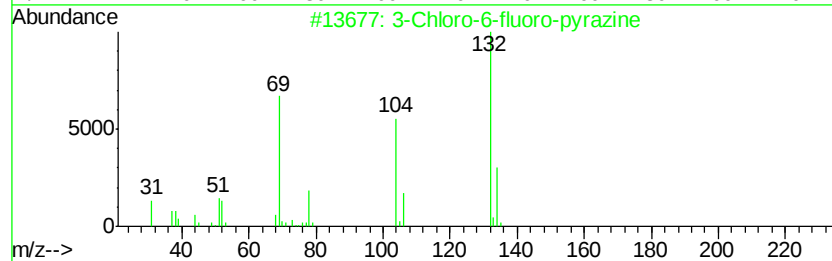
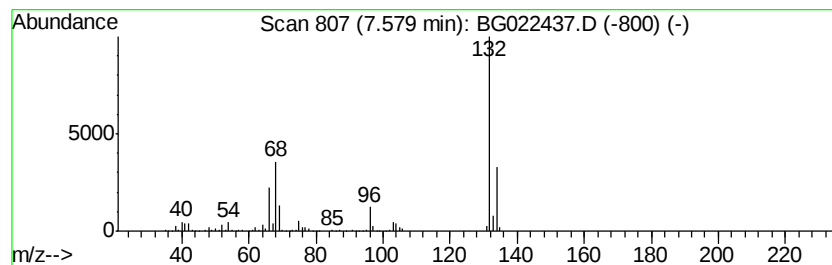
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 9 unknown7.58 Concentration Rank 2

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 7.58 | 60.85 ng | 709974 | 1,4-Dichlorobenzene-d4 | 8.04 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 38   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 17   |
| 3    |    | Amphetaminil                     | 250 | C17H18N2  | 017590-01-1  | 12   |
| 4    |    | Tranylcypromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 12   |
| 5    |    | Benzene, 1,3,5-trifluoro-        | 132 | C6H3F3    | 000372-38-3  | 9    |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061716\  
Data File : BG022437.D  
Acq On : 18 Jun 2016 23:58  
Operator : UM/SJ  
Sample : H3471-25  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
STA-1160-(0-4)

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

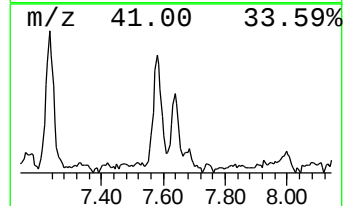
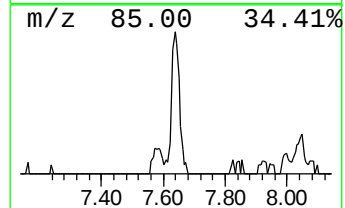
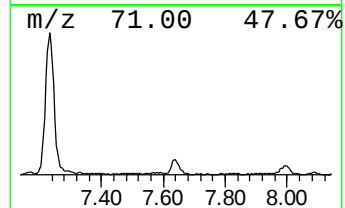
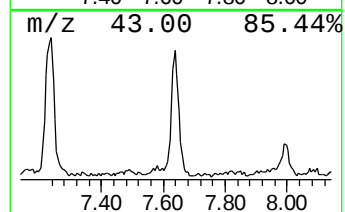
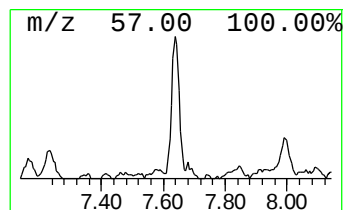
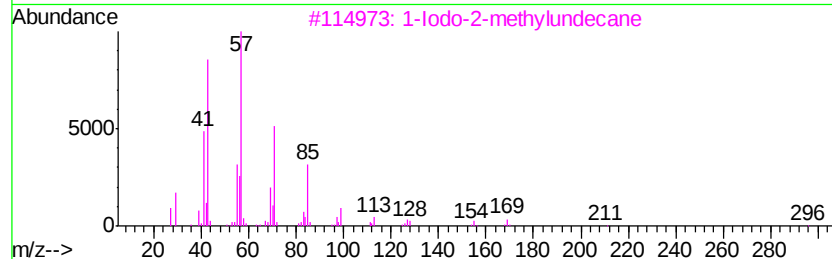
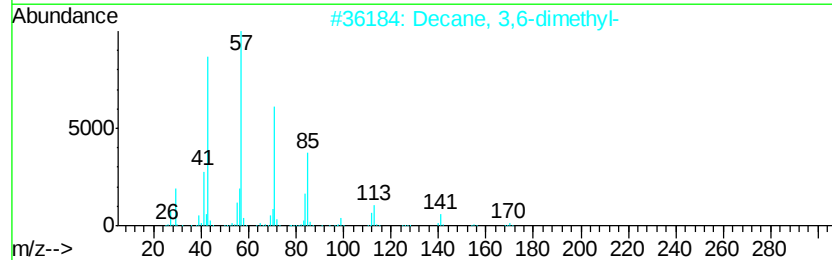
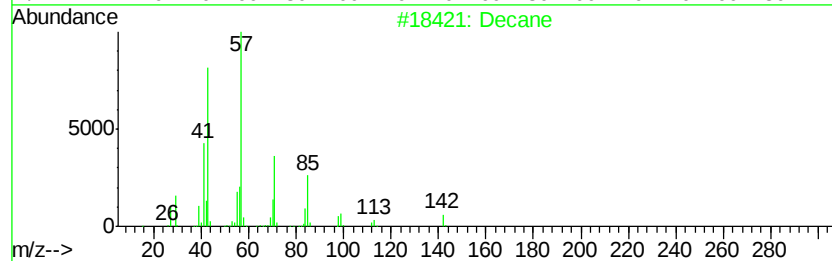
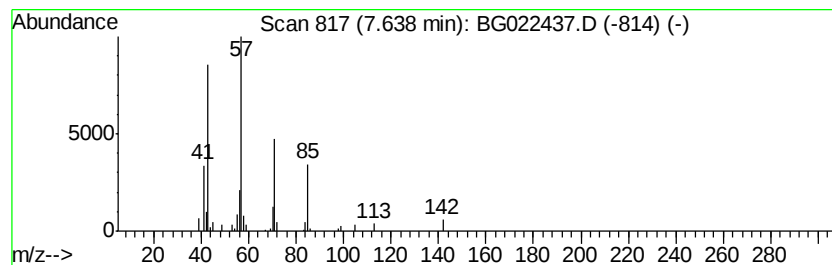
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Decane Concentration Rank 12

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 7.64 | 4.83 ng | 56344 | 1,4-Dichlorobenzene-d4 | 8.04 |

| Hit# | of                          | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|-----------------------------|---|--------------|-----|---------|-------------|------|
| 1    | Decane                      |   |              | 142 | C10H22  | 000124-18-5 | 87   |
| 2    | Decane, 3,6-dimethyl-       |   |              | 170 | C12H26  | 017312-53-7 | 78   |
| 3    | 1-Iodo-2-methylundecane     |   |              | 296 | C12H25I | 073105-67-6 | 78   |
| 4    | Dodecane, 2,6,11-trimethyl- |   |              | 212 | C15H32  | 031295-56-4 | 64   |
| 5    | Dotriacontane               |   |              | 451 | C32H66  | 000544-85-4 | 64   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061716\  
Data File : BG022437.D  
Acq On : 18 Jun 2016 23:58  
Operator : UM/SJ  
Sample : H3471-25  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
STA-1160-(0-4)

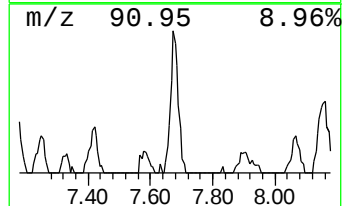
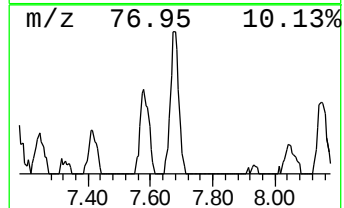
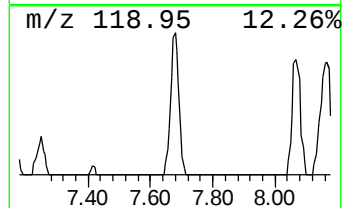
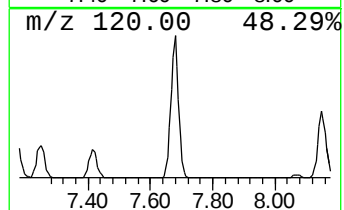
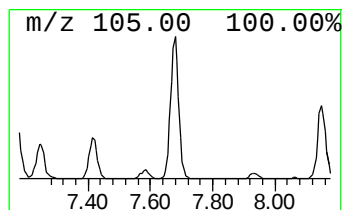
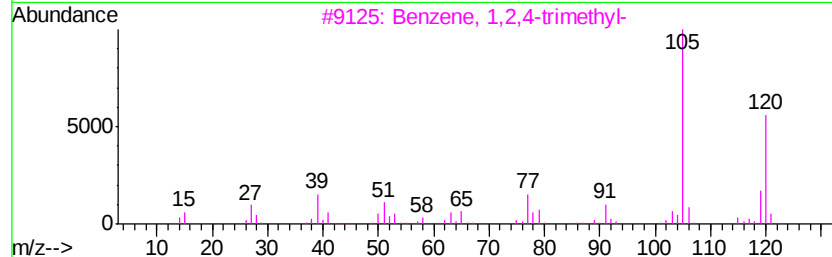
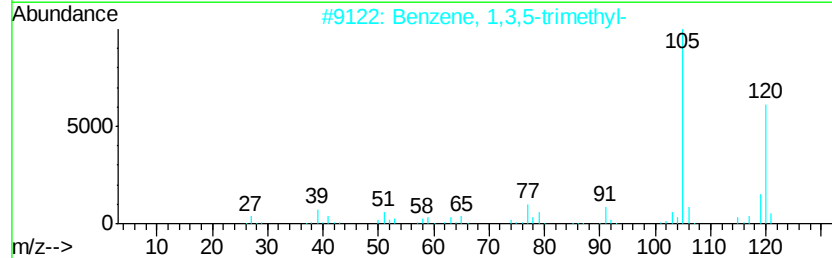
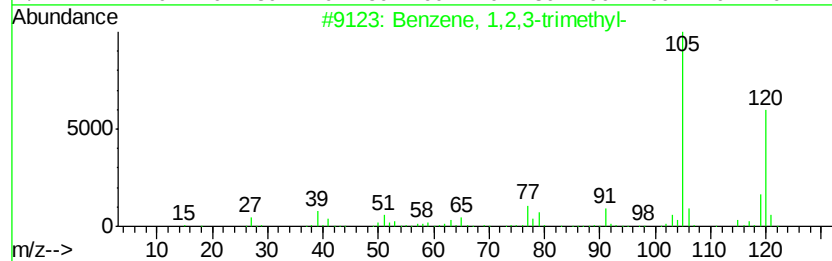
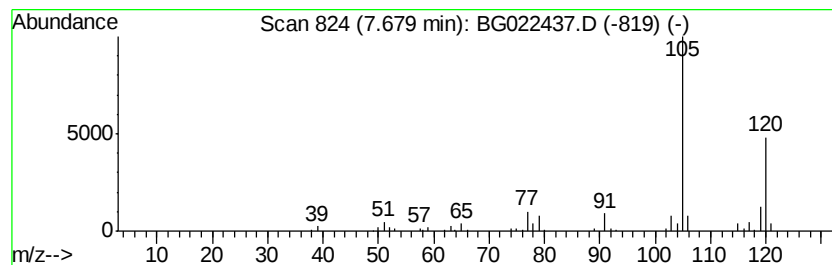
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 11 Benzene, 1,2,3-trimethyl- Concentration Rank 4

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 7.68 | 15.44 ng | 180184 | 1,4-Dichlorobenzene-d4 | 8.04 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3-trimethyl-          | 120 | C9H12   | 000526-73-8 | 97   |
| 2    |    | Benzene, 1,3,5-trimethyl-          | 120 | C9H12   | 000108-67-8 | 93   |
| 3    |    | Benzene, 1,2,4-trimethyl-          | 120 | C9H12   | 000095-63-6 | 91   |
| 4    |    | 2,3-Heptadien-5-yne, 2,4-dimethyl- | 120 | C9H12   | 041898-89-9 | 87   |
| 5    |    | Benzene, 1-ethyl-4-methyl-         | 120 | C9H12   | 000622-96-8 | 87   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061716\  
Data File : BG022437.D  
Acq On : 18 Jun 2016 23:58  
Operator : UM/SJ  
Sample : H3471-25  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
STA-1160-(0-4)

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

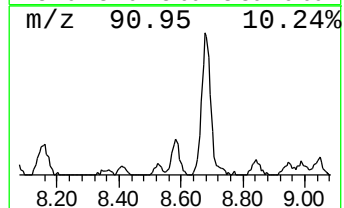
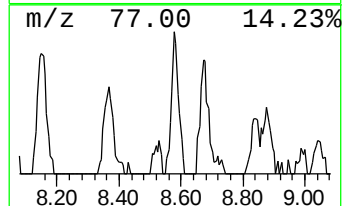
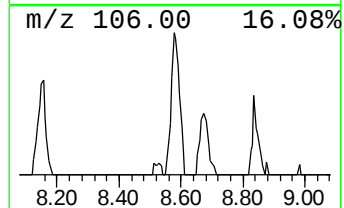
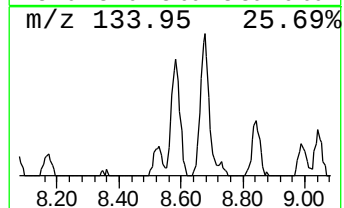
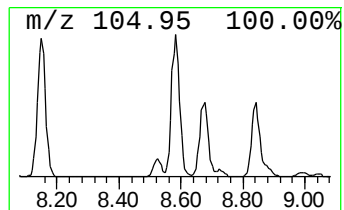
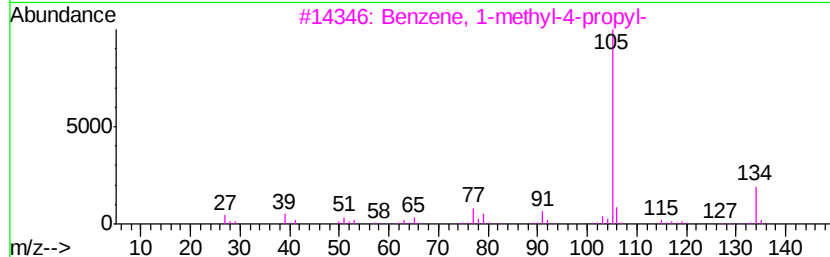
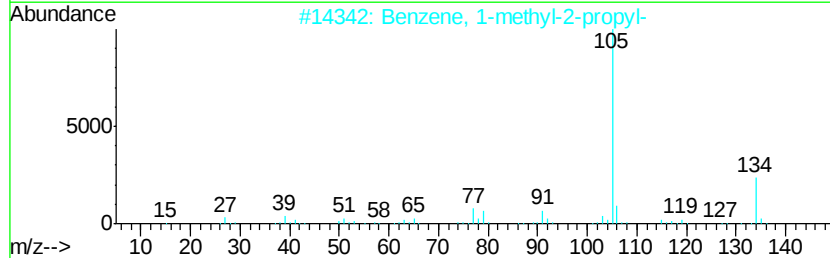
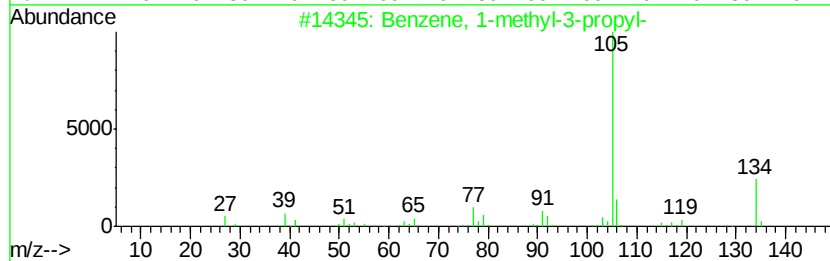
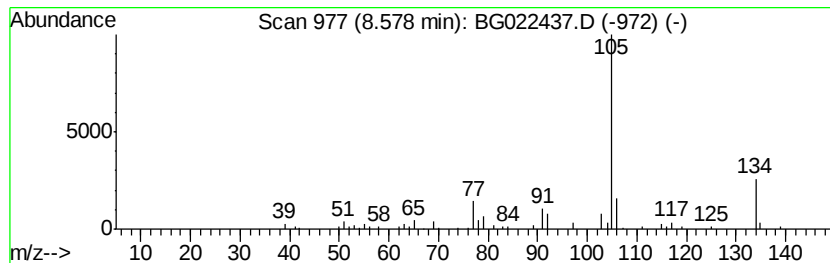
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Benzene, 1-methyl-3-propyl- Concentration Rank 6

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 8.58 | 8.96 ng | 104547 | 1,4-Dichlorobenzene-d4 | 8.04 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-methyl-3-propyl-         | 134 | C10H14  | 001074-43-7 | 91   |
| 2    |    |   | Benzene, 1-methyl-2-propyl-         | 134 | C10H14  | 001074-17-5 | 91   |
| 3    |    |   | Benzene, 1-methyl-4-propyl-         | 134 | C10H14  | 001074-55-1 | 91   |
| 4    |    |   | Benzene, (1-methylpropyl)-          | 134 | C10H14  | 000135-98-8 | 86   |
| 5    |    |   | Benzeneacetaldehyde, .alpha.-met... | 134 | C9H10O  | 000093-53-8 | 72   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061716\  
Data File : BG022437.D  
Acq On : 18 Jun 2016 23:58  
Operator : UM/SJ  
Sample : H3471-25  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
STA-1160-(0-4)

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

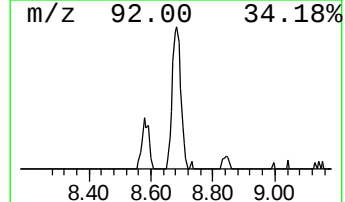
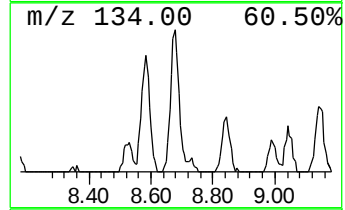
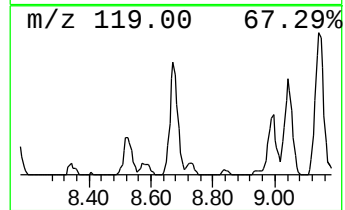
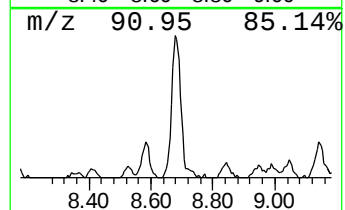
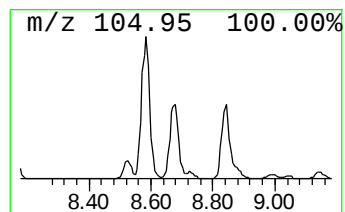
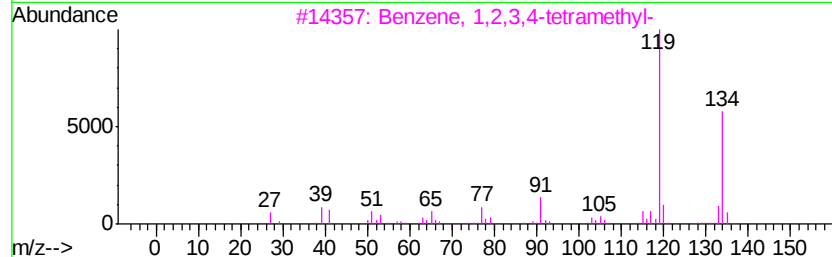
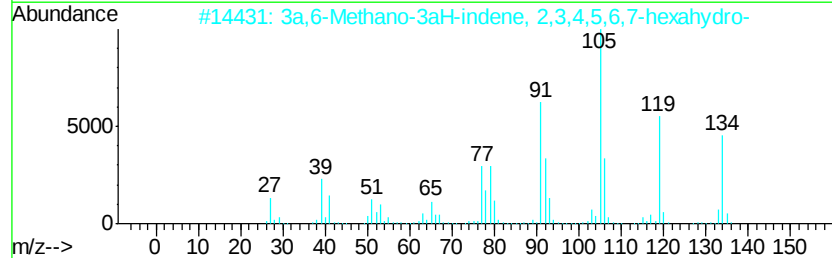
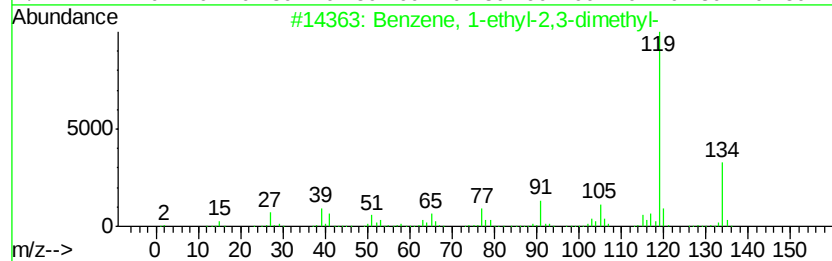
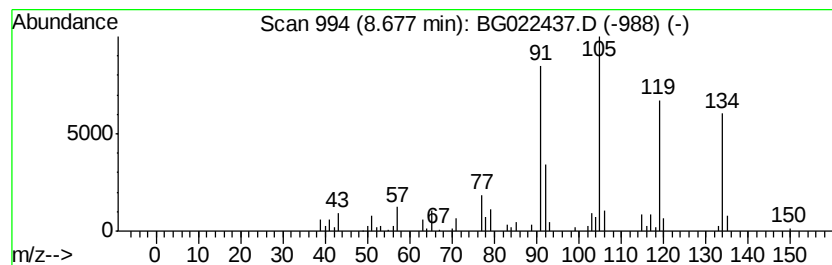
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 5

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 8.68 | 9.35 ng | 109058 | 1,4-Dichlorobenzene-d4 | 8.04 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 87   |
| 2    |    | 3a,6-Methano-3aH-indene, 2,3,4,5... | 134 | C10H14  | 098640-10-9 | 81   |
| 3    |    | Benzene, 1,2,3,4-tetramethyl-       | 134 | C10H14  | 000488-23-3 | 81   |
| 4    |    | Benzene, 2-ethyl-1,3-dimethyl-      | 134 | C10H14  | 002870-04-4 | 81   |
| 5    |    | Benzene, 1,2,4,5-tetramethyl-       | 134 | C10H14  | 000095-93-2 | 81   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061716\  
Data File : BG022437.D  
Acq On : 18 Jun 2016 23:58  
Operator : UM/SJ  
Sample : H3471-25  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
STA-1160-(0-4)

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

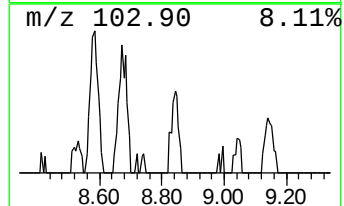
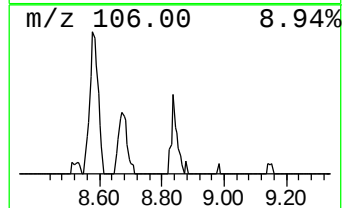
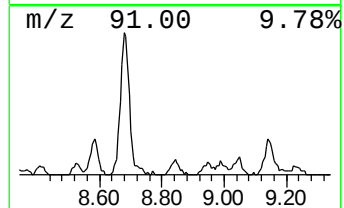
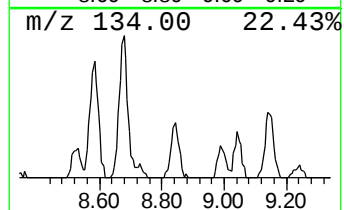
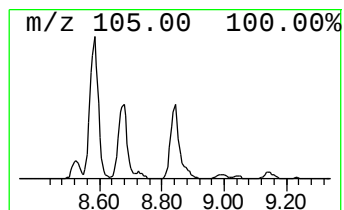
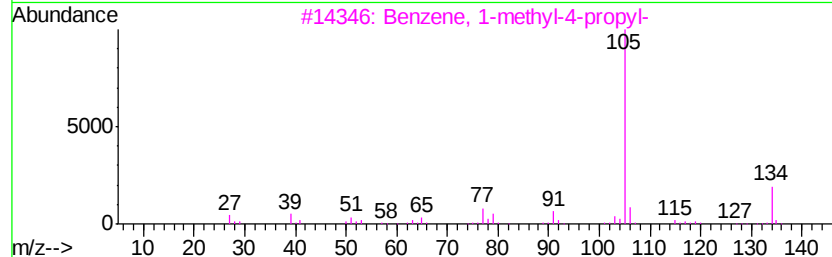
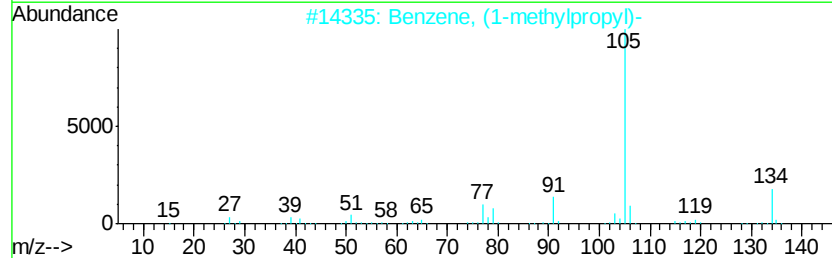
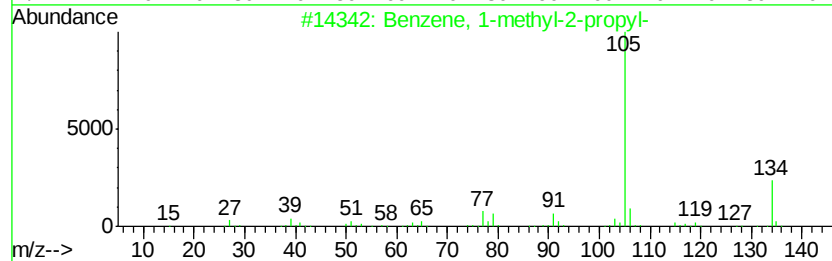
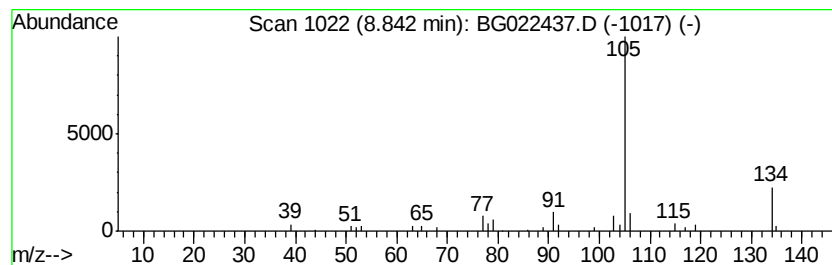
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Benzene, 1-methyl-2-propyl- Concentration Rank 16

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 8.84 | 3.93 ng | 45886 | 1,4-Dichlorobenzene-d4 | 8.04 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-methyl-2-propyl-         | 134 | C10H14  | 001074-17-5 | 91   |
| 2    |    |   | Benzene, (1-methylpropyl)-          | 134 | C10H14  | 000135-98-8 | 91   |
| 3    |    |   | Benzene, 1-methyl-4-propyl-         | 134 | C10H14  | 001074-55-1 | 91   |
| 4    |    |   | 2-Tolyloxirane                      | 134 | C9H10O  | 002783-26-8 | 68   |
| 5    |    |   | Benzeneacetaldehyde, .alpha.-met... | 134 | C9H10O  | 000093-53-8 | 64   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061716\  
Data File : BG022437.D  
Acq On : 18 Jun 2016 23:58  
Operator : UM/SJ  
Sample : H3471-25  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
STA-1160-(0-4)

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

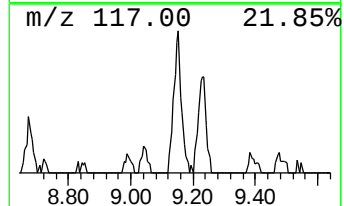
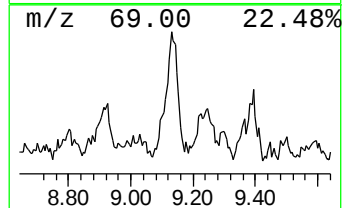
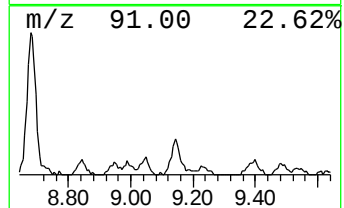
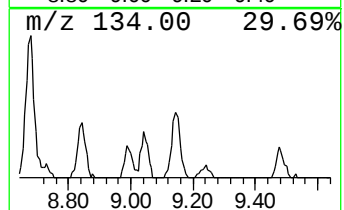
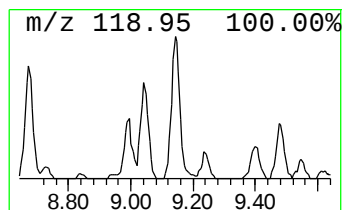
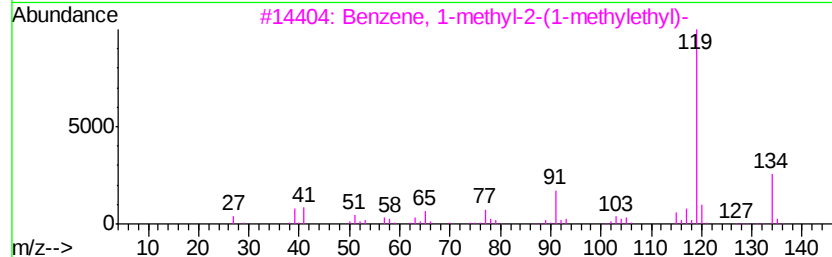
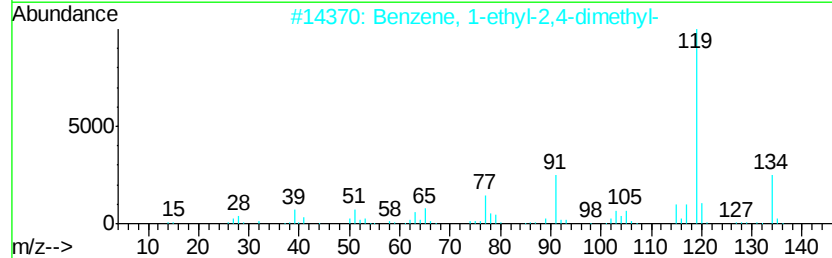
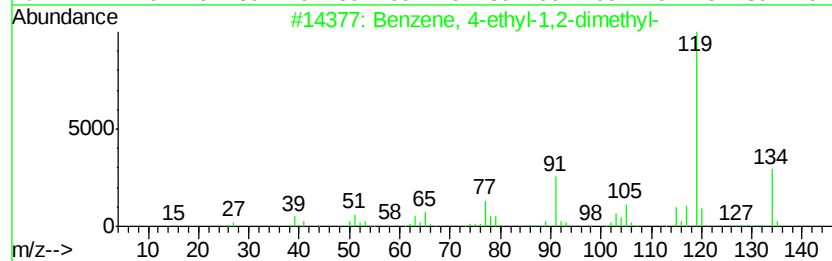
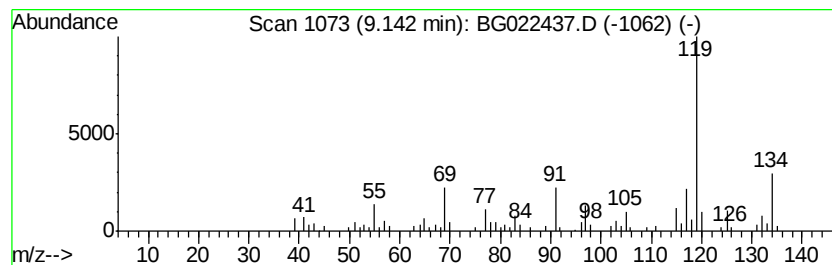
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 8

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 9.14 | 8.43 ng | 98349 | 1,4-Dichlorobenzene-d4 | 8.04 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 4-ethyl-1,2-dimethyl-       | 134 | C10H14  | 000934-80-5 | 92   |
| 2    |    |   | Benzene, 1-ethyl-2,4-dimethyl-       | 134 | C10H14  | 000874-41-9 | 70   |
| 3    |    |   | Benzene, 1-methyl-2-(1-methylethyl)- | 134 | C10H14  | 000527-84-4 | 70   |
| 4    |    |   | Benzene, 1-ethyl-2,3-dimethyl-       | 134 | C10H14  | 000933-98-2 | 70   |
| 5    |    |   | Benzene, 2-ethyl-1,4-dimethyl-       | 134 | C10H14  | 001758-88-9 | 70   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061716\  
Data File : BG022437.D  
Acq On : 18 Jun 2016 23:58  
Operator : UM/SJ  
Sample : H3471-25  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
STA-1160-(0-4)

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

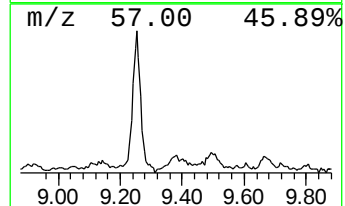
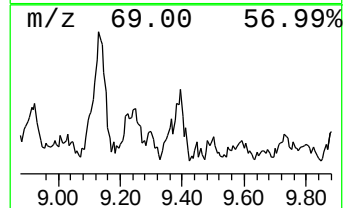
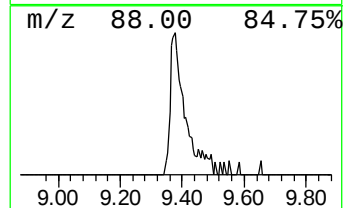
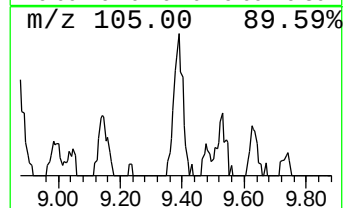
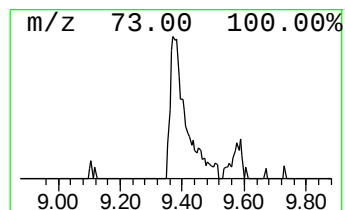
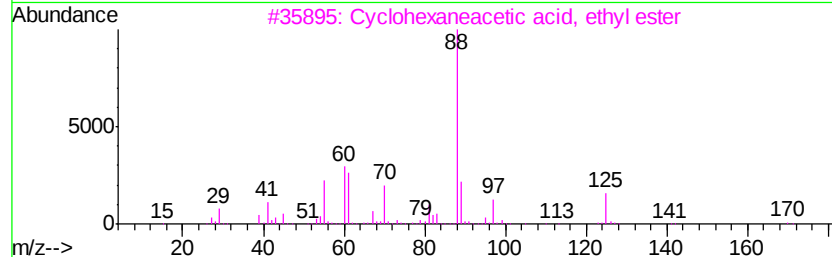
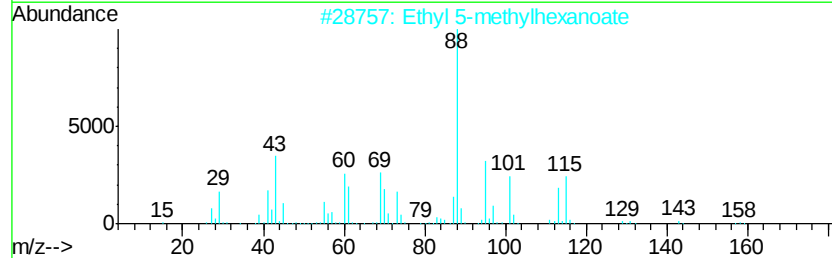
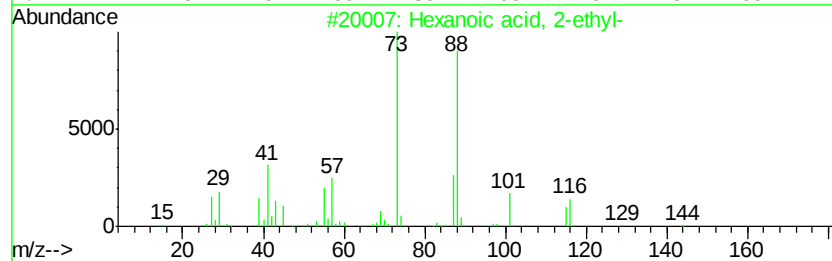
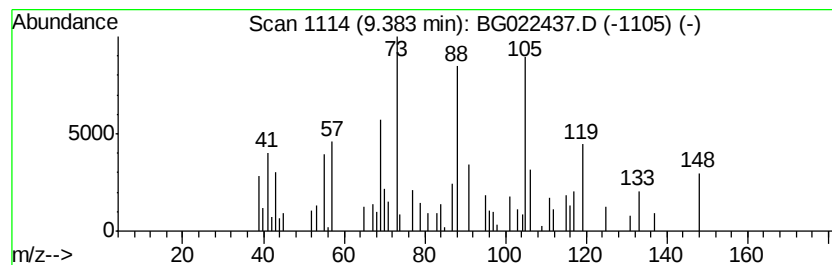
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 unknown9.38 Concentration Rank 11

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 9.38 | 5.59 ng | 65167 | 1,4-Dichlorobenzene-d4 | 8.04 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Hexanoic acid, 2-ethyl-             | 144 | C8H16O2  | 000149-57-5 | 35   |
| 2    |    | Ethyl 5-methylhexanoate             | 158 | C9H18O2  | 010236-10-9 | 12   |
| 3    |    | Cyclohexaneacetic acid, ethyl ester | 170 | C10H18O2 | 005452-75-5 | 12   |
| 4    |    | Pentanoic acid, 3-methyl-, ethyl... | 144 | C8H16O2  | 005870-68-8 | 10   |
| 5    |    | Heptanoic acid, 2,4-dimethyl-, m... | 172 | C10H20O2 | 018524-86-2 | 10   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061716\  
Data File : BG022437.D  
Acq On : 18 Jun 2016 23:58  
Operator : UM/SJ  
Sample : H3471-25  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
STA-1160-(0-4)

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

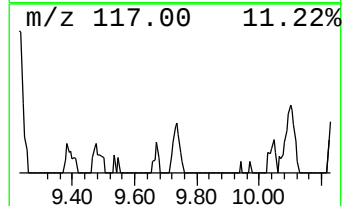
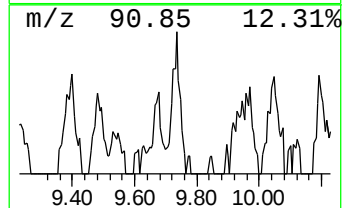
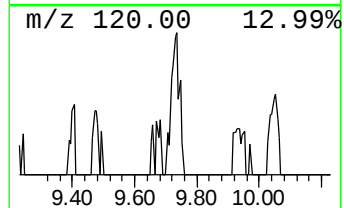
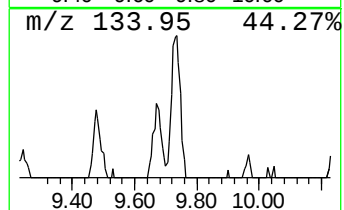
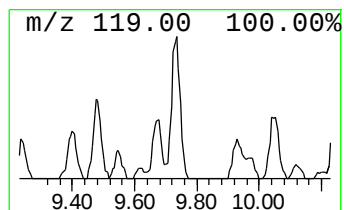
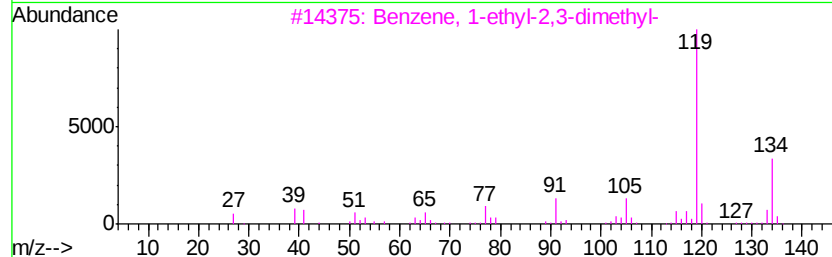
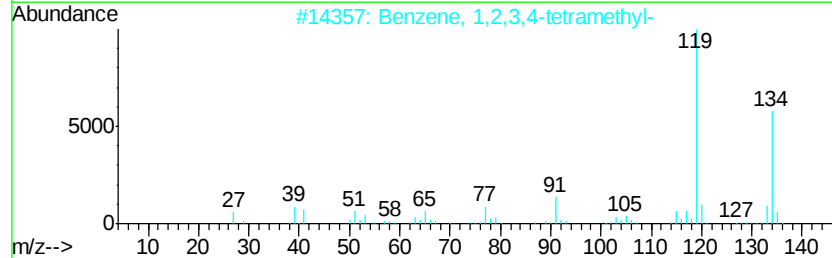
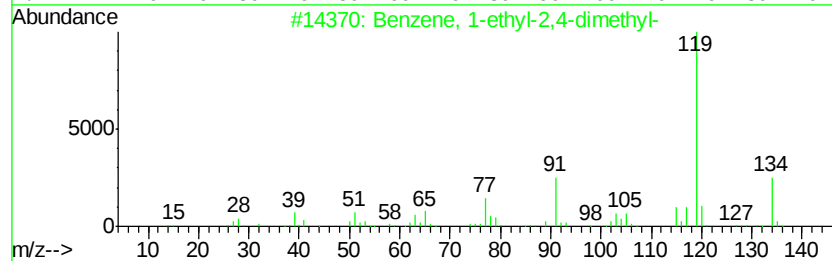
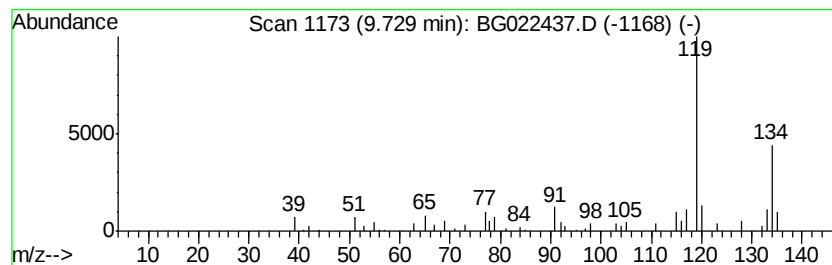
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 20

| R.T. | EstConc | Area  | Relative to ISTD | R.T.  |
|------|---------|-------|------------------|-------|
| 9.73 | 2.56 ng | 46076 | Naphthalene-d8   | 10.86 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-2,4-dimethyl- | 134 | C10H14  | 000874-41-9 | 93   |
| 2    |    |   | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 91   |
| 3    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 91   |
| 4    |    |   | Benzene, 1,2,4,5-tetramethyl-  | 134 | C10H14  | 000095-93-2 | 91   |
| 5    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 90   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061716\  
Data File : BG022437.D  
Acq On : 18 Jun 2016 23:58  
Operator : UM/SJ  
Sample : H3471-25  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
STA-1160-(0-4)

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

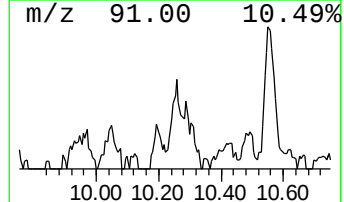
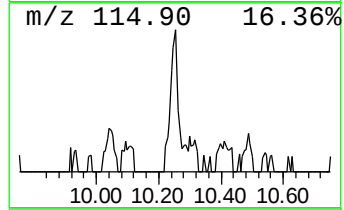
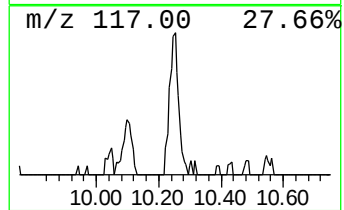
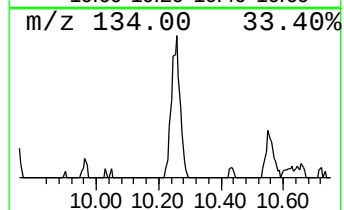
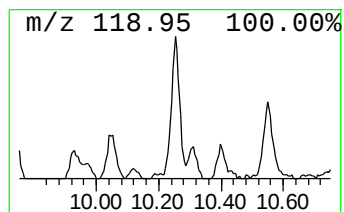
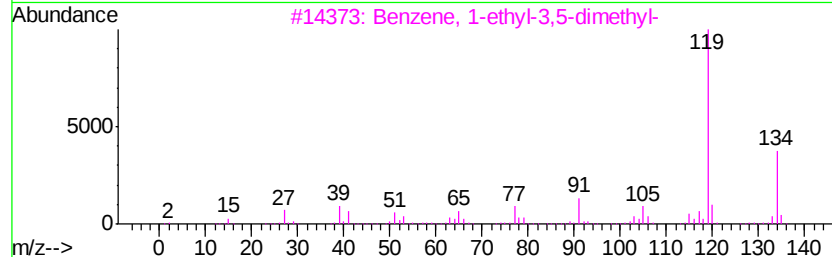
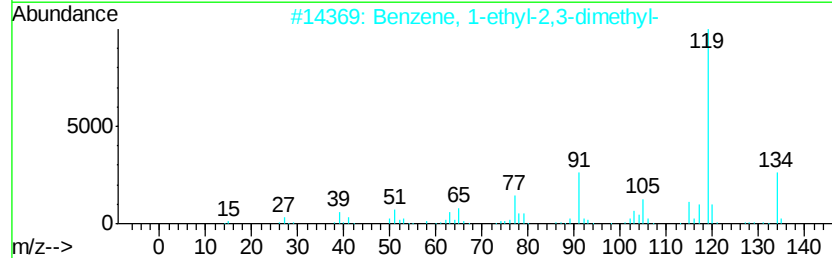
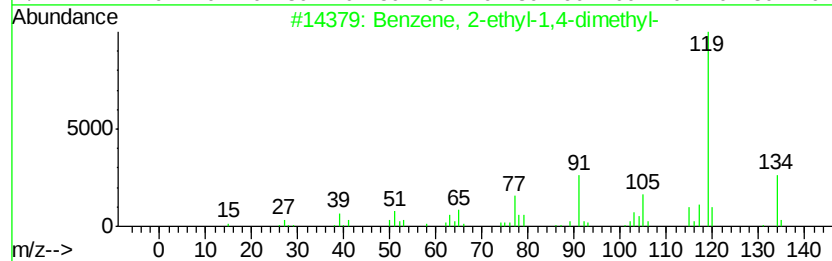
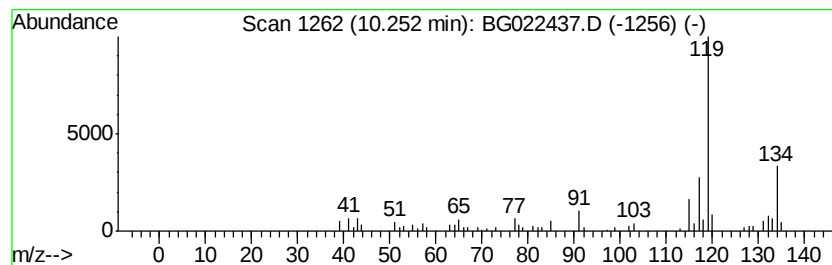
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 17

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 10.25 | 3.93 ng | 70789 | Naphthalene-d8   | 10.86 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 76   |
| 2    |    | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 76   |
| 3    |    | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 72   |
| 4    |    | Benzene, 1,2,3,5-tetramethyl-  | 134 | C10H14  | 000527-53-7 | 72   |
| 5    |    | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 72   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG061716\  
Data File : BG022437.D  
Acq On : 18 Jun 2016 23:58  
Operator : UM/SJ  
Sample : H3471-25  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
STA-1160-(0-4)

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060616.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 |
| unknown2.87          | 2.87  | 111.3   | ng    | 1298780  | 1                     | 8.04  | 233348 | 20.0 |
| 2-Pentanol, 2,4-d... | 5.11  | 3.8     | ng    | 44109    | 1                     | 8.04  | 233348 | 20.0 |
| N-Benzyl-2-phenet... | 7.00  | 3.6     | ng    | 42000    | 1                     | 8.04  | 233348 | 20.0 |
| Benzene, 1-ethyl-... | 7.11  | 8.4     | ng    | 97500    | 1                     | 8.04  | 233348 | 20.0 |
| Benzene, 1-ethyl-... | 7.17  | 6.8     | ng    | 79964    | 1                     | 8.04  | 233348 | 20.0 |
| unknown7.58          | 7.58  | 60.9    | ng    | 709974   | 1                     | 8.04  | 233348 | 20.0 |
| Decane               | 7.64  | 4.8     | ng    | 56344    | 1                     | 8.04  | 233348 | 20.0 |
| Benzene, 1,2,3-tr... | 7.68  | 15.4    | ng    | 180184   | 1                     | 8.04  | 233348 | 20.0 |
| Benzene, 1-methyl... | 8.58  | 9.0     | ng    | 104547   | 1                     | 8.04  | 233348 | 20.0 |
| Benzene, 1-ethyl-... | 8.68  | 9.3     | ng    | 109058   | 1                     | 8.04  | 233348 | 20.0 |
| Benzene, 1-methyl... | 8.84  | 3.9     | ng    | 45886    | 1                     | 8.04  | 233348 | 20.0 |
| Benzene, 4-ethyl-... | 9.14  | 8.4     | ng    | 98349    | 1                     | 8.04  | 233348 | 20.0 |
| unknown9.38          | 9.38  | 5.6     | ng    | 65167    | 1                     | 8.04  | 233348 | 20.0 |
| Benzene, 1-ethyl-... | 9.73  | 2.6     | ng    | 46076    | 2                     | 10.86 | 360183 | 20.0 |
| Benzene, 2-ethyl-... | 10.25 | 3.9     | ng    | 70789    | 2                     | 10.86 | 360183 | 20.0 |