

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF012818\  
 Data File : BF102577.D  
 Acq On : 29 Jan 2018 00:12  
 Operator : SJ/JU  
 Sample : J1257-13  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 WC-17(0-5)

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\_F\METHODS\8270-BF012218.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         | 4.904       | 474           | 479         | 487          | rBV      | 981085         | 1087467       | 29.33%          | 2.471%        |
| 2         | 5.181       | 521           | 526         | 528          | rBV2     | 41588          | 52294         | 1.41%           | 0.119%        |
| 3         | 5.281       | 540           | 543         | 548          | rBV      | 115099         | 126219        | 3.40%           | 0.287%        |
| 4         | 5.328       | 548           | 551         | 559          | rBV      | 1038465        | 1016850       | 27.42%          | 2.310%        |
| 5         | 5.545       | 585           | 588         | 590          | rBV      | 79380          | 78693         | 2.12%           | 0.179%        |
| 6         | 5.604       | 595           | 598         | 603          | rVB      | 98592          | 89917         | 2.43%           | 0.204%        |
| 7         | 5.763       | 621           | 625         | 628          | rBV2     | 39307          | 58209         | 1.57%           | 0.132%        |
| 8         | 5.851       | 633           | 640         | 643          | rBV2     | 60169          | 84370         | 2.28%           | 0.192%        |
| 9         | 5.945       | 654           | 656         | 663          | rBV2     | 72596          | 87520         | 2.36%           | 0.199%        |
| 10        | 6.139       | 684           | 689         | 692          | rBV2     | 88279          | 116112        | 3.13%           | 0.264%        |
| 11        | 6.192       | 695           | 698         | 702          | rVB4     | 163759         | 236417        | 6.38%           | 0.537%        |
| 12        | 6.234       | 702           | 705         | 708          | rBV      | 341858         | 296425        | 7.99%           | 0.673%        |
| 13        | 6.263       | 708           | 710         | 713          | rVV      | 114705         | 102145        | 2.75%           | 0.232%        |
| 14        | 6.292       | 713           | 715         | 716          | rVV      | 152927         | 117022        | 3.16%           | 0.266%        |
| 15        | 6.310       | 716           | 718         | 723          | rVB      | 262504         | 243248        | 6.56%           | 0.553%        |
| 16        | 6.369       | 723           | 728         | 737          | rBV      | 2958120        | 3707765       | 100.00%         | 8.424%        |
| 17        | 6.486       | 744           | 748         | 753          | rVB      | 2544139        | 2226286       | 60.04%          | 5.058%        |
| 18        | 6.539       | 753           | 757         | 764          | rVB2     | 455581         | 491070        | 13.24%          | 1.116%        |
| 19        | 6.645       | 770           | 775         | 777          | rBV2     | 70396          | 79991         | 2.16%           | 0.182%        |
| 20        | 6.716       | 784           | 787         | 793          | rVB      | 1243732        | 1067228       | 28.78%          | 2.425%        |
| 21        | 6.769       | 793           | 796         | 803          | rVB2     | 242311         | 270661        | 7.30%           | 0.615%        |
| 22        | 6.828       | 803           | 806         | 808          | rBV      | 121739         | 112461        | 3.03%           | 0.256%        |
| 23        | 6.869       | 808           | 813         | 819          | rVB      | 3032194        | 3054678       | 82.39%          | 6.940%        |
| 24        | 6.957       | 826           | 828         | 830          | rBV      | 171557         | 151512        | 4.09%           | 0.344%        |
| 25        | 6.986       | 830           | 833         | 836          | rVV2     | 365369         | 346177        | 9.34%           | 0.786%        |
| 26        | 7.033       | 836           | 841         | 842          | rVV2     | 589885         | 597979        | 16.13%          | 1.359%        |
| 27        | 7.045       | 842           | 843         | 849          | rVB      | 257271         | 211498        | 5.70%           | 0.481%        |
| 28        | 7.098       | 849           | 852         | 856          | rBV3     | 185730         | 216954        | 5.85%           | 0.493%        |
| 29        | 7.175       | 863           | 865         | 867          | rVB      | 109959         | 76573         | 2.07%           | 0.174%        |
| 30        | 7.198       | 867           | 869         | 873          | rVB      | 113409         | 86226         | 2.33%           | 0.196%        |
| 31        | 7.245       | 873           | 877         | 880          | rBV      | 355991         | 363456        | 9.80%           | 0.826%        |
| 32        | 7.286       | 880           | 884         | 886          | rVV      | 2440892        | 2310655       | 62.32%          | 5.250%        |
| 33        | 7.304       | 886           | 887         | 891          | rVB      | 303919         | 214242        | 5.78%           | 0.487%        |
| 34        | 7.357       | 891           | 896         | 899          | rVB4     | 133727         | 171854        | 4.63%           | 0.390%        |

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 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 WC-17(0-5)

## 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\_F\METHODS\8270-BF012218.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |       |      |      |      |      |         |         |        |        |
|----|-------|------|------|------|------|---------|---------|--------|--------|
| 35 | 7.398 | 899  | 903  | 905  | rBV2 | 135521  | 177986  | 4.80%  | 0.404% |
| 36 | 7.463 | 910  | 914  | 915  | rBV2 | 110994  | 119692  | 3.23%  | 0.272% |
| 37 | 7.480 | 915  | 917  | 920  | rVV  | 483115  | 451333  | 12.17% | 1.025% |
| 38 | 7.510 | 920  | 922  | 926  | rVB  | 392835  | 306246  | 8.26%  | 0.696% |
| 39 | 7.598 | 933  | 937  | 939  | rBV3 | 167238  | 201416  | 5.43%  | 0.458% |
| 40 | 7.651 | 942  | 946  | 947  | rVV4 | 140962  | 188850  | 5.09%  | 0.429% |
| 41 | 7.669 | 947  | 949  | 953  | rVV2 | 209867  | 297864  | 8.03%  | 0.677% |
| 42 | 7.710 | 953  | 956  | 958  | rVV2 | 164760  | 187018  | 5.04%  | 0.425% |
| 43 | 7.733 | 958  | 960  | 963  | rVV3 | 278168  | 359244  | 9.69%  | 0.816% |
| 44 | 7.763 | 963  | 965  | 967  | rVV  | 209016  | 176257  | 4.75%  | 0.400% |
| 45 | 7.786 | 967  | 969  | 971  | rVV  | 136452  | 130479  | 3.52%  | 0.296% |
| 46 | 7.810 | 971  | 973  | 976  | rVV2 | 119467  | 133799  | 3.61%  | 0.304% |
| 47 | 7.863 | 979  | 982  | 985  | rBV  | 117179  | 109636  | 2.96%  | 0.249% |
| 48 | 7.945 | 990  | 996  | 999  | rBV5 | 103230  | 220751  | 5.95%  | 0.502% |
| 49 | 7.975 | 999  | 1001 | 1002 | rVV  | 354767  | 306672  | 8.27%  | 0.697% |
| 50 | 7.998 | 1002 | 1005 | 1011 | rVV2 | 1558960 | 1901712 | 51.29% | 4.321% |
| 51 | 8.051 | 1011 | 1014 | 1017 | rVB2 | 235189  | 284420  | 7.67%  | 0.646% |
| 52 | 8.080 | 1017 | 1019 | 1022 | rBV  | 139211  | 124257  | 3.35%  | 0.282% |
| 53 | 8.151 | 1027 | 1031 | 1032 | rBV3 | 59416   | 64953   | 1.75%  | 0.148% |
| 54 | 8.192 | 1036 | 1038 | 1041 | rVB2 | 89680   | 74970   | 2.02%  | 0.170% |
| 55 | 8.233 | 1041 | 1045 | 1046 | rBV4 | 43307   | 65007   | 1.75%  | 0.148% |
| 56 | 8.286 | 1049 | 1054 | 1060 | rVB5 | 163065  | 256265  | 6.91%  | 0.582% |
| 57 | 8.339 | 1060 | 1063 | 1066 | rBV4 | 84546   | 99241   | 2.68%  | 0.225% |
| 58 | 8.380 | 1066 | 1070 | 1072 | rBV3 | 151863  | 175129  | 4.72%  | 0.398% |
| 59 | 8.410 | 1072 | 1075 | 1077 | rVV3 | 184107  | 170984  | 4.61%  | 0.388% |
| 60 | 8.463 | 1081 | 1084 | 1086 | rVB2 | 124131  | 130044  | 3.51%  | 0.295% |
| 61 | 8.498 | 1088 | 1090 | 1092 | rBV2 | 85186   | 75436   | 2.03%  | 0.171% |
| 62 | 8.557 | 1097 | 1100 | 1102 | rVB  | 122938  | 84614   | 2.28%  | 0.192% |
| 63 | 8.586 | 1102 | 1105 | 1108 | rBV2 | 259925  | 246162  | 6.64%  | 0.559% |
| 64 | 8.680 | 1118 | 1121 | 1124 | rVB4 | 107114  | 134255  | 3.62%  | 0.305% |
| 65 | 8.710 | 1124 | 1126 | 1129 | rBV  | 406222  | 321282  | 8.67%  | 0.730% |
| 66 | 8.769 | 1134 | 1136 | 1138 | rBV3 | 44886   | 55122   | 1.49%  | 0.125% |
| 67 | 8.810 | 1138 | 1143 | 1146 | rVV  | 239304  | 237030  | 6.39%  | 0.539% |
| 68 | 8.869 | 1150 | 1153 | 1155 | rBV3 | 91088   | 106001  | 2.86%  | 0.241% |
| 69 | 8.898 | 1156 | 1158 | 1163 | rVB3 | 105513  | 103576  | 2.79%  | 0.235% |
| 70 | 8.945 | 1163 | 1166 | 1170 | rBV4 | 83486   | 92882   | 2.51%  | 0.211% |
| 71 | 8.986 | 1170 | 1173 | 1175 | rBV3 | 90526   | 86525   | 2.33%  | 0.197% |

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Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

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\_F\METHODS\8270-BF012218.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|     |        |      |      |      |      |         |         |        |        |
|-----|--------|------|------|------|------|---------|---------|--------|--------|
| 72  | 9.010  | 1175 | 1177 | 1180 | rBV  | 225866  | 198177  | 5.34%  | 0.450% |
| 73  | 9.075  | 1183 | 1188 | 1191 | rBV  | 3979305 | 3564738 | 96.14% | 8.099% |
| 74  | 9.145  | 1196 | 1200 | 1203 | rBV  | 317431  | 379685  | 10.24% | 0.863% |
| 75  | 9.222  | 1210 | 1213 | 1216 | rVB4 | 70875   | 80174   | 2.16%  | 0.182% |
| 76  | 9.269  | 1219 | 1221 | 1227 | rVB3 | 65767   | 66512   | 1.79%  | 0.151% |
| 77  | 9.339  | 1230 | 1233 | 1236 | rBV2 | 176608  | 207912  | 5.61%  | 0.472% |
| 78  | 9.410  | 1242 | 1245 | 1247 | rBV2 | 173287  | 207368  | 5.59%  | 0.471% |
| 79  | 9.469  | 1250 | 1255 | 1257 | rVB2 | 668772  | 667669  | 18.01% | 1.517% |
| 80  | 9.527  | 1263 | 1265 | 1267 | rBV  | 74007   | 51710   | 1.39%  | 0.117% |
| 81  | 9.557  | 1268 | 1270 | 1274 | rVB4 | 149897  | 163157  | 4.40%  | 0.371% |
| 82  | 9.610  | 1276 | 1279 | 1284 | rBV4 | 247773  | 307582  | 8.30%  | 0.699% |
| 83  | 9.651  | 1284 | 1286 | 1289 | rBV  | 268788  | 247223  | 6.67%  | 0.562% |
| 84  | 9.680  | 1289 | 1291 | 1295 | rBV2 | 193280  | 189609  | 5.11%  | 0.431% |
| 85  | 9.722  | 1295 | 1298 | 1299 | rVV3 | 136592  | 145614  | 3.93%  | 0.331% |
| 86  | 9.751  | 1300 | 1303 | 1306 | rVB  | 1300987 | 1086618 | 29.31% | 2.469% |
| 87  | 9.786  | 1306 | 1309 | 1312 | rBV3 | 109695  | 124655  | 3.36%  | 0.283% |
| 88  | 9.851  | 1318 | 1320 | 1325 | rBV3 | 126011  | 142087  | 3.83%  | 0.323% |
| 89  | 9.939  | 1333 | 1335 | 1337 | rBV  | 112169  | 98314   | 2.65%  | 0.223% |
| 90  | 9.998  | 1342 | 1345 | 1349 | rBV4 | 270338  | 357140  | 9.63%  | 0.811% |
| 91  | 10.069 | 1354 | 1357 | 1361 | rBV3 | 214023  | 315356  | 8.51%  | 0.716% |
| 92  | 10.151 | 1368 | 1371 | 1373 | rBV3 | 373367  | 357812  | 9.65%  | 0.813% |
| 93  | 10.174 | 1373 | 1375 | 1378 | rVB  | 233121  | 224413  | 6.05%  | 0.510% |
| 94  | 10.398 | 1409 | 1413 | 1416 | rBV  | 856701  | 890038  | 24.00% | 2.022% |
| 95  | 10.516 | 1431 | 1433 | 1435 | rBV2 | 186662  | 138565  | 3.74%  | 0.315% |
| 96  | 10.539 | 1435 | 1437 | 1438 | rBV  | 170109  | 138775  | 3.74%  | 0.315% |
| 97  | 10.669 | 1454 | 1459 | 1461 | rBV  | 1821953 | 2251753 | 60.73% | 5.116% |
| 98  | 11.133 | 1535 | 1538 | 1545 | rVB  | 1886025 | 1850924 | 49.92% | 4.205% |
| 99  | 11.245 | 1554 | 1557 | 1559 | rBV  | 985597  | 969216  | 26.14% | 2.202% |
| 100 | 11.486 | 1595 | 1598 | 1601 | rBV  | 749740  | 785463  | 21.18% | 1.785% |

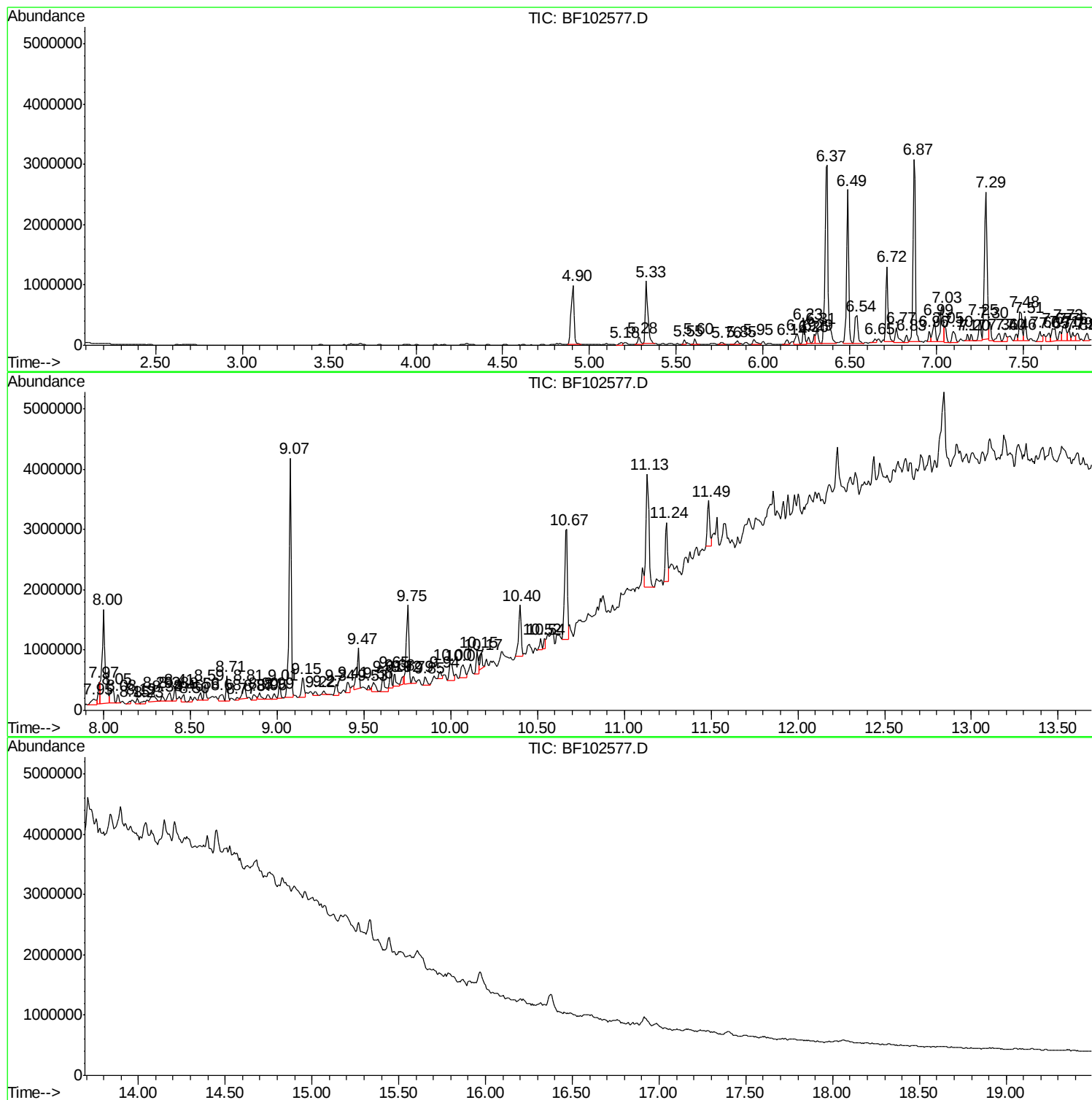
Sum of corrected areas: 44015543

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Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF012218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012818\  
Data File : BF102577.D  
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WC-17(0-5)

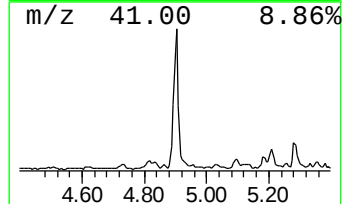
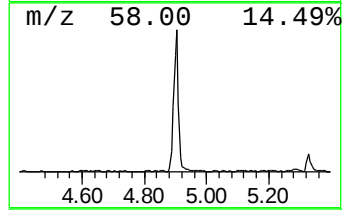
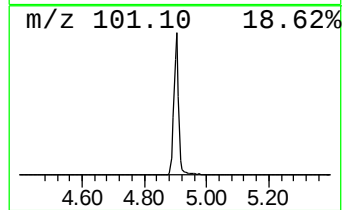
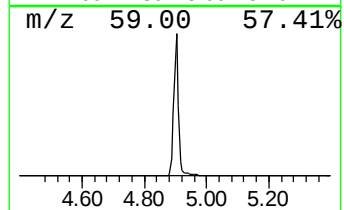
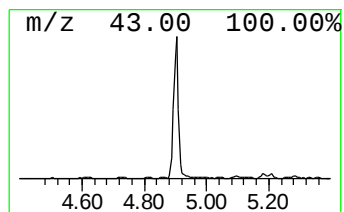
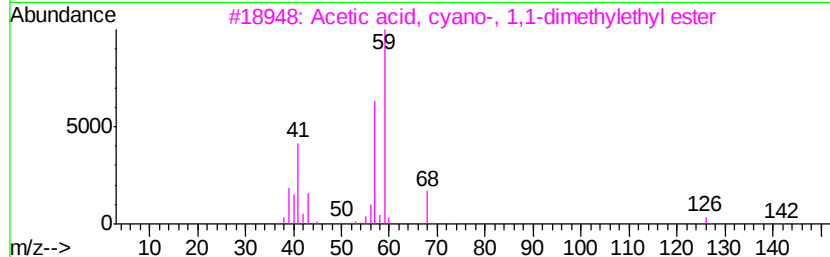
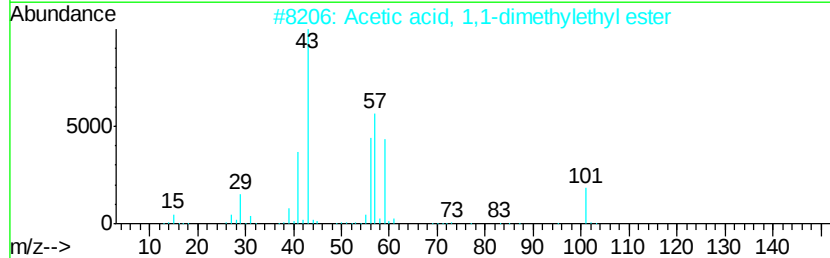
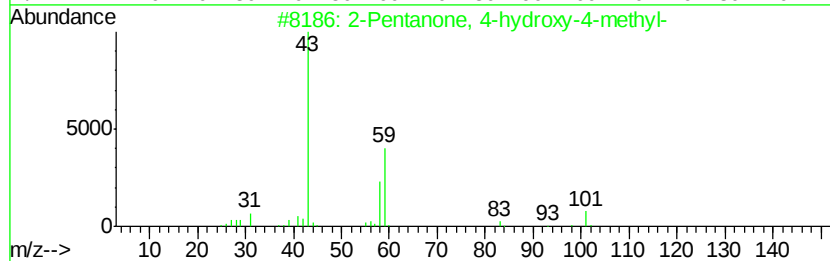
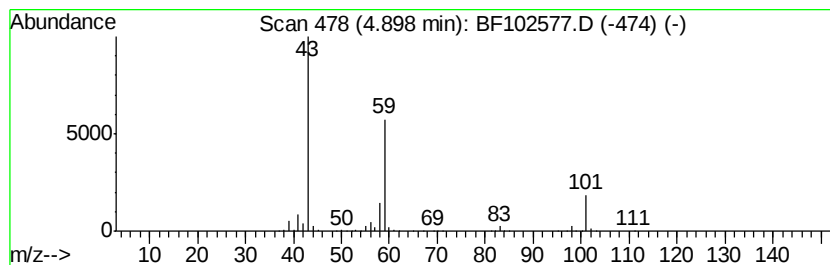
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 4.90 | 20.38 ng | 1087470 | 1,4-Dichlorobenzene-d4 | 6.72 |

| 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, 1,1-dimethylethyl e... | 116 | C6H12O2  | 000540-88-5 | 39   |
| 3    |    | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 17   |
| 4    |    | 2,3-Butanedione, monooxime          | 101 | C4H7NO2  | 000057-71-6 | 9    |
| 5    |    | Morpholine, 4-methyl-               | 101 | C5H11NO  | 000109-02-4 | 9    |



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WC-17(0-5)

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

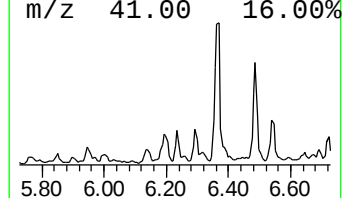
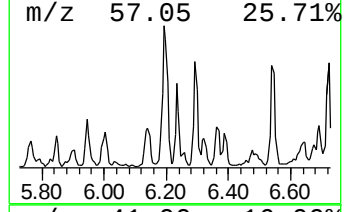
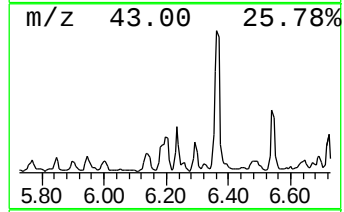
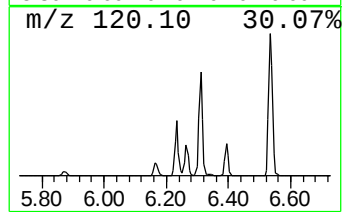
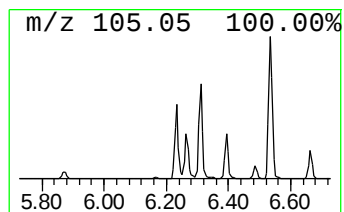
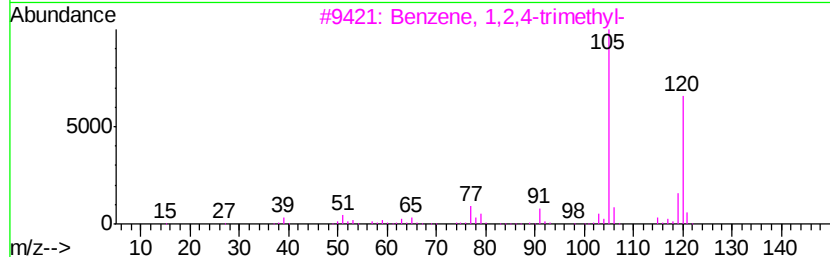
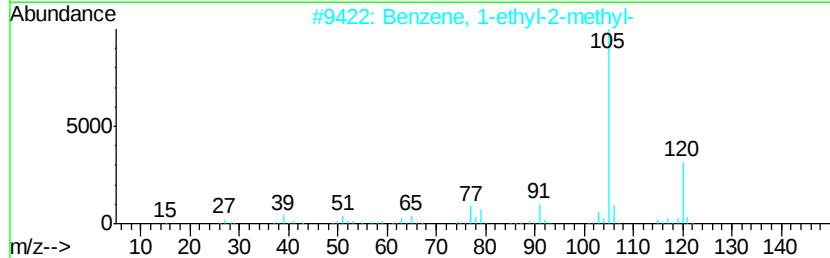
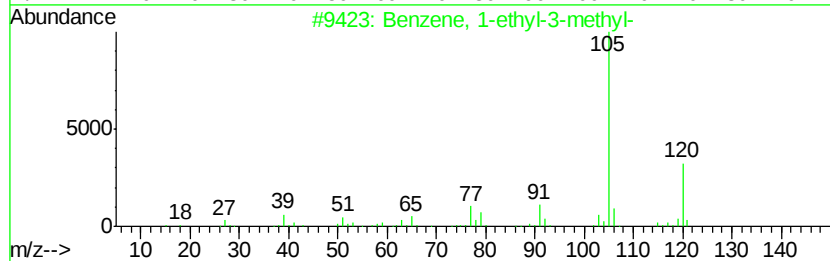
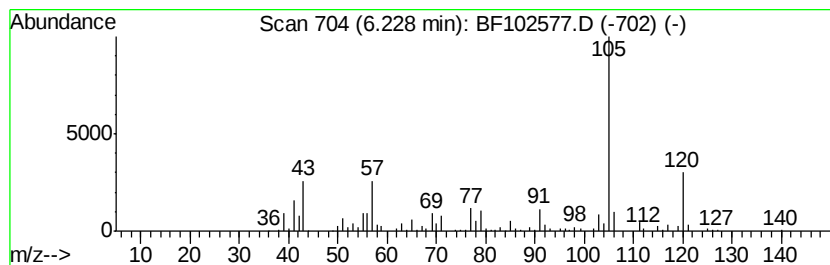
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 16

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.23 | 5.56 ng | 296425 | 1,4-Dichlorobenzene-d4 | 6.72 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 92   |
| 2    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 92   |
| 3    |    | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 64   |
| 4    |    | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 64   |
| 5    |    | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 64   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF012818\  
Data File : BF102577.D  
Acq On : 29 Jan 2018 00:12  
Operator : SJ/JU  
Sample : J1257-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF012218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

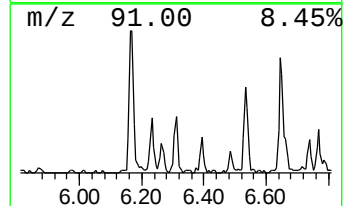
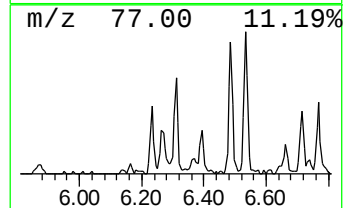
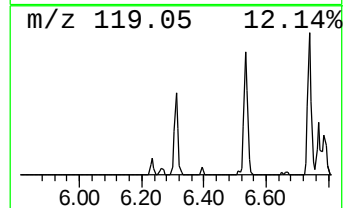
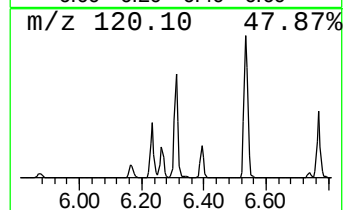
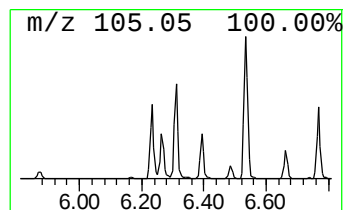
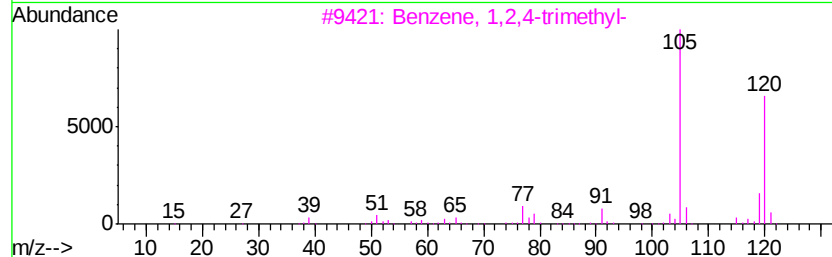
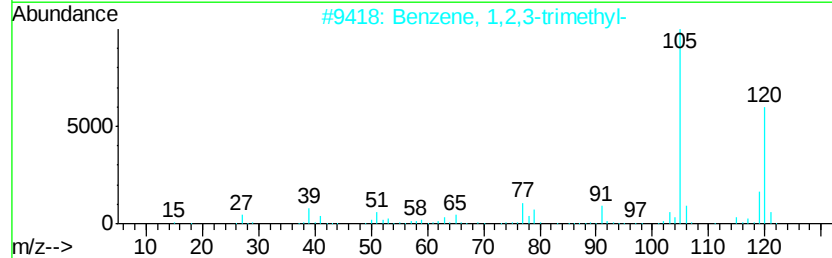
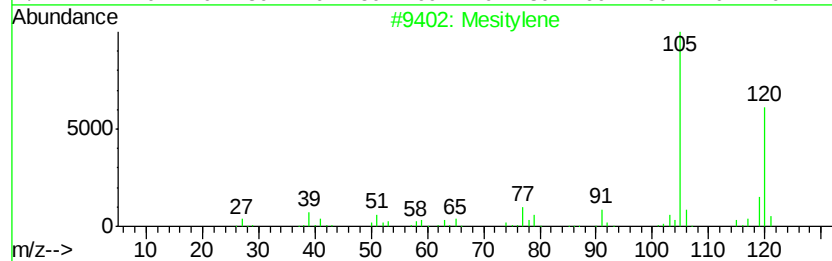
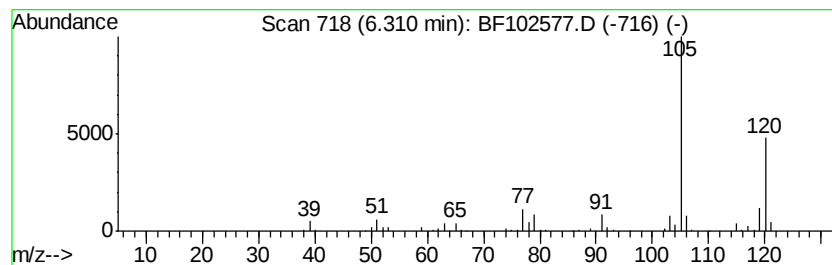
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Mesitylene Concentration Rank 19

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.31 | 4.56 ng | 243248 | 1,4-Dichlorobenzene-d4 | 6.72 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 97   |
| 2    |    | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 97   |
| 3    |    | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 95   |
| 4    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 90   |
| 5    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 87   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012818\  
Data File : BF102577.D  
Acq On : 29 Jan 2018 00:12  
Operator : SJ/JU  
Sample : J1257-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

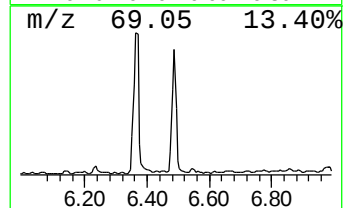
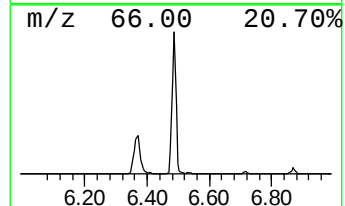
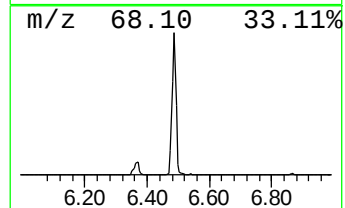
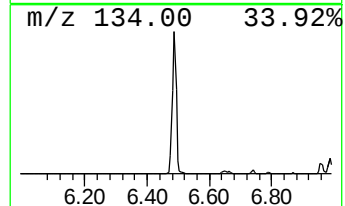
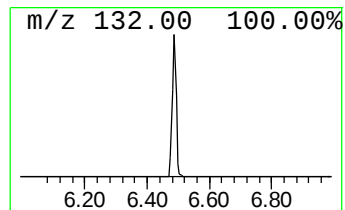
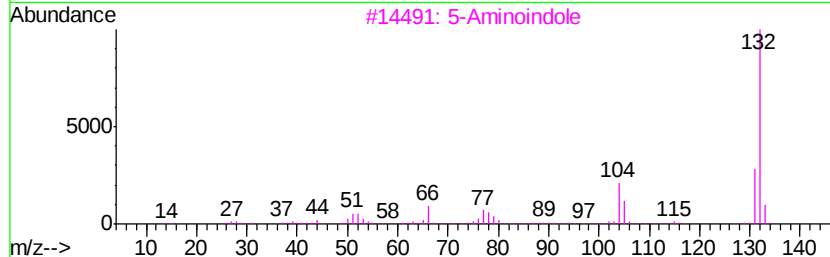
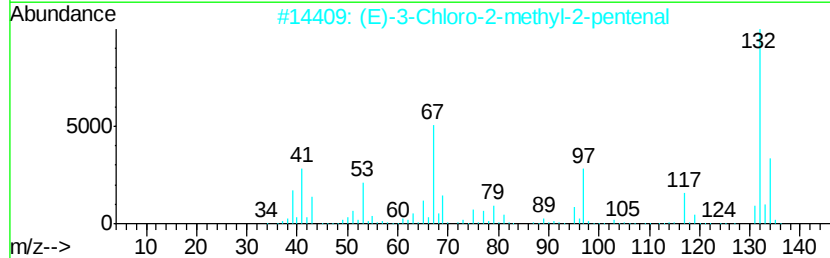
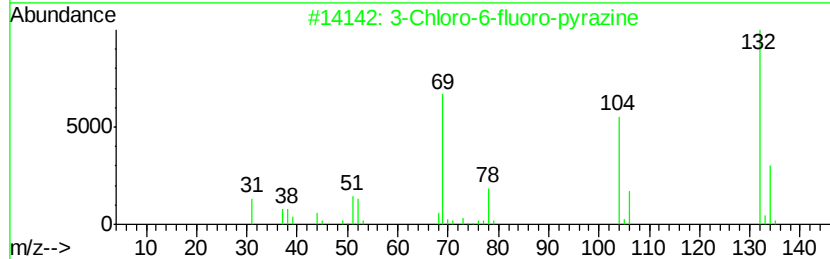
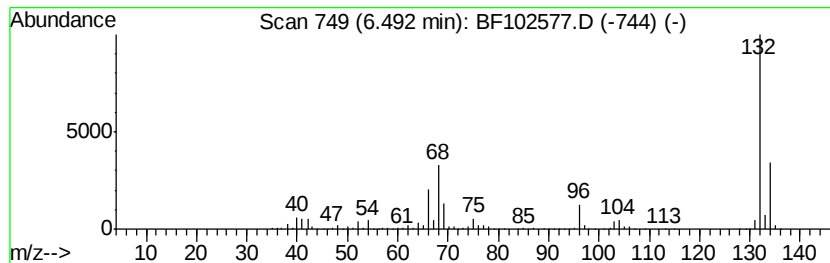
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 unknown6.49 Concentration Rank 3

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.49 | 41.72 ng | 2226290 | 1,4-Dichlorobenzene-d4 | 6.72 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 12   |
| 3    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 12   |
| 4    |    | Tranylcypromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 10   |
| 5    |    | 1,3,5-Trifluorobenzene           | 132 | C6H3F3    | 000372-38-3  | 9    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF012818\  
Data File : BF102577.D  
Acq On : 29 Jan 2018 00:12  
Operator : SJ/JU  
Sample : J1257-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

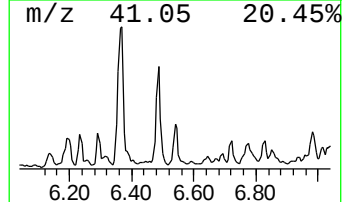
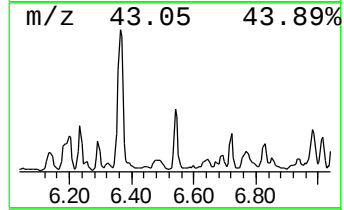
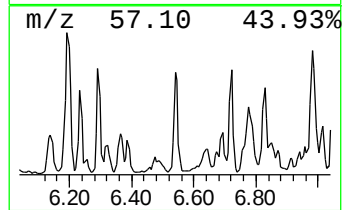
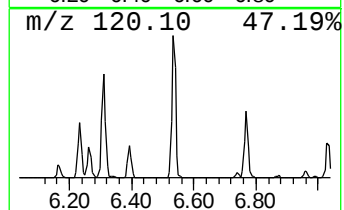
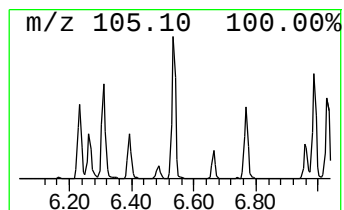
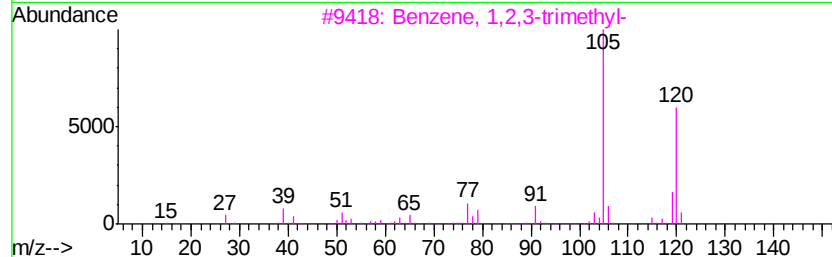
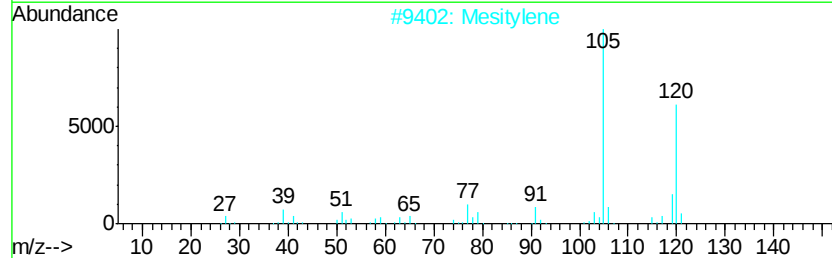
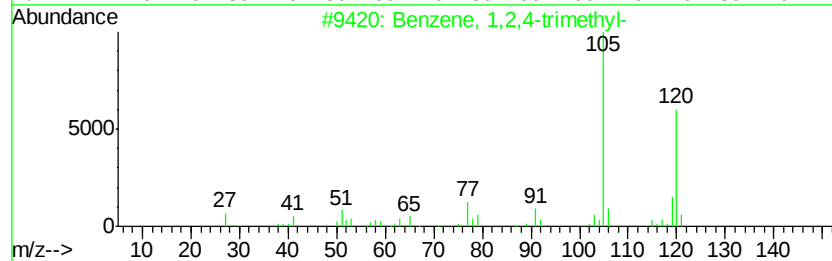
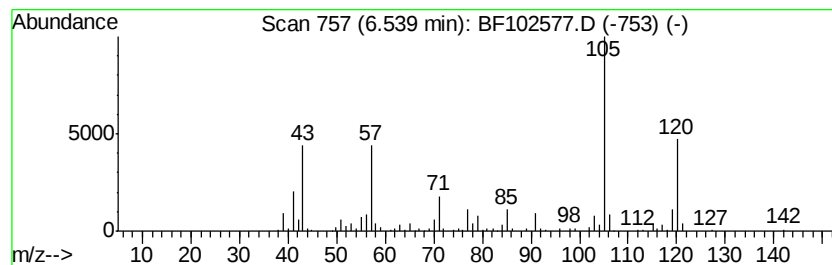
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 5 Benzene, 1,2,4-trimethyl- Concentration Rank 9

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.54 | 9.20 ng | 491070 | 1,4-Dichlorobenzene-d4 | 6.72 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 94   |
| 2    |    | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 94   |
| 3    |    | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 94   |
| 4    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 87   |
| 5    |    | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 87   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012818\  
Data File : BF102577.D  
Acq On : 29 Jan 2018 00:12  
Operator : SJ/JU  
Sample : J1257-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

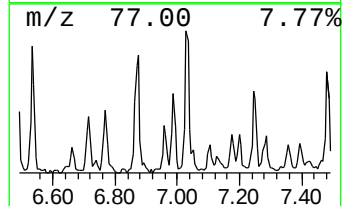
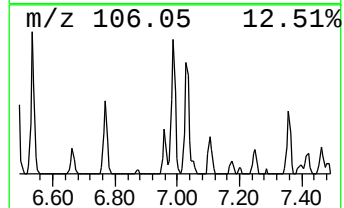
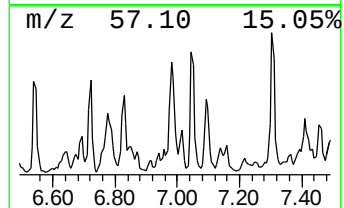
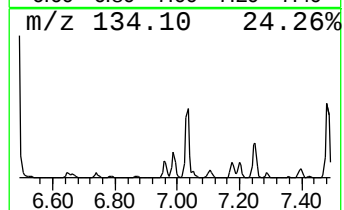
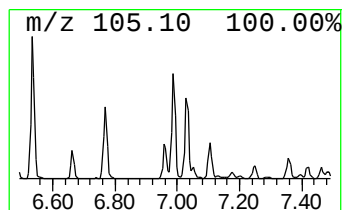
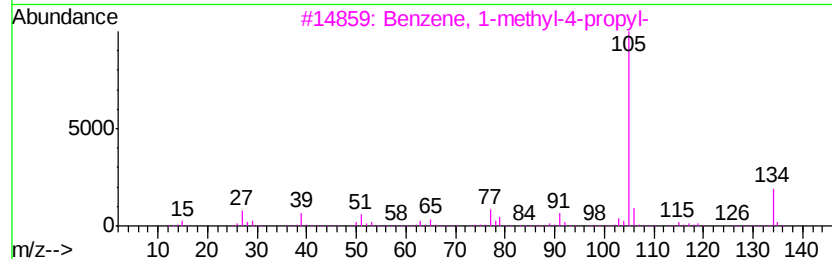
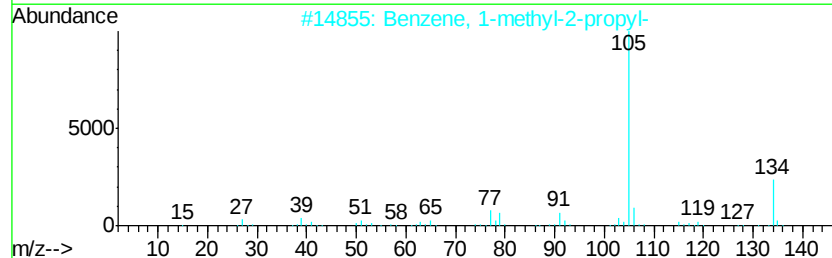
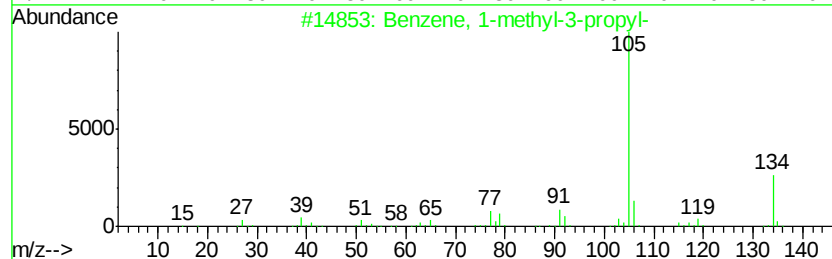
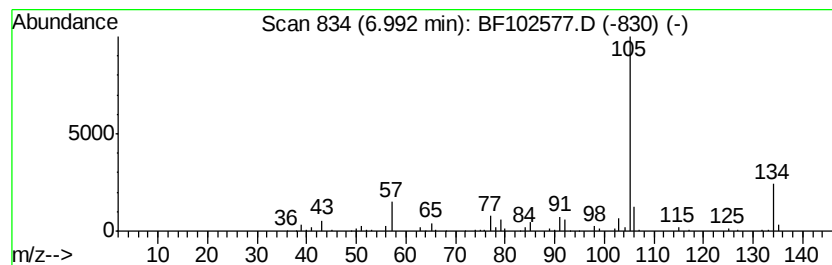
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Benzene, 1-methyl-3-propyl- Concentration Rank 13

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.99 | 6.49 ng | 346177 | 1,4-Dichlorobenzene-d4 | 6.72 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-3-propyl-         | 134 | C10H14  | 001074-43-7 | 76   |
| 2    |    | Benzene, 1-methyl-2-propyl-         | 134 | C10H14  | 001074-17-5 | 76   |
| 3    |    | Benzene, 1-methyl-4-propyl-         | 134 | C10H14  | 001074-55-1 | 70   |
| 4    |    | Benzeneacetaldehyde, .alpha.-met... | 134 | C9H10O  | 000093-53-8 | 58   |
| 5    |    | Benzene, (1-methylpropyl)-          | 134 | C10H14  | 000135-98-8 | 53   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF012818\  
Data File : BF102577.D  
Acq On : 29 Jan 2018 00:12  
Operator : SJ/JU  
Sample : J1257-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF012218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

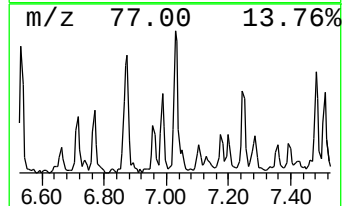
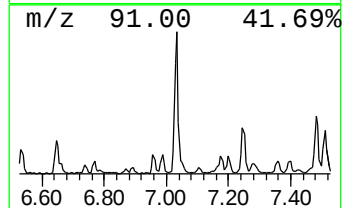
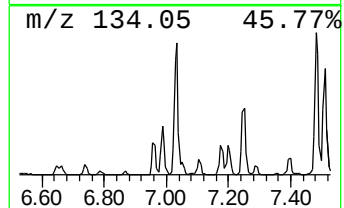
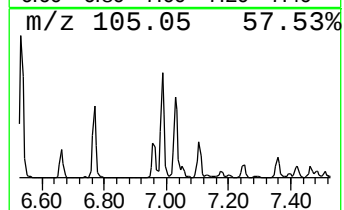
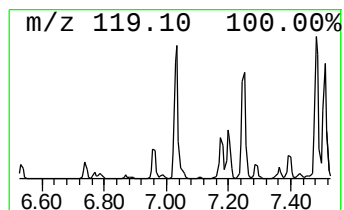
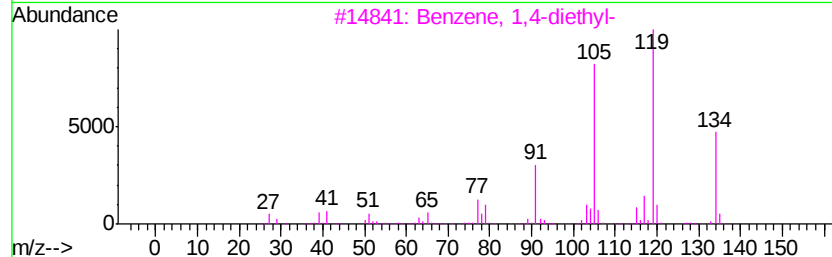
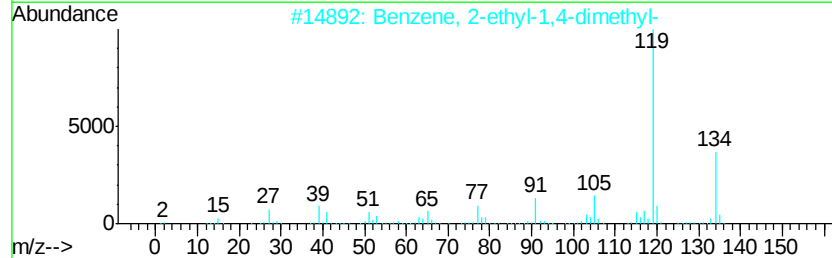
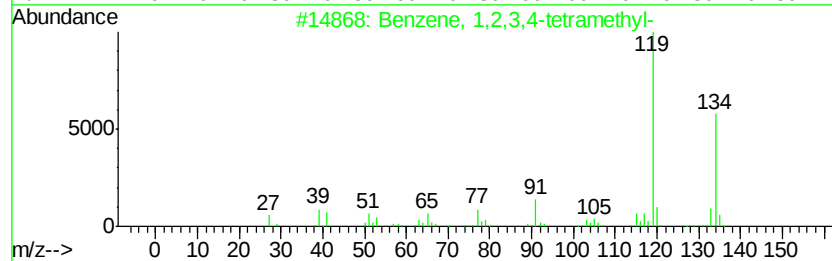
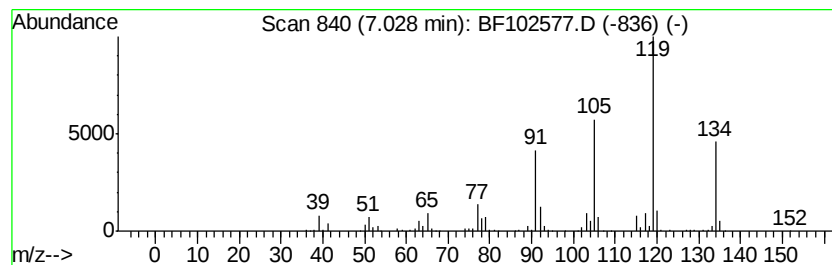
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 8

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 7.03 | 11.21 ng | 597979 | 1,4-Dichlorobenzene-d4 | 6.72 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 91   |
| 2    |    | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 91   |
| 3    |    | Benzene, 1,4-diethyl-          | 134 | C10H14  | 000105-05-5 | 91   |
| 4    |    | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 91   |
| 5    |    | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 91   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012818\  
Data File : BF102577.D  
Acq On : 29 Jan 2018 00:12  
Operator : SJ/JU  
Sample : J1257-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

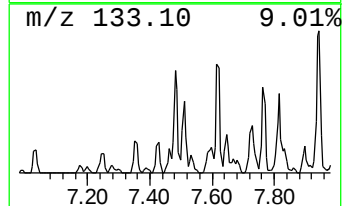
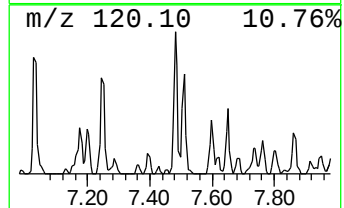
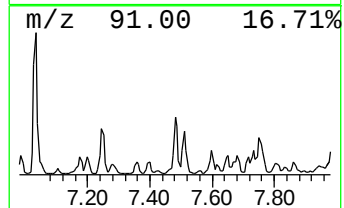
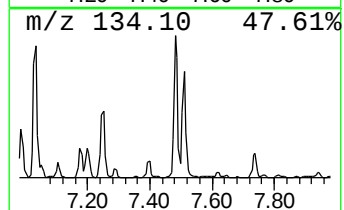
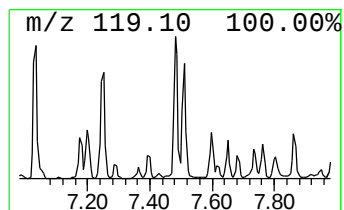
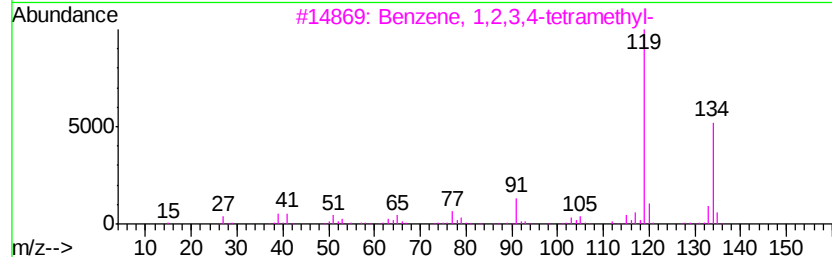
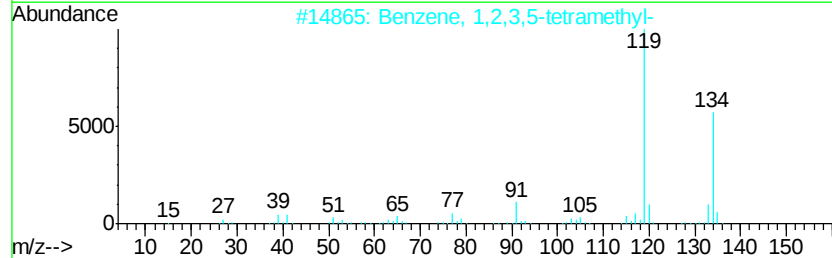
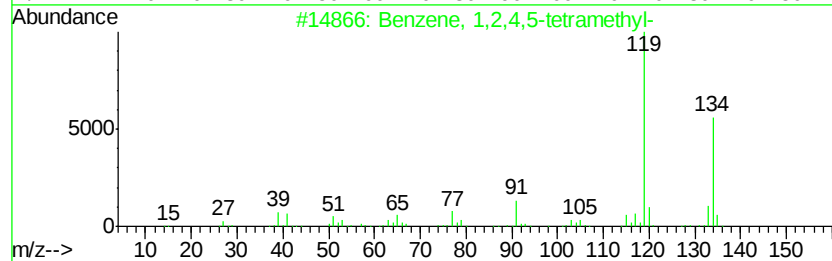
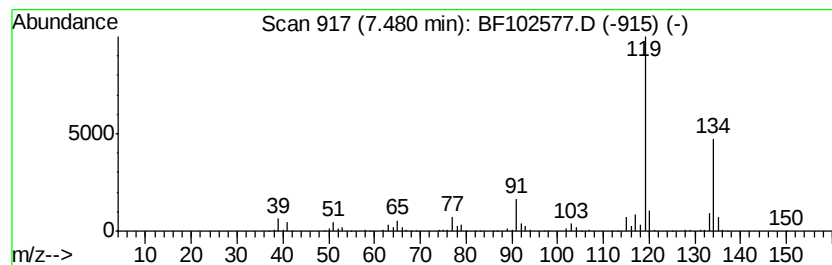
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 18

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 7.48 | 4.75 ng | 451333 | Naphthalene-d8   | 8.00 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,4,5-tetramethyl-       | 134 | C10H14  | 000095-93-2 | 95   |
| 2    |    |   | Benzene, 1,2,3,5-tetramethyl-       | 134 | C10H14  | 000527-53-7 | 94   |
| 3    |    |   | Benzene, 1,2,3,4-tetramethyl-       | 134 | C10H14  | 000488-23-3 | 94   |
| 4    |    |   | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 94   |
| 5    |    |   | Benzene, 1-ethyl-3,5-dimethyl-      | 134 | C10H14  | 000934-74-7 | 91   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF012818\  
Data File : BF102577.D  
Acq On : 29 Jan 2018 00:12  
Operator : SJ/JU  
Sample : J1257-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF012218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

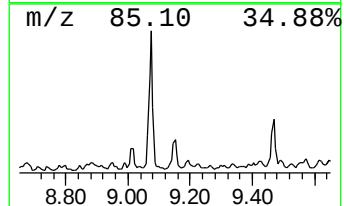
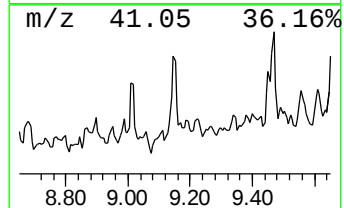
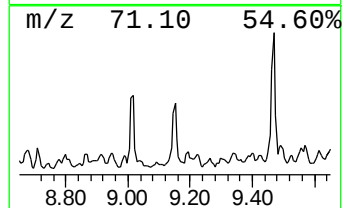
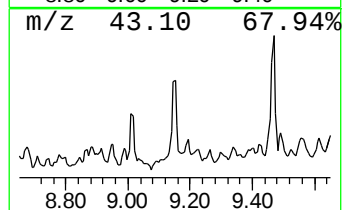
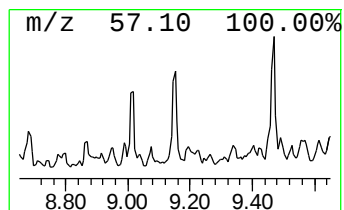
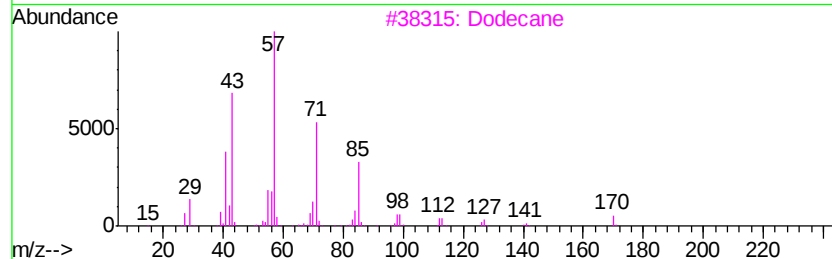
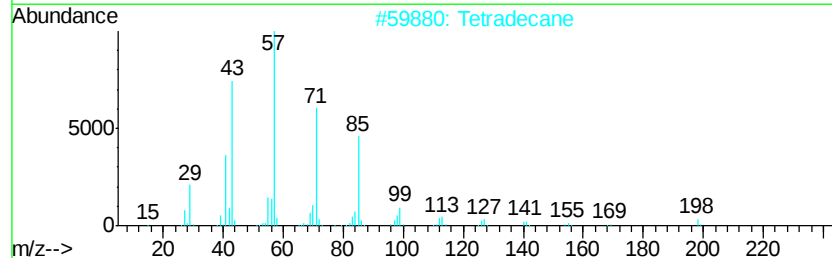
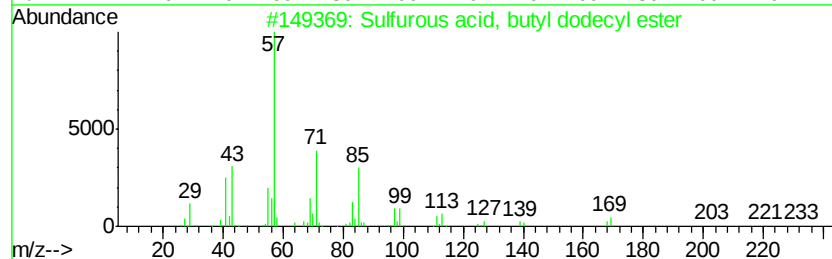
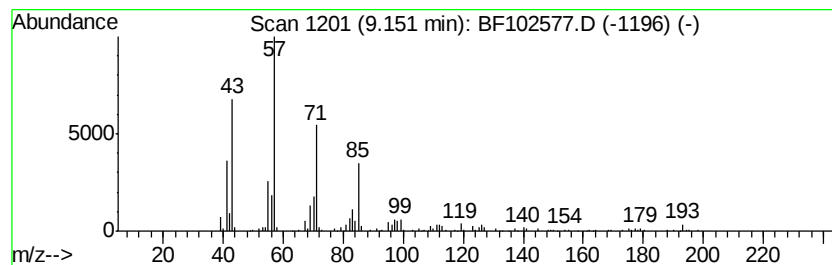
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Sulfurous acid, butyl dodec... Concentration Rank 10

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.15 | 6.99 ng | 379685 | Acenaphthene-d10 | 9.75 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Sulfurous acid, butyl dodecyl ester | 306 | C16H34O3S | 1000309-17-9 | 53   |
| 2    |    | Tetradecane                         | 198 | C14H30    | 000629-59-4  | 52   |
| 3    |    | Dodecane                            | 170 | C12H26    | 000112-40-3  | 50   |
| 4    |    | Hexatriacontane                     | 507 | C36H74    | 000630-06-8  | 50   |
| 5    |    | Tridecane                           | 184 | C13H28    | 000629-50-5  | 50   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF012818\  
Data File : BF102577.D  
Acq On : 29 Jan 2018 00:12  
Operator : SJ/JU  
Sample : J1257-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF012218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

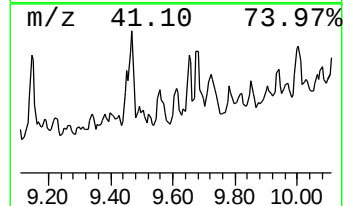
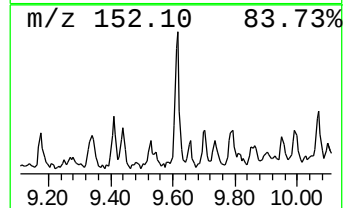
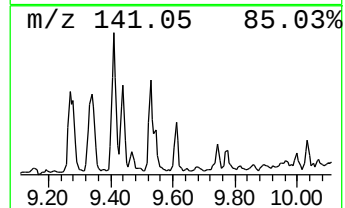
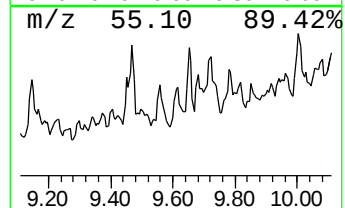
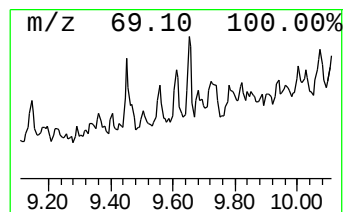
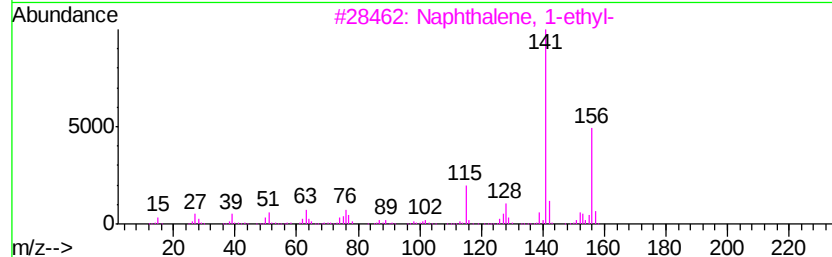
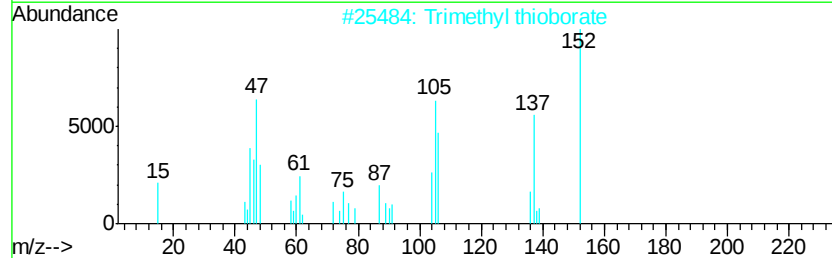
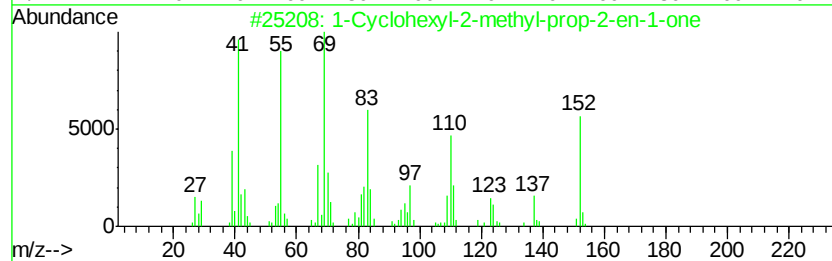
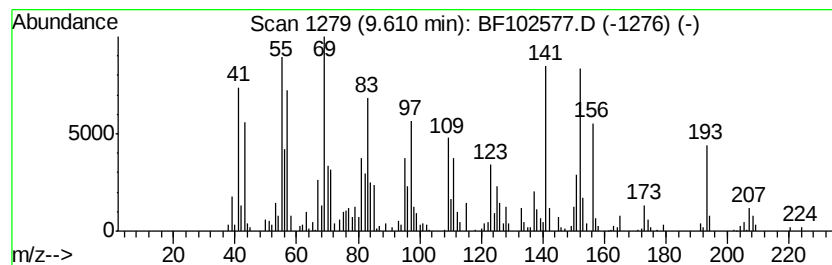
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 unknown9.61 Concentration Rank 15

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.61 | 5.66 ng | 307582 | Acenaphthene-d10 | 9.75 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1-Cyclohexyl-2-methyl-prop-2-en-... | 152 | C10H16O | 025183-82-8 | 41   |
| 2    |    |   | Trimethyl thioborate                | 152 | C3H9BS3 | 000997-49-9 | 25   |
| 3    |    |   | Naphthalene, 1-ethyl-               | 156 | C12H12  | 001127-76-0 | 25   |
| 4    |    |   | 1-Cyclohexyl-1-(4-methylcyclohex... | 208 | C15H28  | 002027-12-5 | 25   |
| 5    |    |   | Naphthalene, 2-ethyl-               | 156 | C12H12  | 000939-27-5 | 25   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF012818\  
Data File : BF102577.D  
Acq On : 29 Jan 2018 00:12  
Operator : SJ/JU  
Sample : J1257-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF012218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

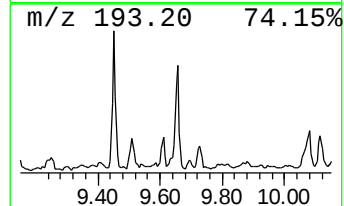
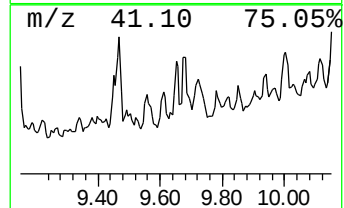
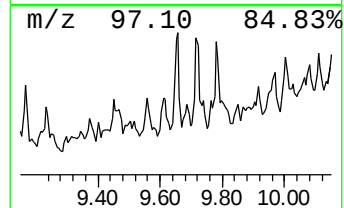
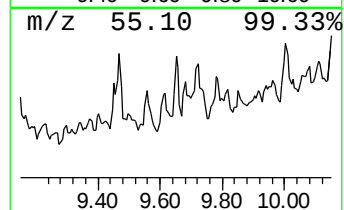
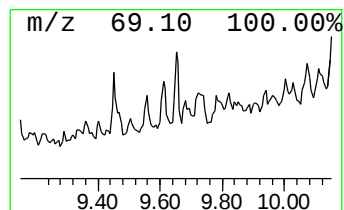
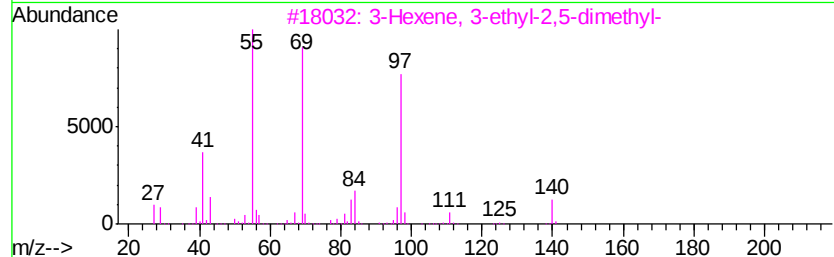
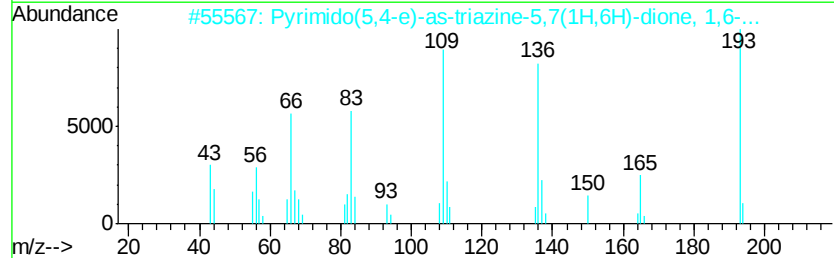
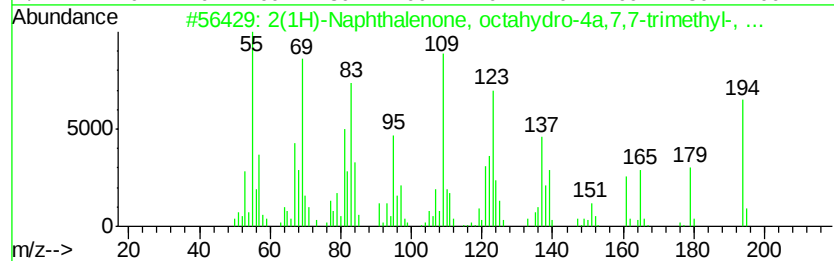
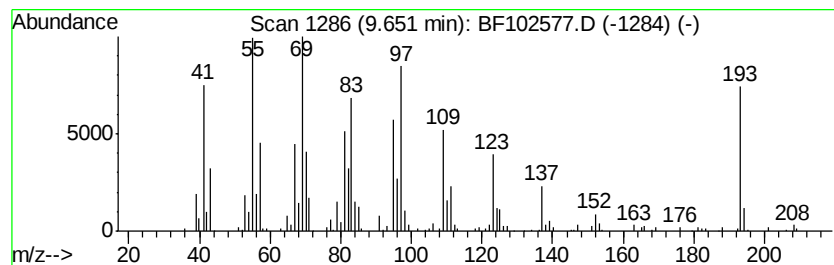
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 unknown9.65 Concentration Rank 20

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.65 | 4.55 ng | 247223 | Acenaphthene-d10 | 9.75 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2(1H)-Naphthalenone, octahydro-4... | 194 | C13H22O  | 054699-31-9 | 43   |
| 2    |    |   | Pyrimido(5,4-e)-as-triazine-5,7(... | 193 | C7H7N5O2 | 000084-82-2 | 35   |
| 3    |    |   | 3-Hexene, 3-ethyl-2,5-dimethyl-     | 140 | C10H20   | 062338-08-3 | 27   |
| 4    |    |   | Cyclohexanone, 2-(1-mercapto-1-m... | 186 | C10H18OS | 033281-91-3 | 22   |
| 5    |    |   | 2-Butenamide, 2,3-dimethyl-N-phe... | 189 | C12H15NO | 024735-54-4 | 22   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012818\  
Data File : BF102577.D  
Acq On : 29 Jan 2018 00:12  
Operator : SJ/JU  
Sample : J1257-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

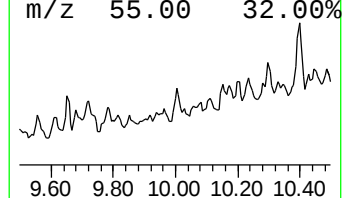
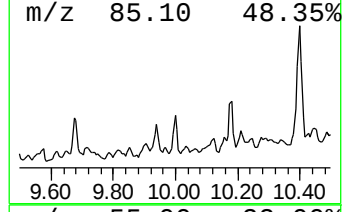
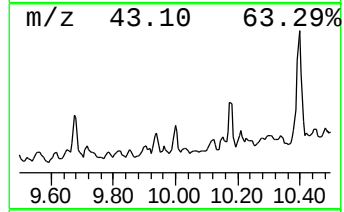
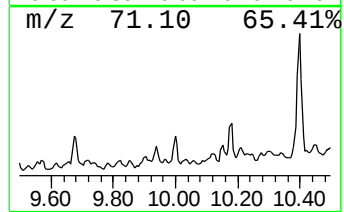
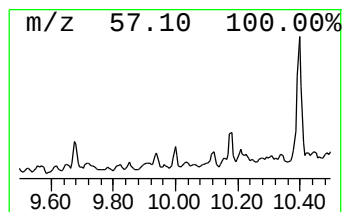
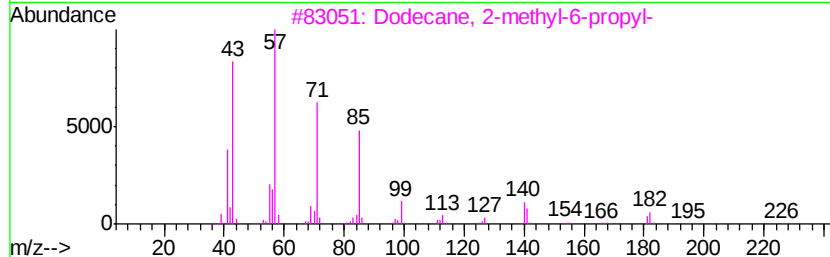
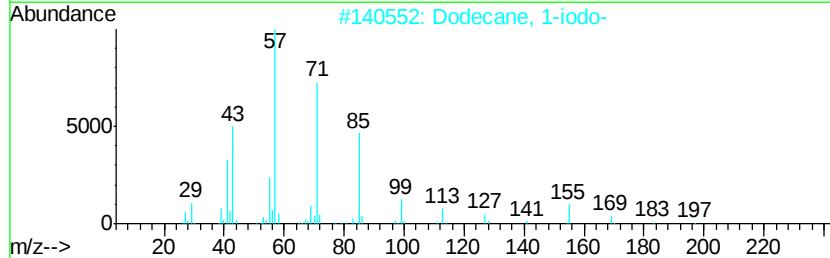
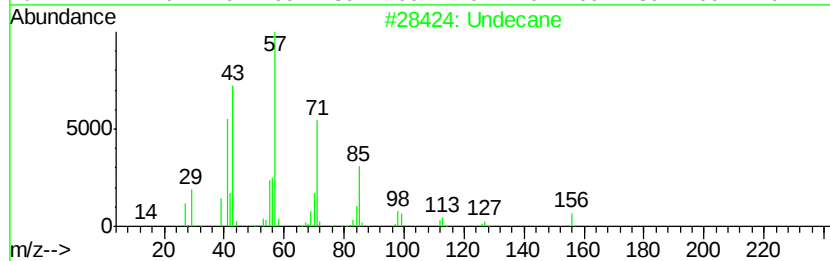
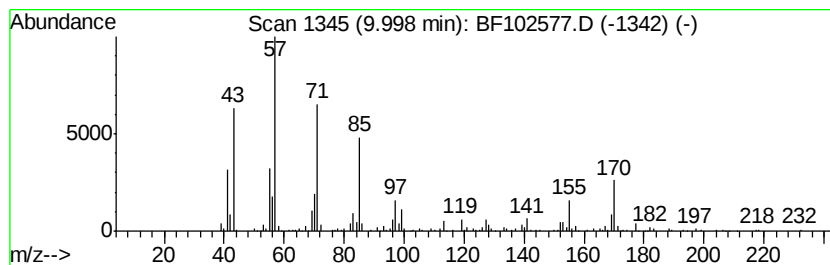
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Undecane Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T. |
|-------|---------|--------|------------------|------|
| 10.00 | 6.57 ng | 357140 | Acenaphthene-d10 | 9.75 |

| Hit# | of                                  | 5 | Tentative ID | MW  | MolForm   | CAS#         | Qual |
|------|-------------------------------------|---|--------------|-----|-----------|--------------|------|
| 1    | Undecane                            |   |              | 156 | C11H24    | 001120-21-4  | 58   |
| 2    | Dodecane, 1-iodo-                   |   |              | 296 | C12H25I   | 004292-19-7  | 53   |
| 3    | Dodecane, 2-methyl-6-propyl-        |   |              | 226 | C16H34    | 055045-08-4  | 49   |
| 4    | Sulfurous acid, 2-ethylhexyl iso... |   |              | 278 | C14H30O3S | 1000309-19-0 | 47   |
| 5    | Undecane, 2,10-dimethyl-            |   |              | 184 | C13H28    | 017301-27-8  | 47   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012818\  
Data File : BF102577.D  
Acq On : 29 Jan 2018 00:12  
Operator : SJ/JU  
Sample : J1257-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

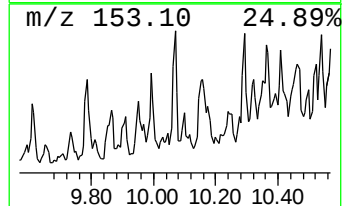
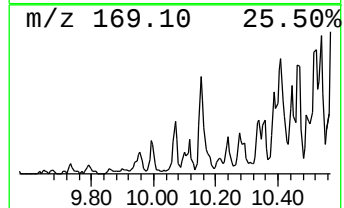
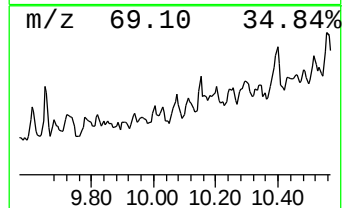
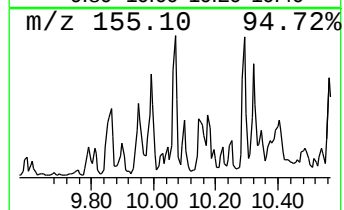
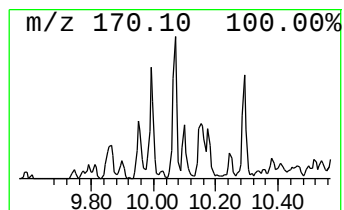
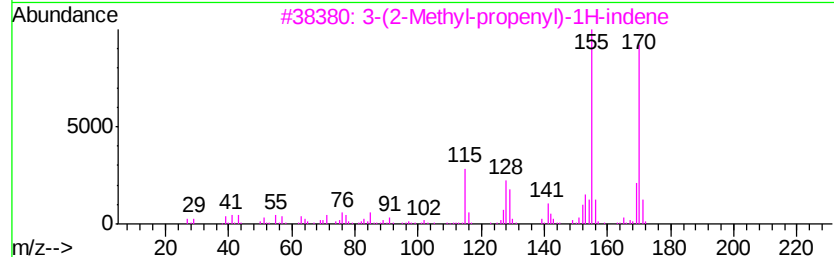
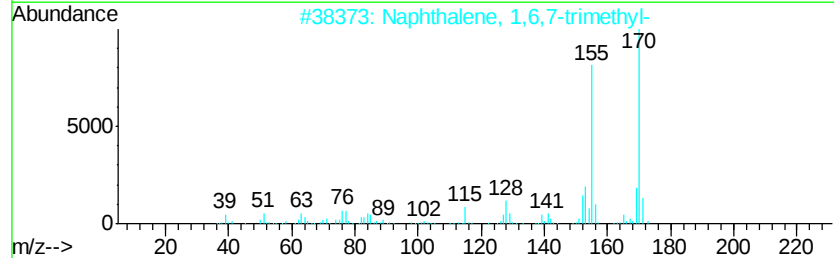
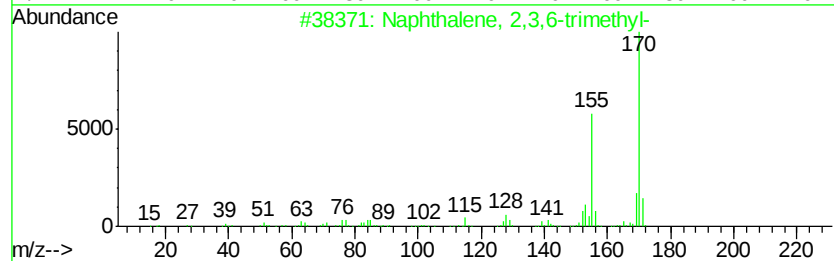
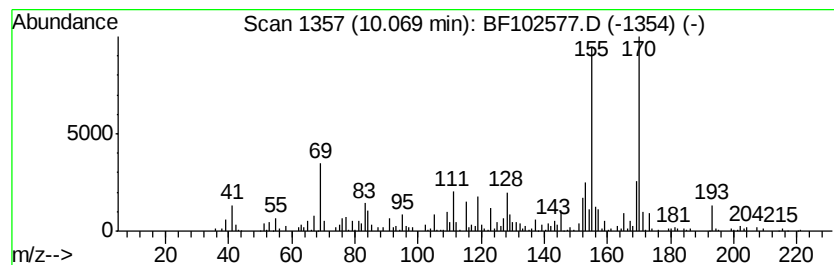
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 Naphthalene, 2,3,6-trimethyl- Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T. |
|-------|---------|--------|------------------|------|
| 10.07 | 5.80 ng | 315356 | Acenaphthene-d10 | 9.75 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Naphthalene, 2,3,6-trimethyl-       | 170 | C13H14  | 000829-26-5  | 90   |
| 2    |    | Naphthalene, 1,6,7-trimethyl-       | 170 | C13H14  | 002245-38-7  | 81   |
| 3    |    | 3-(2-Methyl-propenyl)-1H-indene     | 170 | C13H14  | 1000187-78-5 | 76   |
| 4    |    | Naphthalene, 1,4,6-trimethyl-       | 170 | C13H14  | 002131-42-2  | 68   |
| 5    |    | 1-(2,2-Dimethylcyclopropyl)-2-ph... | 170 | C13H14  | 1000294-91-1 | 64   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012818\  
Data File : BF102577.D  
Acq On : 29 Jan 2018 00:12  
Operator : SJ/JU  
Sample : J1257-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

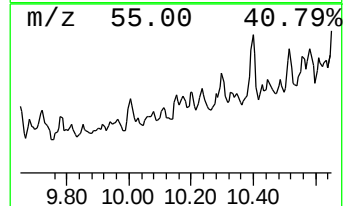
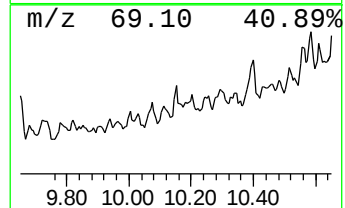
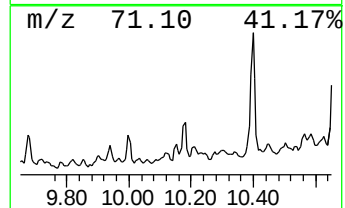
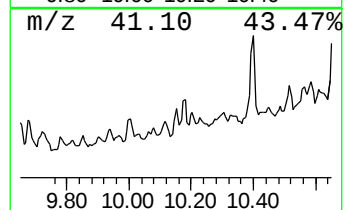
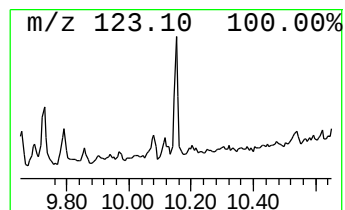
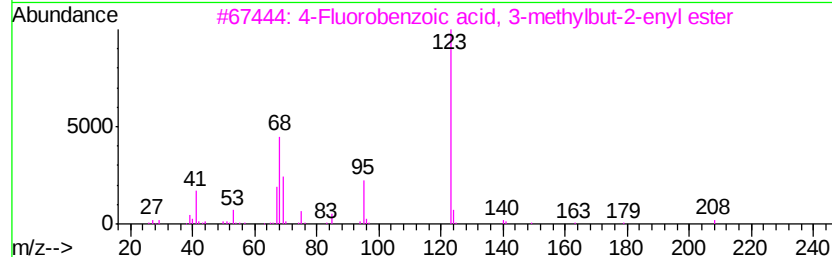
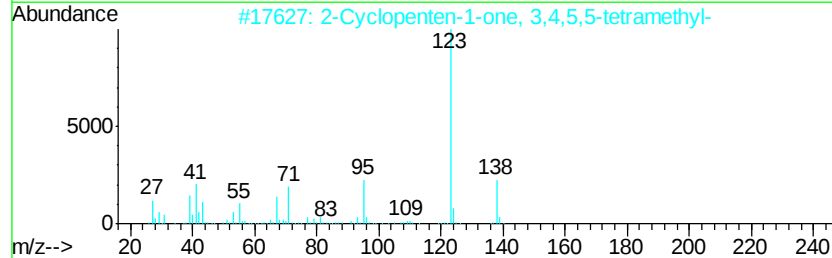
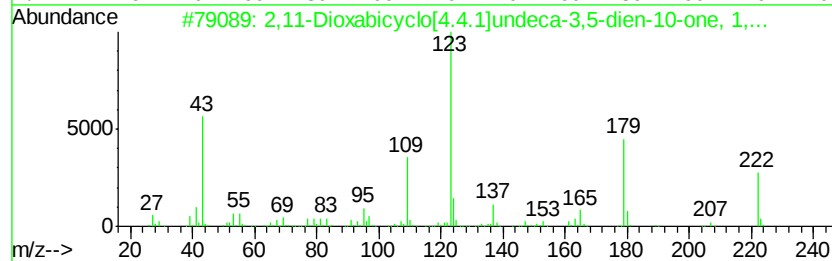
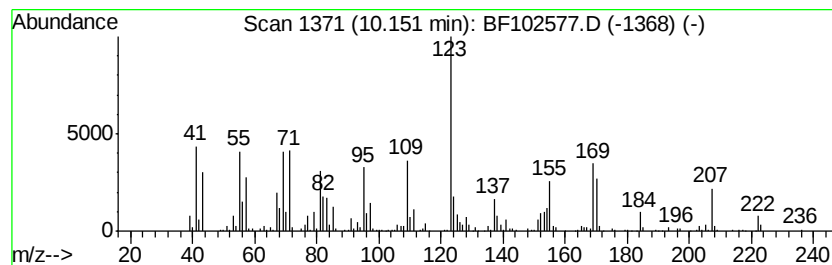
Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 16 unknown10.15 Concentration Rank 11

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 10.15     | 6.59 ng                             | 357812 | Acenaphthene-d10 | 9.75         |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 2,11-Dioxabicyclo[4.4.1]undeca-3... | 222    | C13H18O3         | 070412-52-1  | 35   |
| 2         | 2-Cyclopenten-1-one, 3,4,5,5-tet... | 138    | C9H14O           | 090611-02-2  | 35   |
| 3         | 4-Fluorobenzoic acid, 3-methylbu... | 208    | C12H13FO2        | 1000299-15-1 | 30   |
| 4         | 1-Propanone, 1-(4-fluorophenyl)-    | 152    | C9H9FO           | 000456-03-1  | 30   |
| 5         | 4,4-Dimethoxy-2,5-cyclohexadien...  | 154    | C8H10O3          | 000935-50-2  | 30   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012818\  
Data File : BF102577.D  
Acq On : 29 Jan 2018 00:12  
Operator : SJ/JU  
Sample : J1257-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

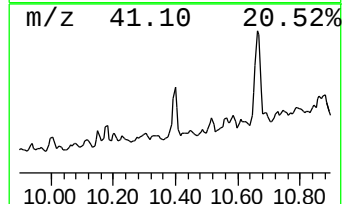
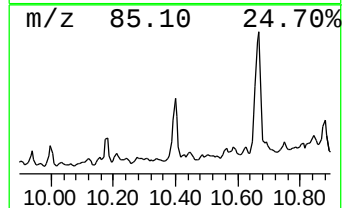
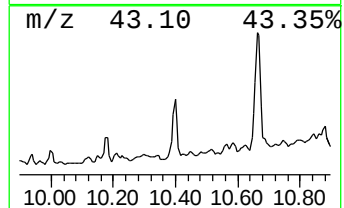
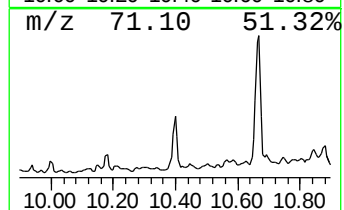
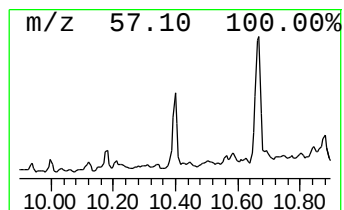
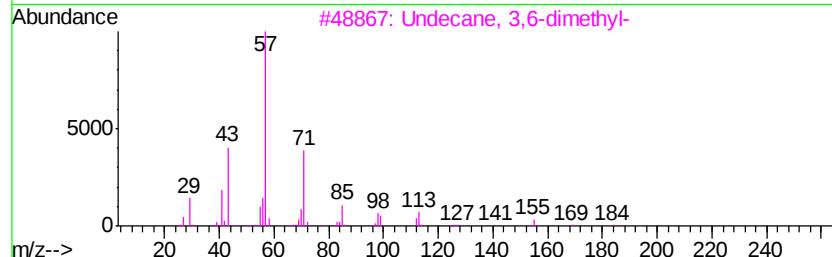
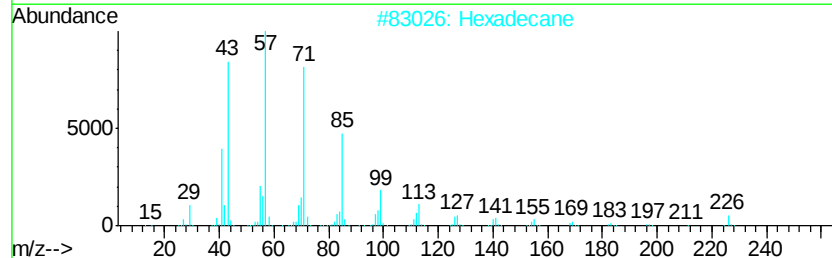
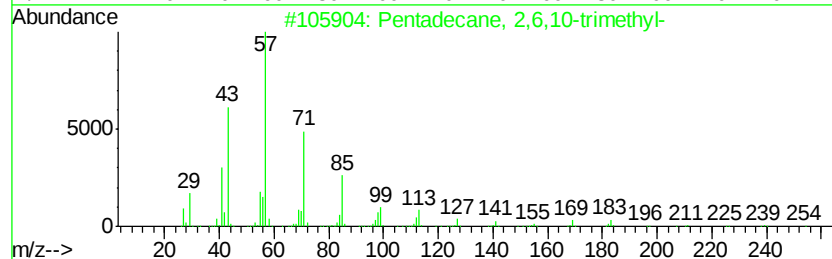
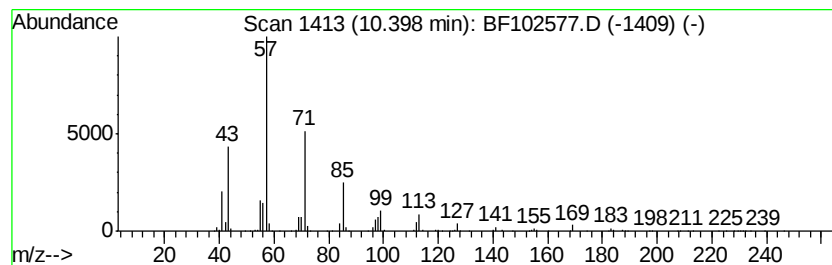
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 Pentadecane, 2,6,10-trimethyl- Concentration Rank 6

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T. |
|-------|----------|--------|------------------|------|
| 10.40 | 16.38 ng | 890038 | Acenaphthene-d10 | 9.75 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Pentadecane, 2,6,10-trimethyl- | 254 | C18H38  | 003892-00-0 | 90   |
| 2    |    | Hexadecane                     | 226 | C16H34  | 000544-76-3 | 86   |
| 3    |    | Undecane, 3,6-dimethyl-        | 184 | C13H28  | 017301-28-9 | 76   |
| 4    |    | Dodecane, 2,7,10-trimethyl-    | 212 | C15H32  | 074645-98-0 | 64   |
| 5    |    | Heptacosane                    | 380 | C27H56  | 000593-49-7 | 64   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF012818\  
Data File : BF102577.D  
Acq On : 29 Jan 2018 00:12  
Operator : SJ/JU  
Sample : J1257-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

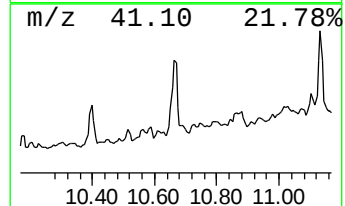
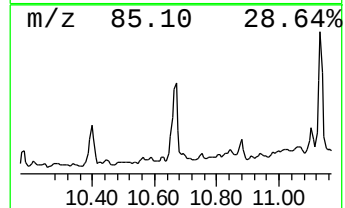
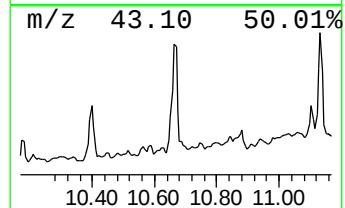
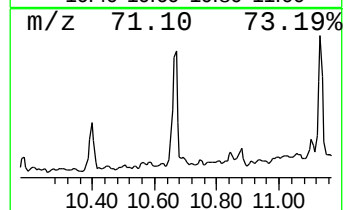
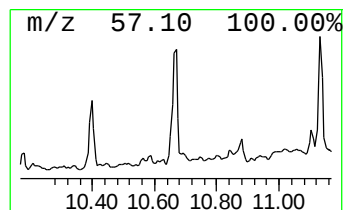
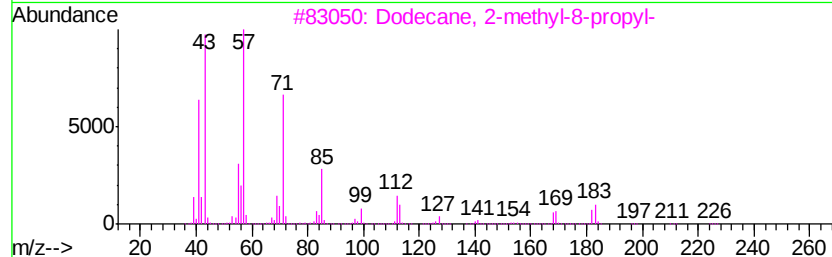
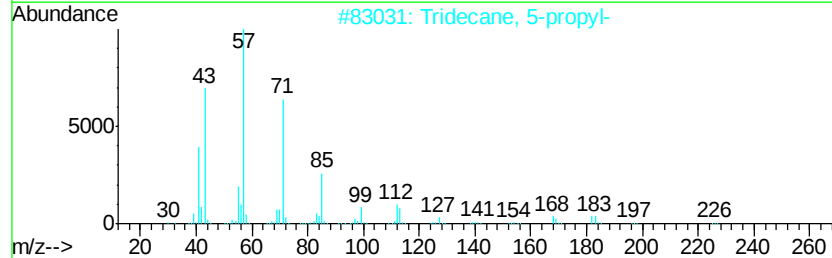
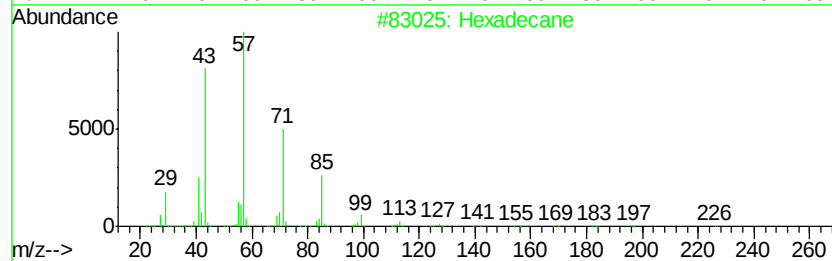
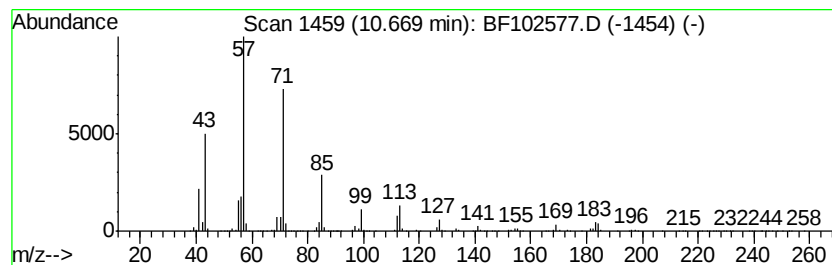
Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF012218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 18 Hexadecane Concentration Rank 2

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 10.67     | 46.47 ng                            | 2251750 | Phenanthrene-d10 | 11.24       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | Hexadecane                          | 226     | C16H34           | 000544-76-3 | 93   |
| 2         | Tridecane, 5-propyl-                | 226     | C16H34           | 055045-11-9 | 86   |
| 3         | Dodecane, 2-methyl-8-propyl-        | 226     | C16H34           | 055045-07-3 | 83   |
| 4         | Pentadecane, 2,6,10,14-tetramethyl- | 268     | C19H40           | 001921-70-6 | 80   |
| 5         | Tridecane, 3-methyl-                | 198     | C14H30           | 006418-41-3 | 68   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF012818\  
Data File : BF102577.D  
Acq On : 29 Jan 2018 00:12  
Operator : SJ/JU  
Sample : J1257-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF012218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

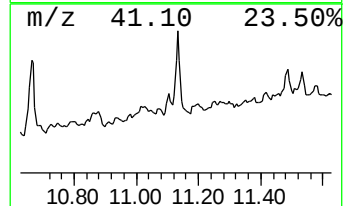
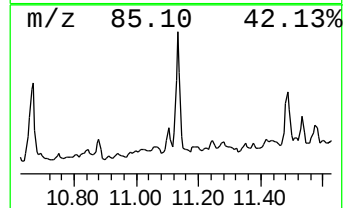
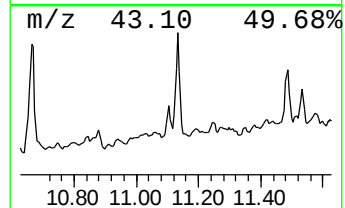
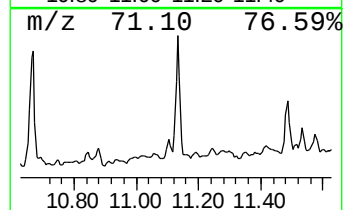
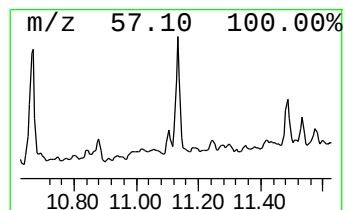
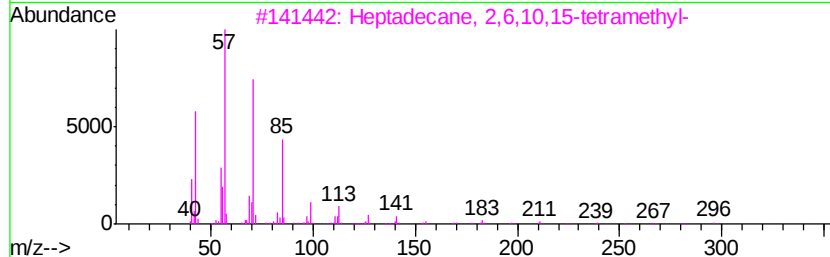
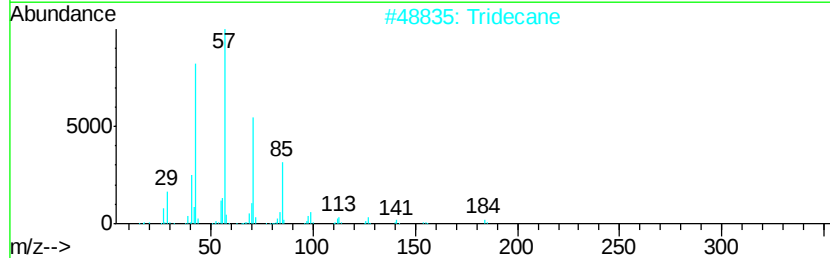
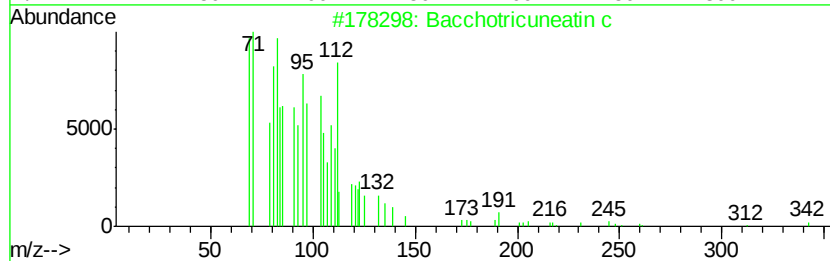
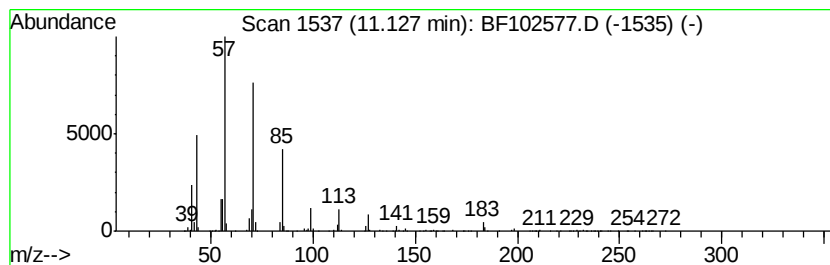
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Bacchotricuneatin c Concentration Rank 4

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 11.13 | 38.19 ng | 1850920 | Phenanthrene-d10 | 11.24 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Bacchotricuneatin c                 | 342 | C20H22O5 | 066563-30-2 | 92   |
| 2    |    | Tridecane                           | 184 | C13H28   | 000629-50-5 | 91   |
| 3    |    | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44   | 054833-48-6 | 86   |
| 4    |    | Hexacosane                          | 366 | C26H54   | 000630-01-3 | 83   |
| 5    |    | Tritetracontane                     | 605 | C43H88   | 007098-21-7 | 78   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012818\  
Data File : BF102577.D  
Acq On : 29 Jan 2018 00:12  
Operator : SJ/JU  
Sample : J1257-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC-17(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

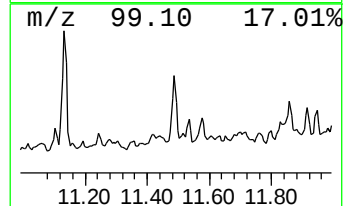
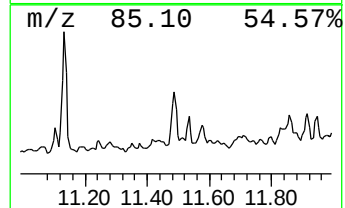
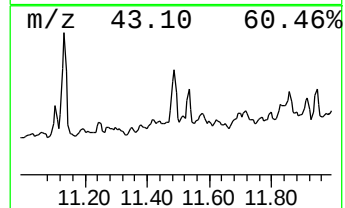
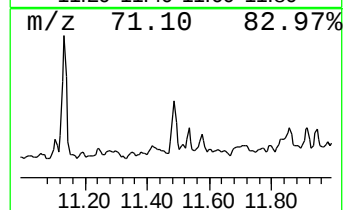
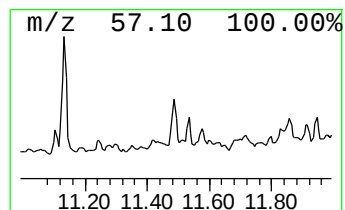
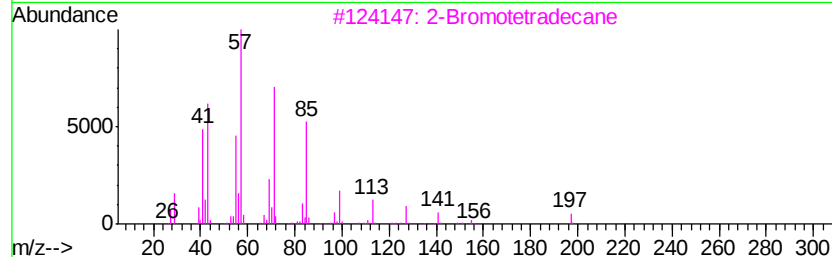
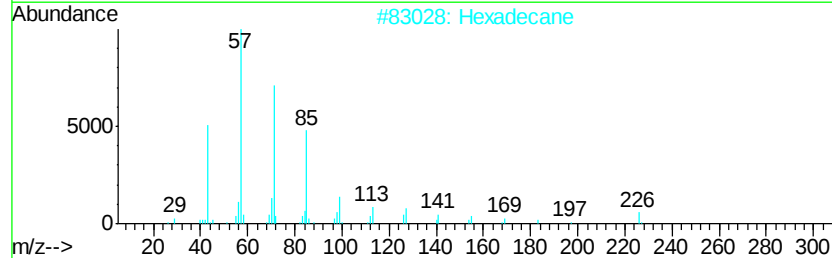
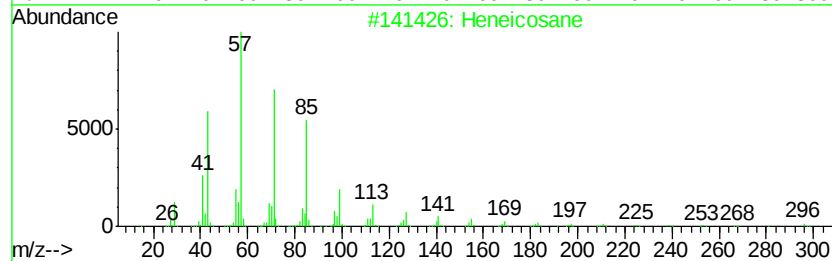
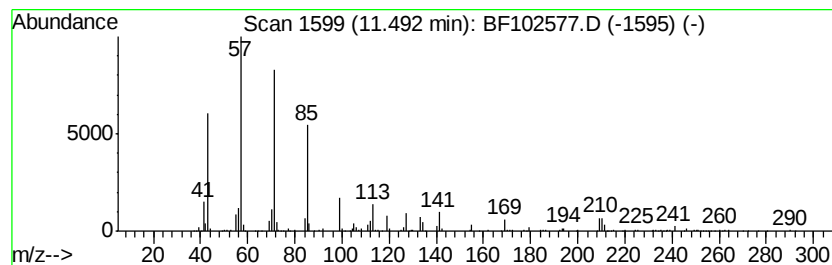
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Heneicosane Concentration Rank 7

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 11.49 | 16.21 ng | 785463 | Phenanthrene-d10 | 11.24 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Heneicosane                         | 296 | C21H44   | 000629-94-7 | 87   |
| 2    |    | Hexadecane                          | 226 | C16H34   | 000544-76-3 | 87   |
| 3    |    | 2-Bromotetradecane                  | 276 | C14H29Br | 074036-95-6 | 87   |
| 4    |    | Octadecane                          | 254 | C18H38   | 000593-45-3 | 86   |
| 5    |    | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44   | 054833-48-6 | 86   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF012818\  
 Data File : BF102577.D  
 Acq On : 29 Jan 2018 00:12  
 Operator : SJ/JU  
 Sample : J1257-13  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 WC-17(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF012218.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| 2-Pentanone, 4-hy... | 4.90  | 20.4    | ng    | 1087470  | 1                     | 6.72  | 1067230 | 20.0 |
| Benzene, 1-ethyl-... | 6.23  | 5.6     | ng    | 296425   | 1                     | 6.72  | 1067230 | 20.0 |
| Mesitylene           | 6.31  | 4.6     | ng    | 243248   | 1                     | 6.72  | 1067230 | 20.0 |
| unknown6.49          | 6.49  | 41.7    | ng    | 2226290  | 1                     | 6.72  | 1067230 | 20.0 |
| Benzene, 1,2,4-tr... | 6.54  | 9.2     | ng    | 491070   | 1                     | 6.72  | 1067230 | 20.0 |
| Benzene, 1-methyl... | 6.99  | 6.5     | ng    | 346177   | 1                     | 6.72  | 1067230 | 20.0 |
| Benzene, 1,2,3,4-... | 7.03  | 11.2    | ng    | 597979   | 1                     | 6.72  | 1067230 | 20.0 |
| Benzene, 1,2,4,5-... | 7.48  | 4.8     | ng    | 451333   | 2                     | 8.00  | 1901710 | 20.0 |
| Sulfurous acid, b... | 9.15  | 7.0     | ng    | 379685   | 3                     | 9.75  | 1086620 | 20.0 |
| unknown9.61          | 9.61  | 5.7     | ng    | 307582   | 3                     | 9.75  | 1086620 | 20.0 |
| unknown9.65          | 9.65  | 4.5     | ng    | 247223   | 3                     | 9.75  | 1086620 | 20.0 |
| Undecane             | 10.00 | 6.6     | ng    | 357140   | 3                     | 9.75  | 1086620 | 20.0 |
| Naphthalene, 2,3,... | 10.07 | 5.8     | ng    | 315356   | 3                     | 9.75  | 1086620 | 20.0 |
| unknown10.15         | 10.15 | 6.6     | ng    | 357812   | 3                     | 9.75  | 1086620 | 20.0 |
| Pentadecane, 2,6,... | 10.40 | 16.4    | ng    | 890038   | 3                     | 9.75  | 1086620 | 20.0 |
| Hexadecane           | 10.67 | 46.5    | ng    | 2251750  | 4                     | 11.24 | 969216  | 20.0 |
| Bacchotricuneatin c  | 11.13 | 38.2    | ng    | 1850920  | 4                     | 11.24 | 969216  | 20.0 |
| Heneicosane          | 11.49 | 16.2    | ng    | 785463   | 4                     | 11.24 | 969216  | 20.0 |