

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
 Data File : BP007961.D  
 Acq On : 12 Nov 2021 00:01  
 Operator : CG/JU  
 Sample : M4630-07  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SB-13N-16.0-16.5

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:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP007961.D\data.ms

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 4.799       | 247           | 257         | 262          | rBV      | 191913         | 295263        | 2.61%           | 0.170%        |
| 2         | 4.893       | 268           | 273         | 277          | rBV      | 139706         | 224549        | 1.98%           | 0.130%        |
| 3         | 4.952       | 278           | 283         | 287          | rVV5     | 98215          | 205529        | 1.82%           | 0.119%        |
| 4         | 4.999       | 287           | 291         | 296          | rVV      | 257376         | 369736        | 3.27%           | 0.213%        |
| 5         | 5.052       | 296           | 300         | 306          | rVV      | 447617         | 723876        | 6.40%           | 0.418%        |
| 6         | 5.216       | 324           | 328         | 332          | rVV      | 241701         | 337325        | 2.98%           | 0.195%        |
| 7         | 5.305       | 337           | 343         | 346          | rVV3     | 562504         | 1202421       | 10.62%          | 0.694%        |
| 8         | 5.334       | 346           | 348         | 356          | rVB2     | 406299         | 509104        | 4.50%           | 0.294%        |
| 9         | 5.434       | 356           | 365         | 372          | rBV2     | 2010227        | 3321473       | 29.34%          | 1.917%        |
| 10        | 5.546       | 381           | 384         | 394          | rVB3     | 140295         | 304022        | 2.69%           | 0.175%        |
| 11        | 5.657       | 394           | 403         | 409          | rBV2     | 291051         | 529883        | 4.68%           | 0.306%        |
| 12        | 5.734       | 409           | 416         | 420          | rBV3     | 518477         | 1016111       | 8.98%           | 0.586%        |
| 13        | 5.810       | 424           | 429         | 434          | rVV      | 951201         | 1469840       | 12.99%          | 0.848%        |
| 14        | 5.875       | 434           | 440         | 448          | rVB      | 3362834        | 5322086       | 47.02%          | 3.071%        |
| 15        | 5.952       | 448           | 453         | 462          | rBV2     | 359530         | 594886        | 5.26%           | 0.343%        |
| 16        | 6.034       | 462           | 467         | 475          | rVB2     | 158326         | 276918        | 2.45%           | 0.160%        |
| 17        | 6.128       | 475           | 483         | 490          | rBV4     | 1261647        | 2881578       | 25.46%          | 1.663%        |
| 18        | 6.252       | 497           | 504         | 510          | rVB2     | 651701         | 1385954       | 12.24%          | 0.800%        |
| 19        | 6.346       | 514           | 520         | 527          | rBV3     | 1202699        | 2687827       | 23.75%          | 1.551%        |
| 20        | 6.416       | 527           | 532         | 536          | rBV      | 2222534        | 3276418       | 28.95%          | 1.891%        |
| 21        | 6.493       | 536           | 545         | 548          | rVV3     | 2062267        | 4027648       | 35.58%          | 2.324%        |
| 22        | 6.528       | 548           | 551         | 557          | rVB      | 1499197        | 2195739       | 19.40%          | 1.267%        |
| 23        | 6.610       | 561           | 565         | 570          | rVB3     | 446269         | 716721        | 6.33%           | 0.414%        |
| 24        | 6.675       | 570           | 576         | 582          | rVB4     | 507487         | 918515        | 8.11%           | 0.530%        |
| 25        | 6.757       | 582           | 590         | 595          | rVB2     | 973839         | 2387216       | 21.09%          | 1.378%        |
| 26        | 6.822       | 595           | 601         | 602          | rBV      | 911486         | 1459412       | 12.89%          | 0.842%        |
| 27        | 6.852       | 602           | 606         | 611          | rVB      | 2229591        | 3474572       | 30.70%          | 2.005%        |
| 28        | 6.904       | 611           | 615         | 619          | rBV      | 1343021        | 1984902       | 17.54%          | 1.145%        |
| 29        | 6.946       | 619           | 622         | 626          | rVB2     | 826552         | 1097572       | 9.70%           | 0.633%        |
| 30        | 7.016       | 626           | 634         | 640          | rVB4     | 2642320        | 6399990       | 56.54%          | 3.693%        |
| 31        | 7.075       | 640           | 644         | 650          | rVB      | 1072505        | 1519787       | 13.43%          | 0.877%        |
| 32        | 7.134       | 650           | 654         | 657          | rBV      | 194015         | 279726        | 2.47%           | 0.161%        |
| 33        | 7.181       | 657           | 662         | 667          | rVB3     | 526309         | 1009081       | 8.92%           | 0.582%        |
| 34        | 7.240       | 667           | 672         | 676          | rBV3     | 945733         | 1471964       | 13.00%          | 0.849%        |
| 35        | 7.275       | 676           | 678         | 680          | rBV2     | 209180         | 220919        | 1.95%           | 0.127%        |
| 36        | 7.310       | 680           | 684         | 686          | rBV      | 474649         | 633654        | 5.60%           | 0.366%        |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
 Data File : BP007961.D  
 Acq On : 12 Nov 2021 00:01  
 Operator : CG/JU  
 Sample : M4630-07  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SB-13N-16.0-16.5

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:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |          |         |        |
|----|--------|------|------|------|------|---------|----------|---------|--------|
| 37 | 7.340  | 686  | 689  | 694  | rVV  | 1150057 | 1943701  | 17.17%  | 1.122% |
| 38 | 7.416  | 698  | 702  | 708  | rVB3 | 651471  | 1230406  | 10.87%  | 0.710% |
| 39 | 7.487  | 708  | 714  | 721  | rBV2 | 6009099 | 11318792 | 100.00% | 6.532% |
| 40 | 7.575  | 726  | 729  | 735  | rVB3 | 558249  | 920439   | 8.13%   | 0.531% |
| 41 | 7.628  | 735  | 738  | 741  | rBV  | 226207  | 254792   | 2.25%   | 0.147% |
| 42 | 7.675  | 741  | 746  | 747  | rBV3 | 491134  | 768066   | 6.79%   | 0.443% |
| 43 | 7.757  | 755  | 760  | 761  | rBV2 | 710950  | 974261   | 8.61%   | 0.562% |
| 44 | 7.840  | 768  | 774  | 779  | rVB  | 4004666 | 6748029  | 59.62%  | 3.894% |
| 45 | 7.893  | 779  | 783  | 784  | rBV  | 347458  | 395652   | 3.50%   | 0.228% |
| 46 | 7.928  | 784  | 789  | 792  | rVV3 | 753667  | 1495423  | 13.21%  | 0.863% |
| 47 | 7.969  | 792  | 796  | 803  | rVV3 | 1074286 | 2156527  | 19.05%  | 1.245% |
| 48 | 8.046  | 804  | 809  | 813  | rVV2 | 1207796 | 1936192  | 17.11%  | 1.117% |
| 49 | 8.087  | 813  | 816  | 818  | rVV2 | 281630  | 372334   | 3.29%   | 0.215% |
| 50 | 8.146  | 818  | 826  | 832  | rVV3 | 1988056 | 4505099  | 39.80%  | 2.600% |
| 51 | 8.204  | 832  | 836  | 844  | rVB3 | 673363  | 1220214  | 10.78%  | 0.704% |
| 52 | 8.293  | 847  | 851  | 857  | rBV3 | 494607  | 832062   | 7.35%   | 0.480% |
| 53 | 8.393  | 857  | 868  | 874  | rBV5 | 1836514 | 4479709  | 39.58%  | 2.585% |
| 54 | 8.451  | 874  | 878  | 881  | rVB  | 1052723 | 1280534  | 11.31%  | 0.739% |
| 55 | 8.493  | 881  | 885  | 887  | rBV2 | 1111094 | 1515466  | 13.39%  | 0.875% |
| 56 | 8.522  | 887  | 890  | 893  | rVB  | 1096430 | 1355850  | 11.98%  | 0.782% |
| 57 | 8.628  | 902  | 908  | 911  | rBV  | 1396040 | 2326793  | 20.56%  | 1.343% |
| 58 | 8.681  | 914  | 917  | 922  | rVB  | 881796  | 1360660  | 12.02%  | 0.785% |
| 59 | 8.804  | 935  | 938  | 942  | rBV  | 454904  | 708954   | 6.26%   | 0.409% |
| 60 | 8.857  | 942  | 947  | 954  | rVB  | 702265  | 1361143  | 12.03%  | 0.785% |
| 61 | 8.963  | 954  | 965  | 969  | rBV4 | 2180709 | 5315294  | 46.96%  | 3.067% |
| 62 | 9.004  | 969  | 972  | 976  | rVV  | 966236  | 1747167  | 15.44%  | 1.008% |
| 63 | 9.057  | 976  | 981  | 983  | rVV4 | 1048000 | 1990037  | 17.58%  | 1.148% |
| 64 | 9.093  | 983  | 987  | 996  | rVB  | 4975671 | 8286978  | 73.21%  | 4.782% |
| 65 | 9.181  | 996  | 1002 | 1003 | rBV3 | 803679  | 1331485  | 11.76%  | 0.768% |
| 66 | 9.210  | 1003 | 1007 | 1013 | rVB3 | 1323177 | 2725993  | 24.08%  | 1.573% |
| 67 | 9.293  | 1013 | 1021 | 1023 | rBV3 | 550393  | 1116773  | 9.87%   | 0.644% |
| 68 | 9.346  | 1024 | 1030 | 1036 | rVB4 | 824245  | 1779323  | 15.72%  | 1.027% |
| 69 | 9.504  | 1050 | 1057 | 1060 | rBV2 | 688174  | 1404832  | 12.41%  | 0.811% |
| 70 | 9.540  | 1060 | 1063 | 1066 | rVV  | 385413  | 487610   | 4.31%   | 0.281% |
| 71 | 9.745  | 1091 | 1098 | 1100 | rBV3 | 383423  | 722642   | 6.38%   | 0.417% |
| 72 | 9.781  | 1100 | 1104 | 1109 | rVV2 | 995783  | 1781967  | 15.74%  | 1.028% |
| 73 | 9.834  | 1109 | 1113 | 1121 | rVB6 | 875255  | 1840428  | 16.26%  | 1.062% |
| 74 | 10.004 | 1135 | 1142 | 1146 | rBV3 | 663331  | 1057497  | 9.34%   | 0.610% |
| 75 | 10.057 | 1146 | 1151 | 1154 | rBV3 | 781904  | 1357711  | 12.00%  | 0.784% |
| 76 | 10.193 | 1169 | 1174 | 1178 | rBV2 | 396166  | 698321   | 6.17%   | 0.403% |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
 Data File : BP007961.D  
 Acq On : 12 Nov 2021 00:01  
 Operator : CG/JU  
 Sample : M4630-07  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SB-13N-16.0-16.5

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:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|     |        |      |      |      |      |         |         |        |        |
|-----|--------|------|------|------|------|---------|---------|--------|--------|
| 77  | 10.240 | 1179 | 1182 | 1186 | rVV5 | 272574  | 469103  | 4.14%  | 0.271% |
| 78  | 10.281 | 1187 | 1189 | 1195 | rVB  | 244551  | 300112  | 2.65%  | 0.173% |
| 79  | 10.357 | 1196 | 1202 | 1205 | rBV  | 274410  | 461629  | 4.08%  | 0.266% |
| 80  | 10.434 | 1211 | 1215 | 1220 | rVB4 | 187973  | 322556  | 2.85%  | 0.186% |
| 81  | 10.522 | 1225 | 1230 | 1232 | rVV3 | 212421  | 364070  | 3.22%  | 0.210% |
| 82  | 10.557 | 1232 | 1236 | 1241 | rVV4 | 250834  | 530886  | 4.69%  | 0.306% |
| 83  | 10.634 | 1242 | 1249 | 1255 | rVV2 | 1908712 | 3547278 | 31.34% | 2.047% |
| 84  | 10.698 | 1255 | 1260 | 1268 | rVV  | 600306  | 1276797 | 11.28% | 0.737% |
| 85  | 10.769 | 1268 | 1272 | 1277 | rVV4 | 202798  | 412647  | 3.65%  | 0.238% |
| 86  | 10.822 | 1277 | 1281 | 1286 | rVV  | 561308  | 811622  | 7.17%  | 0.468% |
| 87  | 10.875 | 1286 | 1290 | 1298 | rVV4 | 130954  | 259113  | 2.29%  | 0.150% |
| 88  | 10.951 | 1298 | 1303 | 1308 | rVB4 | 100472  | 206319  | 1.82%  | 0.119% |
| 89  | 11.357 | 1361 | 1372 | 1380 | rBV7 | 135083  | 373605  | 3.30%  | 0.216% |
| 90  | 11.687 | 1422 | 1428 | 1437 | rBV4 | 134636  | 277864  | 2.45%  | 0.160% |
| 91  | 12.081 | 1489 | 1495 | 1505 | rBV  | 147810  | 230193  | 2.03%  | 0.133% |
| 92  | 12.304 | 1524 | 1533 | 1541 | rVB2 | 114172  | 225469  | 1.99%  | 0.130% |
| 93  | 13.110 | 1663 | 1670 | 1682 | rBV  | 2108800 | 3108653 | 27.46% | 1.794% |
| 94  | 14.481 | 1896 | 1903 | 1911 | rVB  | 683318  | 1015617 | 8.97%  | 0.586% |
| 95  | 15.975 | 2151 | 2157 | 2173 | rBV  | 2142289 | 3062380 | 27.06% | 1.767% |
| 96  | 17.239 | 2365 | 2372 | 2378 | rBV  | 1007098 | 1412214 | 12.48% | 0.815% |
| 97  | 19.804 | 2802 | 2808 | 2814 | rBV  | 5712856 | 6648867 | 58.74% | 3.837% |
| 98  | 21.045 | 3015 | 3019 | 3025 | rBV  | 430243  | 679523  | 6.00%  | 0.392% |
| 99  | 21.333 | 3062 | 3068 | 3072 | rBV  | 1795036 | 2390763 | 21.12% | 1.380% |
| 100 | 23.739 | 3470 | 3477 | 3487 | rVB2 | 1448592 | 3059104 | 27.03% | 1.765% |

Sum of corrected areas: 173283757



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

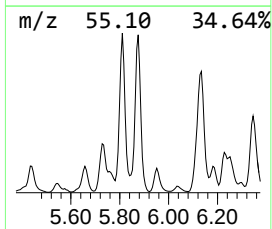
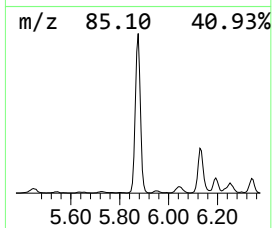
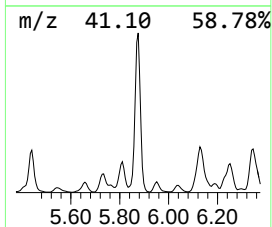
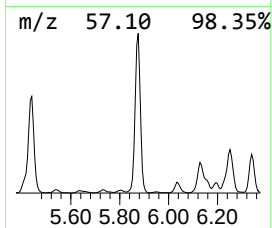
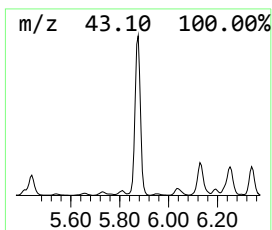
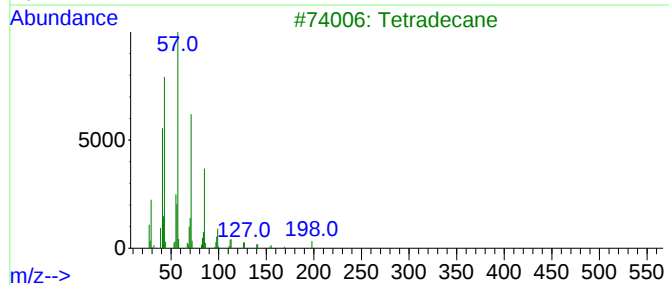
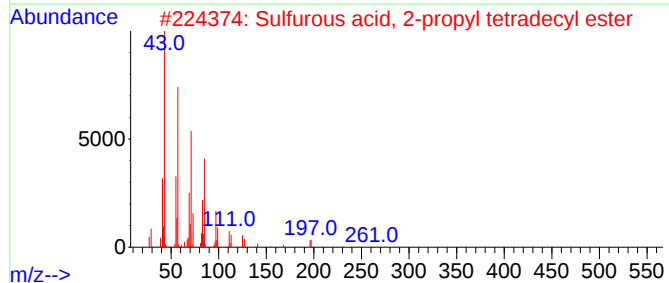
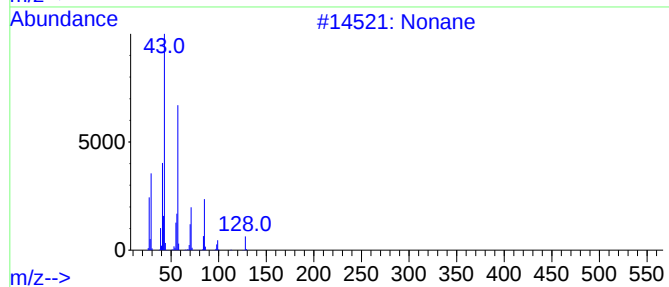
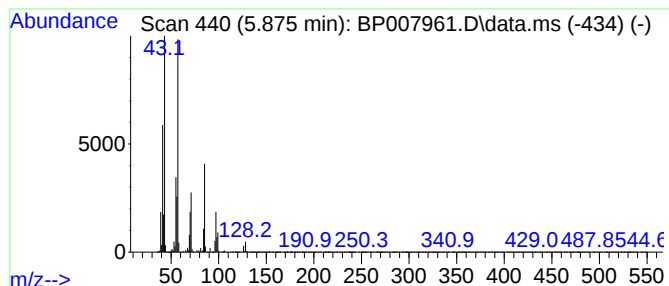
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 Nonane Concentration Rank 2

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 5.875 | 15.77 ng | 5322090 | 1,4-Dichlorobenzene-d4 | 7.851 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm   | CAS#         | Qual |
|------|----|---|--------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Nonane                               | 128 | C9H20     | 000111-84-2  | 90   |
| 2    |    |   | Sulfurous acid, 2-propyl tetradec... | 320 | C17H36O3S | 1010309-12-5 | 59   |
| 3    |    |   | Tetradecane                          | 198 | C14H30    | 000629-59-4  | 53   |
| 4    |    |   | Carbonic acid, nonyl vinyl ester     | 214 | C12H22O3  | 1000383-25-6 | 53   |
| 5    |    |   | 1-Decanol, 2-ethyl-                  | 186 | C12H26O   | 021078-65-9  | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

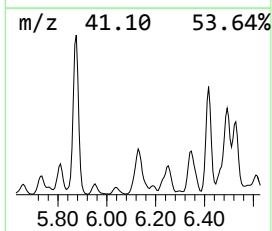
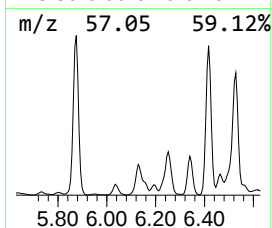
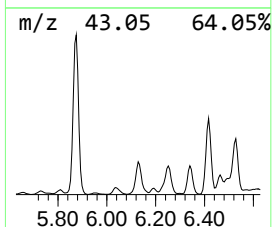
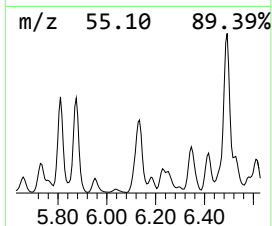
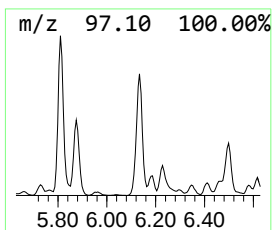
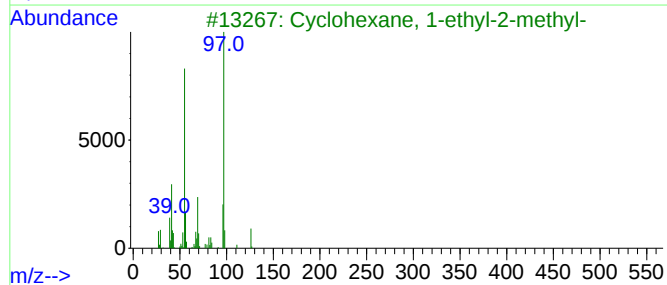
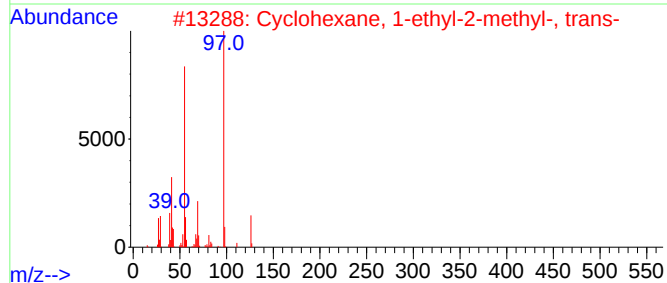
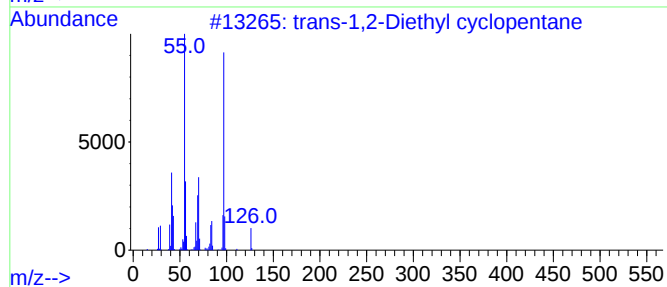
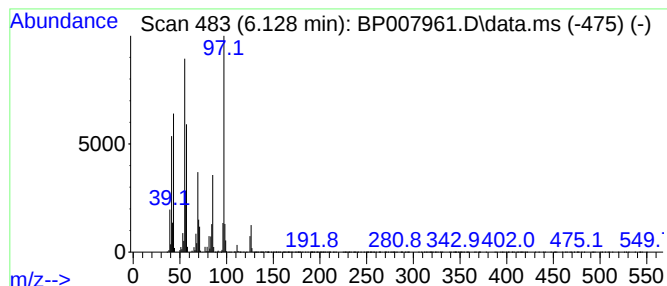
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 2 trans-1,2-Diethyl cyclopentane Concentration Rank 11

| R.T.  | EstConc | Area    | Relative to ISTD       | R.T.  |
|-------|---------|---------|------------------------|-------|
| 6.128 | 8.54 ng | 2881580 | 1,4-Dichlorobenzene-d4 | 7.851 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | trans-1,2-Diethyl cyclopentane      | 126 | C9H18   | 000932-40-1 | 87   |
| 2    |    |   | Cyclohexane, 1-ethyl-2-methyl-, ... | 126 | C9H18   | 004923-78-8 | 62   |
| 3    |    |   | Cyclohexane, 1-ethyl-2-methyl-      | 126 | C9H18   | 003728-54-9 | 58   |
| 4    |    |   | Cyclohexane, 1-methyl-2-propyl-     | 140 | C10H20  | 004291-79-6 | 53   |
| 5    |    |   | Cyclohexane, 1,3-dimethyl-, cis-    | 112 | C8H16   | 000638-04-0 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

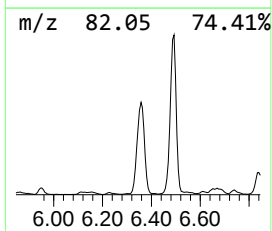
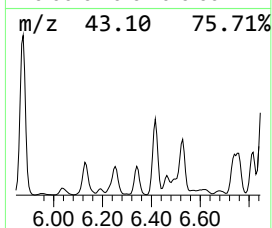
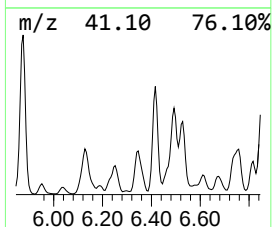
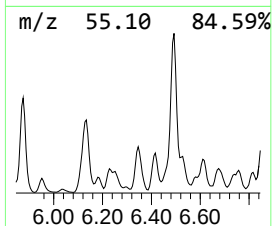
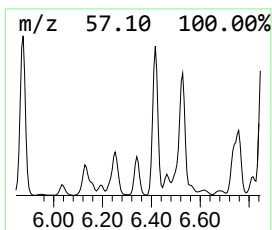
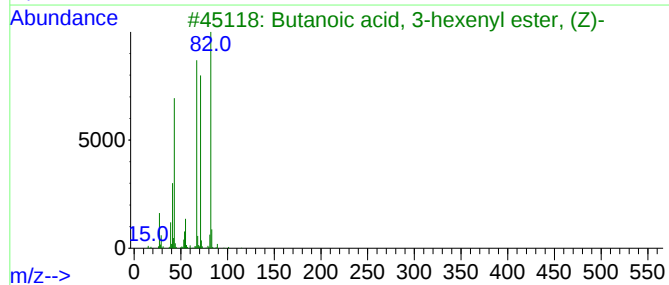
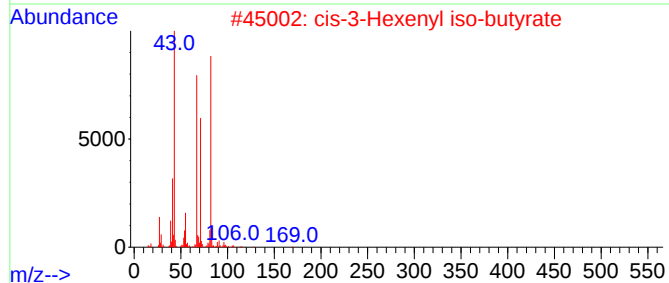
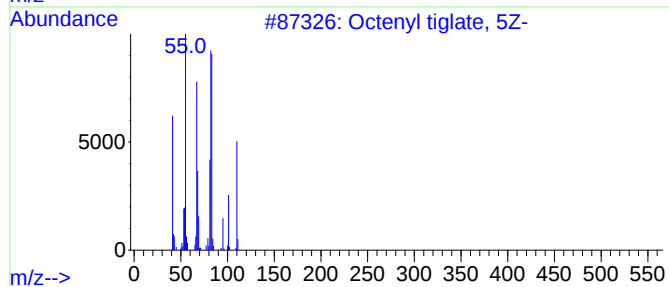
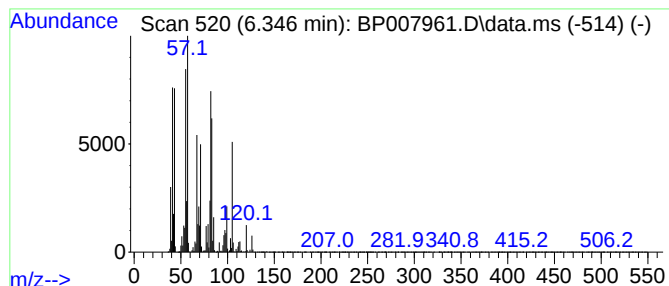
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 3 unknown6.346 Concentration Rank 13

| R.T.  | EstConc | Area    | Relative to ISTD       | R.T.  |
|-------|---------|---------|------------------------|-------|
| 6.346 | 7.97 ng | 2687830 | 1,4-Dichlorobenzene-d4 | 7.851 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Octenyl tiglate, 5Z-                | 210 | C13H22O2 | 1000383-60-1 | 43   |
| 2    |    |   | cis-3-Hexenyl iso-butyrate          | 170 | C10H18O2 | 041519-23-7  | 43   |
| 3    |    |   | Butanoic acid, 3-hexenyl ester, ... | 170 | C10H18O2 | 016491-36-4  | 43   |
| 4    |    |   | Glutaric acid, cyclohexylmethyl ... | 310 | C18H30O4 | 1000394-02-8 | 35   |
| 5    |    |   | 2,4-Hexadiene, (E,Z)-               | 82  | C6H10    | 005194-50-3  | 30   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

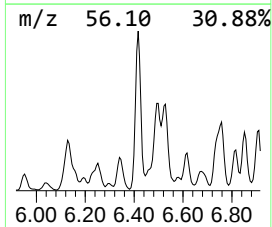
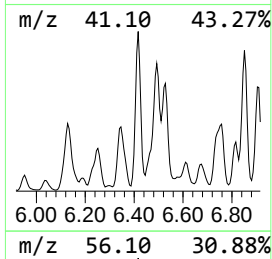
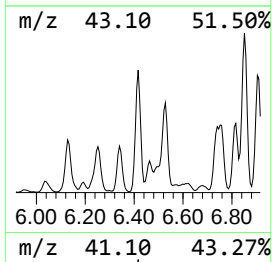
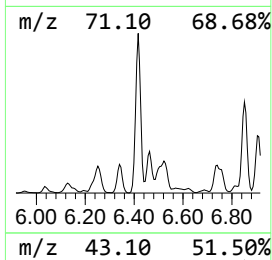
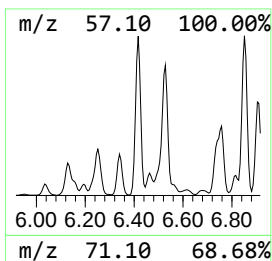
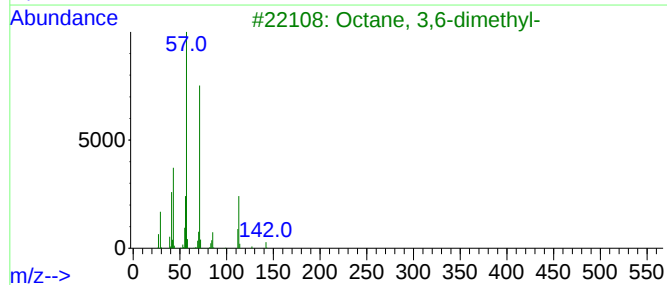
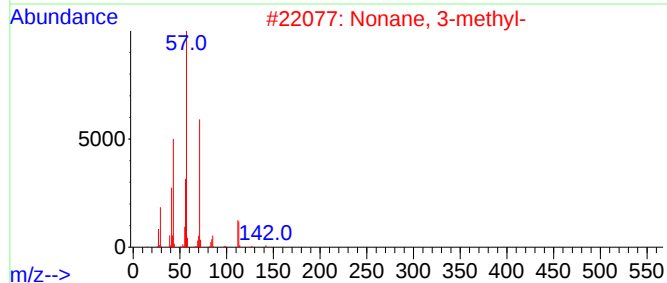
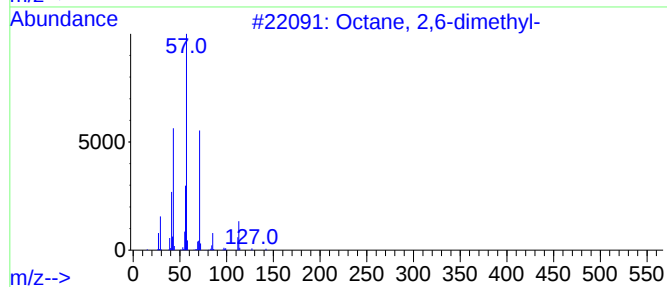
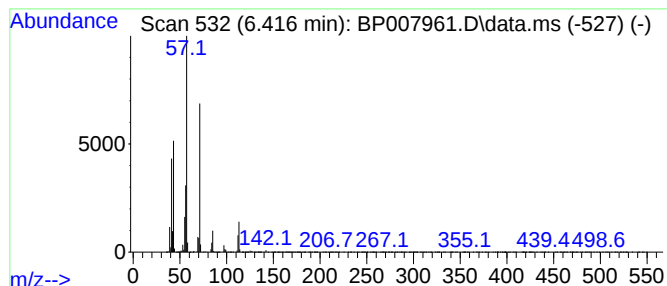
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 Octane, 2,6-dimethyl- Concentration Rank 10

| R.T.  | EstConc | Area    | Relative to ISTD       | R.T.  |
|-------|---------|---------|------------------------|-------|
| 6.416 | 9.71 ng | 3276420 | 1,4-Dichlorobenzene-d4 | 7.851 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Octane, 2,6-dimethyl-               | 142 | C10H22    | 002051-30-1  | 93   |
| 2    |    |   | Nonane, 3-methyl-                   | 142 | C10H22    | 005911-04-6  | 91   |
| 3    |    |   | Octane, 3,6-dimethyl-               | 142 | C10H22    | 015869-94-0  | 91   |
| 4    |    |   | Sulfurous acid, di(2-ethylhexyl)... | 306 | C16H34O3S | 1000309-19-1 | 59   |
| 5    |    |   | Octane, 2-iodo-                     | 240 | C8H17I    | 000557-36-8  | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

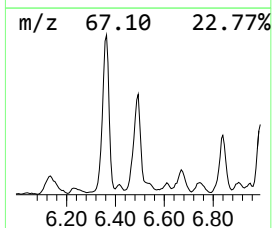
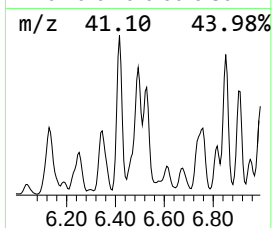
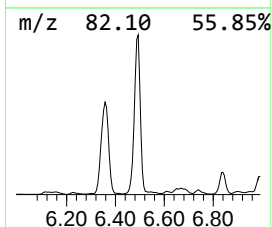
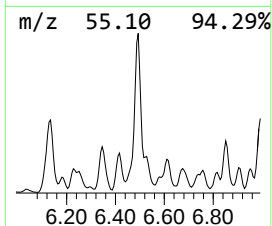
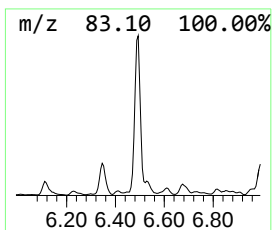
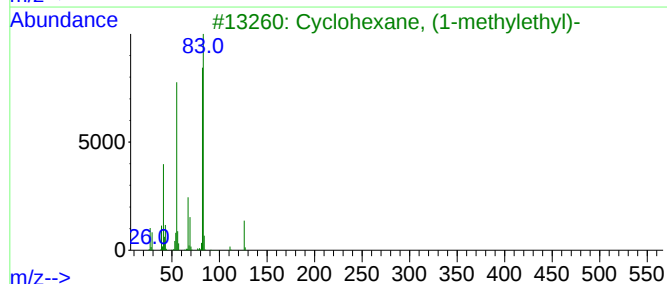
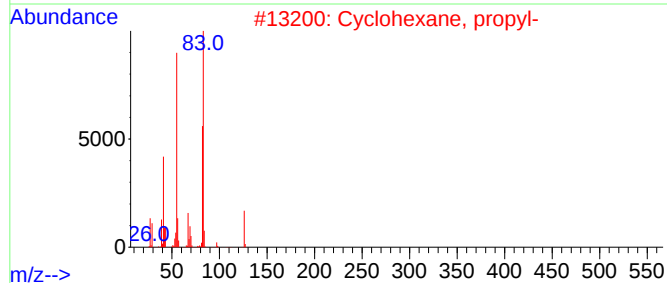
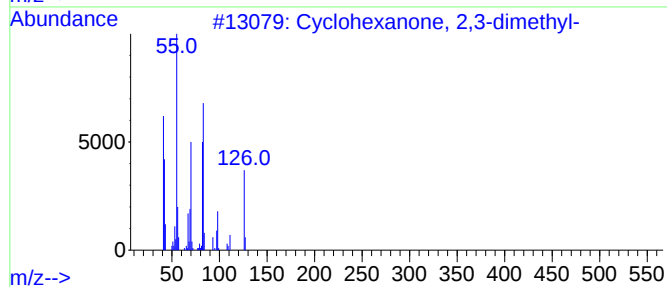
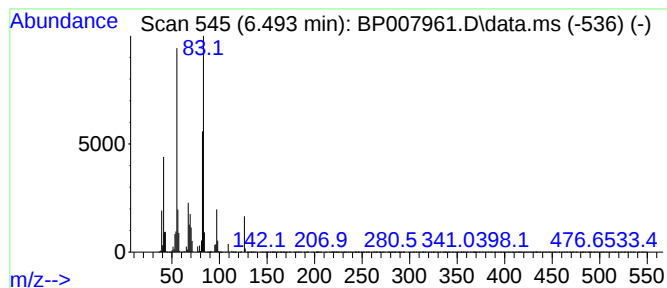
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Cyclohexanone, 2,3-dimethyl- Concentration Rank 5

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 6.493 | 11.94 ng | 4027650 | 1,4-Dichlorobenzene-d4 | 7.851 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexanone, 2,3-dimethyl-       | 126 | C8H14O  | 013395-76-1 | 87   |
| 2    |    |   | Cyclohexane, propyl-               | 126 | C9H18   | 001678-92-8 | 87   |
| 3    |    |   | Cyclohexane, (1-methylethyl)-      | 126 | C9H18   | 000696-29-7 | 60   |
| 4    |    |   | Cyclohexane, ethyl-                | 112 | C8H16   | 001678-91-7 | 59   |
| 5    |    |   | Ethanone, 1-(1-methylcyclopentyl)- | 126 | C8H14O  | 013388-93-7 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

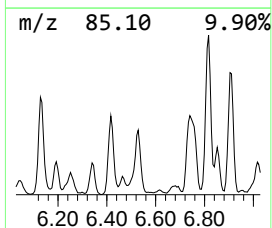
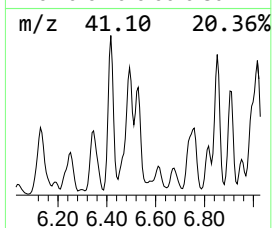
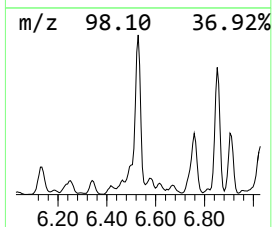
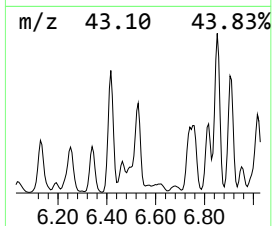
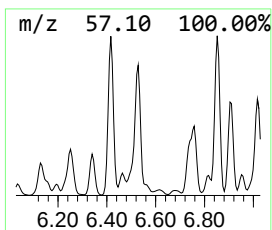
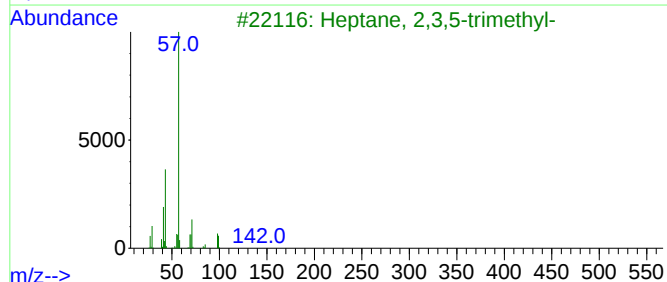
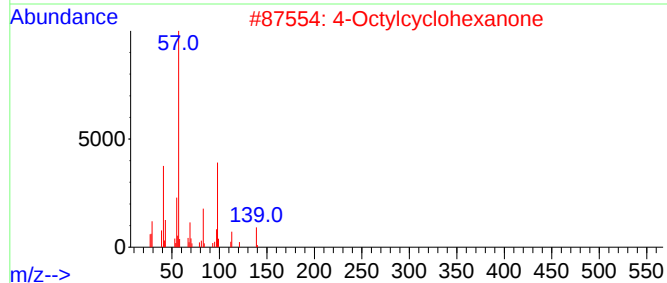
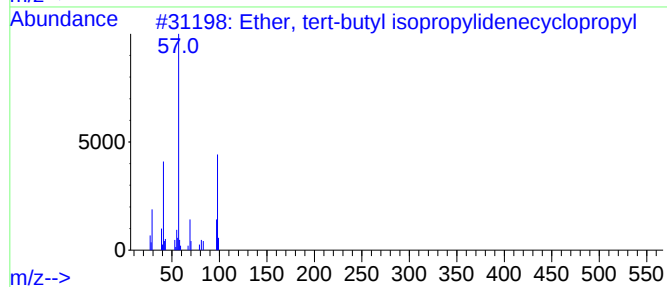
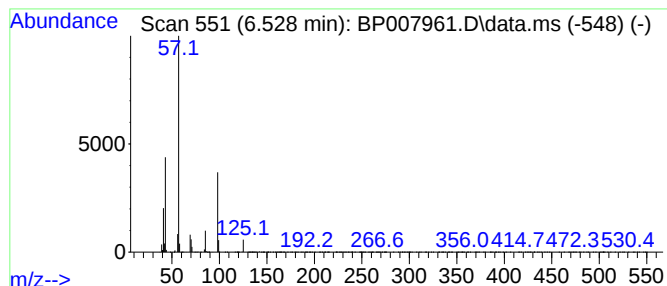
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 6 Ether, tert-butyl isopropyl... Concentration Rank 17

| R.T.  | EstConc | Area    | Relative to ISTD       | R.T.  |
|-------|---------|---------|------------------------|-------|
| 6.528 | 6.51 ng | 2195740 | 1,4-Dichlorobenzene-d4 | 7.851 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Ether, tert-butyl isopropylidene... | 154 | C10H18O | 024524-56-9  | 53   |
| 2    |    |   | 4-Octylcyclohexanone                | 210 | C14H26O | 1000428-94-1 | 50   |
| 3    |    |   | Heptane, 2,3,5-trimethyl-           | 142 | C10H22  | 020278-85-7  | 47   |
| 4    |    |   | Hexane, 3-ethyl-2,5-dimethyl-       | 142 | C10H22  | 052897-04-8  | 47   |
| 5    |    |   | Octane, 2,3-dimethyl-               | 142 | C10H22  | 007146-60-3  | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

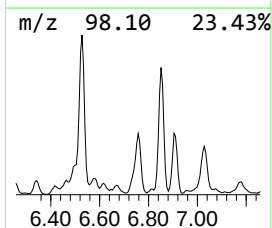
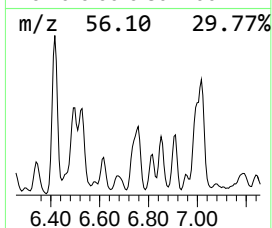
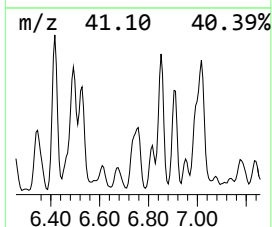
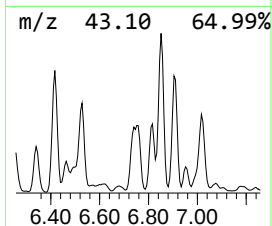
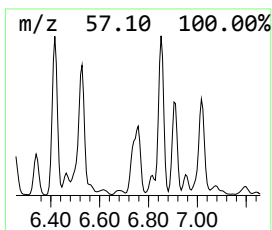
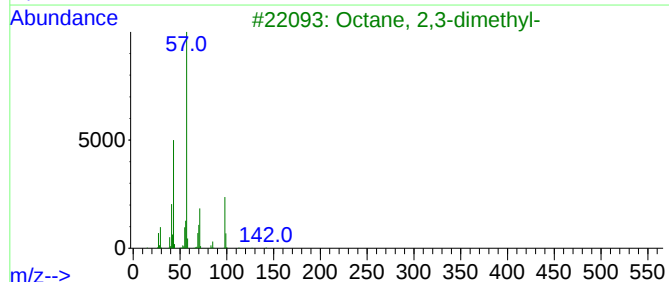
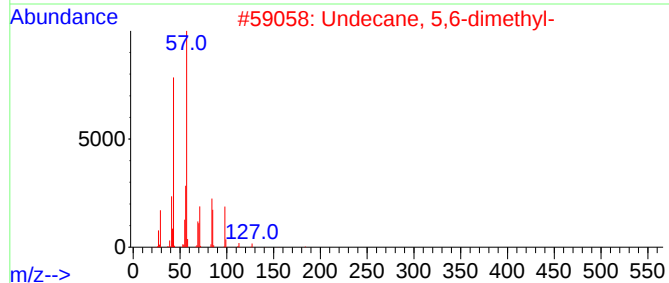
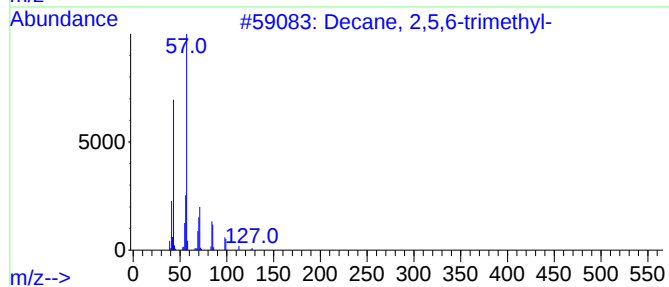
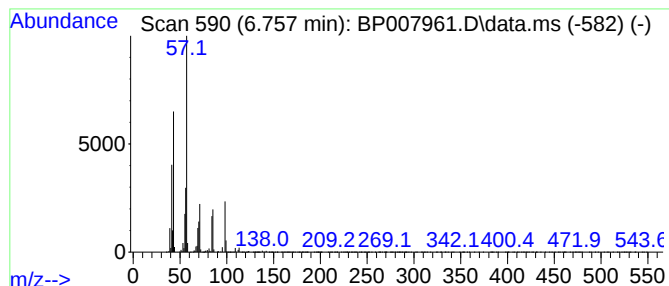
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Decane, 2,5,6-trimethyl- Concentration Rank 16

| R.T.  | EstConc | Area    | Relative to ISTD       | R.T.  |
|-------|---------|---------|------------------------|-------|
| 6.757 | 7.08 ng | 2387220 | 1,4-Dichlorobenzene-d4 | 7.851 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 2,5,6-trimethyl-   | 184 | C13H28  | 062108-23-0 | 72   |
| 2    |    |   | Undecane, 5,6-dimethyl-    | 184 | C13H28  | 017615-91-7 | 64   |
| 3    |    |   | Octane, 2,3-dimethyl-      | 142 | C10H22  | 007146-60-3 | 58   |
| 4    |    |   | Heptane, 3-ethyl-2-methyl- | 142 | C10H22  | 014676-29-0 | 58   |
| 5    |    |   | Nonane                     | 128 | C9H20   | 000111-84-2 | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

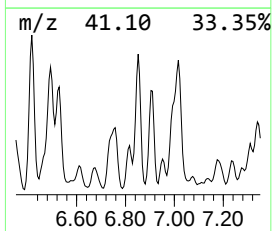
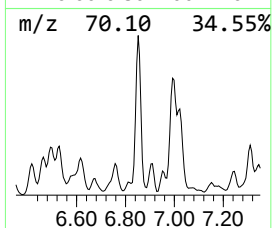
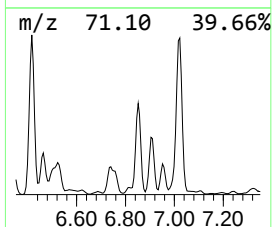
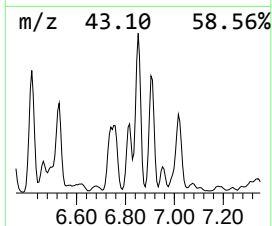
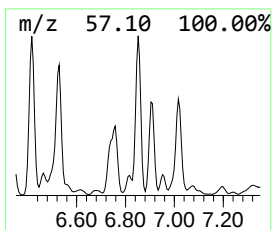
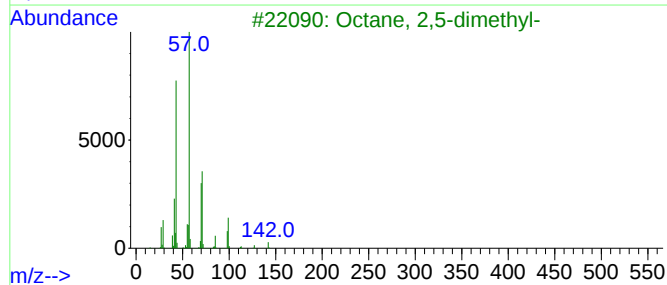
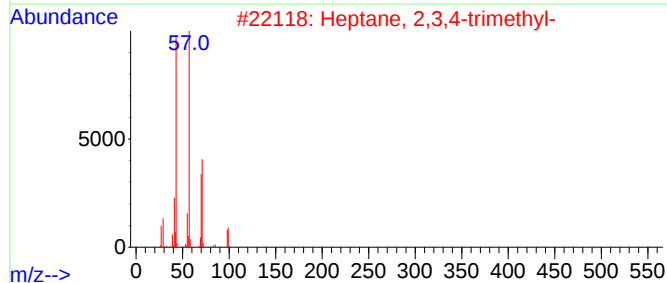
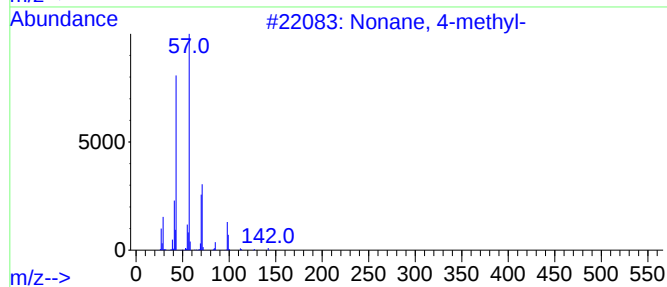
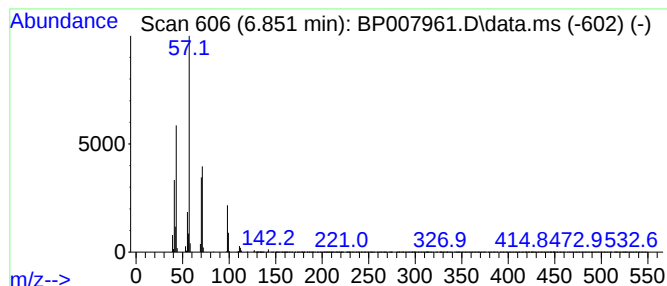
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 Nonane, 4-methyl- Concentration Rank 7

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 6.851 | 10.30 ng | 3474570 | 1,4-Dichlorobenzene-d4 | 7.851 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Nonane, 4-methyl-             | 142 | C10H22  | 017301-94-9 | 87   |
| 2    |    |   | Heptane, 2,3,4-trimethyl-     | 142 | C10H22  | 052896-95-4 | 59   |
| 3    |    |   | Octane, 2,5-dimethyl-         | 142 | C10H22  | 015869-89-3 | 59   |
| 4    |    |   | Pentane, 2,2,3,3-tetramethyl- | 128 | C9H20   | 007154-79-2 | 53   |
| 5    |    |   | Undecane, 4,5-dimethyl-       | 184 | C13H28  | 017312-79-7 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

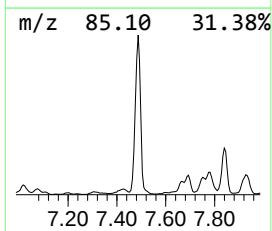
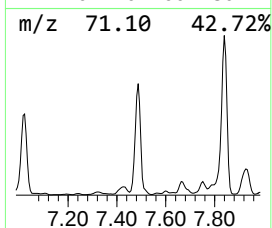
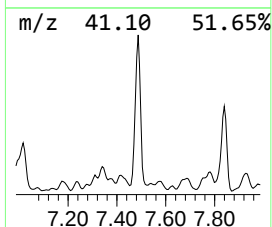
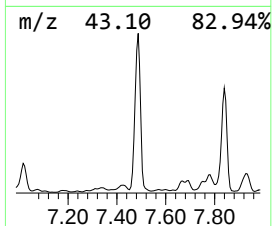
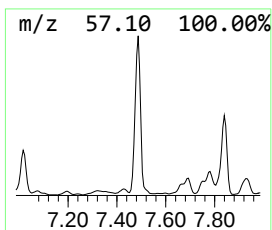
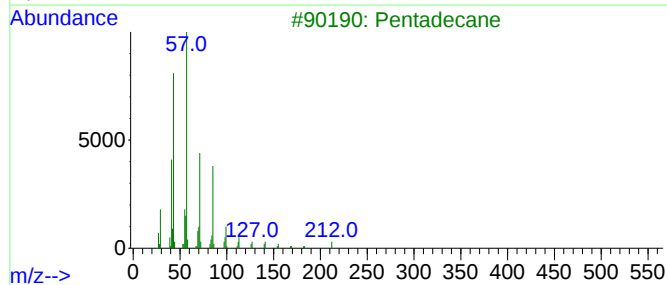
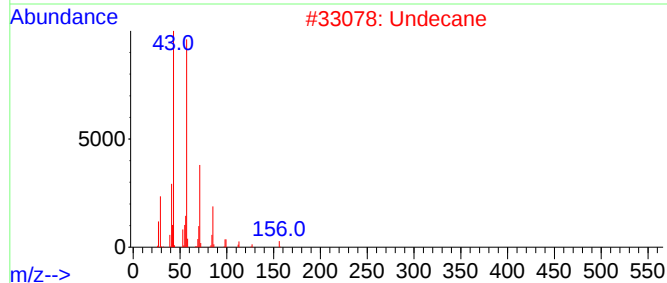
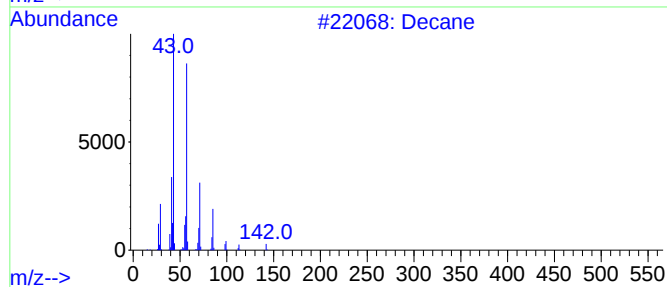
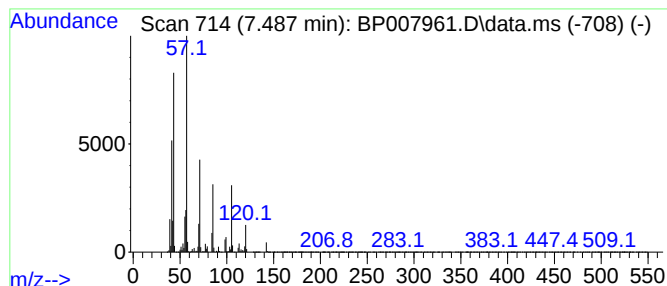
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Decane Concentration Rank 1

| R.T.  | EstConc  | Area     | Relative to ISTD       | R.T.  |
|-------|----------|----------|------------------------|-------|
| 7.487 | 33.55 ng | 11318800 | 1,4-Dichlorobenzene-d4 | 7.851 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Decane                              | 142 | C10H22   | 000124-18-5  | 90   |
| 2    |    |   | Undecane                            | 156 | C11H24   | 001120-21-4  | 59   |
| 3    |    |   | Pentadecane                         | 212 | C15H32   | 000629-62-9  | 59   |
| 4    |    |   | Carbonic acid, prop-1-en-2-yl tr... | 284 | C17H32O3 | 1000382-90-8 | 58   |
| 5    |    |   | Tridecane                           | 184 | C13H28   | 000629-50-5  | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

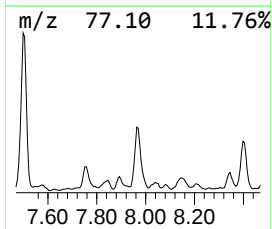
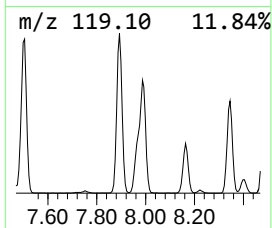
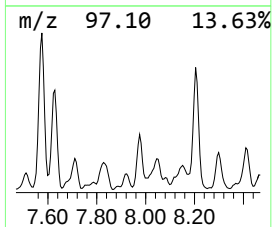
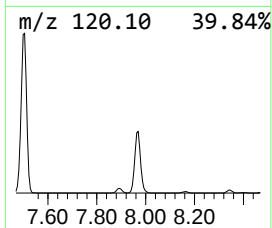
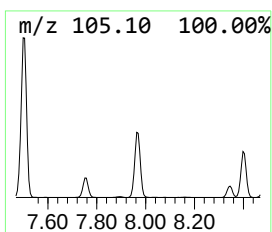
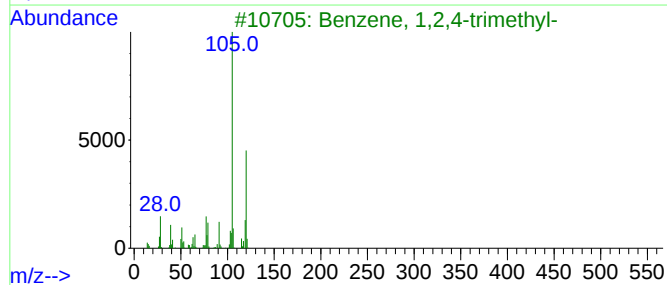
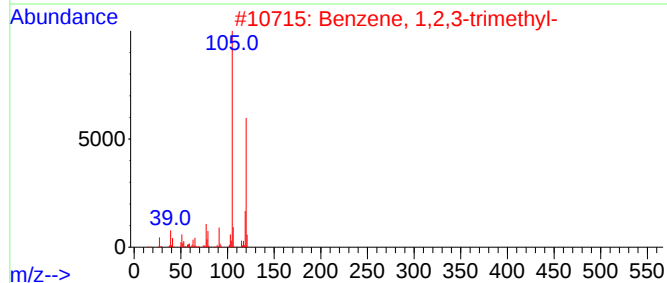
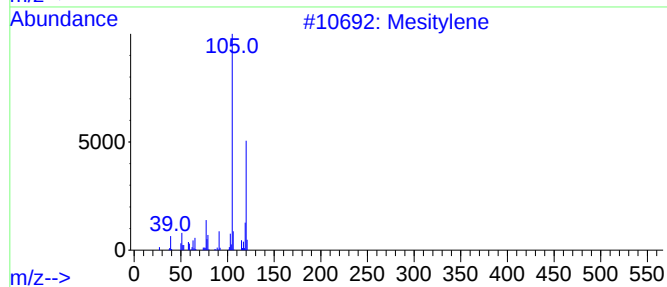
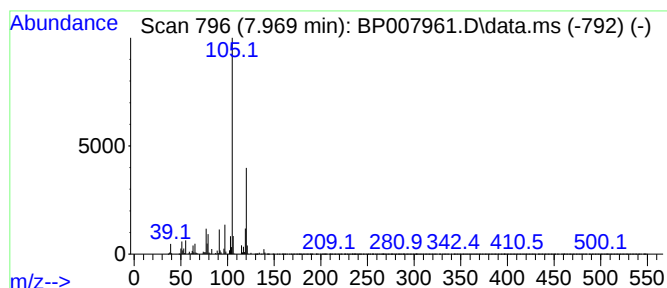
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Mesitylene Concentration Rank 18

| R.T.  | EstConc | Area    | Relative to ISTD       | R.T.  |
|-------|---------|---------|------------------------|-------|
| 7.969 | 6.39 ng | 2156530 | 1,4-Dichlorobenzene-d4 | 7.851 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 95   |
| 2    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 94   |
| 3    |    |   | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 93   |
| 4    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 87   |
| 5    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

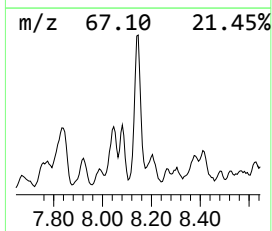
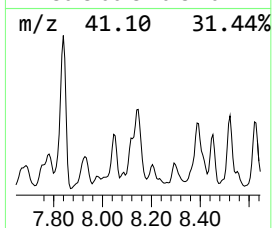
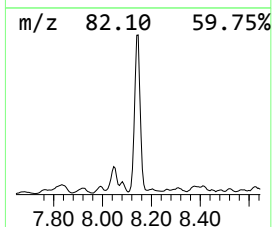
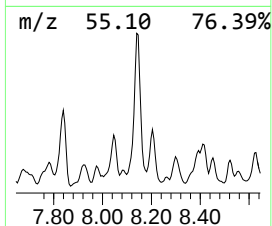
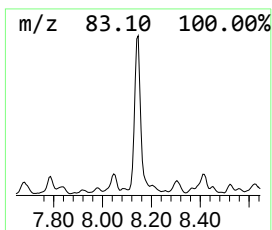
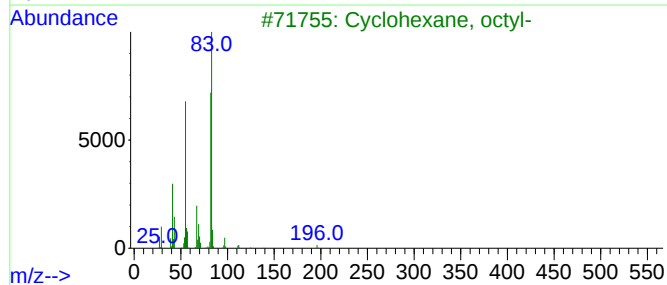
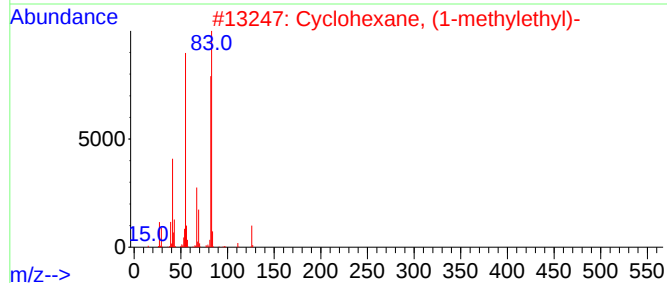
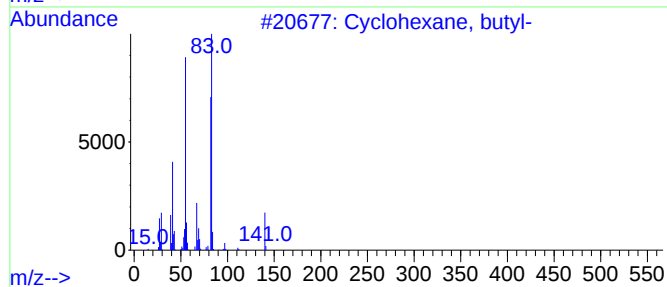
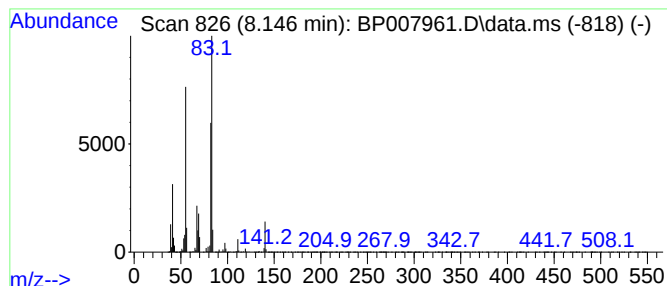
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Cyclohexane, butyl- Concentration Rank 3

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 8.146 | 13.35 ng | 4505100 | 1,4-Dichlorobenzene-d4 | 7.851 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, butyl-           | 140 | C10H20  | 001678-93-9 | 90   |
| 2    |    |   | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 80   |
| 3    |    |   | Cyclohexane, octyl-           | 196 | C14H28  | 001795-15-9 | 78   |
| 4    |    |   | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 72   |
| 5    |    |   | Cyclohexane, decyl-           | 224 | C16H32  | 001795-16-0 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

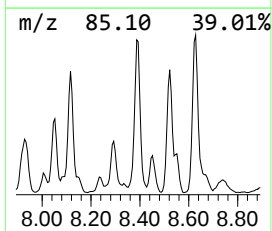
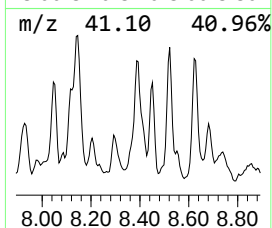
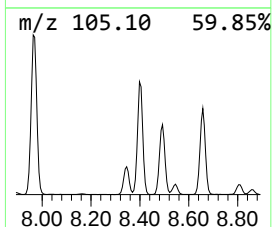
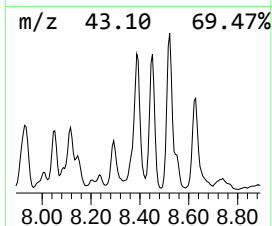
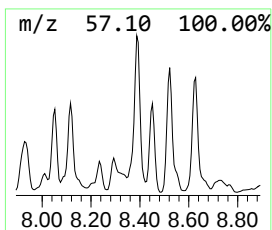
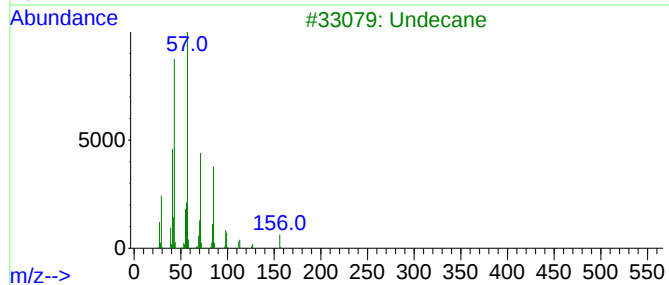
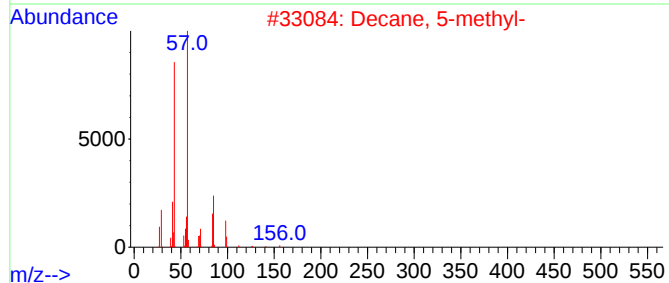
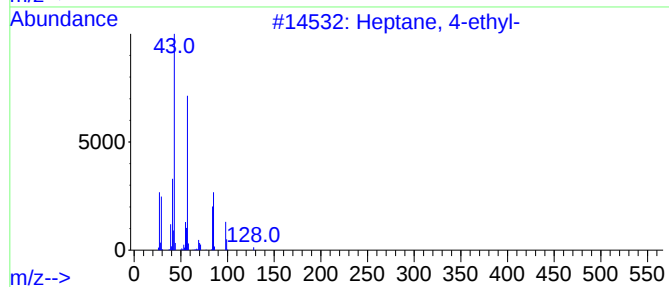
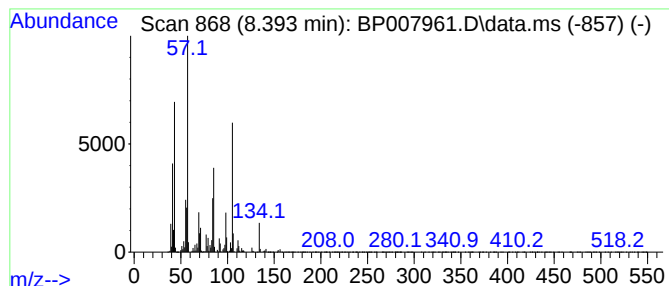
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Heptane, 4-ethyl- Concentration Rank 4

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 8.393 | 13.28 ng | 4479710 | 1,4-Dichlorobenzene-d4 | 7.851 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Heptane, 4-ethyl-                   | 128 | C9H20    | 002216-32-2 | 58   |
| 2    |    |   | Decane, 5-methyl-                   | 156 | C11H24   | 013151-35-4 | 52   |
| 3    |    |   | Undecane                            | 156 | C11H24   | 001120-21-4 | 50   |
| 4    |    |   | 2H-Pyran, 2-(1,1-dimethylethoxy)... | 158 | C9H18O2  | 001927-69-1 | 38   |
| 5    |    |   | 2,2-Dimethylpropionic acid, 4-me... | 186 | C11H22O2 | 150588-62-8 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

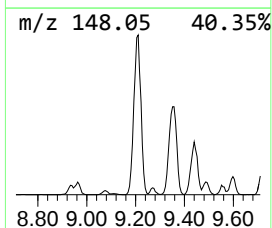
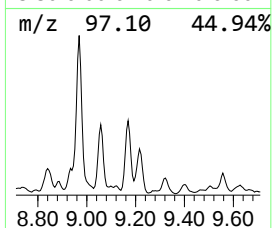
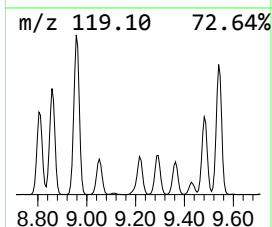
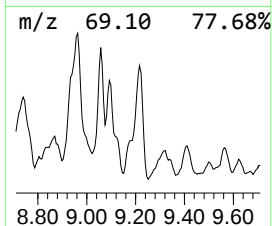
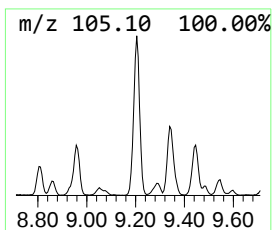
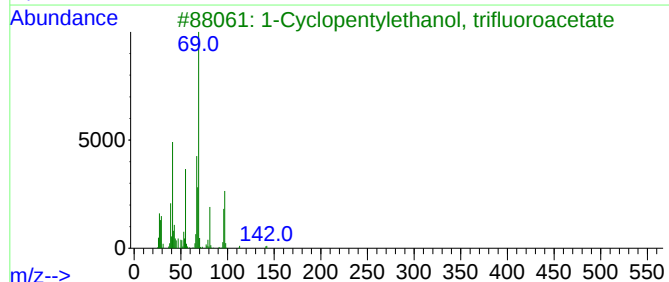
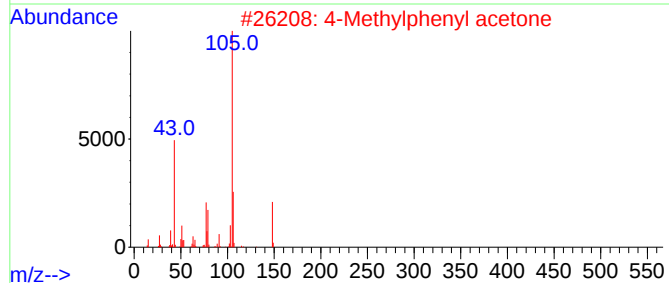
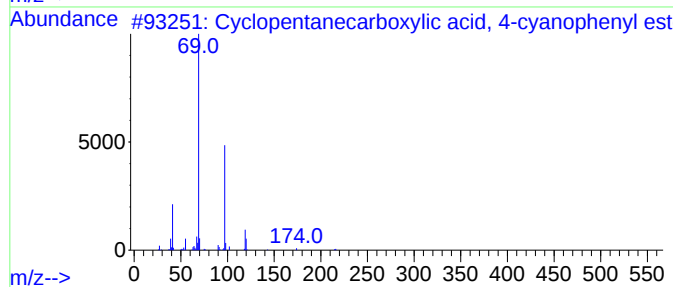
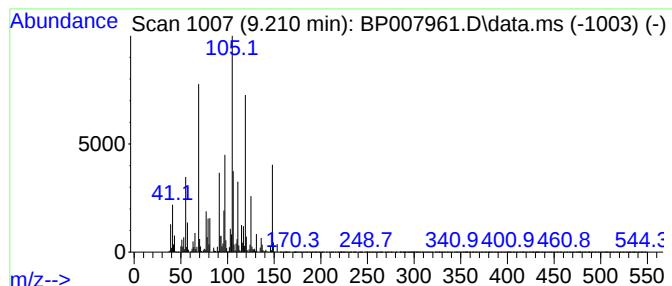
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 13 unknown9.210 Concentration Rank 12

| R.T.  | EstConc | Area    | Relative to ISTD       | R.T.  |
|-------|---------|---------|------------------------|-------|
| 9.210 | 8.08 ng | 2725990 | 1,4-Dichlorobenzene-d4 | 7.851 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Cyclopentanecarboxylic acid, 4-c... | 215 | C13H13NO2 | 1000307-57-8 | 27   |
| 2    |    |   | 4-Methylphenyl acetone              | 148 | C10H12O   | 002096-86-8  | 25   |
| 3    |    |   | 1-Cyclopentylethanol, trifluoroa... | 210 | C9H13F3O2 | 1000364-11-9 | 22   |
| 4    |    |   | Benzene, 1-methyl-4-(2-methylpro... | 148 | C11H16    | 005161-04-6  | 15   |
| 5    |    |   | 1-[1-(1-Hydroxy-butyl)-cyclopent... | 260 | C17H24O2  | 097234-38-3  | 14   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

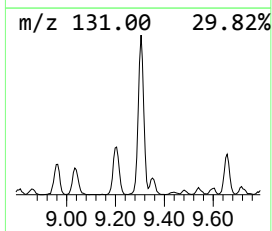
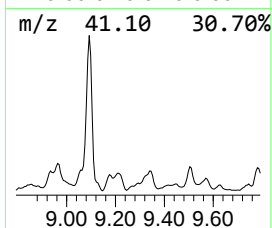
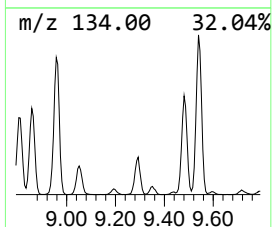
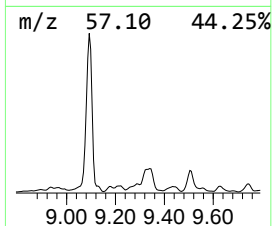
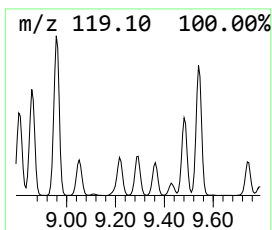
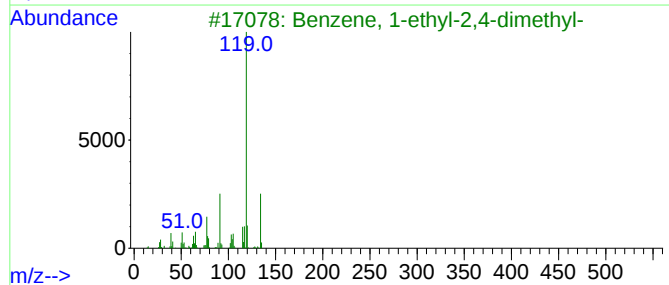
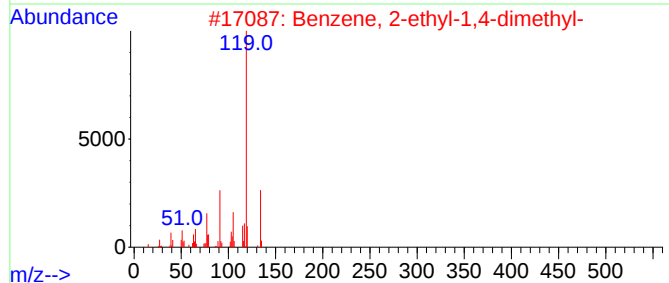
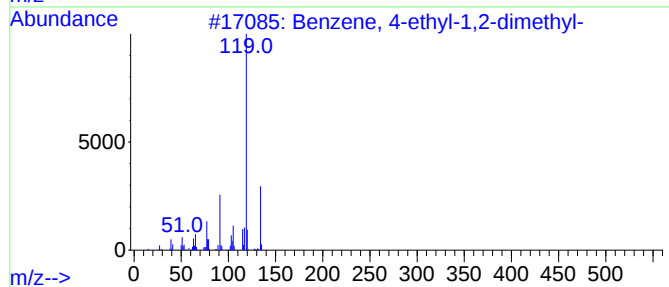
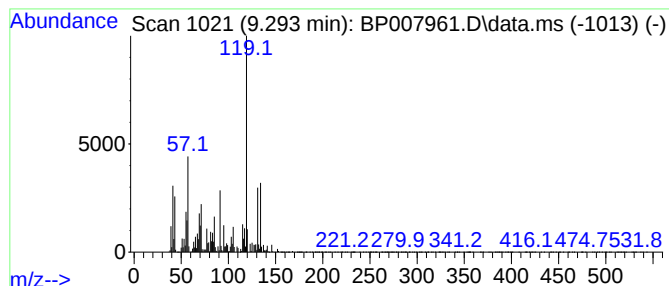
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 19

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.   |
|-------|---------|---------|------------------|--------|
| 9.293 | 6.30 ng | 1116770 | Naphthalene-d8   | 10.645 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 92   |
| 2    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 78   |
| 3    |    |   | Benzene, 1-ethyl-2,4-dimethyl- | 134 | C10H14  | 000874-41-9 | 78   |
| 4    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 78   |
| 5    |    |   | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 78   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

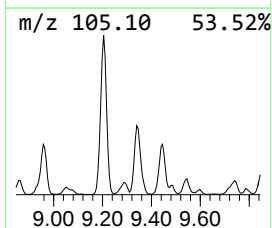
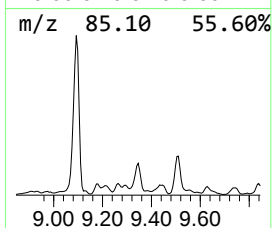
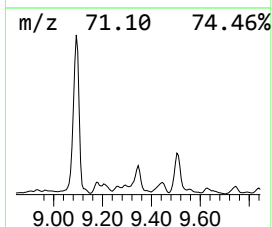
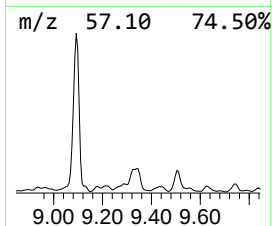
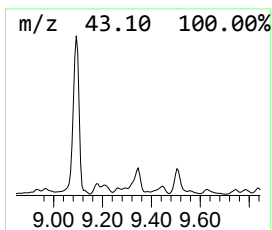
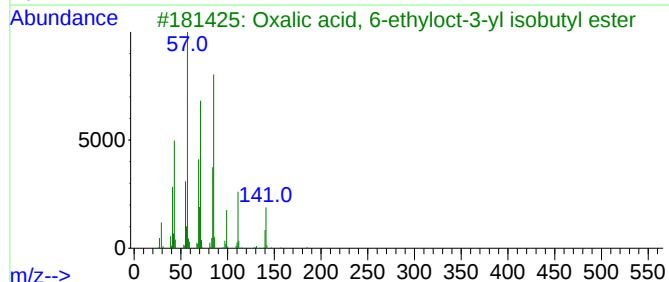
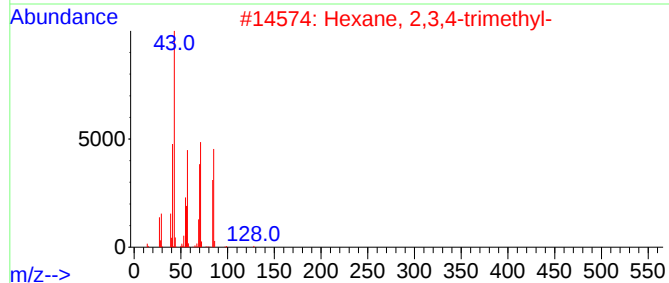
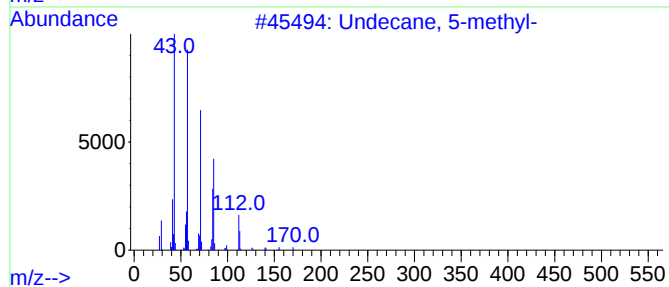
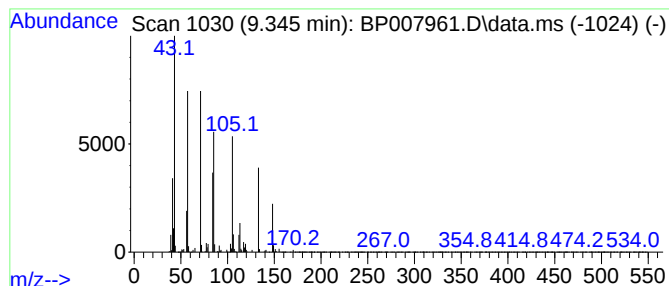
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Undecane, 5-methyl- Concentration Rank 9

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.   |
|-------|----------|---------|------------------|--------|
| 9.345 | 10.03 ng | 1779320 | Naphthalene-d8   | 10.645 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Undecane, 5-methyl-                 | 170 | C12H26   | 001632-70-8  | 58   |
| 2    |    |   | Hexane, 2,3,4-trimethyl-            | 128 | C9H20    | 000921-47-1  | 50   |
| 3    |    |   | Oxalic acid, 6-ethyloct-3-yl iso... | 286 | C16H30O4 | 1010309-34-1 | 50   |
| 4    |    |   | Octane, 4-methyl-                   | 128 | C9H20    | 002216-34-4  | 50   |
| 5    |    |   | 2-Bromo dodecane                    | 248 | C12H25Br | 013187-99-0  | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

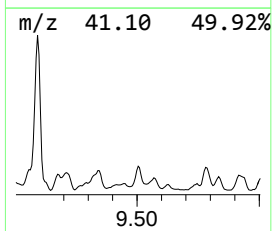
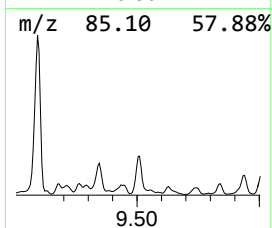
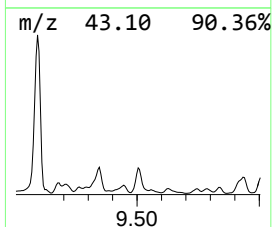
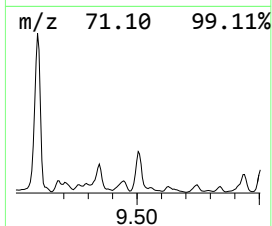
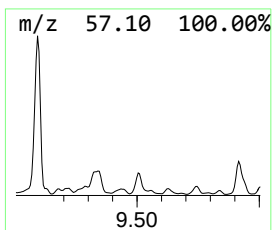
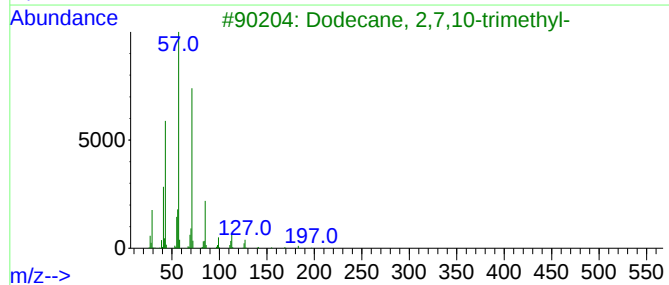
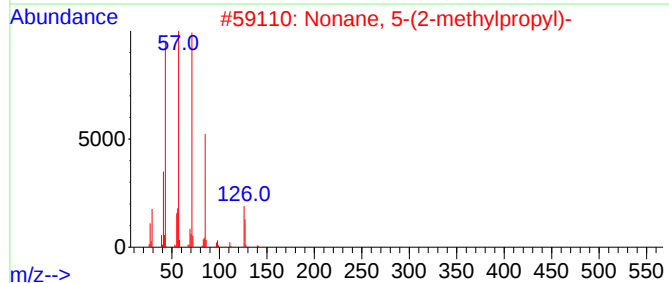
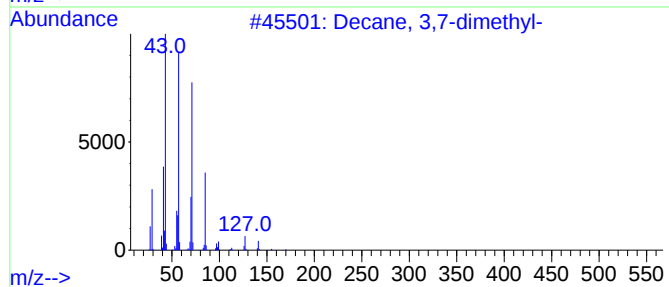
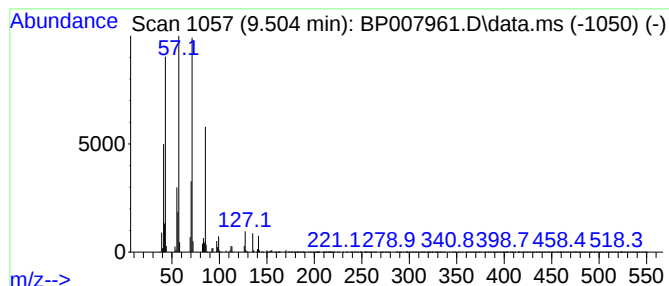
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Decane, 3,7-dimethyl- Concentration Rank 14

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.   |
|-------|---------|---------|------------------|--------|
| 9.504 | 7.92 ng | 1404830 | Naphthalene-d8   | 10.645 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 3,7-dimethyl-       | 170 | C12H26  | 017312-54-8 | 93   |
| 2    |    |   | Nonane, 5-(2-methylpropyl)- | 184 | C13H28  | 062185-53-9 | 72   |
| 3    |    |   | Dodecane, 2,7,10-trimethyl- | 212 | C15H32  | 074645-98-0 | 59   |
| 4    |    |   | Undecane, 2-methyl-         | 170 | C12H26  | 007045-71-8 | 58   |
| 5    |    |   | Octane, 3-ethyl-            | 142 | C10H22  | 005881-17-4 | 52   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

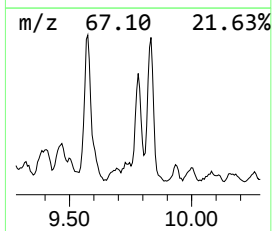
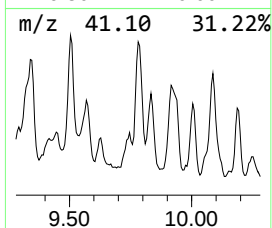
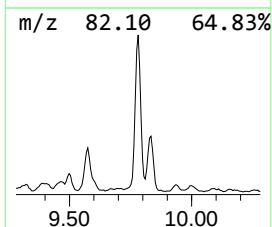
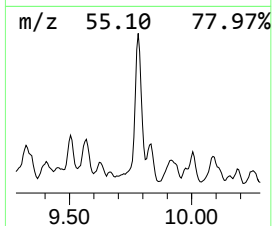
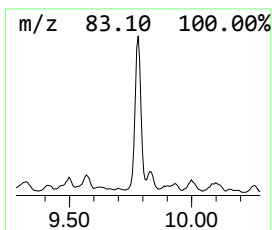
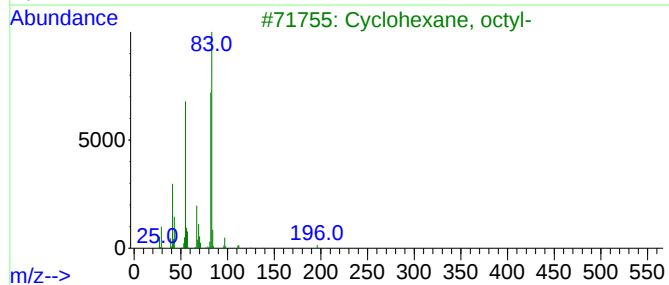
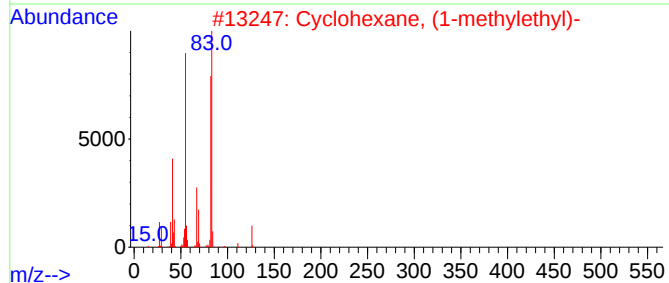
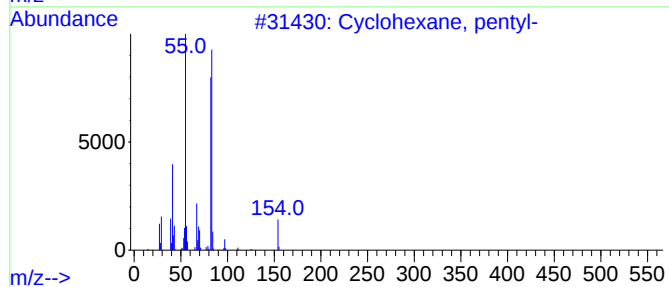
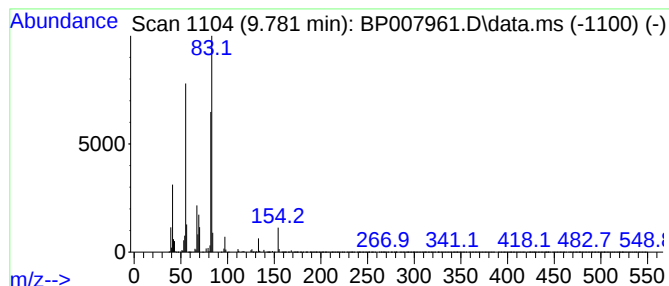
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Cyclohexane, pentyl- Concentration Rank 8

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.   |
|-------|----------|---------|------------------|--------|
| 9.781 | 10.05 ng | 1781970 | Naphthalene-d8   | 10.645 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, pentyl-          | 154 | C11H22  | 004292-92-6 | 95   |
| 2    |    |   | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 80   |
| 3    |    |   | Cyclohexane, octyl-           | 196 | C14H28  | 001795-15-9 | 78   |
| 4    |    |   | Cyclohexane, hexyl-           | 168 | C12H24  | 004292-75-5 | 78   |
| 5    |    |   | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 78   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

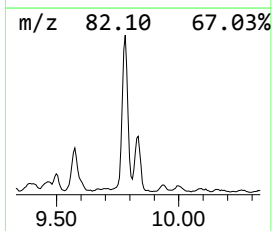
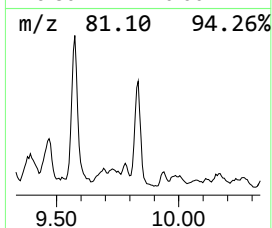
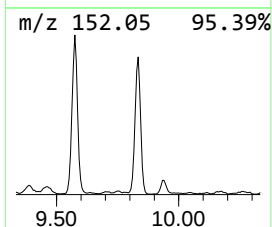
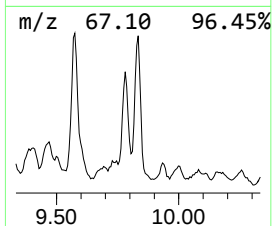
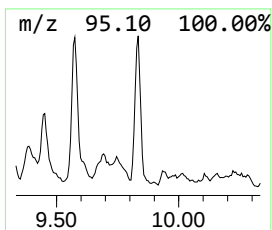
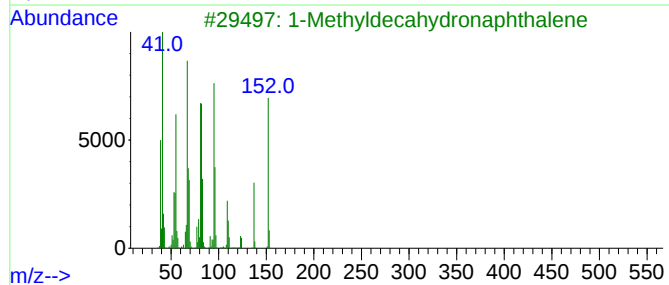
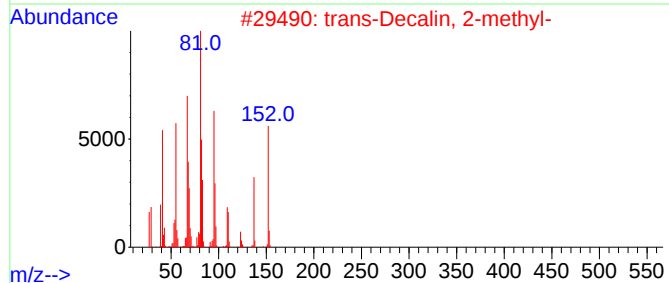
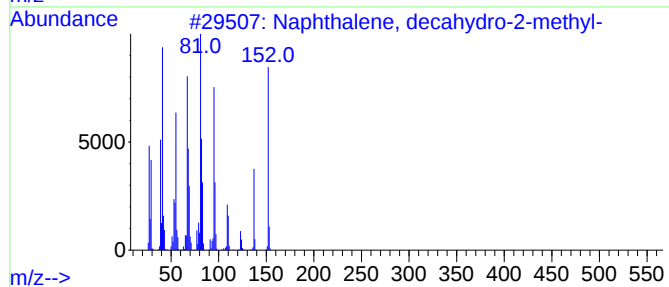
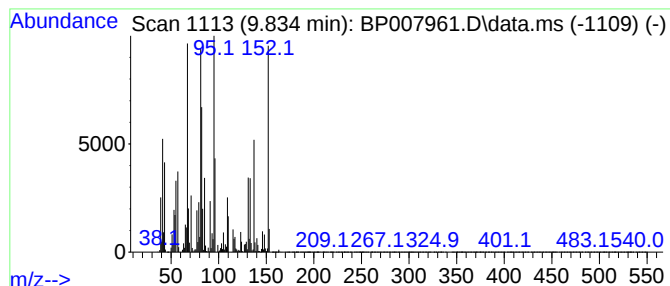
\*\*\*\*\*

Peak Number 18 Naphthalene, decahydro-2-me... Concentration Rank 6

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.   |
|-------|----------|---------|------------------|--------|
| 9.834 | 10.38 ng | 1840430 | Naphthalene-d8   | 10.645 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|------|-------------------------------------|-----|---------|--------------|------|
| 1    |      | Naphthalene, decahydro-2-methyl-    | 152 | C11H20  | 002958-76-1  | 91   |
| 2    |      | trans-Decalin, 2-methyl-            | 152 | C11H20  | 1000152-47-3 | 86   |
| 3    |      | 1-Methyldecahydronaphthalene        | 152 | C11H20  | 002958-75-0  | 86   |
| 4    |      | cis-Decalin, 2-syn-methyl-          | 152 | C11H20  | 1000155-85-6 | 72   |
| 5    |      | trans-4a-Methyl-decahydronaphtha... | 152 | C11H20  | 002547-27-5  | 70   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

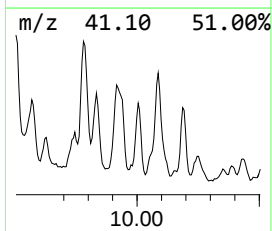
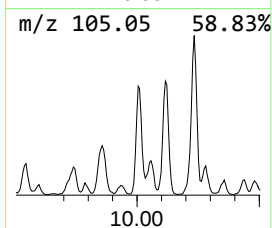
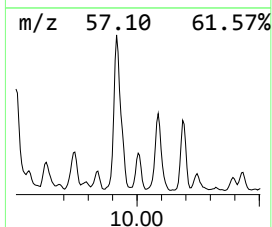
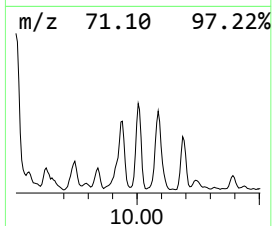
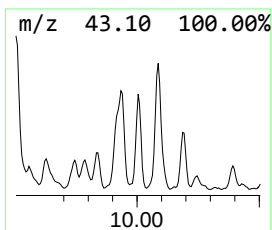
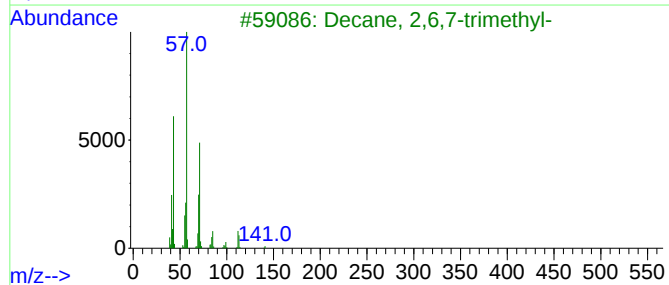
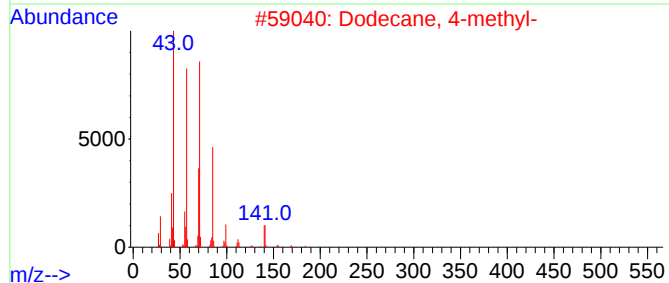
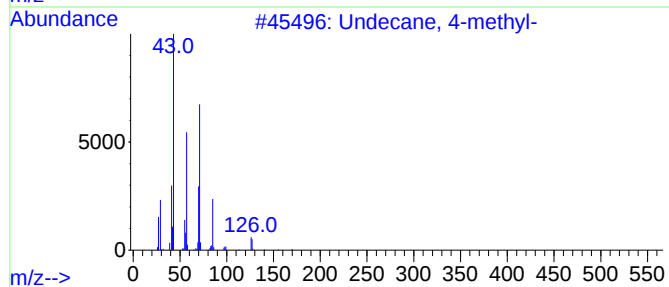
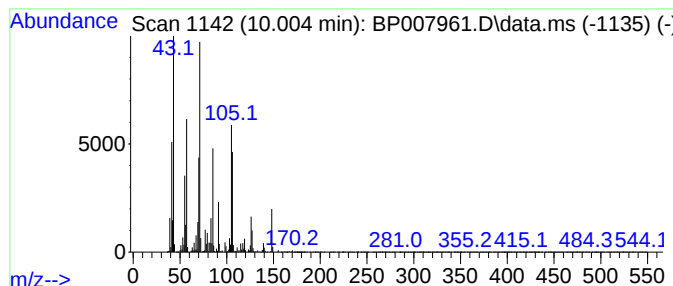
Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 Undecane, 4-methyl- Concentration Rank 20

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|--------------|------|
| 10.004    | 5.96 ng                             | 1057500 | Naphthalene-d8   | 10.645       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#         | Qual |
| 1         | Undecane, 4-methyl-                 | 170     | C12H26           | 002980-69-0  | 53   |
| 2         | Dodecane, 4-methyl-                 | 184     | C13H28           | 006117-97-1  | 47   |
| 3         | Decane, 2,6,7-trimethyl-            | 184     | C13H28           | 062108-25-2  | 38   |
| 4         | Decane, 4-methyl-                   | 156     | C11H24           | 002847-72-5  | 38   |
| 5         | 4-Methyl-3-heptanol, trifluoroac... | 226     | C10H17F3O2       | 1000365-51-8 | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

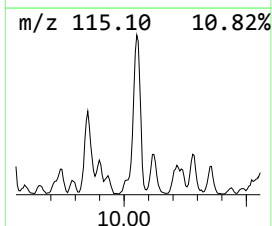
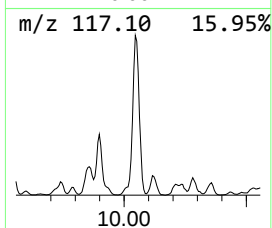
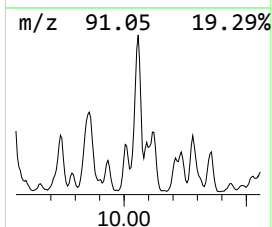
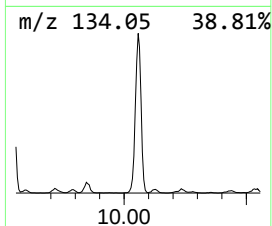
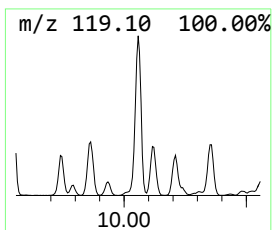
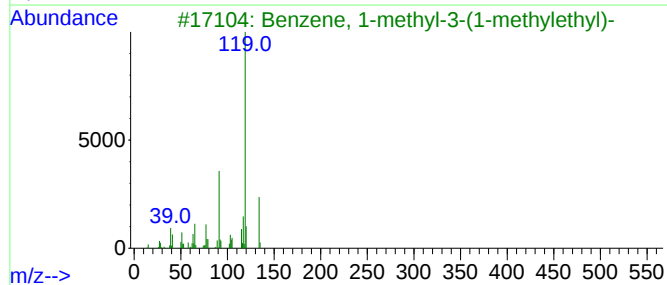
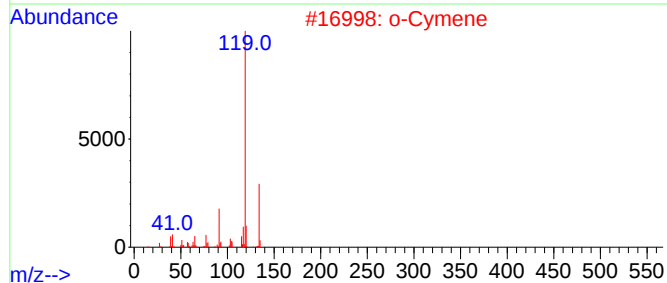
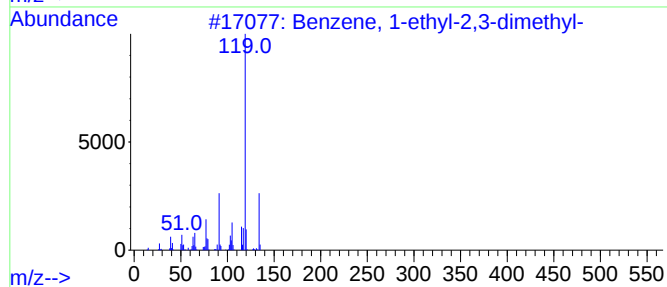
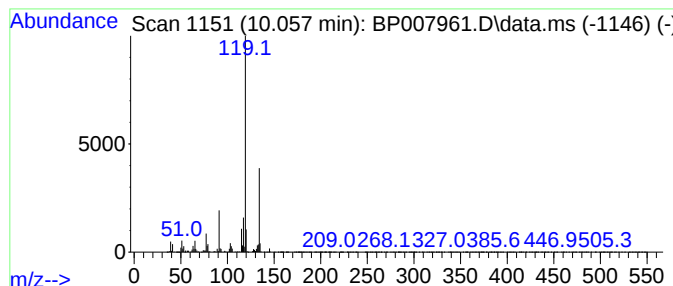
\*\*\*\*\*

Peak Number 20 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 15

| R.T.   | EstConc | Area    | Relative to ISTD | R.T.   |
|--------|---------|---------|------------------|--------|
| 10.057 | 7.65 ng | 1357710 | Naphthalene-d8   | 10.645 |

| Hit# | of 5 | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|------|--------------------------------------|-----|---------|-------------|------|
| 1    |      | Benzene, 1-ethyl-2,3-dimethyl-       | 134 | C10H14  | 000933-98-2 | 94   |
| 2    |      | o-Cymene                             | 134 | C10H14  | 000527-84-4 | 93   |
| 3    |      | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 91   |
| 4    |      | Benzene, 4-ethyl-1,2-dimethyl-       | 134 | C10H14  | 000934-80-5 | 91   |
| 5    |      | p-Cymene                             | 134 | C10H14  | 000099-87-6 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111121\  
Data File : BP007961.D  
Acq On : 12 Nov 2021 00:01  
Operator : CG/JU  
Sample : M4630-07  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SB-13N-16.0-16.5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L  
TIC Integration Parameters: LSCINT.P

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
|                    |        |         |       |          | #                     | RT     | Resp    | Conc |
| Nonane             | 5.875  | 15.8    | ng    | 5322090  | 1                     | 7.851  | 6748030 | 20.0 |
| trans-1,2-Dieth... | 6.128  | 8.5     | ng    | 2881580  | 1                     | 7.851  | 6748030 | 20.0 |
| unknown6.346       | 6.346  | 8.0     | ng    | 2687830  | 1                     | 7.851  | 6748030 | 20.0 |
| Octane, 2,6-dim... | 6.416  | 9.7     | ng    | 3276420  | 1                     | 7.851  | 6748030 | 20.0 |
| Cyclohexanone, ... | 6.493  | 11.9    | ng    | 4027650  | 1                     | 7.851  | 6748030 | 20.0 |
| Ether, tert-but... | 6.528  | 6.5     | ng    | 2195740  | 1                     | 7.851  | 6748030 | 20.0 |
| Decane, 2,5,6-t... | 6.757  | 7.1     | ng    | 2387220  | 1                     | 7.851  | 6748030 | 20.0 |
| Nonane, 4-methyl-  | 6.851  | 10.3    | ng    | 3474570  | 1                     | 7.851  | 6748030 | 20.0 |
| Decane             | 7.487  | 33.5    | ng    | 11318800 | 1                     | 7.851  | 6748030 | 20.0 |
| Mesitylene         | 7.969  | 6.4     | ng    | 2156530  | 1                     | 7.851  | 6748030 | 20.0 |
| Cyclohexane, bu... | 8.146  | 13.4    | ng    | 4505100  | 1                     | 7.851  | 6748030 | 20.0 |
| Heptane, 4-ethyl-  | 8.393  | 13.3    | ng    | 4479710  | 1                     | 7.851  | 6748030 | 20.0 |
| unknown9.210       | 9.210  | 8.1     | ng    | 2725990  | 1                     | 7.851  | 6748030 | 20.0 |
| Benzene, 4-ethy... | 9.293  | 6.3     | ng    | 1116770  | 2                     | 10.645 | 3547280 | 20.0 |
| Undecane, 5-met... | 9.345  | 10.0    | ng    | 1779320  | 2                     | 10.645 | 3547280 | 20.0 |
| Decane, 3,7-dim... | 9.504  | 7.9     | ng    | 1404830  | 2                     | 10.645 | 3547280 | 20.0 |
| Cyclohexane, pe... | 9.781  | 10.1    | ng    | 1781970  | 2                     | 10.645 | 3547280 | 20.0 |
| Naphthalene, de... | 9.834  | 10.4    | ng    | 1840430  | 2                     | 10.645 | 3547280 | 20.0 |
| Undecane, 4-met... | 10.004 | 6.0     | ng    | 1057500  | 2                     | 10.645 | 3547280 | 20.0 |
| Benzene, 1-ethy... | 10.057 | 7.7     | ng    | 1357710  | 2                     | 10.645 | 3547280 | 20.0 |