

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
 Data File : BP008973.D  
 Acq On : 31 Jan 2022 19:43  
 Operator : CG/JU  
 Sample : N1303-03  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 CC-6

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-BP011422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP008973.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         | 3.040       | 4             | 9           | 15           | rVB      | 354545         | 467770        | 1.27%           | 0.168%        |
| 2         | 3.752       | 125           | 130         | 141          | rVB      | 1023843        | 1362426       | 3.71%           | 0.488%        |
| 3         | 4.369       | 226           | 235         | 240          | rBV      | 17682409       | 36751176      | 100.00%         | 13.170%       |
| 4         | 4.969       | 325           | 337         | 344          | rBV      | 8313835        | 14827978      | 40.35%          | 5.314%        |
| 5         | 5.346       | 396           | 401         | 406          | rVB      | 547201         | 804371        | 2.19%           | 0.288%        |
| 6         | 5.475       | 417           | 423         | 441          | rVB      | 2212730        | 4311634       | 11.73%          | 1.545%        |
| 7         | 5.787       | 471           | 476         | 480          | rBV      | 334929         | 511486        | 1.39%           | 0.183%        |
| 8         | 5.846       | 480           | 486         | 495          | rVB2     | 1660032        | 2742935       | 7.46%           | 0.983%        |
| 9         | 6.105       | 522           | 530         | 538          | rBV3     | 668509         | 1570817       | 4.27%           | 0.563%        |
| 10        | 6.228       | 544           | 551         | 557          | rVB2     | 396934         | 794942        | 2.16%           | 0.285%        |
| 11        | 6.322       | 561           | 567         | 574          | rBV5     | 738835         | 1598299       | 4.35%           | 0.573%        |
| 12        | 6.393       | 574           | 579         | 583          | rBV      | 1230619        | 1871065       | 5.09%           | 0.671%        |
| 13        | 6.469       | 583           | 592         | 595          | rVV3     | 1070357        | 2506185       | 6.82%           | 0.898%        |
| 14        | 6.499       | 595           | 597         | 605          | rVB      | 856011         | 1296365       | 3.53%           | 0.465%        |
| 15        | 6.587       | 607           | 612         | 617          | rVB3     | 292602         | 494393        | 1.35%           | 0.177%        |
| 16        | 6.646       | 617           | 622         | 628          | rVB3     | 311053         | 548888        | 1.49%           | 0.197%        |
| 17        | 6.728       | 628           | 636         | 642          | rVB3     | 754757         | 1744828       | 4.75%           | 0.625%        |
| 18        | 6.822       | 642           | 652         | 658          | rBV3     | 1717309        | 4129572       | 11.24%          | 1.480%        |
| 19        | 6.881       | 658           | 662         | 665          | rBV      | 677550         | 987417        | 2.69%           | 0.354%        |
| 20        | 6.922       | 665           | 669         | 673          | rVV      | 2310163        | 3340557       | 9.09%           | 1.197%        |
| 21        | 6.981       | 673           | 679         | 687          | rVV4     | 2104809        | 5183039       | 14.10%          | 1.857%        |
| 22        | 7.057       | 687           | 692         | 697          | rVB      | 2301923        | 3680659       | 10.02%          | 1.319%        |
| 23        | 7.157       | 704           | 709         | 714          | rVB5     | 427357         | 829065        | 2.26%           | 0.297%        |
| 24        | 7.216       | 714           | 719         | 723          | rBV2     | 1218250        | 1966968       | 5.35%           | 0.705%        |
| 25        | 7.316       | 728           | 736         | 741          | rBV      | 1114625        | 2169550       | 5.90%           | 0.777%        |
| 26        | 7.393       | 745           | 749         | 755          | rVB3     | 653599         | 1161190       | 3.16%           | 0.416%        |
| 27        | 7.481       | 755           | 764         | 772          | rBV2     | 7238831        | 15728964      | 42.80%          | 5.637%        |
| 28        | 7.552       | 772           | 776         | 781          | rVB2     | 548040         | 1005471       | 2.74%           | 0.360%        |
| 29        | 7.663       | 787           | 795         | 801          | rBV4     | 587432         | 1624259       | 4.42%           | 0.582%        |
| 30        | 7.752       | 802           | 810         | 815          | rBV5     | 964589         | 2691141       | 7.32%           | 0.964%        |
| 31        | 7.816       | 815           | 821         | 827          | rVB2     | 4746966        | 7694686       | 20.94%          | 2.758%        |
| 32        | 7.875       | 827           | 831         | 832          | rBV      | 513416         | 696955        | 1.90%           | 0.250%        |
| 33        | 7.904       | 832           | 836         | 839          | rVV2     | 663006         | 1111562       | 3.02%           | 0.398%        |
| 34        | 7.946       | 839           | 843         | 851          | rVB2     | 2904950        | 5138059       | 13.98%          | 1.841%        |
| 35        | 8.022       | 851           | 856         | 860          | rVB2     | 1236577        | 1794453       | 4.88%           | 0.643%        |
| 36        | 8.116       | 865           | 872         | 880          | rVB3     | 1956377        | 4424587       | 12.04%          | 1.586%        |

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 Operator : CG/JU  
 Sample : N1303-03  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 CC-6

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-BP011422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 8.199  | 880  | 886  | 891  | rVB3 | 1134068 | 2395273 | 6.52%  | 0.858% |
| 38 | 8.275  | 894  | 899  | 903  | rBV3 | 486048  | 790560  | 2.15%  | 0.283% |
| 39 | 8.328  | 903  | 908  | 910  | rBV3 | 745289  | 1265480 | 3.44%  | 0.454% |
| 40 | 8.375  | 910  | 916  | 921  | rVB3 | 2272992 | 4738886 | 12.89% | 1.698% |
| 41 | 8.428  | 921  | 925  | 928  | rVB  | 1029311 | 1332009 | 3.62%  | 0.477% |
| 42 | 8.475  | 928  | 933  | 941  | rBV3 | 3044685 | 6715575 | 18.27% | 2.407% |
| 43 | 8.599  | 949  | 954  | 957  | rBV  | 1121962 | 1812074 | 4.93%  | 0.649% |
| 44 | 8.657  | 957  | 964  | 969  | rVB2 | 1306102 | 3099135 | 8.43%  | 1.111% |
| 45 | 8.787  | 981  | 986  | 990  | rBV  | 1000053 | 1602810 | 4.36%  | 0.574% |
| 46 | 8.834  | 990  | 994  | 1001 | rVB  | 1367428 | 2300714 | 6.26%  | 0.825% |
| 47 | 8.940  | 1001 | 1012 | 1016 | rBV2 | 3405372 | 7238184 | 19.70% | 2.594% |
| 48 | 8.987  | 1016 | 1020 | 1024 | rVV  | 1749160 | 3418696 | 9.30%  | 1.225% |
| 49 | 9.034  | 1024 | 1028 | 1029 | rVV3 | 1363717 | 2124728 | 5.78%  | 0.761% |
| 50 | 9.069  | 1029 | 1034 | 1043 | rVB  | 4494786 | 8007796 | 21.79% | 2.870% |
| 51 | 9.151  | 1043 | 1048 | 1050 | rBV3 | 982107  | 1630405 | 4.44%  | 0.584% |
| 52 | 9.187  | 1050 | 1054 | 1060 | rVB4 | 1798914 | 3656956 | 9.95%  | 1.311% |
| 53 | 9.269  | 1060 | 1068 | 1071 | rBV3 | 1110182 | 2283265 | 6.21%  | 0.818% |
| 54 | 9.322  | 1074 | 1077 | 1084 | rVB5 | 1009868 | 1748963 | 4.76%  | 0.627% |
| 55 | 9.463  | 1097 | 1101 | 1107 | rVV2 | 909818  | 2357963 | 6.42%  | 0.845% |
| 56 | 9.522  | 1107 | 1111 | 1114 | rVV  | 1367597 | 2466342 | 6.71%  | 0.884% |
| 57 | 9.551  | 1114 | 1116 | 1122 | rVV2 | 1093585 | 1509879 | 4.11%  | 0.541% |
| 58 | 9.716  | 1137 | 1144 | 1147 | rBV3 | 469262  | 1001553 | 2.73%  | 0.359% |
| 59 | 9.757  | 1147 | 1151 | 1156 | rVB3 | 886345  | 1412741 | 3.84%  | 0.506% |
| 60 | 9.816  | 1156 | 1161 | 1169 | rVB4 | 1081824 | 2646017 | 7.20%  | 0.948% |
| 61 | 9.893  | 1169 | 1174 | 1175 | rBV2 | 413447  | 607552  | 1.65%  | 0.218% |
| 62 | 9.987  | 1185 | 1190 | 1193 | rBV3 | 625768  | 979234  | 2.66%  | 0.351% |
| 63 | 10.034 | 1193 | 1198 | 1205 | rVV3 | 1406800 | 3133780 | 8.53%  | 1.123% |
| 64 | 10.098 | 1205 | 1209 | 1216 | rVB2 | 493405  | 841474  | 2.29%  | 0.302% |
| 65 | 10.169 | 1216 | 1221 | 1226 | rBV2 | 289316  | 694821  | 1.89%  | 0.249% |
| 66 | 10.216 | 1226 | 1229 | 1233 | rVV4 | 312961  | 482857  | 1.31%  | 0.173% |
| 67 | 10.257 | 1233 | 1236 | 1243 | rVB  | 240350  | 405126  | 1.10%  | 0.145% |
| 68 | 10.334 | 1244 | 1249 | 1253 | rBV  | 329730  | 542298  | 1.48%  | 0.194% |
| 69 | 10.498 | 1272 | 1277 | 1280 | rVV4 | 217878  | 435803  | 1.19%  | 0.156% |
| 70 | 10.534 | 1280 | 1283 | 1289 | rVV6 | 268286  | 602946  | 1.64%  | 0.216% |
| 71 | 10.616 | 1289 | 1297 | 1303 | rVV2 | 1791778 | 4221135 | 11.49% | 1.513% |
| 72 | 10.675 | 1303 | 1307 | 1315 | rVV  | 2093261 | 3983990 | 10.84% | 1.428% |
| 73 | 10.745 | 1315 | 1319 | 1324 | rVV5 | 283439  | 599123  | 1.63%  | 0.215% |
| 74 | 10.798 | 1324 | 1328 | 1333 | rVV  | 571874  | 899008  | 2.45%  | 0.322% |
| 75 | 12.287 | 1572 | 1581 | 1589 | rVB  | 413244  | 751011  | 2.04%  | 0.269% |
| 76 | 12.510 | 1613 | 1619 | 1628 | rVB  | 221404  | 404345  | 1.10%  | 0.145% |

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 Sample : N1303-03  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 CC-6

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-BP011422.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|     |        |      |      |      |      |         |         |        |        |
|-----|--------|------|------|------|------|---------|---------|--------|--------|
| 77  | 13.092 | 1710 | 1718 | 1726 | rBV  | 4520275 | 6738043 | 18.33% | 2.415% |
| 78  | 14.469 | 1943 | 1952 | 1958 | rBV  | 1186516 | 1869674 | 5.09%  | 0.670% |
| 79  | 17.222 | 2414 | 2420 | 2424 | rBV  | 1347668 | 2096124 | 5.70%  | 0.751% |
| 80  | 17.269 | 2424 | 2428 | 2435 | rVV  | 1479768 | 2281996 | 6.21%  | 0.818% |
| 81  | 18.092 | 2561 | 2568 | 2572 | rBV  | 273282  | 517796  | 1.41%  | 0.186% |
| 82  | 18.280 | 2595 | 2600 | 2604 | rVV4 | 340851  | 648947  | 1.77%  | 0.233% |
| 83  | 18.404 | 2616 | 2621 | 2630 | rBV3 | 238363  | 467022  | 1.27%  | 0.167% |
| 84  | 18.492 | 2631 | 2636 | 2641 | rBV5 | 186535  | 408348  | 1.11%  | 0.146% |
| 85  | 18.545 | 2641 | 2645 | 2649 | rBV2 | 332380  | 534760  | 1.46%  | 0.192% |
| 86  | 18.674 | 2663 | 2667 | 2675 | rVB3 | 239272  | 419258  | 1.14%  | 0.150% |
| 87  | 18.816 | 2684 | 2691 | 2696 | rBV  | 2509948 | 3355972 | 9.13%  | 1.203% |
| 88  | 18.939 | 2707 | 2712 | 2716 | rBV  | 328425  | 510061  | 1.39%  | 0.183% |
| 89  | 19.239 | 2758 | 2763 | 2768 | rBV  | 2035785 | 2710241 | 7.37%  | 0.971% |
| 90  | 19.292 | 2768 | 2772 | 2776 | rVB  | 1161969 | 1420397 | 3.86%  | 0.509% |
| 91  | 19.586 | 2814 | 2822 | 2831 | rBV  | 1782620 | 3191443 | 8.68%  | 1.144% |
| 92  | 19.786 | 2851 | 2856 | 2861 | rVB  | 8024399 | 9739008 | 26.50% | 3.490% |
| 93  | 19.998 | 2888 | 2892 | 2896 | rVB  | 579574  | 684169  | 1.86%  | 0.245% |
| 94  | 20.074 | 2902 | 2905 | 2912 | rVB2 | 389707  | 710091  | 1.93%  | 0.254% |
| 95  | 21.033 | 3065 | 3068 | 3075 | rVB3 | 612334  | 973045  | 2.65%  | 0.349% |
| 96  | 21.316 | 3110 | 3116 | 3120 | rBV2 | 2720947 | 4860386 | 13.23% | 1.742% |
| 97  | 21.351 | 3120 | 3122 | 3131 | rVB  | 1093265 | 1429554 | 3.89%  | 0.512% |
| 98  | 22.992 | 3396 | 3401 | 3407 | rBV  | 1012633 | 2326252 | 6.33%  | 0.834% |
| 99  | 23.621 | 3503 | 3508 | 3515 | rVB  | 810521  | 1618181 | 4.40%  | 0.580% |
| 100 | 23.727 | 3520 | 3526 | 3532 | rVV2 | 1427490 | 2828047 | 7.70%  | 1.013% |

Sum of corrected areas: 279043034



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Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

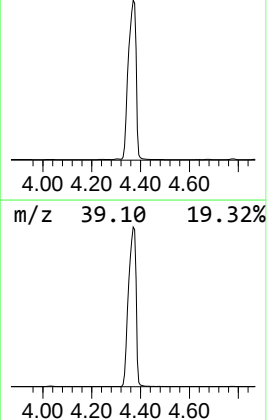
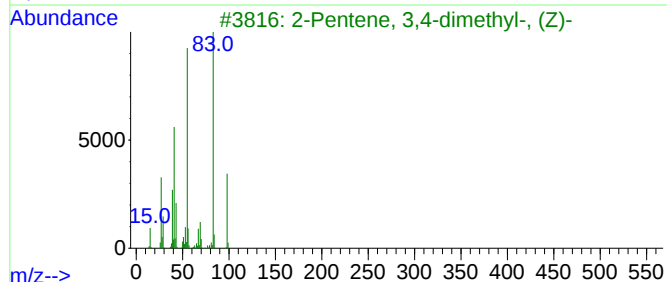
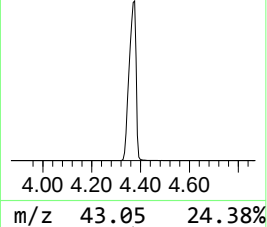
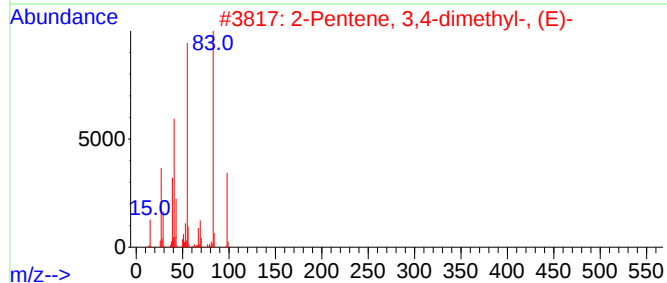
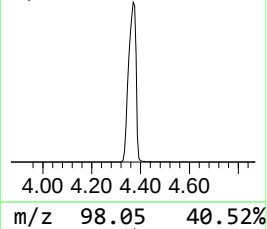
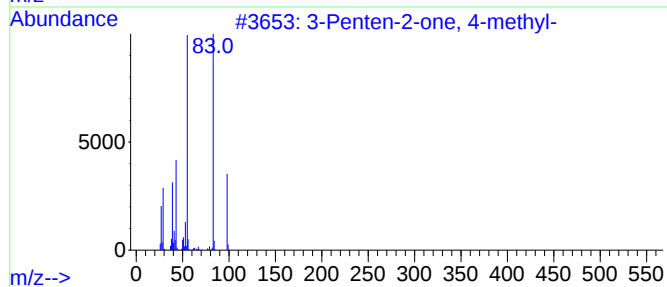
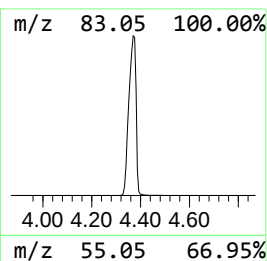
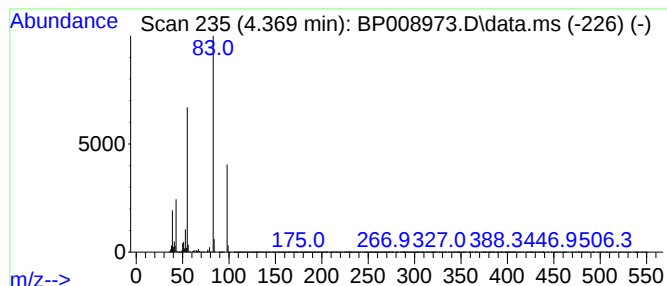
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 1

| R.T.  | EstConc  | Area     | Relative to ISTD       | R.T.  |
|-------|----------|----------|------------------------|-------|
| 4.369 | 95.52 ng | 36751200 | 1,4-Dichlorobenzene-d4 | 7.828 |

| Hit# | of | 5 | Tentative ID                   | MW | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|----|---------|-------------|------|
| 1    |    |   | 3-Penten-2-one, 4-methyl-      | 98 | C6H10O  | 000141-79-7 | 91   |
| 2    |    |   | 2-Pentene, 3,4-dimethyl-, (E)- | 98 | C7H14   | 004914-92-5 | 83   |
| 3    |    |   | 2-Pentene, 3,4-dimethyl-, (Z)- | 98 | C7H14   | 004914-91-4 | 78   |
| 4    |    |   | Furan, 2-methoxy-              | 98 | C5H6O2  | 025414-22-6 | 78   |
| 5    |    |   | 2-Pentene, 2,3-dimethyl-       | 98 | C7H14   | 010574-37-5 | 78   |



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Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

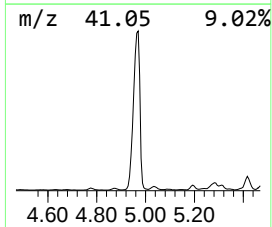
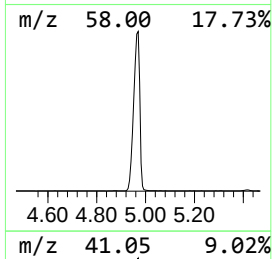
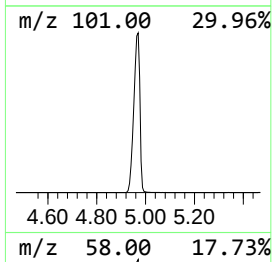
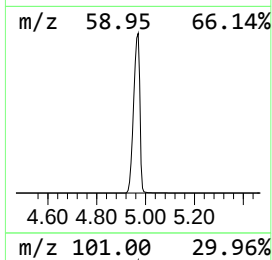
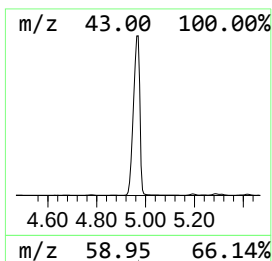
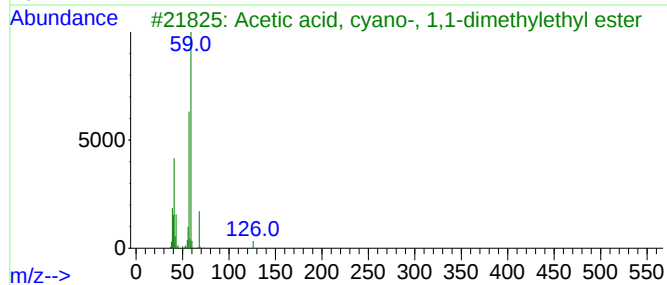
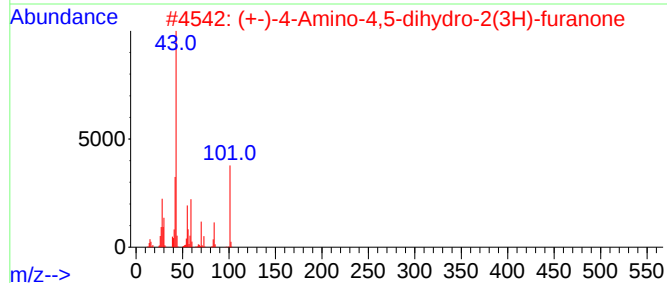
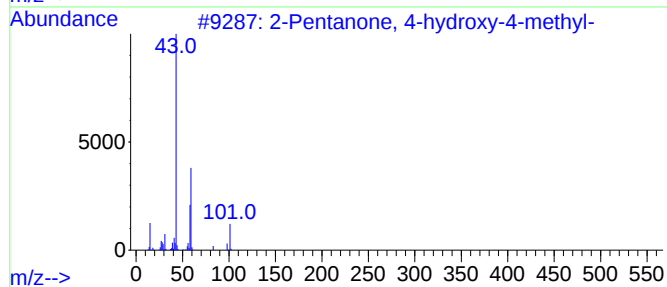
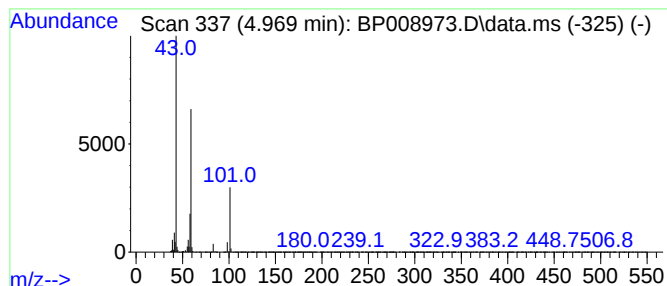
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

| R.T.  | EstConc  | Area     | Relative to ISTD       | R.T.  |
|-------|----------|----------|------------------------|-------|
| 4.969 | 38.54 ng | 14828000 | 1,4-Dichlorobenzene-d4 | 7.828 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 56   |
| 2    |    |   | (+)-4-Amino-4,5-dihydro-2(3H)-f...  | 101 | C4H7NO2  | 016504-58-8 | 47   |
| 3    |    |   | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 17   |
| 4    |    |   | 1-Propen-2-ol, acetate              | 100 | C5H8O2   | 000108-22-5 | 10   |
| 5    |    |   | Morpholine, 4-methyl-               | 101 | C5H11NO  | 000109-02-4 | 9    |



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Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

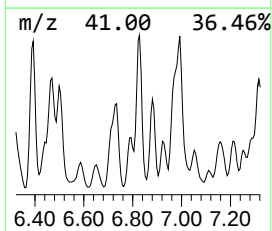
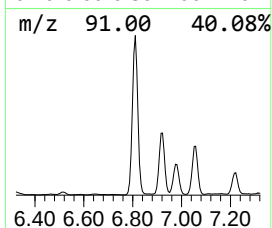
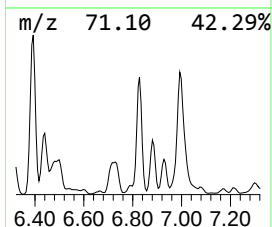
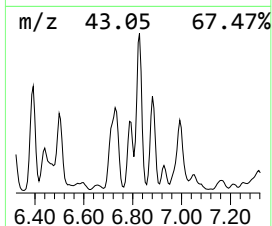
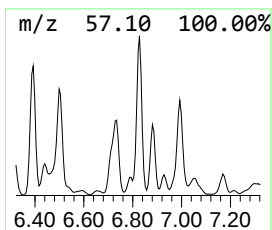
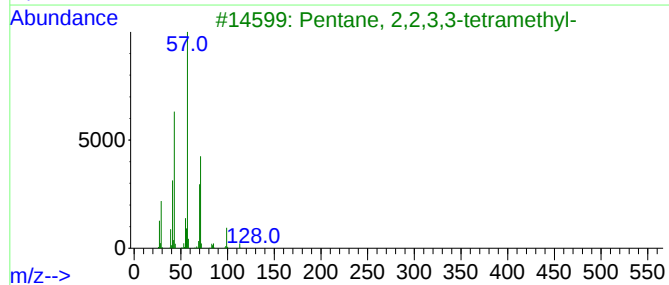
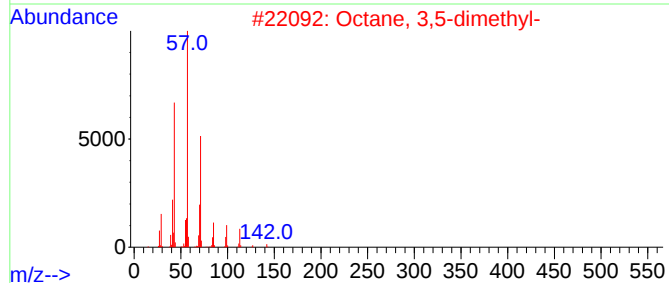
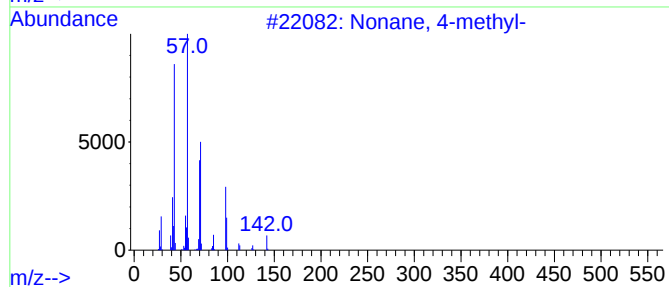
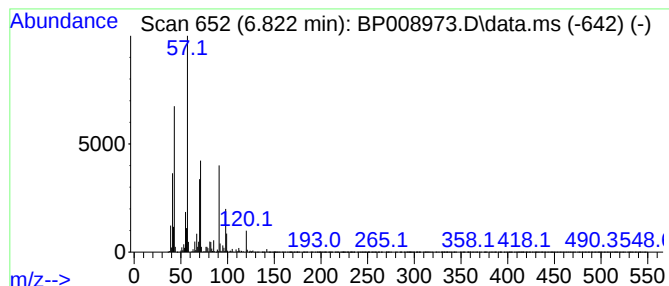
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 Nonane, 4-methyl- Concentration Rank 16

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 6.822 | 10.73 ng | 4129570 | 1,4-Dichlorobenzene-d4 | 7.828 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Nonane, 4-methyl-             | 142 | C10H22  | 017301-94-9 | 64   |
| 2    |    |   | Octane, 3,5-dimethyl-         | 142 | C10H22  | 015869-93-9 | 53   |
| 3    |    |   | Pentane, 2,2,3,3-tetramethyl- | 128 | C9H20   | 007154-79-2 | 50   |
| 4    |    |   | Undecane, 4,5-dimethyl-       | 184 | C13H28  | 017312-79-7 | 50   |
| 5    |    |   | Heptane, 2,3,4-trimethyl-     | 142 | C10H22  | 052896-95-4 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
Data File : BP008973.D  
Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

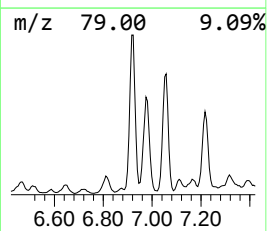
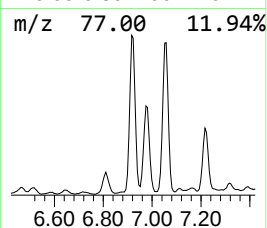
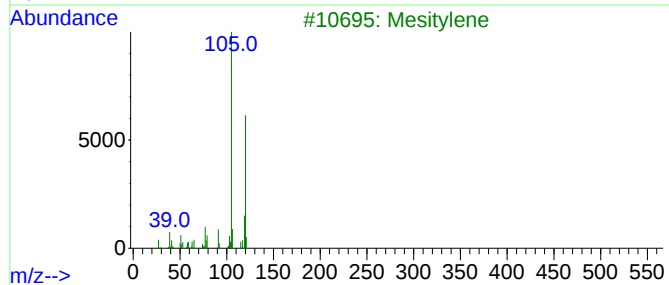
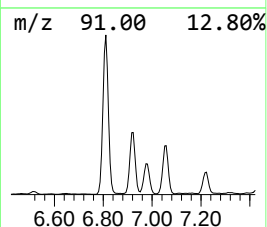
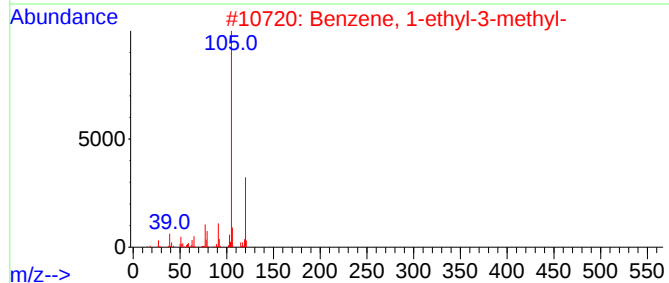
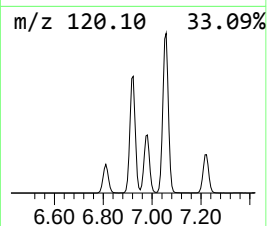
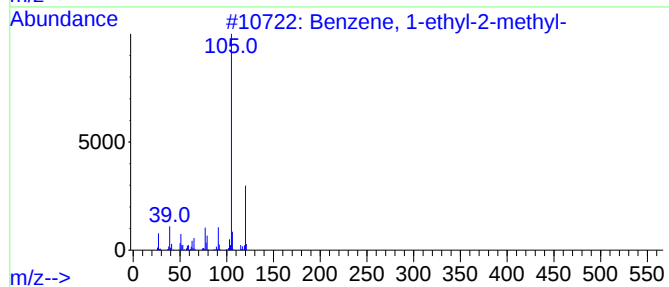
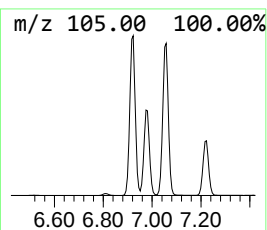
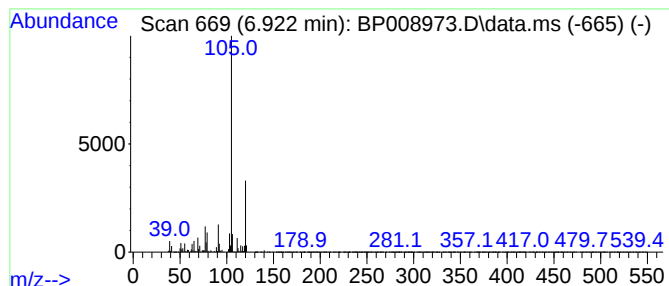
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 18

| R.T.  | EstConc | Area    | Relative to ISTD       | R.T.  |
|-------|---------|---------|------------------------|-------|
| 6.922 | 8.68 ng | 3340560 | 1,4-Dichlorobenzene-d4 | 7.828 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 95   |
| 2    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 95   |
| 3    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 91   |
| 4    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |
| 5    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
Data File : BP008973.D  
Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

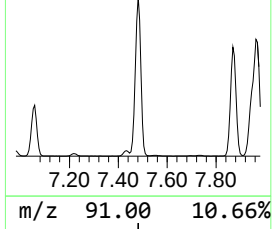
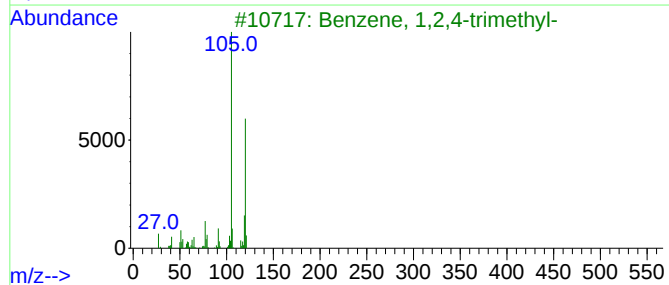
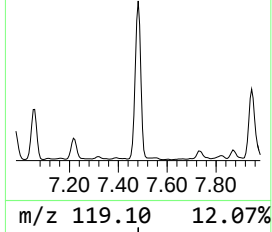
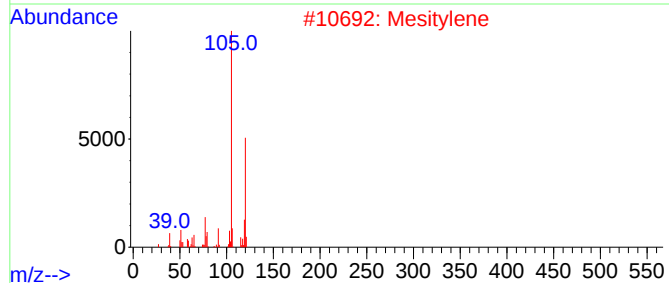
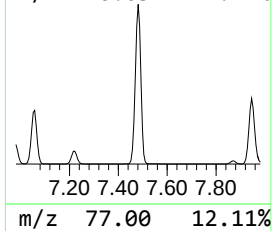
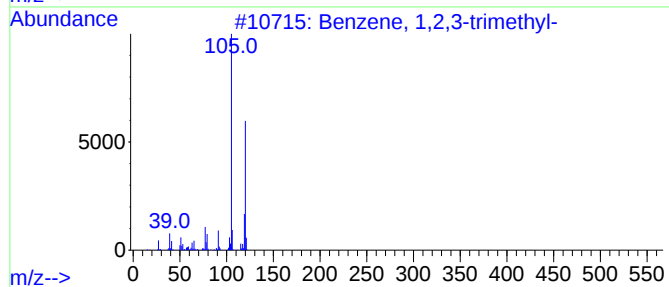
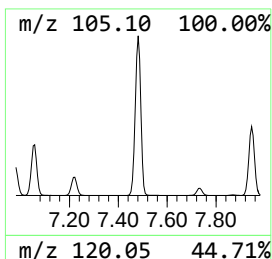
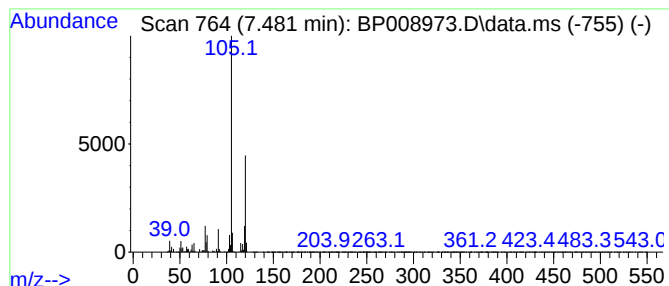
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 2

| R.T.  | EstConc  | Area     | Relative to ISTD       | R.T.  |
|-------|----------|----------|------------------------|-------|
| 7.481 | 40.88 ng | 15729000 | 1,4-Dichlorobenzene-d4 | 7.828 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 97   |
| 2    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 97   |
| 3    |    |   | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 95   |
| 4    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 90   |
| 5    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
Data File : BP008973.D  
Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

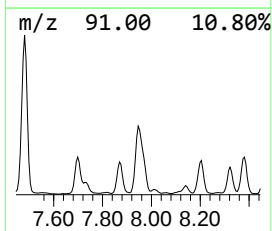
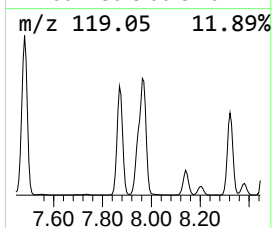
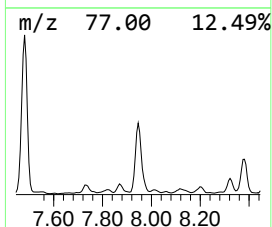
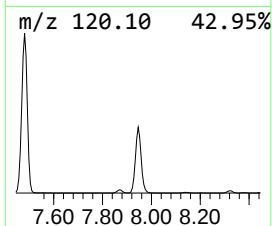
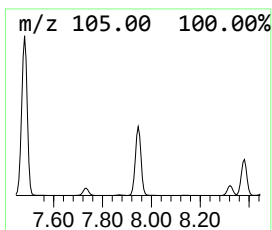
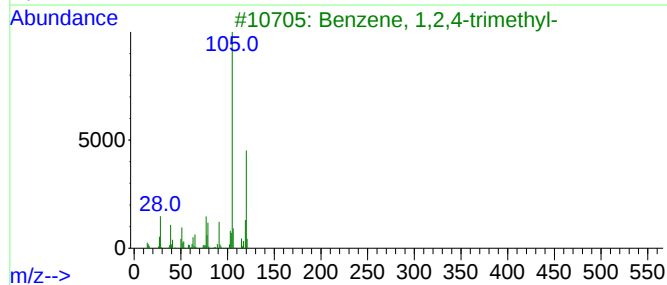
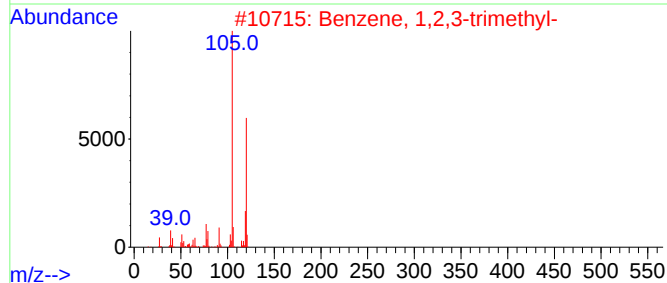
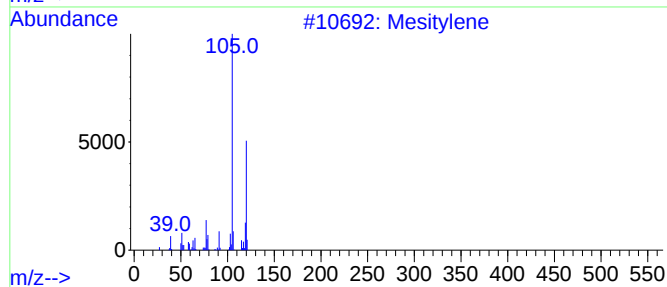
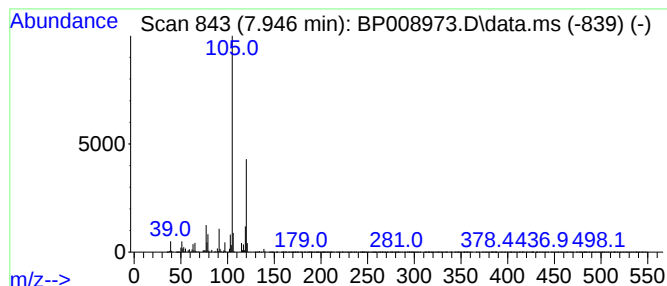
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 Mesitylene Concentration Rank 8

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 7.946 | 13.35 ng | 5138060 | 1,4-Dichlorobenzene-d4 | 7.828 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | 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 | 94   |
| 4    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 90   |
| 5    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
Data File : BP008973.D  
Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

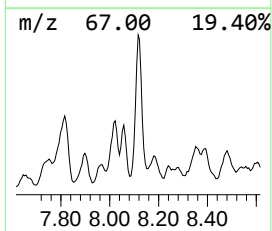
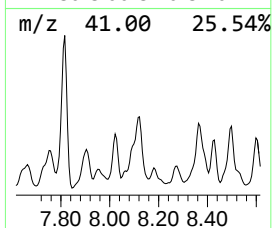
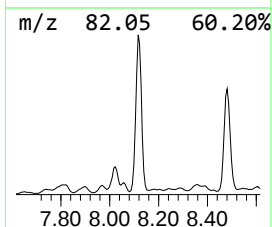
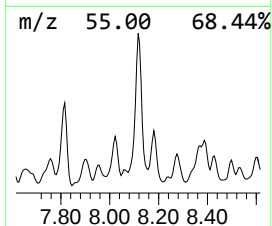
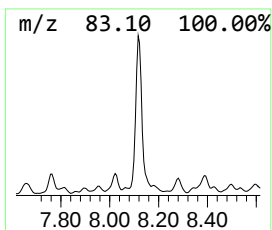
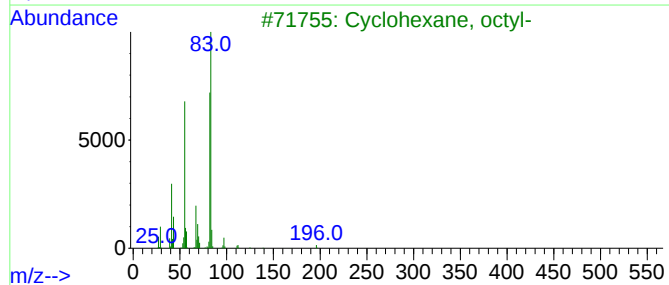
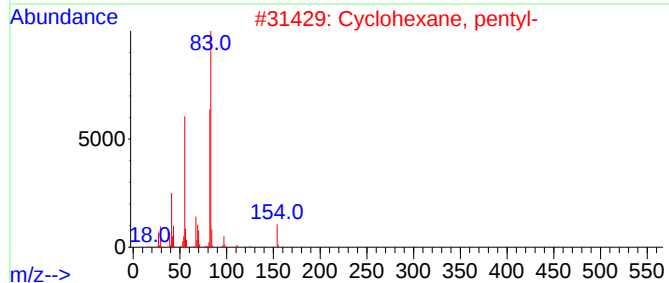
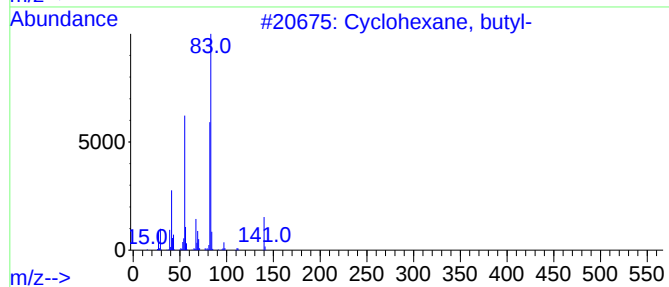
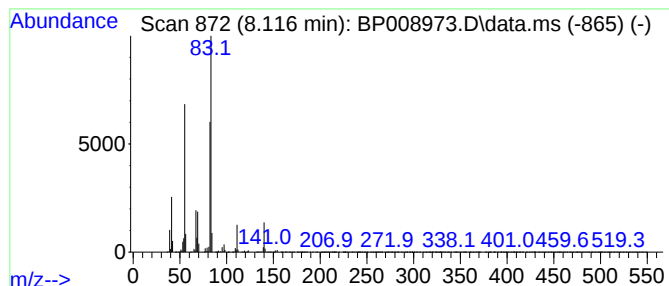
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 Cyclohexane, butyl- Concentration Rank 12

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 8.116 | 11.50 ng | 4424590 | 1,4-Dichlorobenzene-d4 | 7.828 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, butyl-            | 140 | C10H20  | 001678-93-9 | 87   |
| 2    |    |   | Cyclohexane, pentyl-           | 154 | C11H22  | 004292-92-6 | 74   |
| 3    |    |   | Cyclohexane, octyl-            | 196 | C14H28  | 001795-15-9 | 72   |
| 4    |    |   | Cyclohexane, 2-propenyl-       | 124 | C9H16   | 002114-42-3 | 64   |
| 5    |    |   | Cyclohexane, (2-methylpropyl)- | 140 | C10H20  | 001678-98-4 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
Data File : BP008973.D  
Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

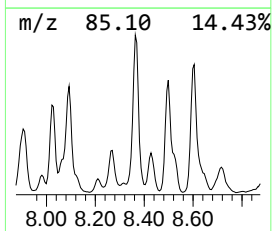
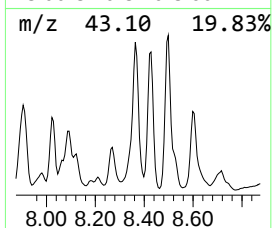
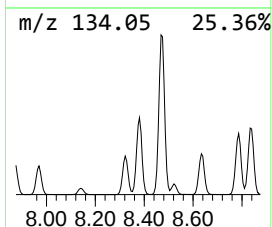
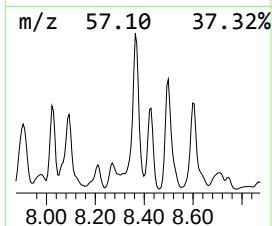
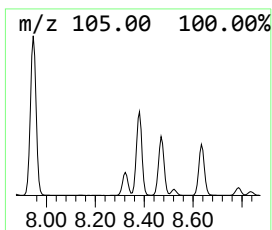
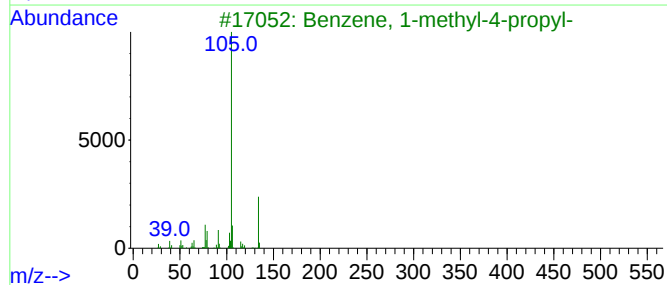
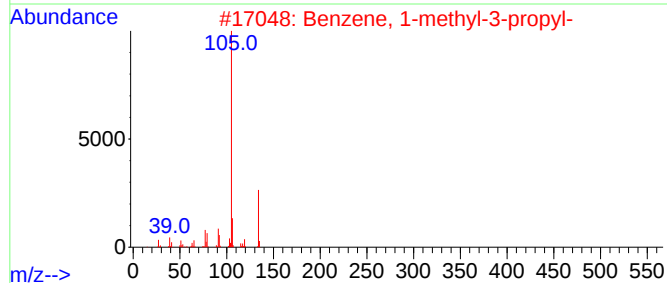
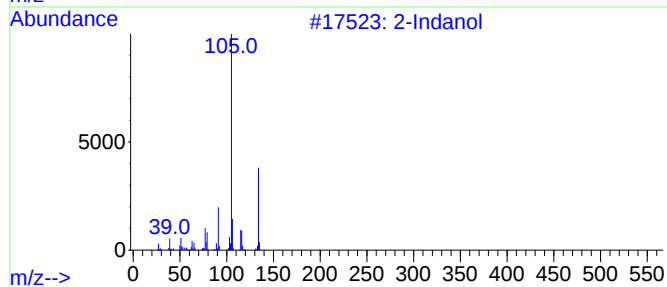
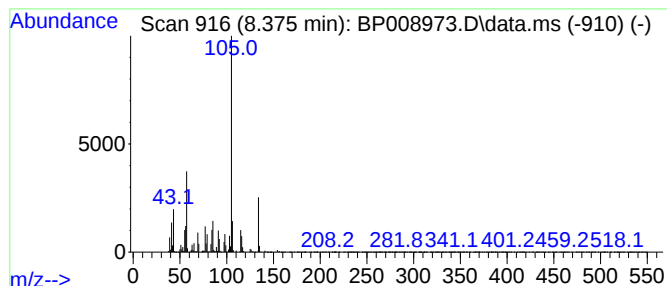
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 2-Indanol Concentration Rank 10

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 8.375 | 12.32 ng | 4738890 | 1,4-Dichlorobenzene-d4 | 7.828 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 2-Indanol                           | 134 | C9H10O  | 004254-29-9 | 62   |
| 2    |    |   | Benzene, 1-methyl-3-propyl-         | 134 | C10H14  | 001074-43-7 | 55   |
| 3    |    |   | Benzene, 1-methyl-4-propyl-         | 134 | C10H14  | 001074-55-1 | 55   |
| 4    |    |   | Benzene, 1-methyl-2-propyl-         | 134 | C10H14  | 001074-17-5 | 49   |
| 5    |    |   | Benzeneacetaldehyde, .alpha.-met... | 134 | C9H10O  | 000093-53-8 | 49   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
Data File : BP008973.D  
Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

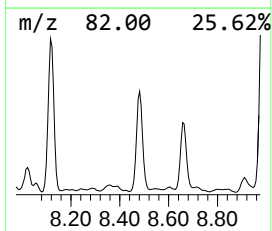
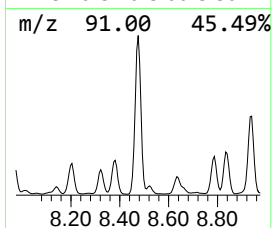
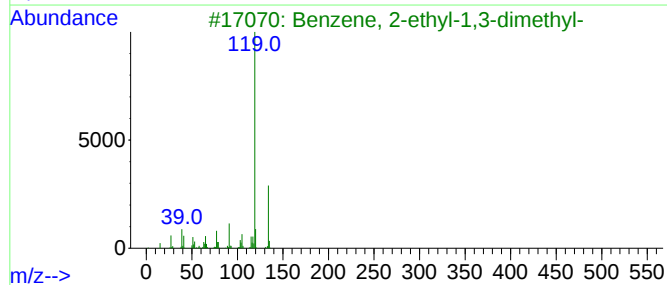
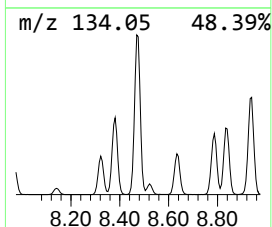
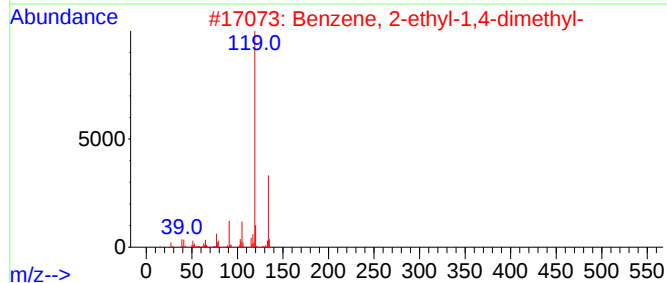
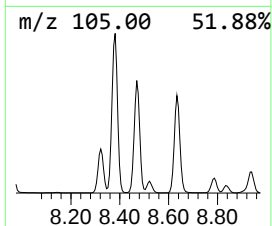
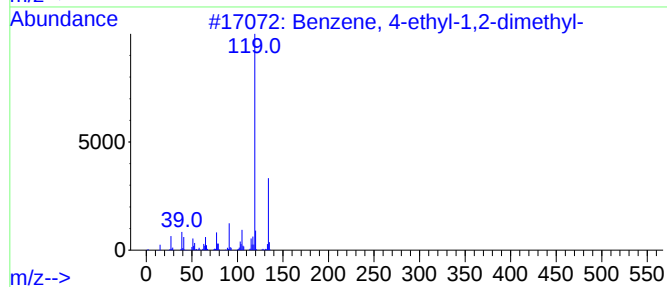
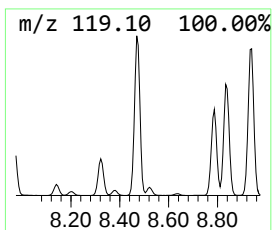
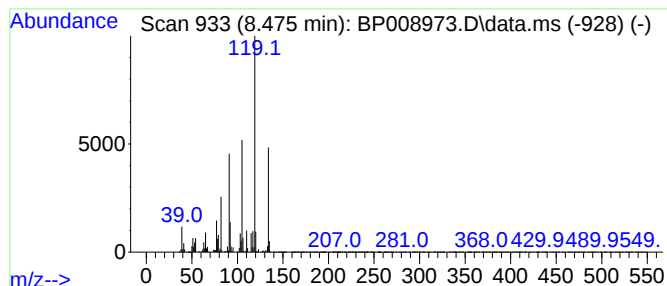
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 6

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 8.475 | 17.46 ng | 6715580 | 1,4-Dichlorobenzene-d4 | 7.828 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 93   |
| 2    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 91   |
| 3    |    |   | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 81   |
| 4    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 81   |
| 5    |    |   | Benzene, 1,4-diethyl-          | 134 | C10H14  | 000105-05-5 | 81   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
Data File : BP008973.D  
Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

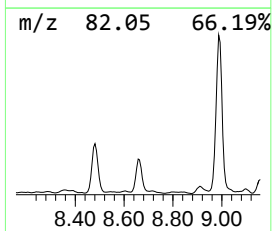
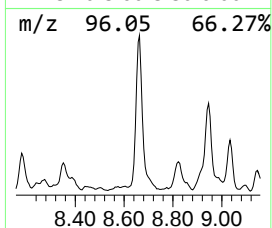
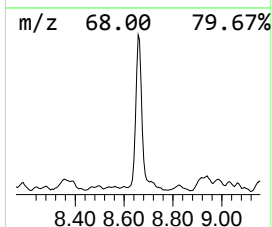
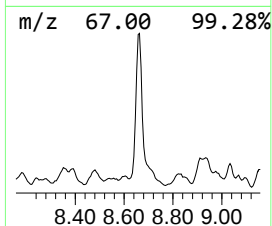
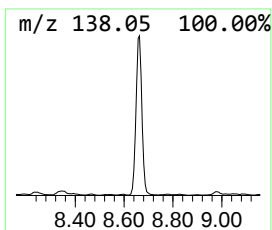
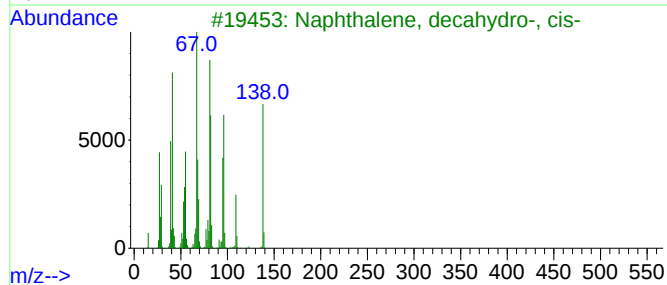
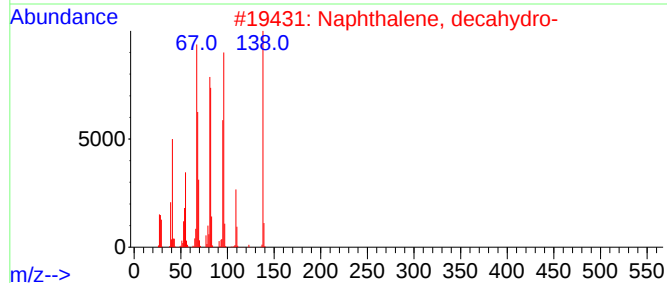
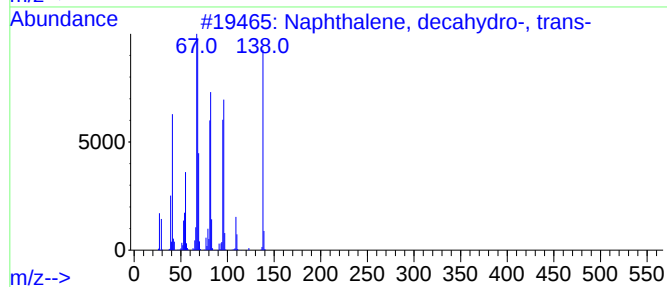
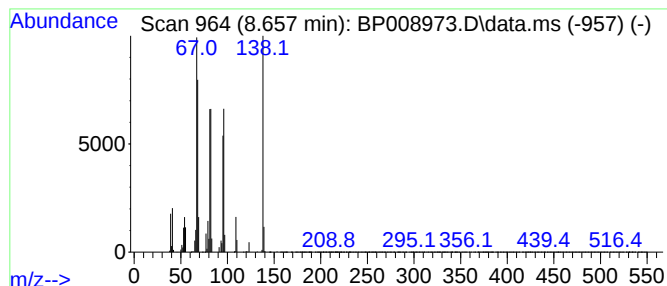
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 11 Naphthalene, decahydro-, tr... Concentration Rank 20

| R.T.  | EstConc | Area    | Relative to ISTD       | R.T.  |
|-------|---------|---------|------------------------|-------|
| 8.657 | 8.06 ng | 3099140 | 1,4-Dichlorobenzene-d4 | 7.828 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, decahydro-, trans-     | 138 | C10H18  | 000493-02-7 | 95   |
| 2    |    |   | Naphthalene, decahydro-             | 138 | C10H18  | 000091-17-8 | 94   |
| 3    |    |   | Naphthalene, decahydro-, cis-       | 138 | C10H18  | 000493-01-6 | 72   |
| 4    |    |   | Cyclopentanone, 2-(2-methylpropy... | 138 | C9H14O  | 016856-72-7 | 70   |
| 5    |    |   | Spiro[4.5]decane                    | 138 | C10H18  | 000176-63-6 | 68   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
Data File : BP008973.D  
Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

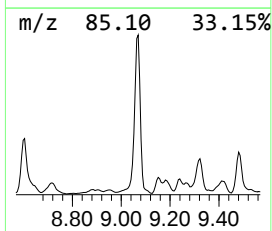
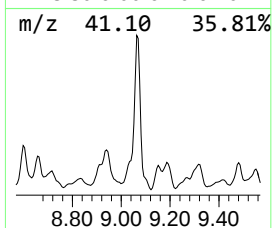
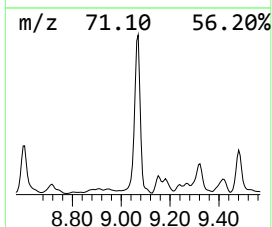
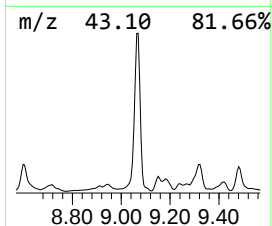
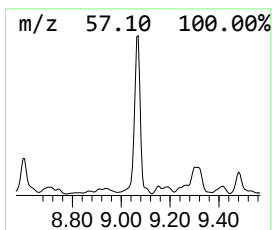
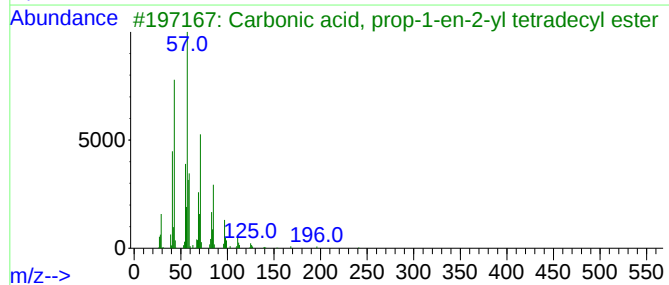
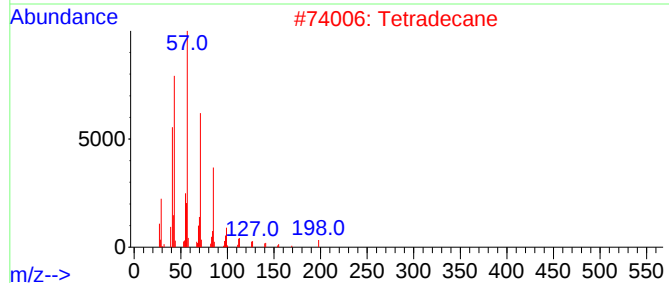
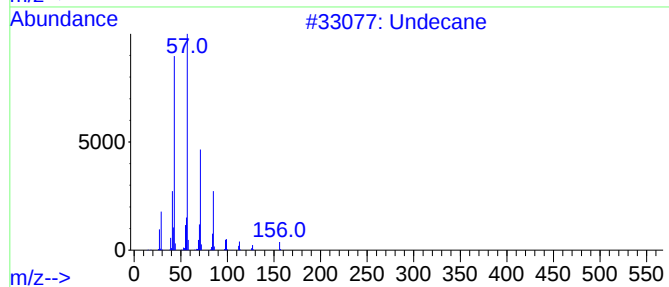
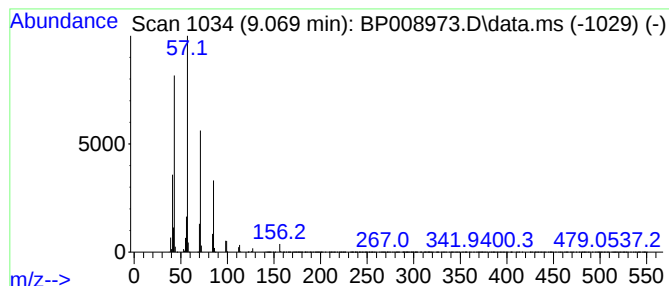
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Undecane Concentration Rank 5

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 9.069 | 20.81 ng | 8007800 | 1,4-Dichlorobenzene-d4 | 7.828 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Undecane                            | 156 | C11H24   | 001120-21-4  | 94   |
| 2    |    |   | Tetradecane                         | 198 | C14H30   | 000629-59-4  | 90   |
| 3    |    |   | Carbonic acid, prop-1-en-2-yl te... | 298 | C18H34O3 | 1000382-90-9 | 90   |
| 4    |    |   | Hexadecane                          | 226 | C16H34   | 000544-76-3  | 90   |
| 5    |    |   | Tridecane                           | 184 | C13H28   | 000629-50-5  | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
Data File : BP008973.D  
Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

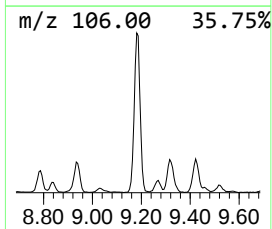
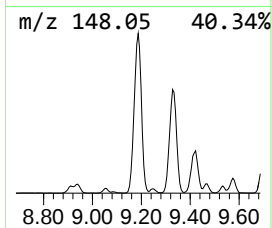
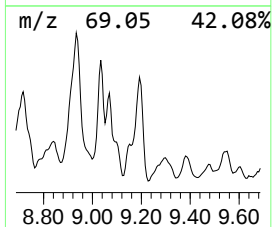
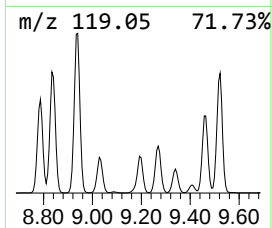
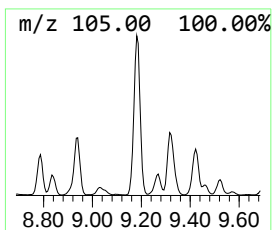
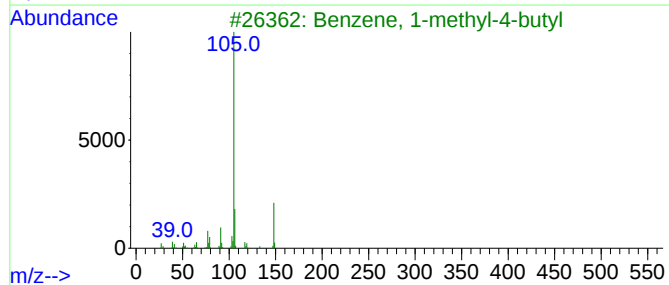
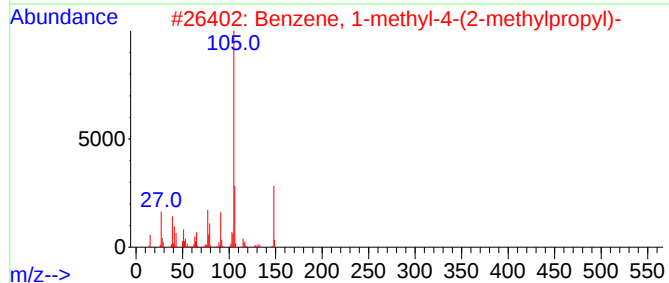
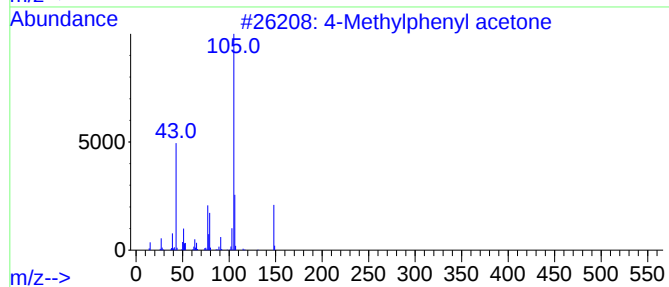
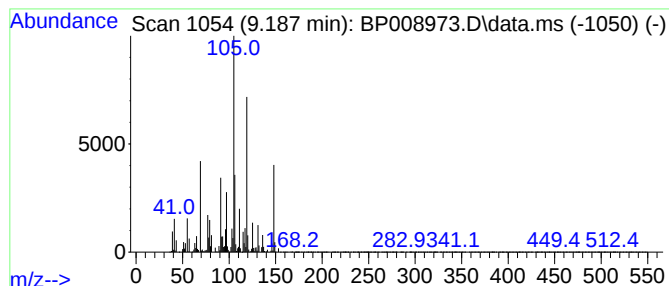
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 unknown9.187 Concentration Rank 17

| R.T.  | EstConc | Area    | Relative to ISTD       | R.T.  |
|-------|---------|---------|------------------------|-------|
| 9.187 | 9.51 ng | 3656960 | 1,4-Dichlorobenzene-d4 | 7.828 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | 4-Methylphenyl acetone              | 148 | C10H12O | 002096-86-8 | 46   |
| 2    |      | Benzene, 1-methyl-4-(2-methylpro... | 148 | C11H16  | 005161-04-6 | 41   |
| 3    |      | Benzene, 1-methyl-4-butyl           | 148 | C11H16  | 001595-05-7 | 38   |
| 4    |      | 3,4-Dihydro-2-[1H]quinoxalinone     | 148 | C8H8N2O | 059564-59-9 | 38   |
| 5    |      | 2-Methylindan-2-ol                  | 148 | C10H12O | 033223-84-6 | 30   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
Data File : BP008973.D  
Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

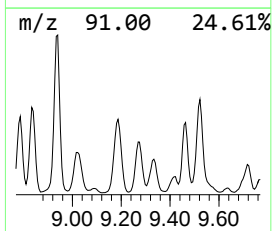
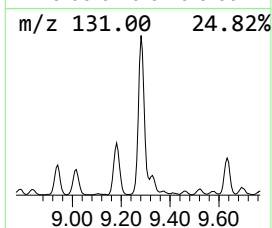
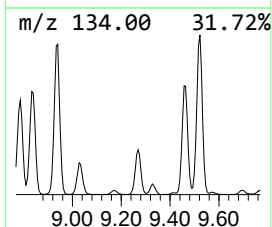
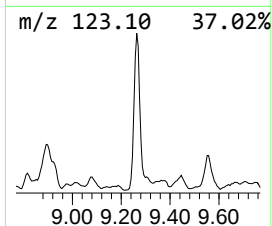
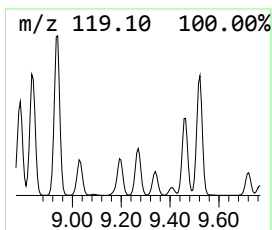
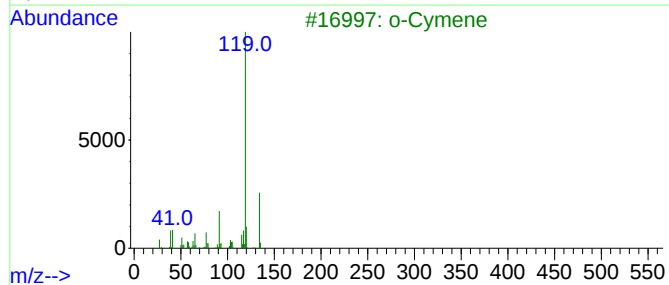
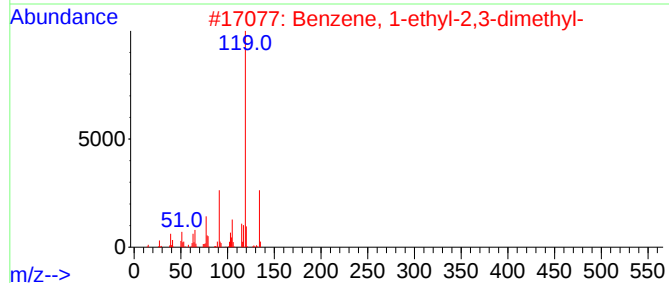
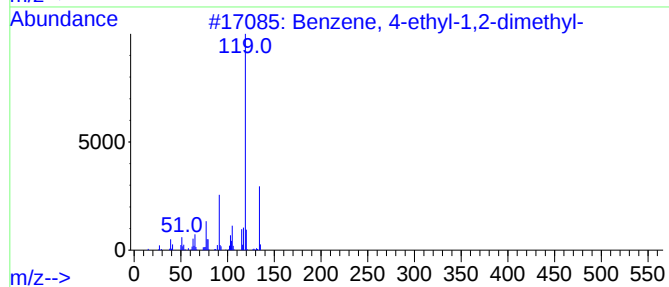
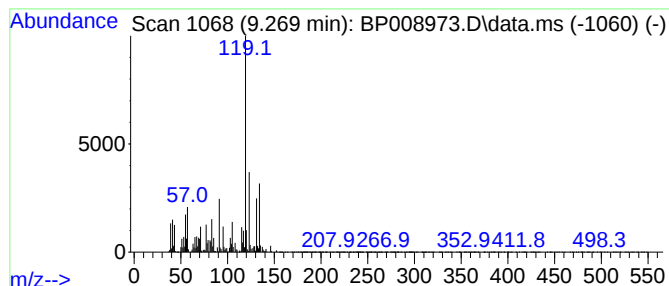
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Peak Number 14 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 15

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.   |
|-------|----------|---------|------------------|--------|
| 9.269 | 10.82 ng | 2283270 | Naphthalene-d8   | 10.628 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 96   |
| 2    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 95   |
| 3    |    |   | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 93   |
| 4    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 93   |
| 5    |    |   | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 92   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
Data File : BP008973.D  
Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

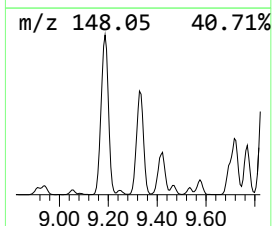
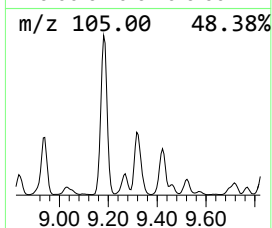
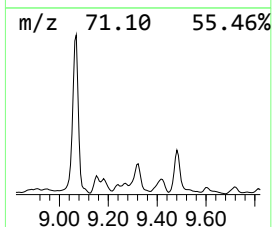
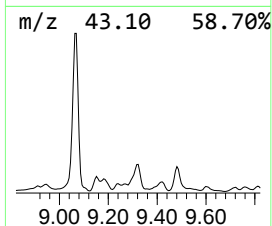
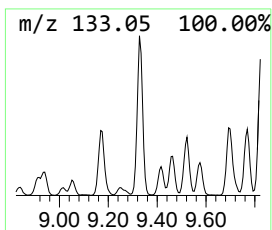
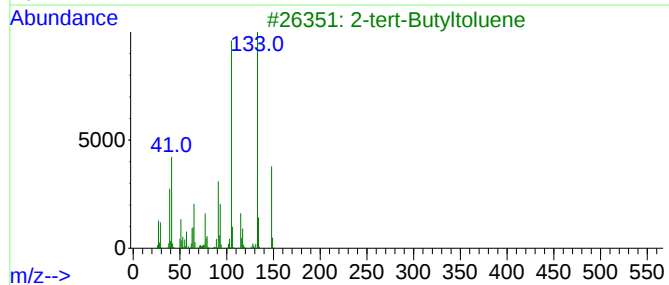
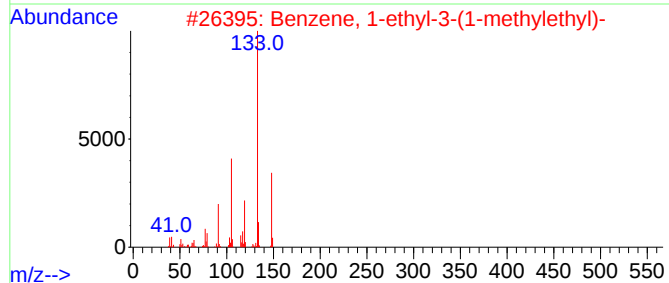
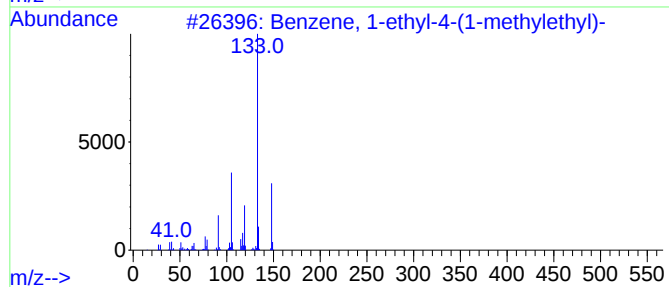
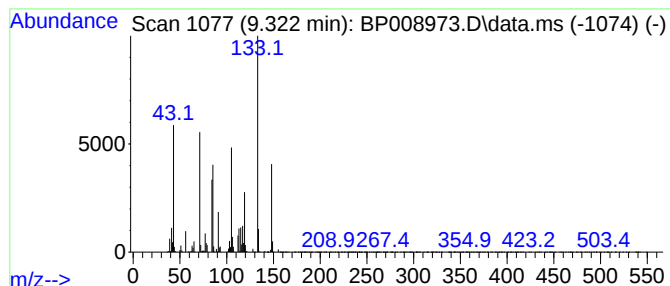
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 15 Benzene, 1-ethyl-4-(1-methy... Concentration Rank 19

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.   |
|-------|---------|---------|------------------|--------|
| 9.322 | 8.29 ng | 1748960 | Naphthalene-d8   | 10.628 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-4-(1-methylethyl)- | 148 | C11H16  | 004218-48-8 | 92   |
| 2    |    |   | Benzene, 1-ethyl-3-(1-methylethyl)- | 148 | C11H16  | 004920-99-4 | 90   |
| 3    |    |   | 2-tert-Butyltoluene                 | 148 | C11H16  | 001074-92-6 | 50   |
| 4    |    |   | Benzene, 1-ethyl-2,4,5-trimethyl-   | 148 | C11H16  | 017851-27-3 | 46   |
| 5    |    |   | Benzene, pentamethyl-               | 148 | C11H16  | 000700-12-9 | 46   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
Data File : BP008973.D  
Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

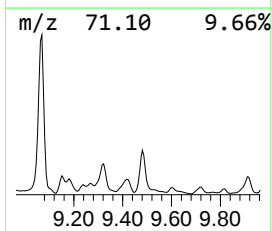
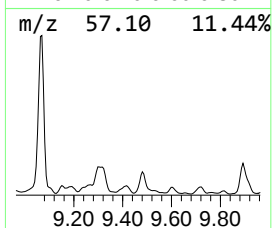
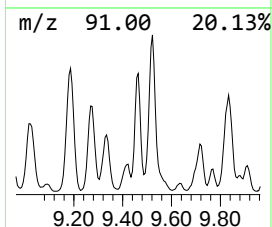
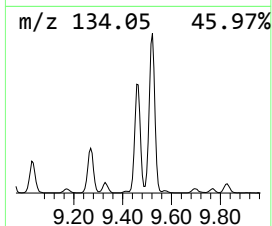
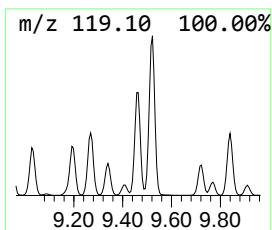
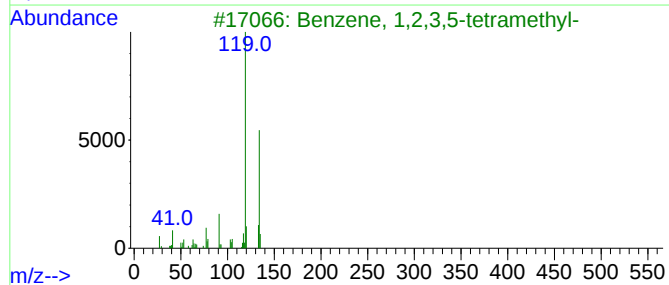
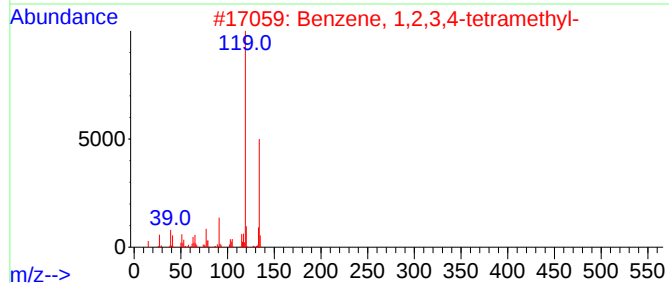
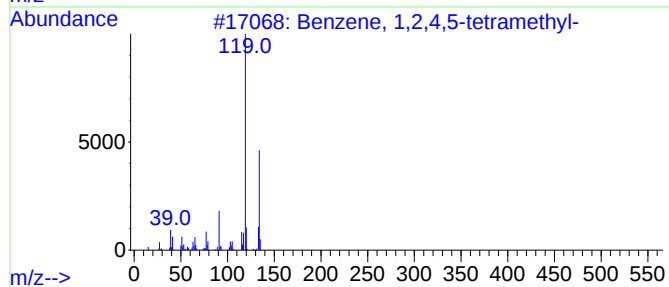
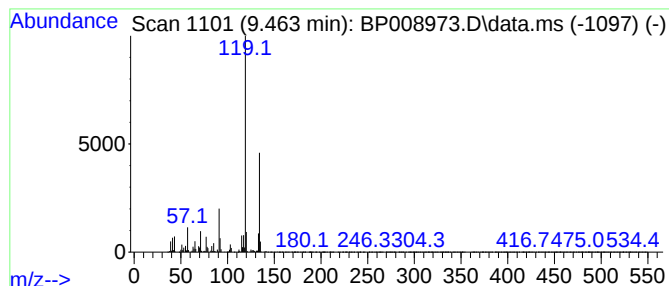
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 14

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.   |
|-------|----------|---------|------------------|--------|
| 9.463 | 11.17 ng | 2357960 | Naphthalene-d8   | 10.628 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,4,5-tetramethyl-       | 134 | C10H14  | 000095-93-2 | 93   |
| 2    |    |   | Benzene, 1,2,3,4-tetramethyl-       | 134 | C10H14  | 000488-23-3 | 90   |
| 3    |    |   | Benzene, 1,2,3,5-tetramethyl-       | 134 | C10H14  | 000527-53-7 | 90   |
| 4    |    |   | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 90   |
| 5    |    |   | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
Data File : BP008973.D  
Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

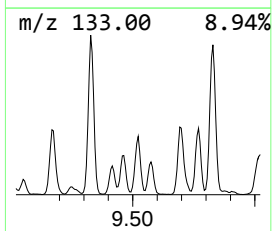
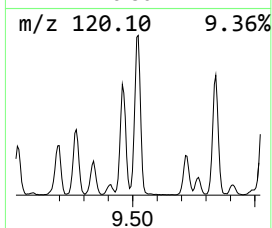
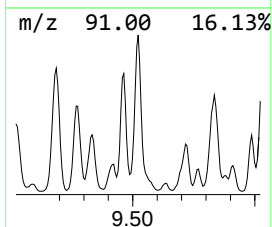
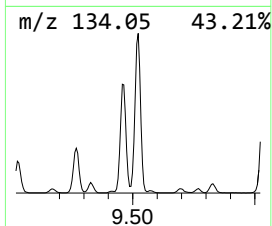
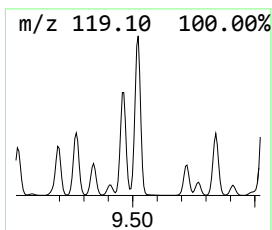
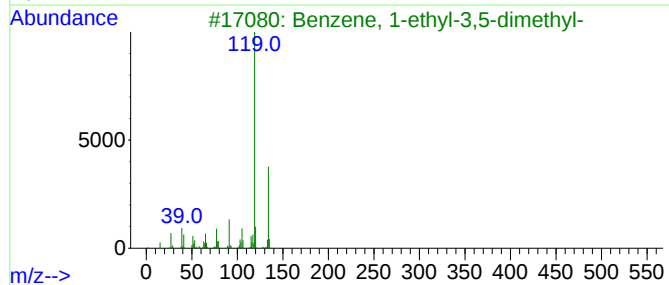
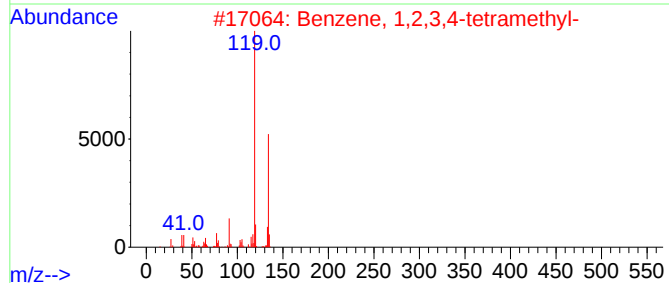
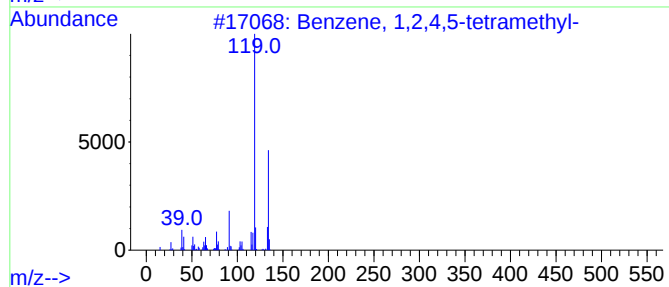
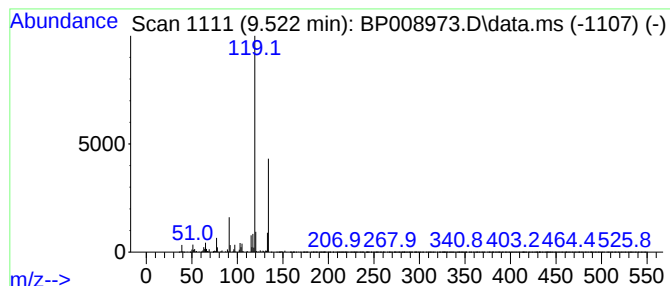
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 11

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.   |
|-------|----------|---------|------------------|--------|
| 9.522 | 11.69 ng | 2466340 | Naphthalene-d8   | 10.628 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,4,5-tetramethyl-  | 134 | C10H14  | 000095-93-2 | 95   |
| 2    |    |   | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 94   |
| 3    |    |   | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 91   |
| 4    |    |   | Benzene, 1,2,3,5-tetramethyl-  | 134 | C10H14  | 000527-53-7 | 91   |
| 5    |    |   | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
Data File : BP008973.D  
Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

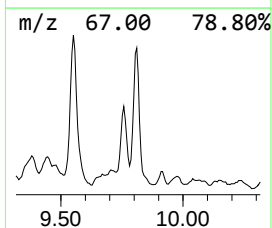
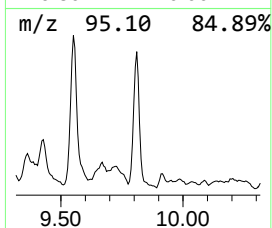
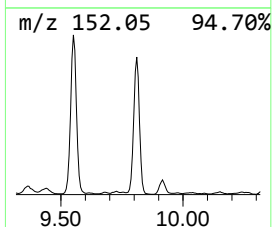
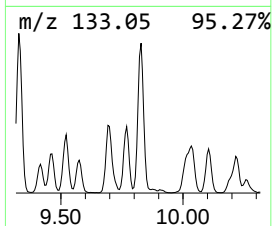
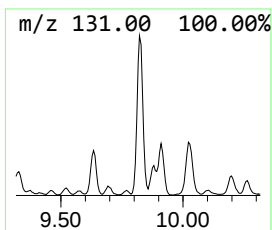
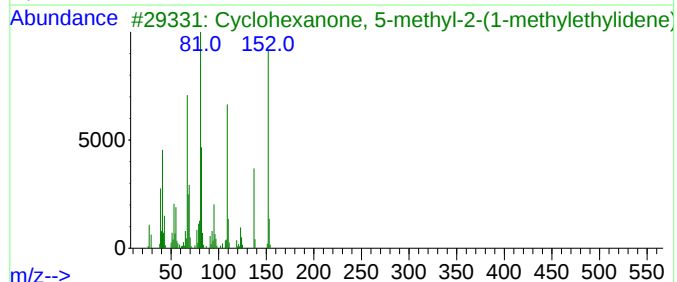
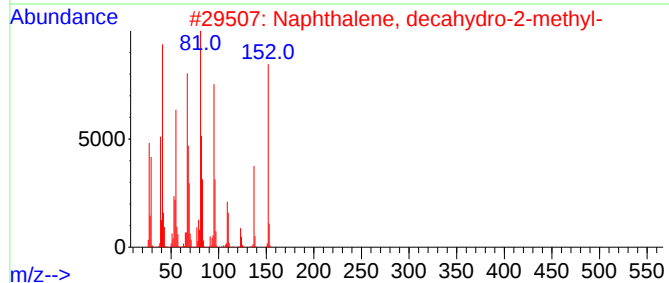
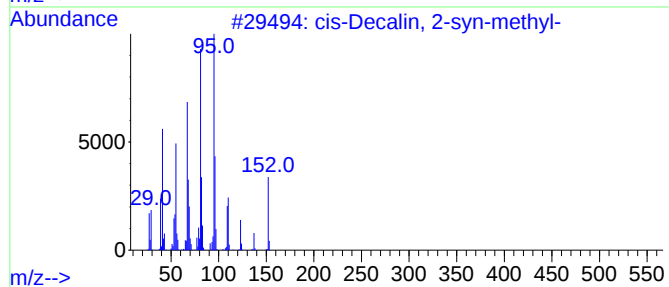
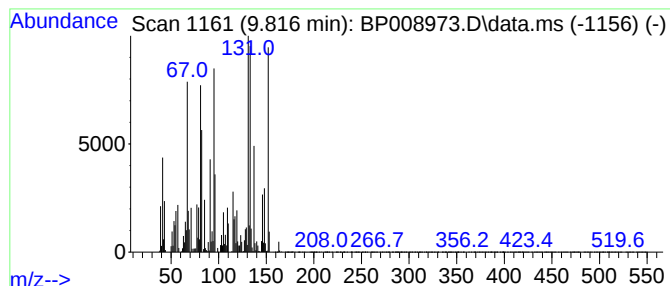
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 18 unknown9.816 Concentration Rank 9

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.   |
|-------|----------|---------|------------------|--------|
| 9.816 | 12.54 ng | 2646020 | Naphthalene-d8   | 10.628 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|------|-------------------------------------|-----|---------|--------------|------|
| 1    |      | cis-Decalin, 2-syn-methyl-          | 152 | C11H20  | 1000155-85-6 | 49   |
| 2    |      | Naphthalene, decahydro-2-methyl-    | 152 | C11H20  | 002958-76-1  | 46   |
| 3    |      | Cyclohexanone, 5-methyl-2-(1-met... | 152 | C10H16O | 015932-80-6  | 38   |
| 4    |      | Bicyclo[4.1.0]heptan-3-one, 4,7,... | 152 | C10H16O | 004176-01-6  | 38   |
| 5    |      | Bicyclo[4.1.0]heptan-3-one, 4,7,... | 152 | C10H16O | 004176-04-9  | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
Data File : BP008973.D  
Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

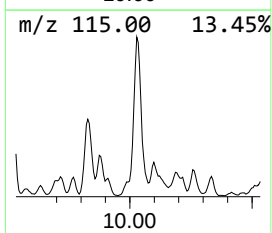
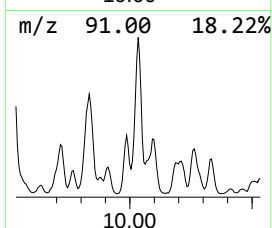
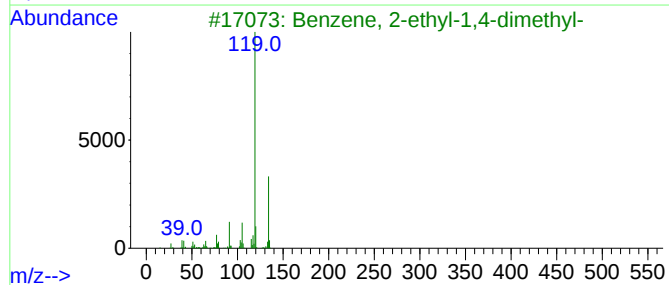
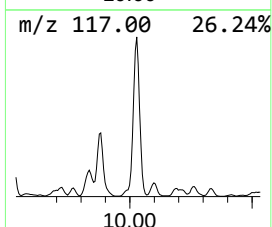
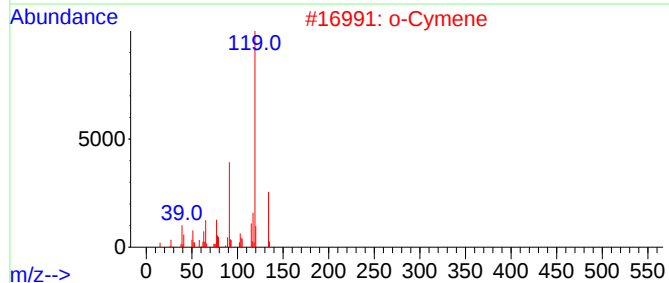
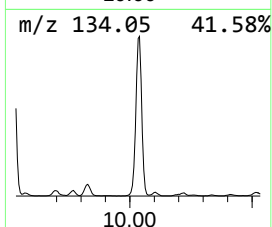
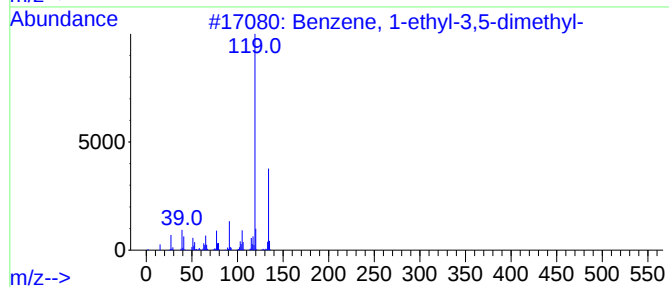
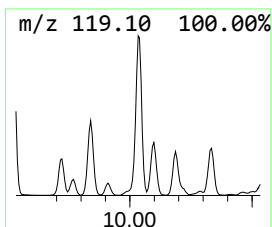
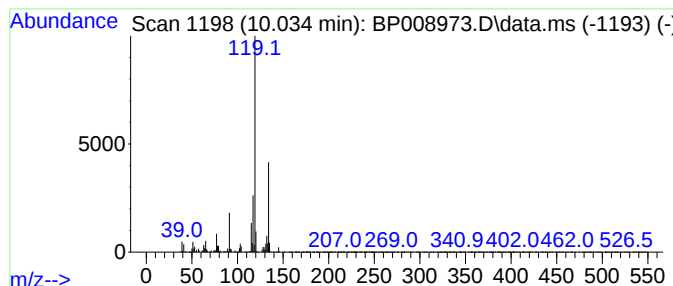
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 7

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 10.034 | 14.85 ng | 3133780 | Naphthalene-d8   | 10.628 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Benzene, 1-ethyl-3,5-dimethyl-      | 134 | C10H14  | 000934-74-7 | 87   |
| 2    |      | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 87   |
| 3    |      | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 83   |
| 4    |      | Benzene, 4-ethyl-1,2-dimethyl-      | 134 | C10H14  | 000934-80-5 | 83   |
| 5    |      | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
Data File : BP008973.D  
Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

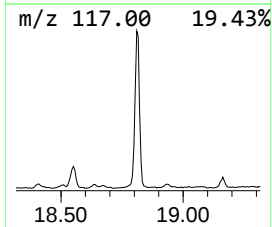
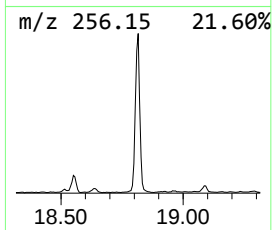
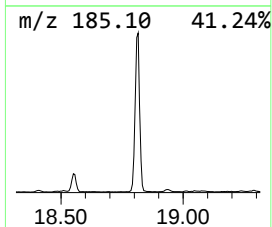
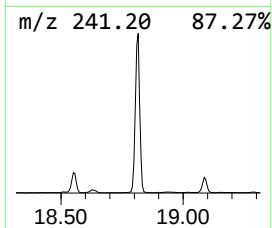
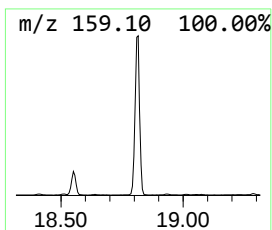
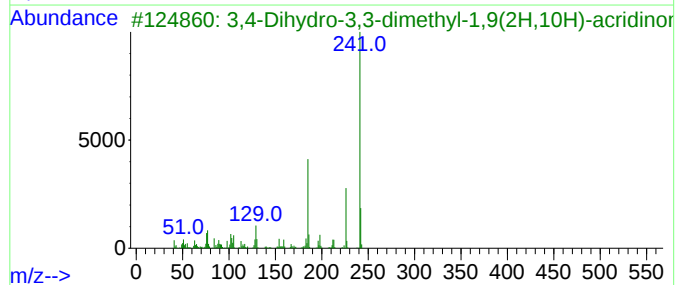
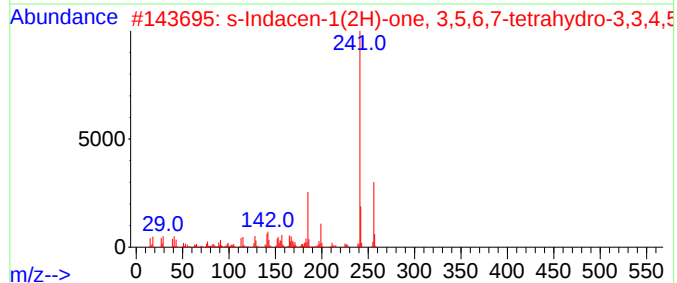
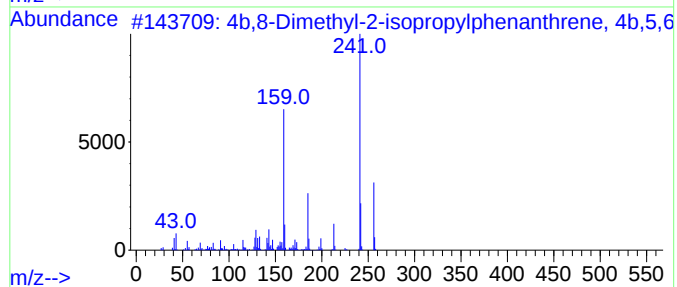
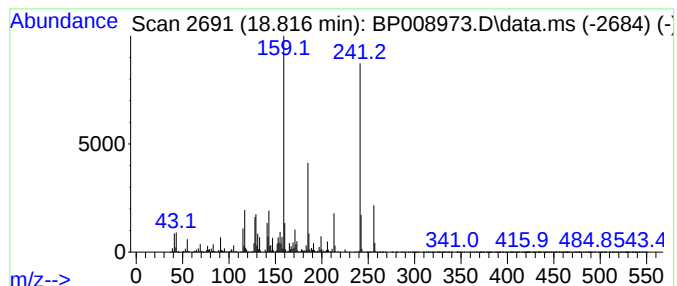
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 4b,8-Dimethyl-2-isopropylph... Concentration Rank 4

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 18.816 | 32.02 ng | 3355970 | Phenanthrene-d10 | 17.222 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 4b,8-Dimethyl-2-isopropylphenant... | 256 | C19H28     | 1000197-14-1 | 97   |
| 2    |    |   | s-Indacen-1(2H)-one, 3,5,6,7-tet... | 256 | C18H24O    | 038754-94-8  | 42   |
| 3    |    |   | 3,4-Dihydro-3,3-dimethyl-1,9(2H,... | 241 | C15H15NO2  | 1010212-95-2 | 38   |
| 4    |    |   | Isoquinoline-4-carbonitrile, 1-e... | 256 | C15H16N2O2 | 1000259-91-1 | 27   |
| 5    |    |   | Isoxazole, 5-methyl-4-phenyl-       | 159 | C10H9NO    | 023253-49-8  | 25   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013122\  
Data File : BP008973.D  
Acq On : 31 Jan 2022 19:43  
Operator : CG/JU  
Sample : N1303-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
CC-6

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.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 |
| 3-Penten-2-one,... | 4.369  | 95.5    | ng    | 36751200 | 1                     | 7.828  | 7694690 | 20.0 |
| 2-Pentanone, 4-... | 4.969  | 38.5    | ng    | 14828000 | 1                     | 7.828  | 7694690 | 20.0 |
| Nonane, 4-methyl-  | 6.822  | 10.7    | ng    | 4129570  | 1                     | 7.828  | 7694690 | 20.0 |
| Benzene, 1-ethy... | 6.922  | 8.7     | ng    | 3340560  | 1                     | 7.828  | 7694690 | 20.0 |
| Benzene, 1,2,3-... | 7.481  | 40.9    | ng    | 15729000 | 1                     | 7.828  | 7694690 | 20.0 |
| Mesitylene         | 7.946  | 13.4    | ng    | 5138060  | 1                     | 7.828  | 7694690 | 20.0 |
| Cyclohexane, bu... | 8.116  | 11.5    | ng    | 4424590  | 1                     | 7.828  | 7694690 | 20.0 |
| 2-Indanol          | 8.375  | 12.3    | ng    | 4738890  | 1                     | 7.828  | 7694690 | 20.0 |
| Benzene, 4-ethy... | 8.475  | 17.5    | ng    | 6715580  | 1                     | 7.828  | 7694690 | 20.0 |
| Naphthalene, de... | 8.657  | 8.1     | ng    | 3099140  | 1                     | 7.828  | 7694690 | 20.0 |
| Undecane           | 9.069  | 20.8    | ng    | 8007800  | 1                     | 7.828  | 7694690 | 20.0 |
| unknown9.187       | 9.187  | 9.5     | ng    | 3656960  | 1                     | 7.828  | 7694690 | 20.0 |
| Benzene, 1-ethy... | 9.269  | 10.8    | ng    | 2283270  | 2                     | 10.628 | 4221140 | 20.0 |
| Benzene, 1-ethy... | 9.322  | 8.3     | ng    | 1748960  | 2                     | 10.628 | 4221140 | 20.0 |
| Benzene, 1,2,4,... | 9.463  | 11.2    | ng    | 2357960  | 2                     | 10.628 | 4221140 | 20.0 |
| Benzene, 1,2,3,... | 9.522  | 11.7    | ng    | 2466340  | 2                     | 10.628 | 4221140 | 20.0 |
| unknown9.816       | 9.816  | 12.5    | ng    | 2646020  | 2                     | 10.628 | 4221140 | 20.0 |
| Benzene, 1-ethy... | 10.034 | 14.9    | ng    | 3133780  | 2                     | 10.628 | 4221140 | 20.0 |
| 4b,8-Dimethyl-2... | 18.816 | 32.0    | ng    | 3355970  | 4                     | 17.222 | 2096120 | 20.0 |