

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120215\  
 Data File : BF083420.D  
 Acq On : 2 Dec 2015 16:12  
 Operator : UM/IZ  
 Sample : G4583-06  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-P3WC-06(0-5)

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0  
 Filtering: 5  
 Min Area: 3 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
 Peak separation: 5

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113015.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 4.647       | 257           | 259         | 275          | rVB      | 791672         | 1312685       | 16.45%          | 1.632%        |
| 2         | 5.127       | 299           | 301         | 315          | rBV      | 3254753        | 3672992       | 46.04%          | 4.568%        |
| 3         | 6.053       | 378           | 382         | 384          | rBV2     | 75005          | 122342        | 1.53%           | 0.152%        |
| 4         | 6.133       | 387           | 389         | 392          | rVB      | 129667         | 120636        | 1.51%           | 0.150%        |
| 5         | 6.213       | 394           | 396         | 398          | rBV      | 4217772        | 4201871       | 52.66%          | 5.226%        |
| 6         | 6.316       | 403           | 405         | 407          | rVV      | 4030257        | 4005707       | 50.21%          | 4.982%        |
| 7         | 6.361       | 407           | 409         | 413          | rVB2     | 217039         | 306423        | 3.84%           | 0.381%        |
| 8         | 6.533       | 423           | 424         | 427          | rBV      | 678437         | 934536        | 11.71%          | 1.162%        |
| 9         | 6.693       | 436           | 438         | 443          | rBV2     | 3090953        | 3556232       | 44.57%          | 4.423%        |
| 10        | 6.876       | 452           | 454         | 455          | rBV      | 131376         | 133008        | 1.67%           | 0.165%        |
| 11        | 7.116       | 472           | 475         | 477          | rVV      | 2291854        | 2860429       | 35.85%          | 3.557%        |
| 12        | 7.253       | 485           | 487         | 491          | rVV2     | 59894          | 156686        | 1.96%           | 0.195%        |
| 13        | 7.356       | 494           | 496         | 498          | rVB      | 138121         | 160023        | 2.01%           | 0.199%        |
| 14        | 7.470       | 503           | 506         | 507          | rVV      | 91685          | 132701        | 1.66%           | 0.165%        |
| 15        | 7.584       | 513           | 516         | 517          | rVV      | 243512         | 416831        | 5.22%           | 0.518%        |
| 16        | 7.619       | 517           | 519         | 521          | rVV2     | 211684         | 417892        | 5.24%           | 0.520%        |
| 17        | 7.847       | 535           | 539         | 542          | rBV      | 4517221        | 7978514       | 100.00%         | 9.922%        |
| 18        | 7.916       | 542           | 545         | 548          | rVB3     | 218843         | 323216        | 4.05%           | 0.402%        |
| 19        | 8.133       | 560           | 564         | 565          | rBV3     | 87671          | 131976        | 1.65%           | 0.164%        |
| 20        | 8.270       | 574           | 576         | 580          | rBV2     | 223561         | 406862        | 5.10%           | 0.506%        |
| 21        | 8.339       | 580           | 582         | 584          | rVB3     | 85959          | 115555        | 1.45%           | 0.144%        |
| 22        | 8.396       | 584           | 587         | 589          | rBV3     | 44954          | 117711        | 1.48%           | 0.146%        |
| 23        | 8.499       | 592           | 596         | 597          | rVB3     | 76332          | 157639        | 1.98%           | 0.196%        |
| 24        | 8.545       | 597           | 600         | 602          | rVB      | 2146303        | 2276511       | 28.53%          | 2.831%        |
| 25        | 8.636       | 606           | 608         | 611          | rVB      | 1769734        | 1624405       | 20.36%          | 2.020%        |
| 26        | 8.739       | 613           | 617         | 619          | rBV4     | 102536         | 227337        | 2.85%           | 0.283%        |
| 27        | 8.876       | 625           | 629         | 630          | rBV      | 181163         | 338374        | 4.24%           | 0.421%        |
| 28        | 8.910       | 630           | 632         | 634          | rVV      | 4453526        | 4325832       | 54.22%          | 5.380%        |
| 29        | 9.002       | 637           | 640         | 643          | rBV2     | 439062         | 550799        | 6.90%           | 0.685%        |
| 30        | 9.093       | 646           | 648         | 652          | rVB      | 613477         | 852686        | 10.69%          | 1.060%        |
| 31        | 9.162       | 652           | 654         | 656          | rBV      | 430288         | 628747        | 7.88%           | 0.782%        |
| 32        | 9.242       | 658           | 661         | 662          | rVV      | 735920         | 877228        | 10.99%          | 1.091%        |
| 33        | 9.310       | 665           | 667         | 669          | rVV2     | 615571         | 753848        | 9.45%           | 0.938%        |
| 34        | 9.356       | 669           | 671         | 675          | rVV2     | 356315         | 513202        | 6.43%           | 0.638%        |

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120215\  
 Data File : BF083420.D  
 Acq On : 2 Dec 2015 16:12  
 Operator : UM/IZ  
 Sample : G4583-06  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-P3WC-06(0-5)

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0

Filtering: 5  
 Min Area: 3 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
 Peak separation: 5

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113015.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |     |     |     |      |         |         |        |        |
|----|--------|-----|-----|-----|------|---------|---------|--------|--------|
| 35 | 9.436  | 675 | 678 | 680 | rVV2 | 341226  | 343508  | 4.31%  | 0.427% |
| 36 | 9.573  | 688 | 690 | 691 | rBV2 | 1555626 | 1538642 | 19.28% | 1.914% |
| 37 | 9.608  | 691 | 693 | 695 | rVB  | 1902601 | 1827620 | 22.91% | 2.273% |
| 38 | 9.676  | 698 | 699 | 702 | rVB  | 132683  | 186953  | 2.34%  | 0.233% |
| 39 | 9.779  | 704 | 708 | 710 | rBV2 | 288118  | 441229  | 5.53%  | 0.549% |
| 40 | 9.825  | 710 | 712 | 713 | rBV  | 137663  | 156150  | 1.96%  | 0.194% |
| 41 | 9.848  | 713 | 714 | 716 | rVB  | 198810  | 202939  | 2.54%  | 0.252% |
| 42 | 9.893  | 716 | 718 | 719 | rBV  | 266105  | 331096  | 4.15%  | 0.412% |
| 43 | 9.985  | 724 | 726 | 728 | rBV3 | 147851  | 275523  | 3.45%  | 0.343% |
| 44 | 10.030 | 728 | 730 | 731 | rVB2 | 228850  | 295489  | 3.70%  | 0.367% |
| 45 | 10.088 | 731 | 735 | 736 | rBV  | 258774  | 348799  | 4.37%  | 0.434% |
| 46 | 10.122 | 736 | 738 | 739 | rBV2 | 527884  | 707571  | 8.87%  | 0.880% |
| 47 | 10.179 | 741 | 743 | 746 | rBV3 | 77788   | 165623  | 2.08%  | 0.206% |
| 48 | 10.248 | 746 | 749 | 750 | rBV2 | 323429  | 475878  | 5.96%  | 0.592% |
| 49 | 10.362 | 755 | 759 | 761 | rBV2 | 2288188 | 2558903 | 32.07% | 3.182% |
| 50 | 10.522 | 769 | 773 | 775 | rBV2 | 325653  | 515449  | 6.46%  | 0.641% |
| 51 | 10.682 | 785 | 787 | 788 | rBV  | 149685  | 205464  | 2.58%  | 0.256% |
| 52 | 10.716 | 788 | 790 | 792 | rVV  | 179871  | 218850  | 2.74%  | 0.272% |
| 53 | 10.773 | 792 | 795 | 797 | rVB4 | 98384   | 210647  | 2.64%  | 0.262% |
| 54 | 10.991 | 812 | 814 | 816 | rVV2 | 188969  | 199483  | 2.50%  | 0.248% |
| 55 | 11.036 | 816 | 818 | 820 | rVB  | 185705  | 233087  | 2.92%  | 0.290% |
| 56 | 11.116 | 822 | 825 | 826 | rBV  | 943616  | 1740067 | 21.81% | 2.164% |
| 57 | 11.139 | 826 | 827 | 830 | rVV  | 2821469 | 2337917 | 29.30% | 2.908% |
| 58 | 11.196 | 830 | 832 | 835 | rVV  | 739591  | 744338  | 9.33%  | 0.926% |
| 59 | 11.471 | 853 | 856 | 857 | rVB2 | 72945   | 110358  | 1.38%  | 0.137% |
| 60 | 11.505 | 857 | 859 | 863 | rVB2 | 119722  | 160579  | 2.01%  | 0.200% |
| 61 | 11.608 | 866 | 868 | 872 | rVB3 | 80356   | 134648  | 1.69%  | 0.167% |
| 62 | 11.722 | 875 | 878 | 879 | rBV  | 369476  | 566699  | 7.10%  | 0.705% |
| 63 | 11.745 | 879 | 880 | 883 | rVB  | 392393  | 568747  | 7.13%  | 0.707% |
| 64 | 11.813 | 883 | 886 | 887 | rBV  | 285642  | 363164  | 4.55%  | 0.452% |
| 65 | 11.848 | 887 | 889 | 891 | rVV  | 677044  | 916009  | 11.48% | 1.139% |
| 66 | 11.882 | 891 | 892 | 895 | rVB  | 345116  | 317612  | 3.98%  | 0.395% |
| 67 | 12.099 | 908 | 911 | 918 | rVB  | 253806  | 531533  | 6.66%  | 0.661% |
| 68 | 12.431 | 937 | 940 | 942 | rBV  | 297274  | 483776  | 6.06%  | 0.602% |
| 69 | 12.476 | 942 | 944 | 946 | rVV2 | 165649  | 277851  | 3.48%  | 0.346% |
| 70 | 12.556 | 949 | 951 | 954 | rVB2 | 99120   | 167169  | 2.10%  | 0.208% |
| 71 | 12.636 | 956 | 958 | 961 | rBV  | 824045  | 1034454 | 12.97% | 1.286% |

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120215\  
 Data File : BF083420.D  
 Acq On : 2 Dec 2015 16:12  
 Operator : UM/IZ  
 Sample : G4583-06  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-P3WC-06(0-5)

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0  
 Filtering: 5  
 Min Area: 3 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
 Peak separation: 5

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113015.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|     |        |      |      |      |      |         |         |        |        |
|-----|--------|------|------|------|------|---------|---------|--------|--------|
| 72  | 12.774 | 968  | 970  | 973  | rVB  | 305190  | 419374  | 5.26%  | 0.522% |
| 73  | 12.876 | 977  | 979  | 982  | rVV  | 95185   | 183562  | 2.30%  | 0.228% |
| 74  | 12.945 | 982  | 985  | 988  | rVV  | 1073727 | 1587676 | 19.90% | 1.974% |
| 75  | 13.014 | 988  | 991  | 994  | rVV5 | 50438   | 134341  | 1.68%  | 0.167% |
| 76  | 13.196 | 1004 | 1007 | 1010 | rVV  | 2605755 | 4184740 | 52.45% | 5.204% |
| 77  | 13.276 | 1011 | 1014 | 1020 | rBV3 | 106847  | 242875  | 3.04%  | 0.302% |
| 78  | 13.402 | 1022 | 1025 | 1026 | rVV3 | 96352   | 186555  | 2.34%  | 0.232% |
| 79  | 13.436 | 1026 | 1028 | 1031 | rVB  | 201707  | 283505  | 3.55%  | 0.353% |
| 80  | 13.528 | 1034 | 1036 | 1038 | rVV  | 148777  | 183745  | 2.30%  | 0.229% |
| 81  | 13.585 | 1038 | 1041 | 1043 | rVB  | 121326  | 179553  | 2.25%  | 0.223% |
| 82  | 13.711 | 1050 | 1052 | 1054 | rVV  | 109121  | 132679  | 1.66%  | 0.165% |
| 83  | 13.757 | 1054 | 1056 | 1060 | rVB  | 107741  | 179291  | 2.25%  | 0.223% |
| 84  | 14.145 | 1087 | 1090 | 1092 | rBV3 | 55562   | 138054  | 1.73%  | 0.172% |
| 85  | 14.202 | 1093 | 1095 | 1099 | rVV2 | 64836   | 111484  | 1.40%  | 0.139% |
| 86  | 14.339 | 1103 | 1107 | 1109 | rBV3 | 61954   | 131467  | 1.65%  | 0.163% |
| 87  | 14.385 | 1109 | 1111 | 1115 | rVV2 | 84341   | 191027  | 2.39%  | 0.238% |
| 88  | 14.499 | 1119 | 1121 | 1126 | rVB2 | 66975   | 143769  | 1.80%  | 0.179% |
| 89  | 14.637 | 1130 | 1133 | 1137 | rVV2 | 178029  | 428088  | 5.37%  | 0.532% |
| 90  | 14.728 | 1137 | 1141 | 1143 | rVV2 | 1044206 | 1915454 | 24.01% | 2.382% |
| 91  | 14.762 | 1143 | 1144 | 1147 | rVV  | 294655  | 362673  | 4.55%  | 0.451% |
| 92  | 14.842 | 1149 | 1151 | 1158 | rVV3 | 67510   | 187309  | 2.35%  | 0.233% |
| 93  | 15.334 | 1191 | 1194 | 1198 | rBV2 | 68794   | 149985  | 1.88%  | 0.187% |
| 94  | 16.282 | 1274 | 1277 | 1279 | rBV  | 149913  | 312546  | 3.92%  | 0.389% |
| 95  | 16.648 | 1307 | 1309 | 1312 | rVB  | 114901  | 156367  | 1.96%  | 0.194% |
| 96  | 16.717 | 1312 | 1315 | 1318 | rBV  | 191242  | 276109  | 3.46%  | 0.343% |
| 97  | 16.785 | 1318 | 1321 | 1334 | rVB  | 544671  | 1000367 | 12.54% | 1.244% |
| 98  | 17.025 | 1338 | 1342 | 1346 | rBV3 | 35463   | 111387  | 1.40%  | 0.139% |
| 99  | 18.168 | 1439 | 1442 | 1447 | rVB2 | 61668   | 134672  | 1.69%  | 0.167% |
| 100 | 18.534 | 1471 | 1474 | 1480 | rVB  | 67775   | 155882  | 1.95%  | 0.194% |

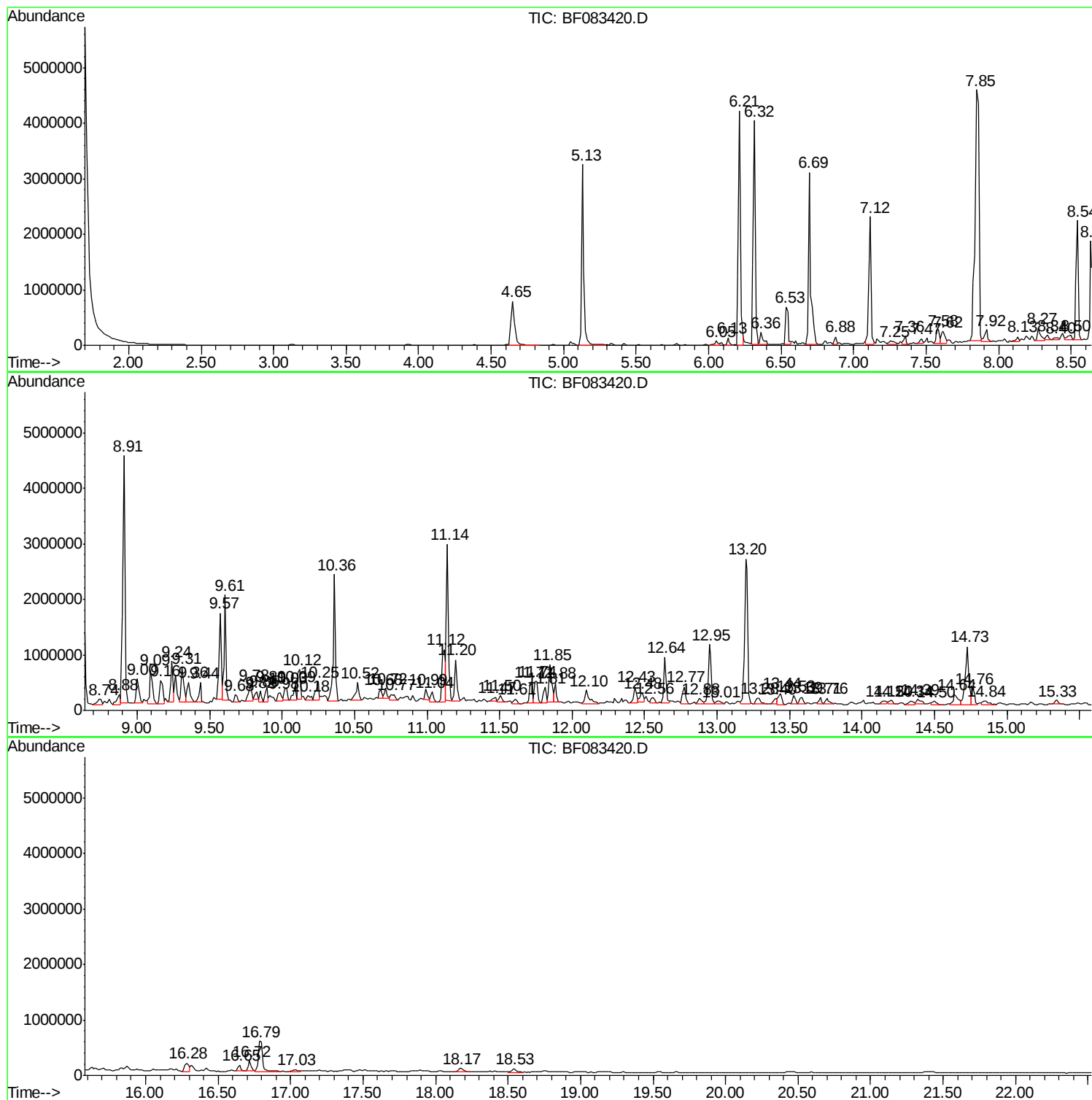
Sum of corrected areas: 80409796

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120215\  
Data File : BF083420.D  
Acq On : 2 Dec 2015 16:12  
Operator : UM/IZ  
Sample : G4583-06  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-P3WC-06(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120215\  
Data File : BF083420.D  
Acq On : 2 Dec 2015 16:12  
Operator : UM/IZ  
Sample : G4583-06  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-P3WC-06(0-5)

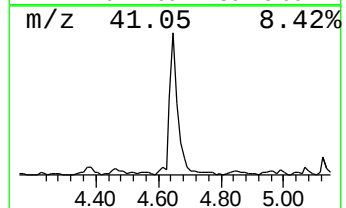
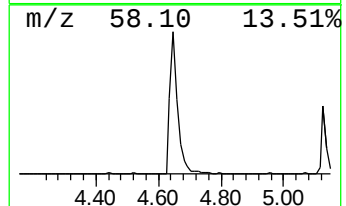
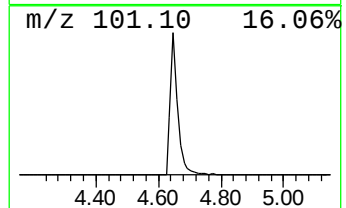
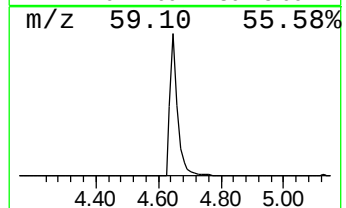
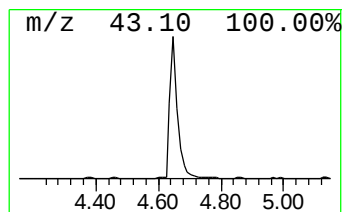
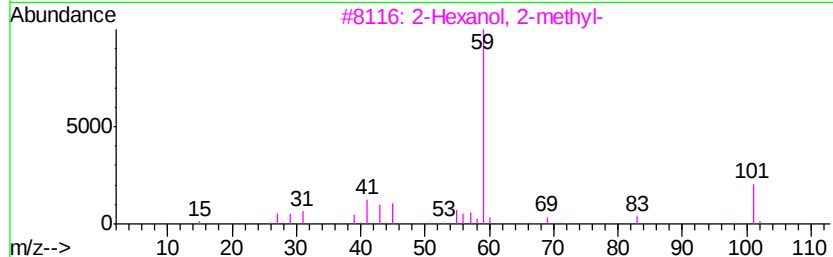
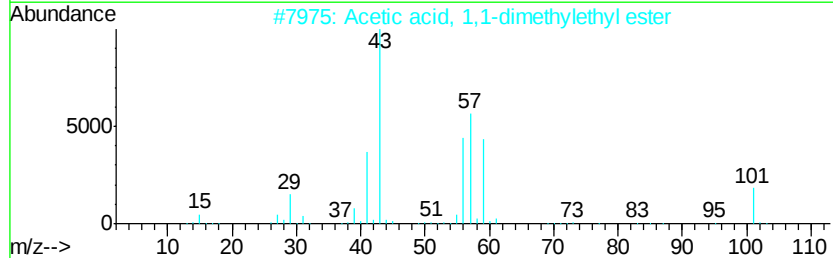
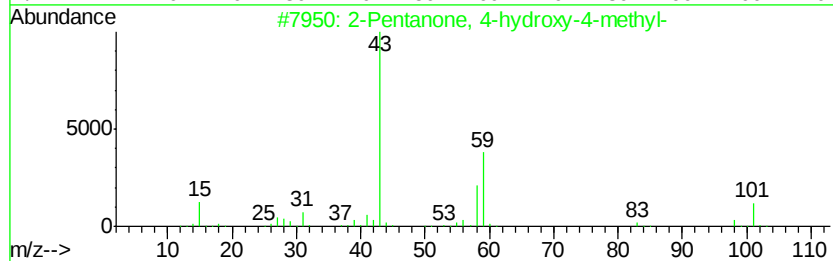
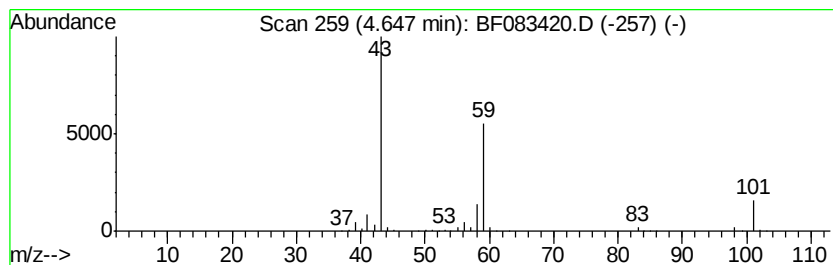
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 4.65 | 28.09 ng | 1312690 | 1,4-Dichlorobenzene-d4 | 6.54 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 50   |
| 2    |    |   | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2  | 000540-88-5 | 39   |
| 3    |    |   | 2-Hexanol, 2-methyl-                | 116 | C7H16O   | 000625-23-0 | 33   |
| 4    |    |   | 3-Hexanol, 4-methyl-                | 116 | C7H16O   | 000615-29-2 | 33   |
| 5    |    |   | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 23   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120215\  
Data File : BF083420.D  
Acq On : 2 Dec 2015 16:12  
Operator : UM/IZ  
Sample : G4583-06  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-P3WC-06(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

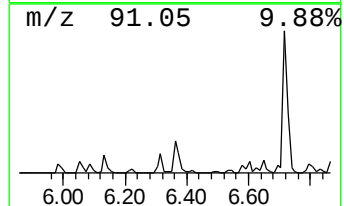
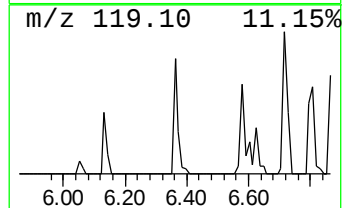
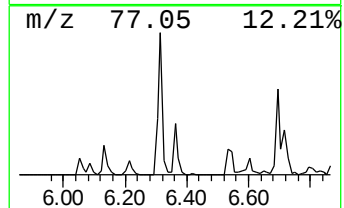
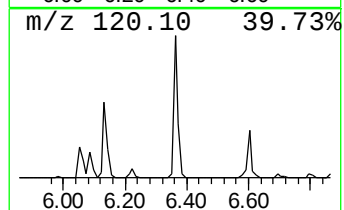
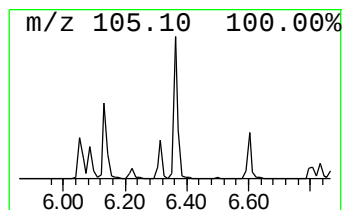
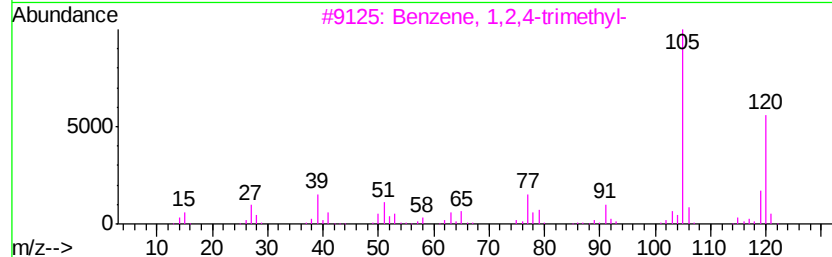
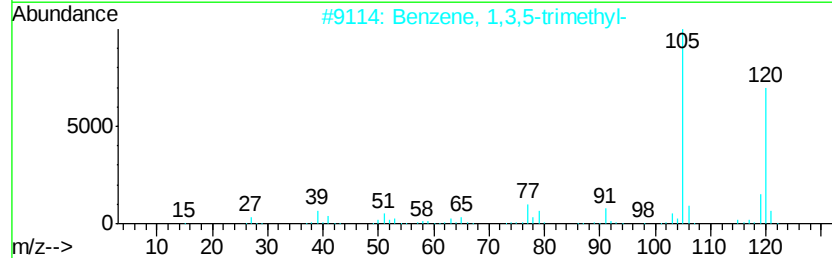
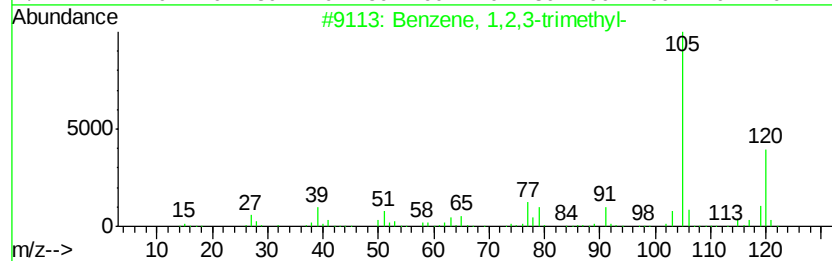
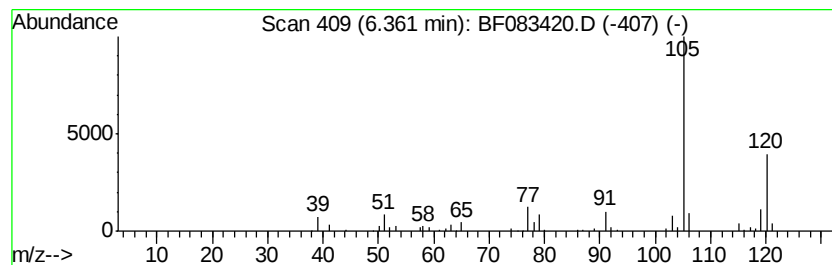
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Benzene, 1,2,3-trimethyl- Concentration Rank 8

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.36 | 6.56 ng | 306423 | 1,4-Dichlorobenzene-d4 | 6.54 |

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



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120215\  
Data File : BF083420.D  
Acq On : 2 Dec 2015 16:12  
Operator : UM/IZ  
Sample : G4583-06  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-P3WC-06(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

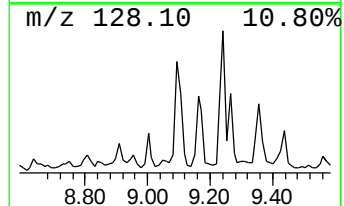
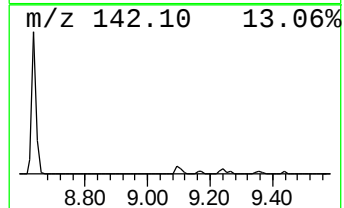
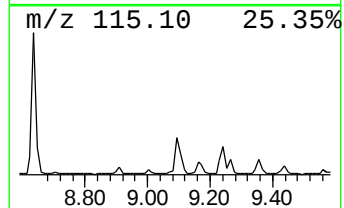
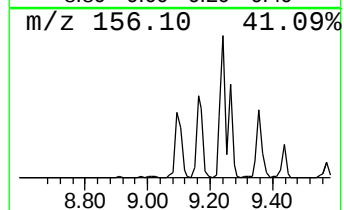
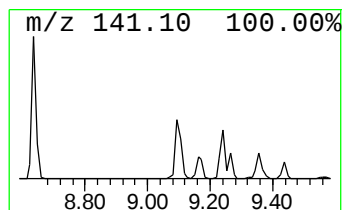
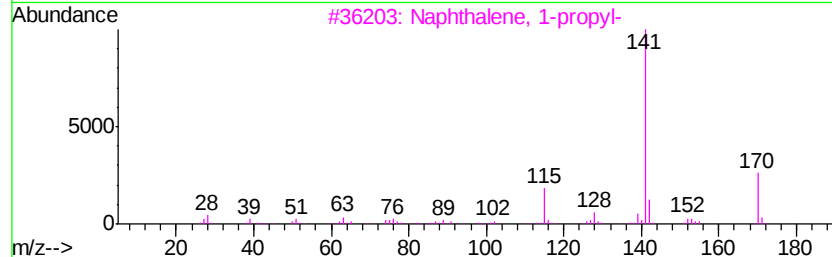
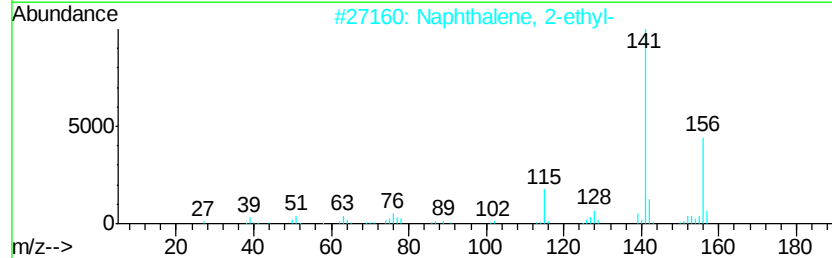
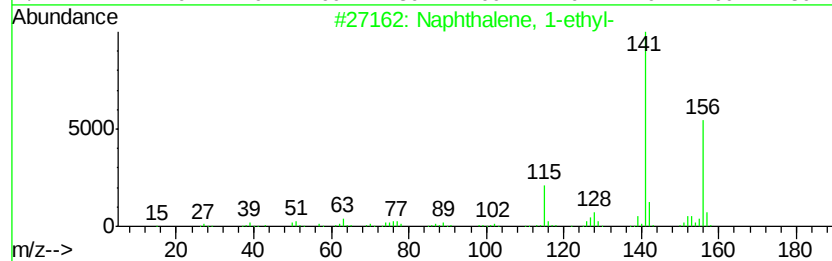
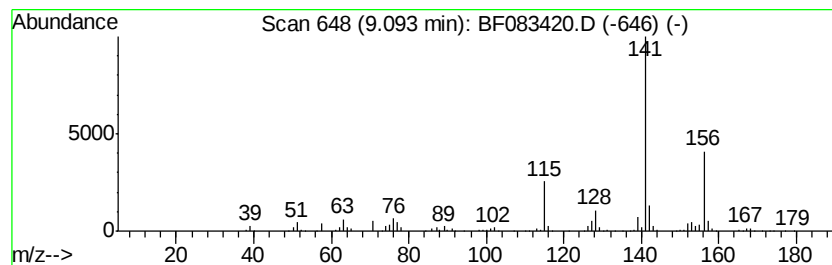
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Naphthalene, 1-ethyl- Concentration Rank 5

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 9.09 | 11.08 ng | 852686 | Acenaphthene-d10 | 9.57 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1-ethyl-        | 156 | C12H12  | 001127-76-0 | 95   |
| 2    |    | Naphthalene, 2-ethyl-        | 156 | C12H12  | 000939-27-5 | 95   |
| 3    |    | Naphthalene, 1-propyl-       | 170 | C13H14  | 002765-18-6 | 59   |
| 4    |    | 3',4'-Difluoroacetophenone   | 156 | C8H6F2O | 000369-33-5 | 53   |
| 5    |    | Methyl phenylphosphinic acid | 156 | C7H9O2P | 004271-13-0 | 53   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120215\  
Data File : BF083420.D  
Acq On : 2 Dec 2015 16:12  
Operator : UM/IZ  
Sample : G4583-06  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-P3WC-06(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

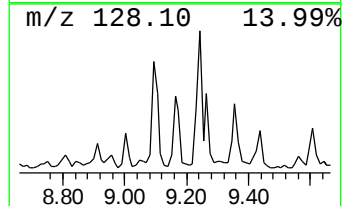
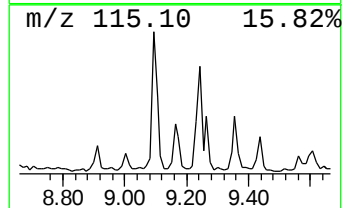
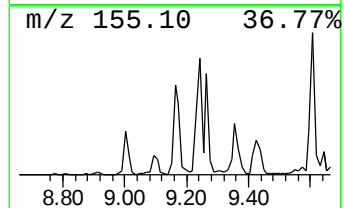
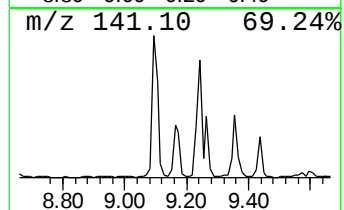
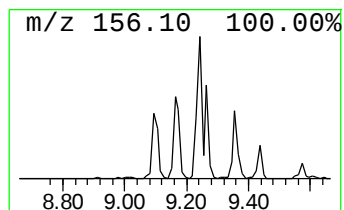
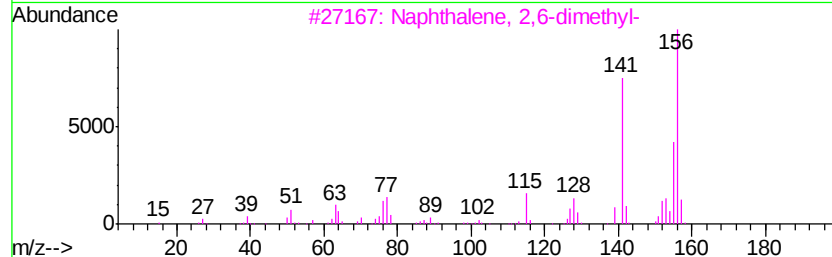
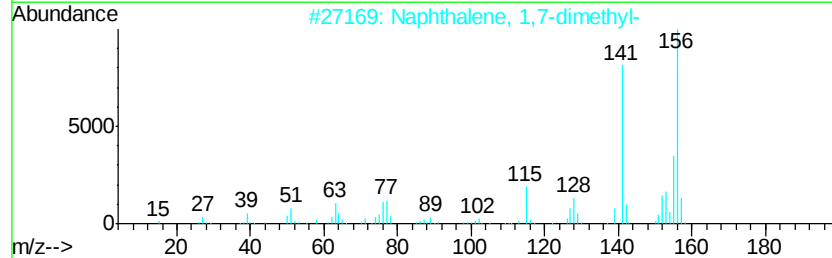
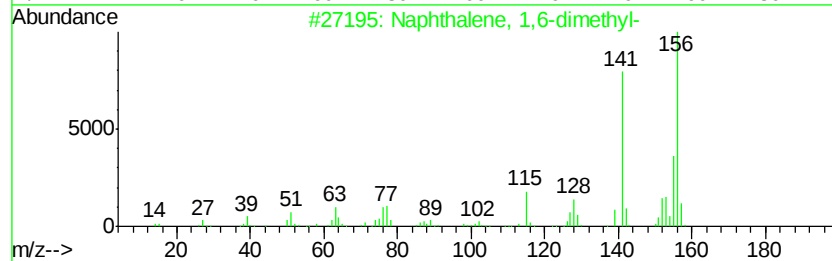
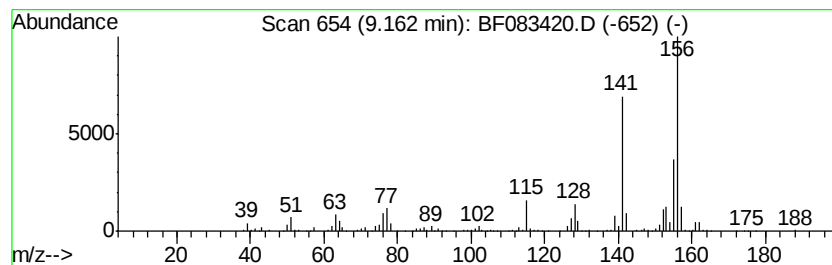
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Naphthalene, 1,6-dimethyl- Concentration Rank 7

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.16 | 8.17 ng | 628747 | Acenaphthene-d10 | 9.57 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 98   |
| 2    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 98   |
| 3    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 97   |
| 4    |    |   | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 97   |
| 5    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 96   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120215\  
Data File : BF083420.D  
Acq On : 2 Dec 2015 16:12  
Operator : UM/IZ  
Sample : G4583-06  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-P3WC-06(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

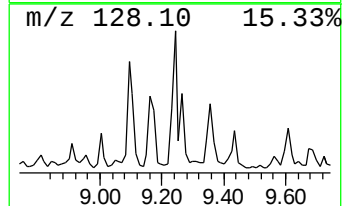
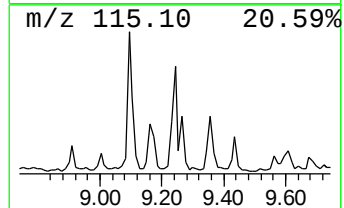
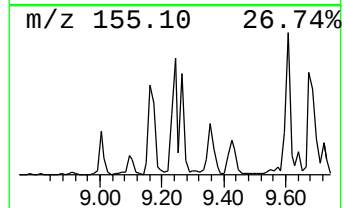
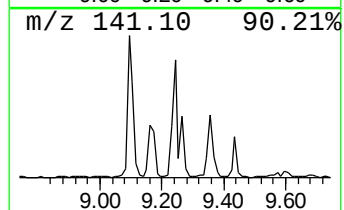
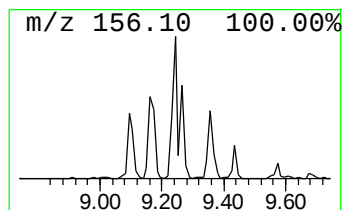
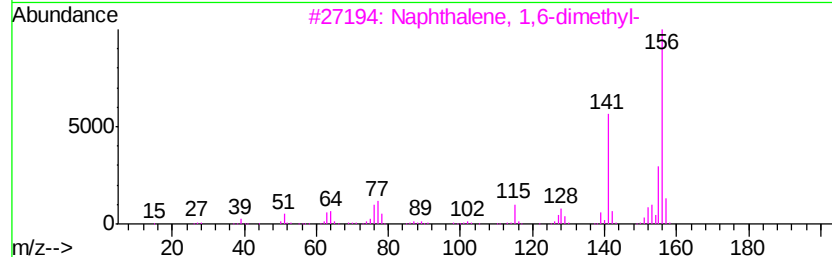
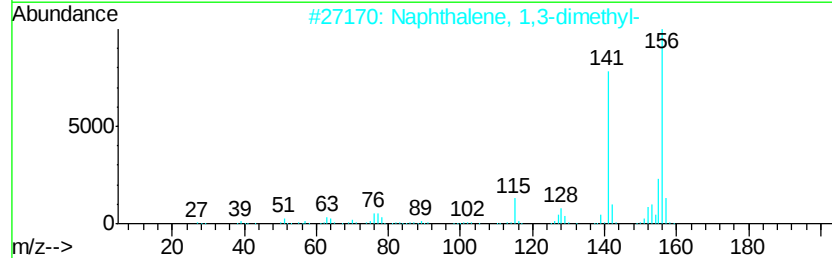
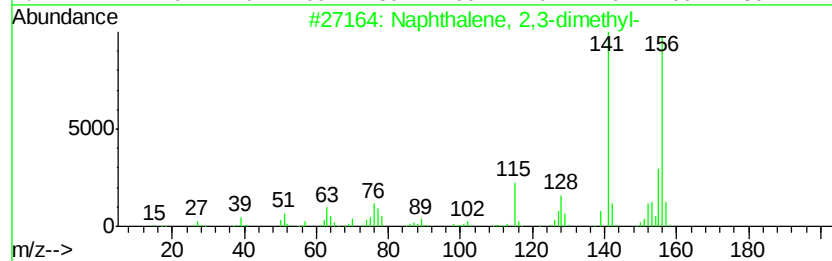
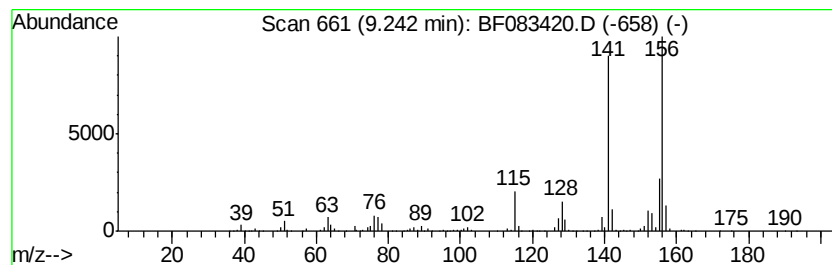
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 Naphthalene, 2,3-dimethyl- Concentration Rank 4

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 9.24 | 11.40 ng | 877228 | Acenaphthene-d10 | 9.57 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 95   |
| 2    |    | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 94   |
| 3    |    | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 94   |
| 4    |    | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 94   |
| 5    |    | Naphthalene, 1,8-dimethyl- | 156 | C12H12  | 000569-41-5 | 93   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120215\  
Data File : BF083420.D  
Acq On : 2 Dec 2015 16:12  
Operator : UM/IZ  
Sample : G4583-06  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-P3WC-06(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

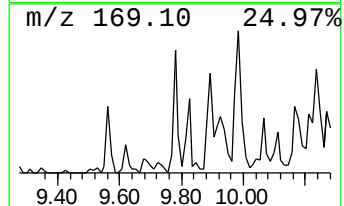
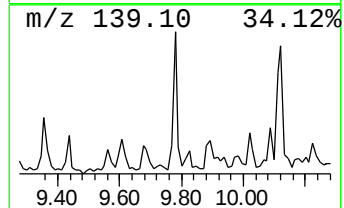
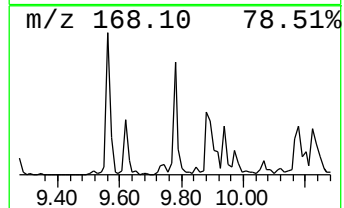
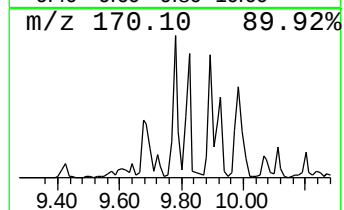
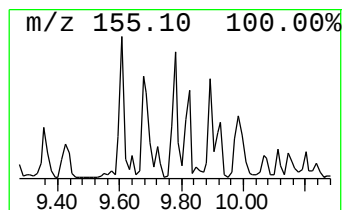
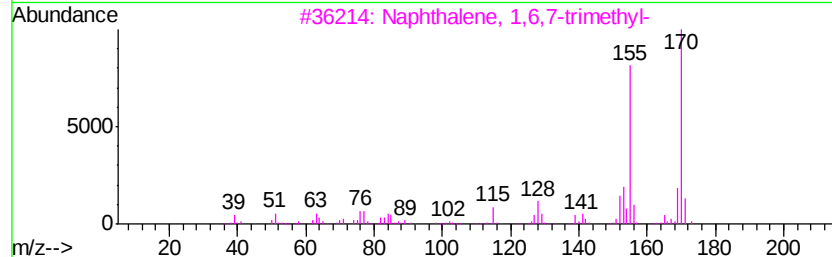
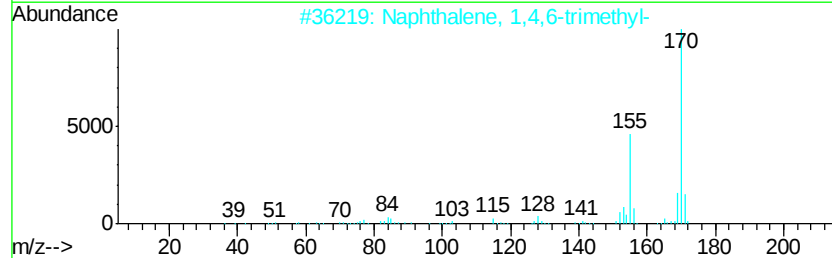
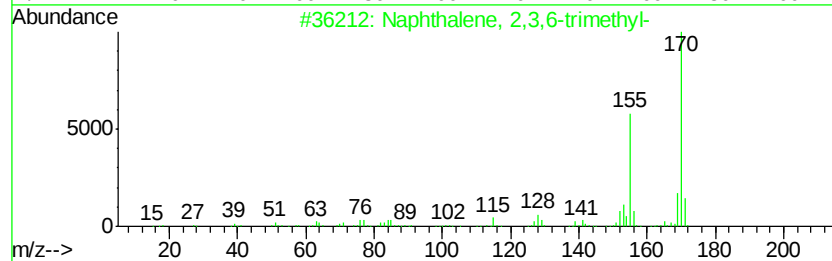
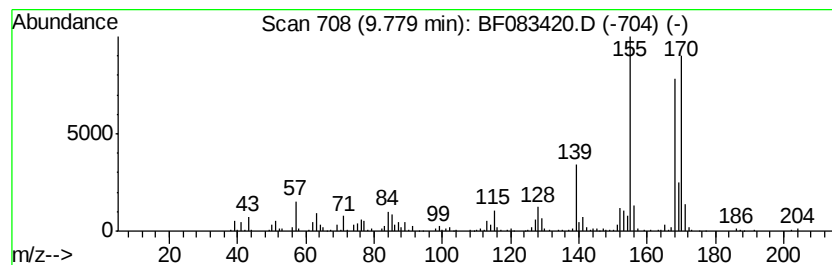
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Naphthalene, 2,3,6-trimethyl- Concentration Rank 14

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.78 | 5.74 ng | 441229 | Acenaphthene-d10 | 9.57 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Naphthalene, 2,3,6-trimethyl-       | 170 | C13H14  | 000829-26-5  | 83   |
| 2    |    |   | Naphthalene, 1,4,6-trimethyl-       | 170 | C13H14  | 002131-42-2  | 83   |
| 3    |    |   | Naphthalene, 1,6,7-trimethyl-       | 170 | C13H14  | 002245-38-7  | 78   |
| 4    |    |   | 1-(2,2-Dimethylcyclopropyl)-2-ph... | 170 | C13H14  | 1000294-91-1 | 58   |
| 5    |    |   | 3-(2-Methyl-propenyl)-1H-indene     | 170 | C13H14  | 1000187-78-5 | 53   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120215\  
Data File : BF083420.D  
Acq On : 2 Dec 2015 16:12  
Operator : UM/IZ  
Sample : G4583-06  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-P3WC-06(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

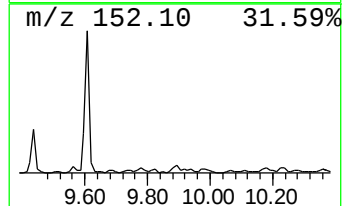
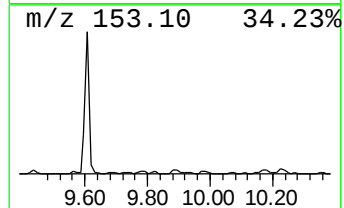
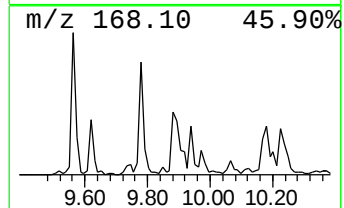
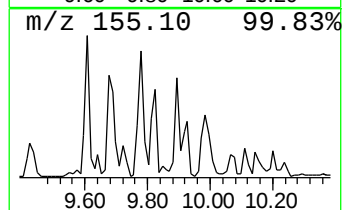
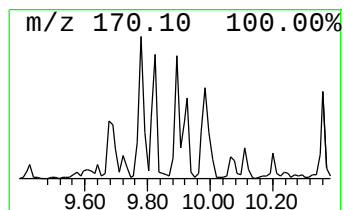
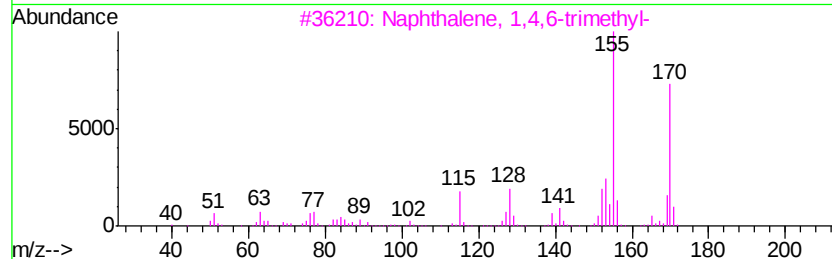
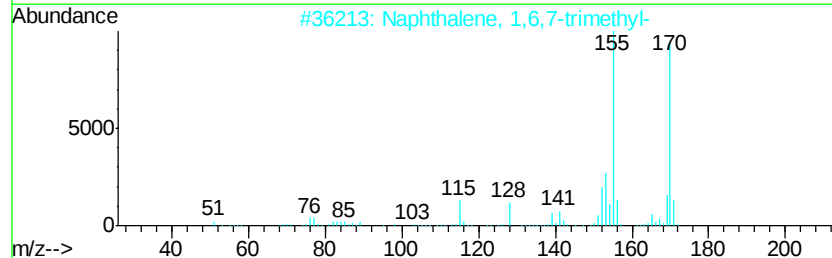
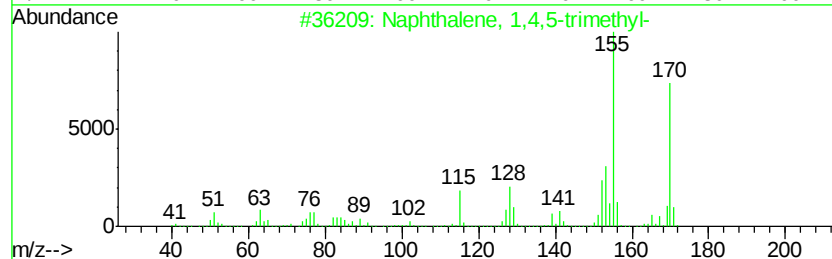
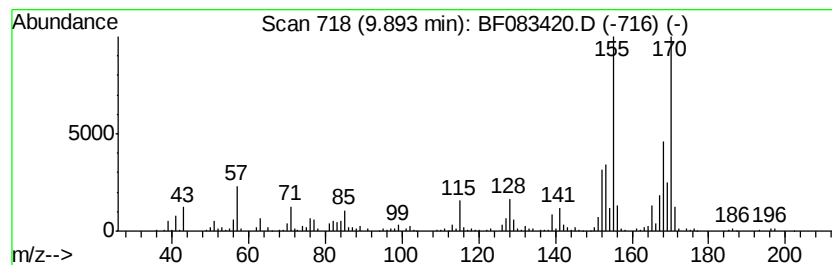
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Naphthalene, 1,4,5-trimethyl- Concentration Rank 18

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.89 | 4.30 ng | 331096 | Acenaphthene-d10 | 9.57 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,4,5-trimethyl- | 170 | C13H14  | 002131-41-1 | 94   |
| 2    |    | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 94   |
| 3    |    | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 94   |
| 4    |    | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 93   |
| 5    |    | Azulene, 4,6,8-trimethyl-     | 170 | C13H14  | 000941-81-1 | 89   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120215\  
Data File : BF083420.D  
Acq On : 2 Dec 2015 16:12  
Operator : UM/IZ  
Sample : G4583-06  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-P3WC-06(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

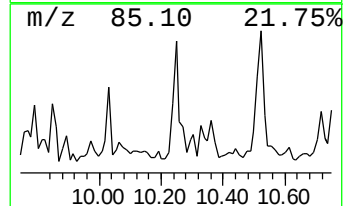
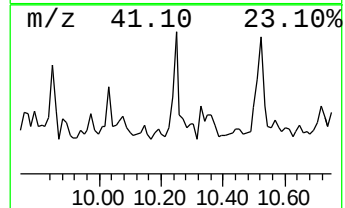
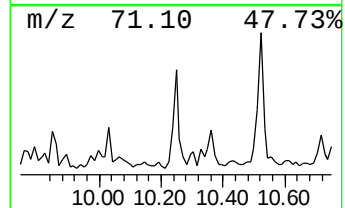
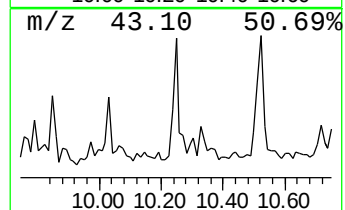
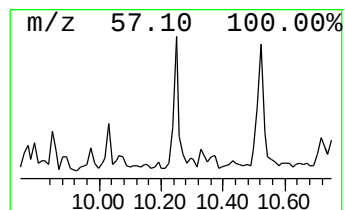
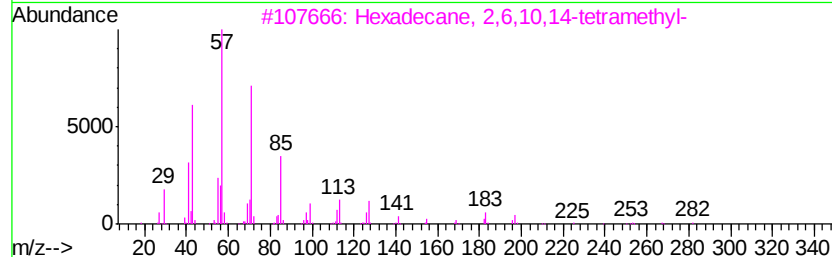
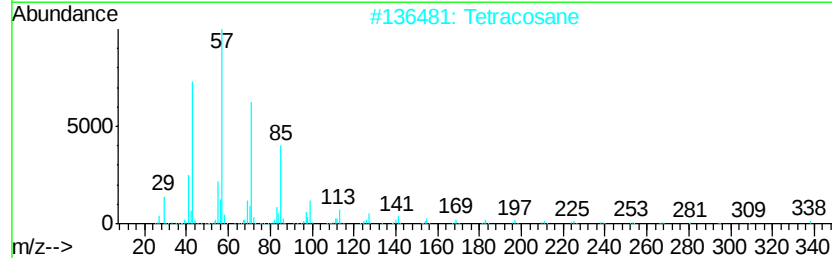
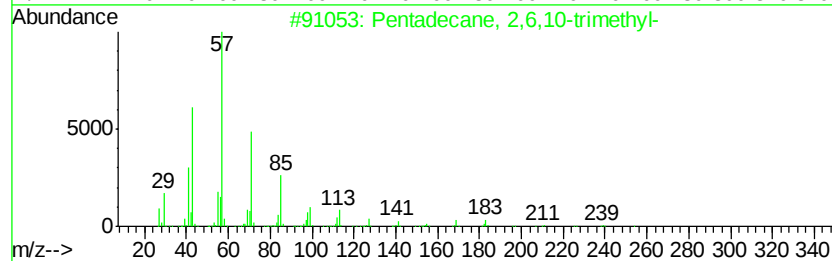
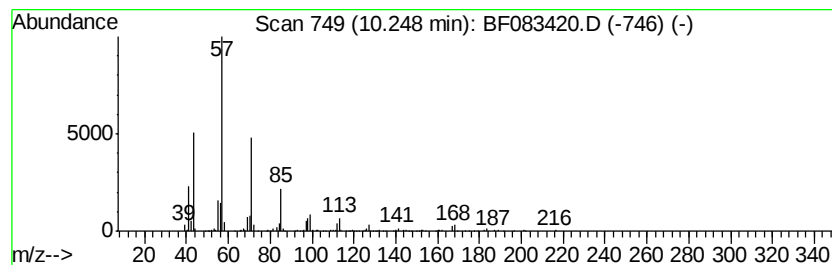
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Pentadecane, 2,6,10-trimethyl- Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T. |
|-------|---------|--------|------------------|------|
| 10.25 | 6.19 ng | 475878 | Acenaphthene-d10 | 9.57 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Pentadecane, 2,6,10-trimethyl-     | 254 | C18H38  | 003892-00-0 | 90   |
| 2    |    | Tetracosane                        | 338 | C24H50  | 000646-31-1 | 80   |
| 3    |    | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8 | 72   |
| 4    |    | Heptacosane                        | 380 | C27H56  | 000593-49-7 | 72   |
| 5    |    | Eicosane                           | 282 | C20H42  | 000112-95-8 | 64   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120215\  
Data File : BF083420.D  
Acq On : 2 Dec 2015 16:12  
Operator : UM/IZ  
Sample : G4583-06  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

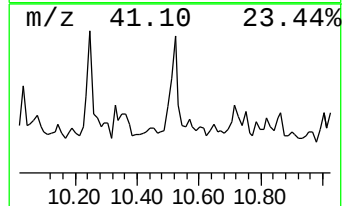
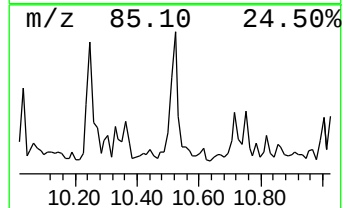
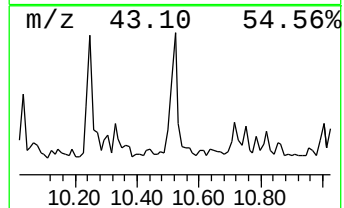
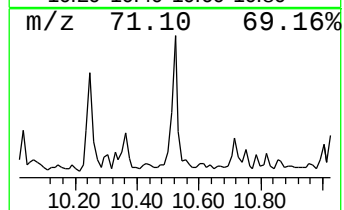
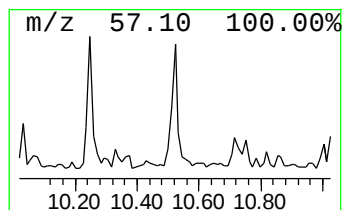
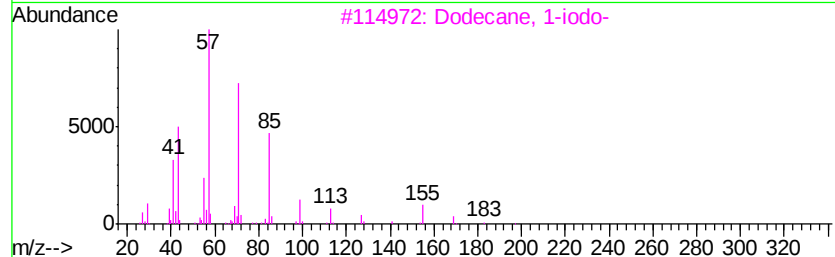
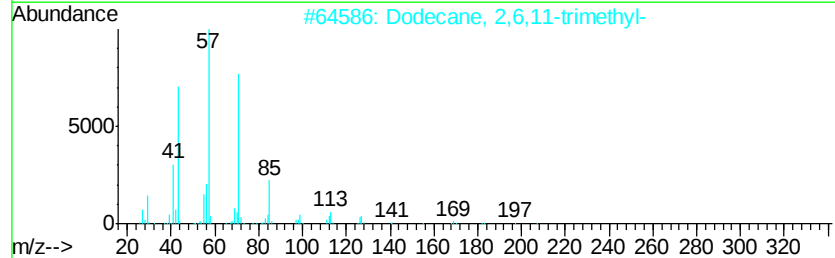
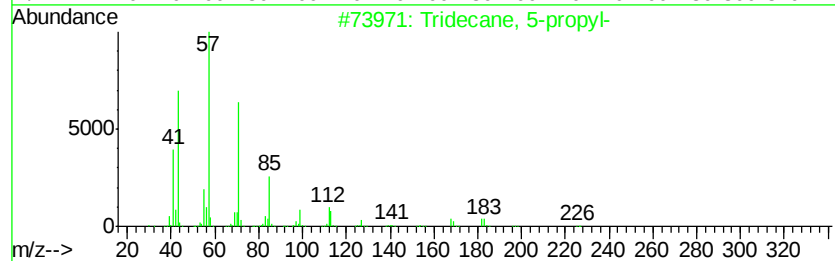
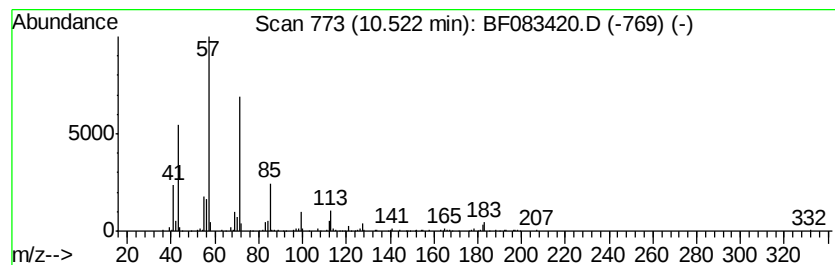
Instrument :  
BNA\_F  
ClientSampleId :  
SB-P3WC-06(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 12 Tridecane, 5-propyl- Concentration Rank 13

| R.T.    | EstConc | Area                        | Relative to ISTD |         | R.T.        |      |
|---------|---------|-----------------------------|------------------|---------|-------------|------|
| 10.52   | 5.92 ng | 515449                      | Phenanthrene-d10 |         | 11.12       |      |
| Hit# of | 5       | Tentative ID                | MW               | MolForm | CAS#        | Qual |
| 1       |         | Tridecane, 5-propyl-        | 226              | C16H34  | 055045-11-9 | 78   |
| 2       |         | Dodecane, 2,6,11-trimethyl- | 212              | C15H32  | 031295-56-4 | 72   |
| 3       |         | Dodecane, 1-iodo-           | 296              | C12H25I | 004292-19-7 | 58   |
| 4       |         | Tridecane, 7-methyl-        | 198              | C14H30  | 026730-14-3 | 53   |
| 5       |         | Dodecane, 4-methyl-         | 184              | C13H28  | 006117-97-1 | 50   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120215\  
Data File : BF083420.D  
Acq On : 2 Dec 2015 16:12  
Operator : UM/IZ  
Sample : G4583-06  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-P3WC-06(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

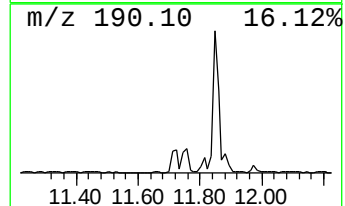
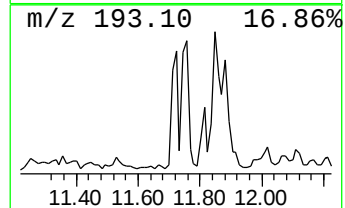
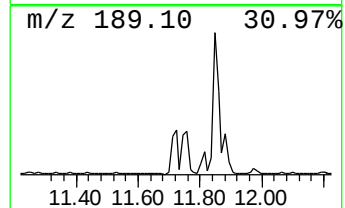
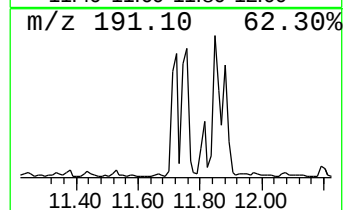
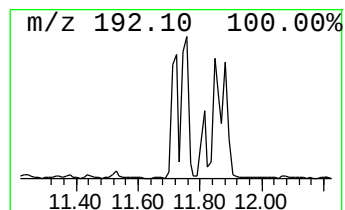
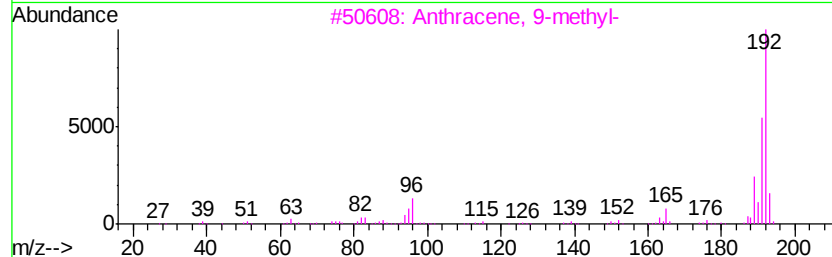
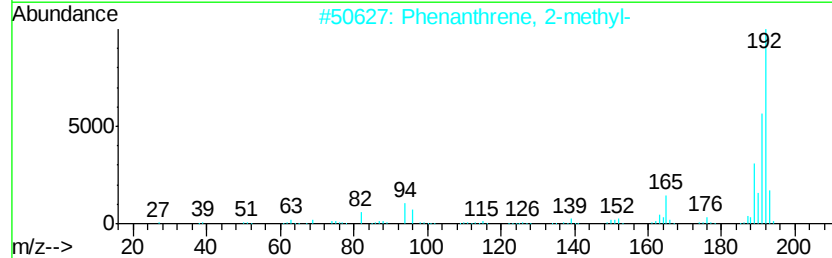
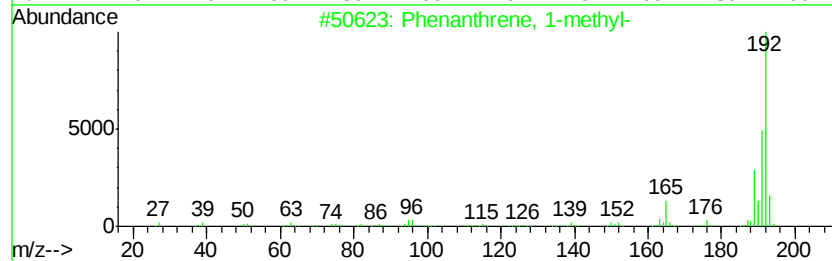
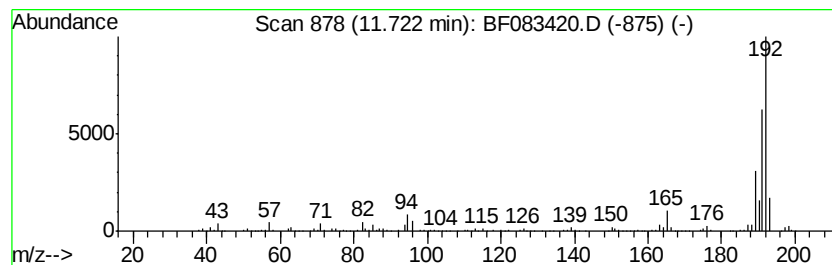
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Phenanthrene, 1-methyl- Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.72 | 6.51 ng | 566699 | Phenanthrene-d10 | 11.12 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 93   |
| 2    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 93   |
| 3    |    |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 91   |
| 4    |    |   | Phenanthrene, 3-methyl- | 192 | C15H12  | 000832-71-3 | 87   |
| 5    |    |   | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 87   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120215\  
Data File : BF083420.D  
Acq On : 2 Dec 2015 16:12  
Operator : UM/IZ  
Sample : G4583-06  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-P3WC-06(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

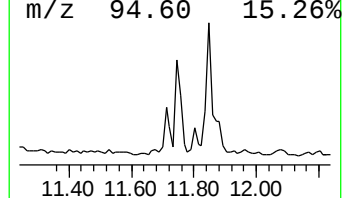
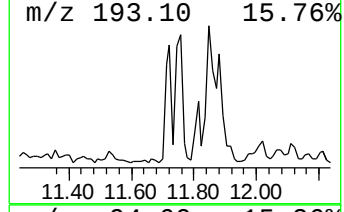
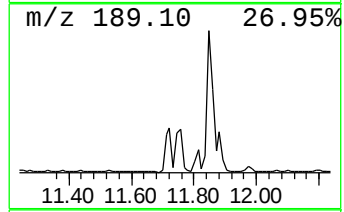
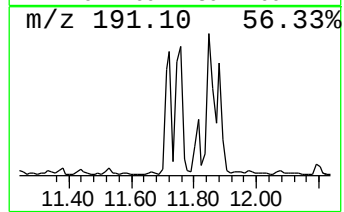
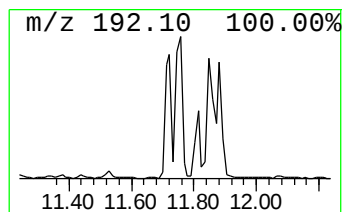
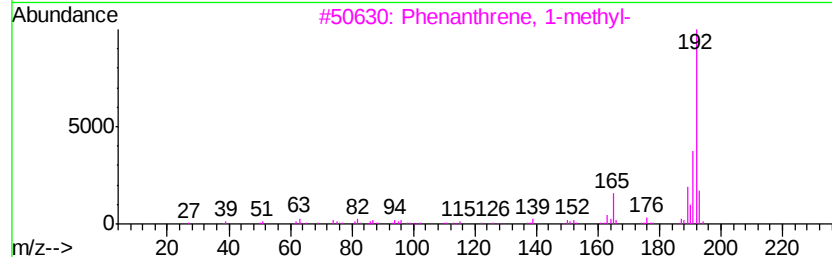
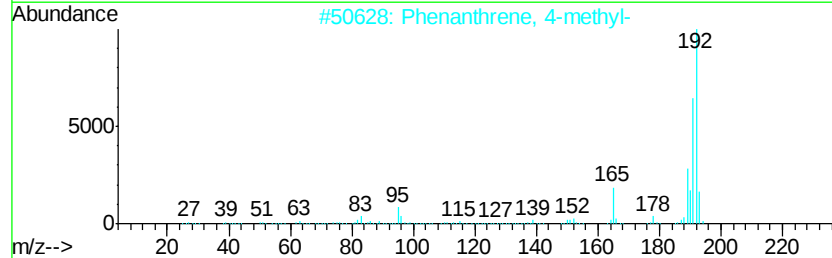
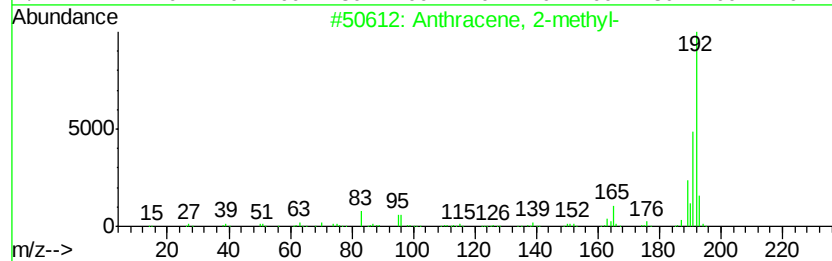
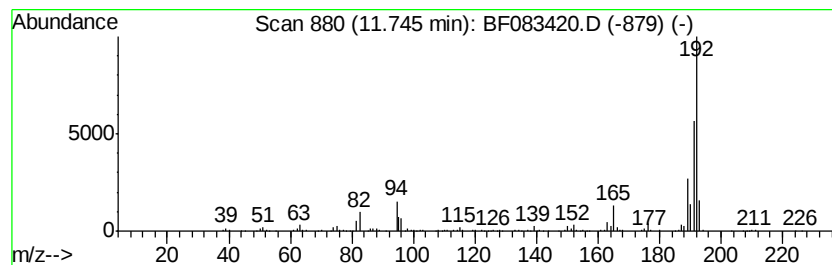
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Anthracene, 2-methyl- Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.74 | 6.54 ng | 568747 | Phenanthrene-d10 | 11.12 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 96   |
| 2    |    | Phenanthrene, 4-methyl- | 192 | C15H12  | 000832-64-4 | 95   |
| 3    |    | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 95   |
| 4    |    | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 95   |
| 5    |    | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 91   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120215\  
Data File : BF083420.D  
Acq On : 2 Dec 2015 16:12  
Operator : UM/IZ  
Sample : G4583-06  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-P3WC-06(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

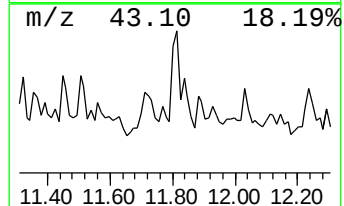
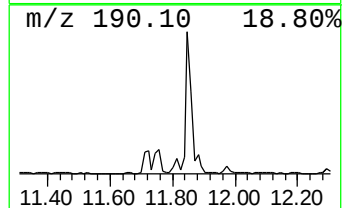
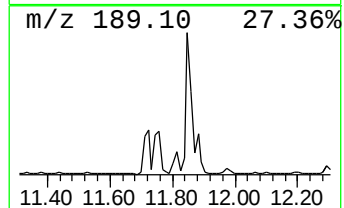
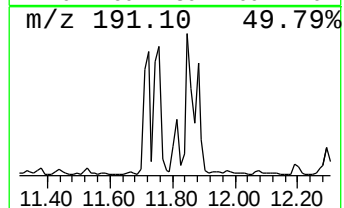
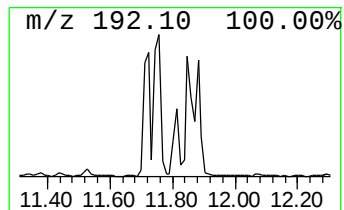
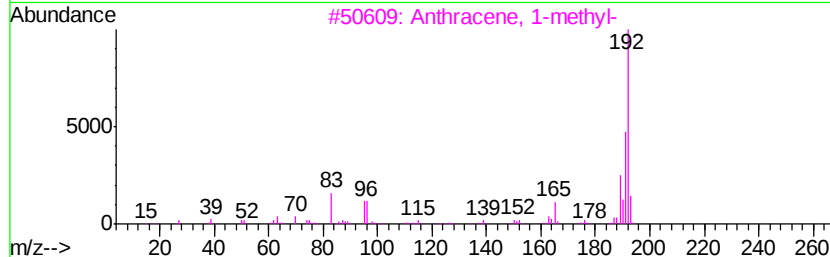
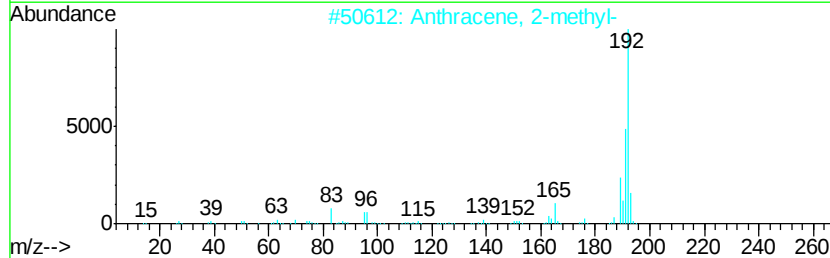
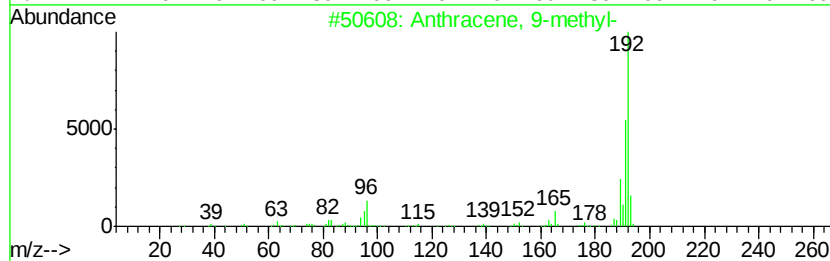
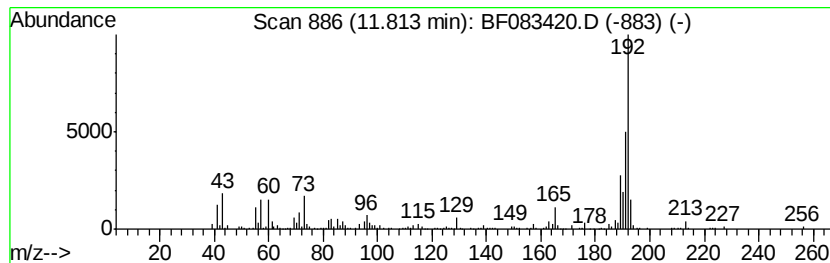
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Anthracene, 9-methyl- Concentration Rank 19

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.81 | 4.17 ng | 363164 | Phenanthrene-d10 | 11.12 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 89   |
| 2    |    |   | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 89   |
| 3    |    |   | Anthracene, 1-methyl-   | 192 | C15H12  | 000610-48-0 | 89   |
| 4    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 89   |
| 5    |    |   | Phenanthrene, 3-methyl- | 192 | C15H12  | 000832-71-3 | 89   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120215\  
Data File : BF083420.D  
Acq On : 2 Dec 2015 16:12  
Operator : UM/IZ  
Sample : G4583-06  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-P3WC-06(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

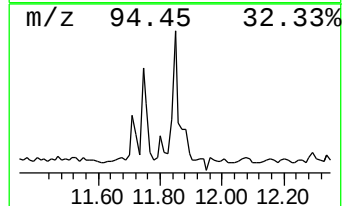
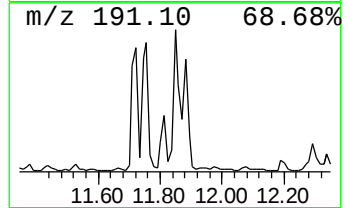
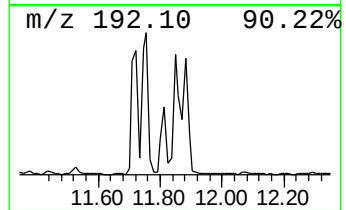
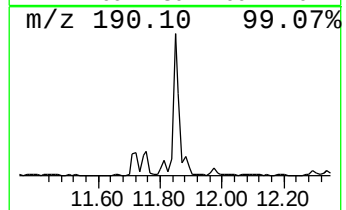
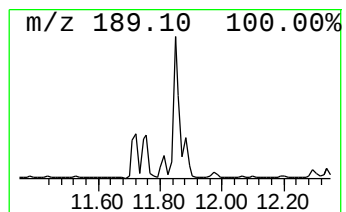
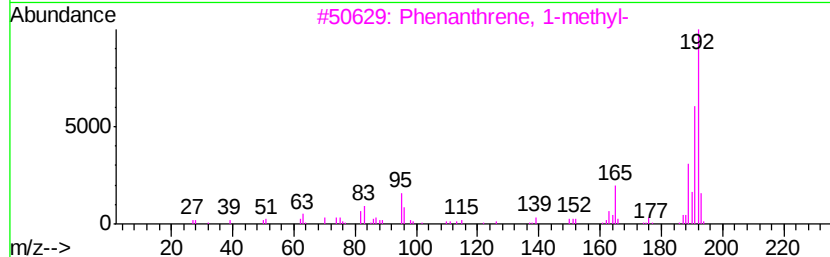
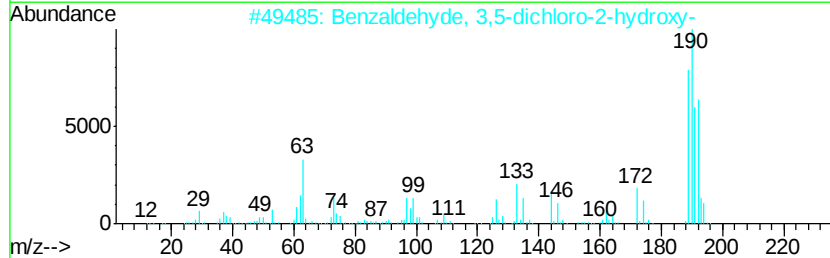
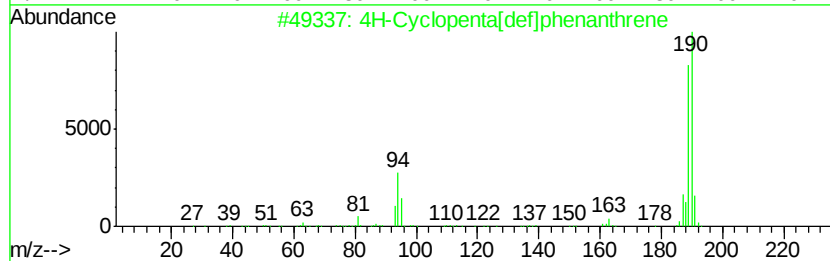
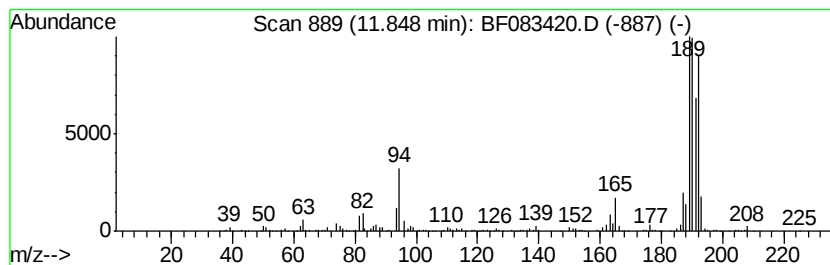
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 11.85 | 10.53 ng | 916009 | Phenanthrene-d10 | 11.12 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10    | 000203-64-5 | 90   |
| 2    |    | Benzaldehyde, 3,5-dichloro-2-hyd... | 190 | C7H4Cl2O2 | 000090-60-8 | 72   |
| 3    |    | Phenanthrene, 1-methyl-             | 192 | C15H12    | 000832-69-9 | 42   |
| 4    |    | Anthracene, 2-methyl-               | 192 | C15H12    | 000613-12-7 | 42   |
| 5    |    | 5H-Dibenzo[a,d]cycloheptene         | 192 | C15H12    | 000256-81-5 | 42   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120215\  
Data File : BF083420.D  
Acq On : 2 Dec 2015 16:12  
Operator : UM/IZ  
Sample : G4583-06  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-P3WC-06(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

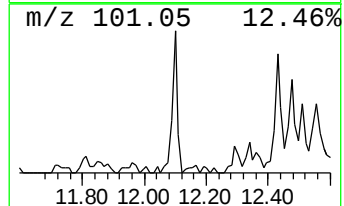
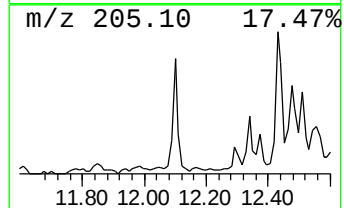
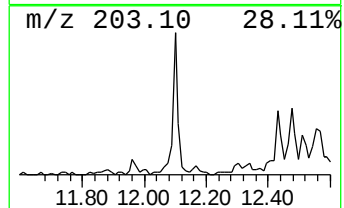
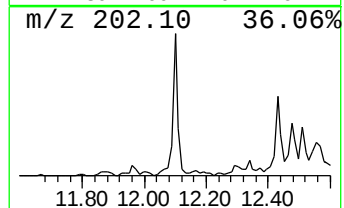
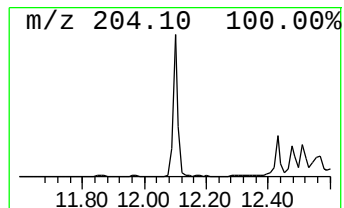
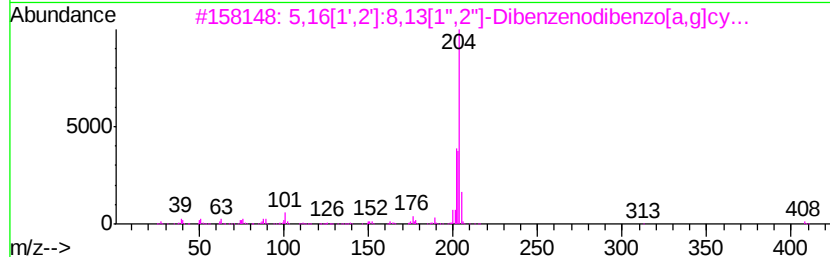
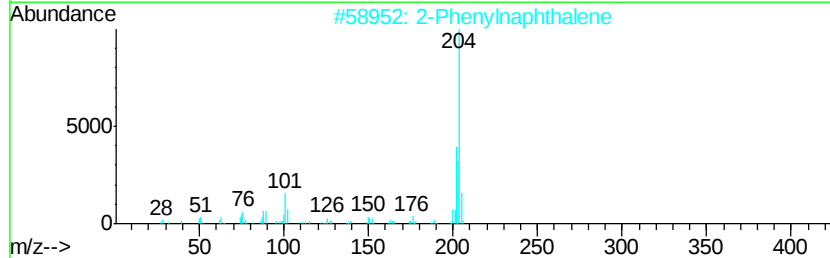
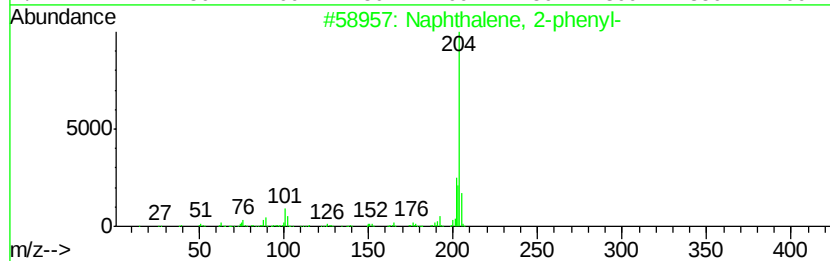
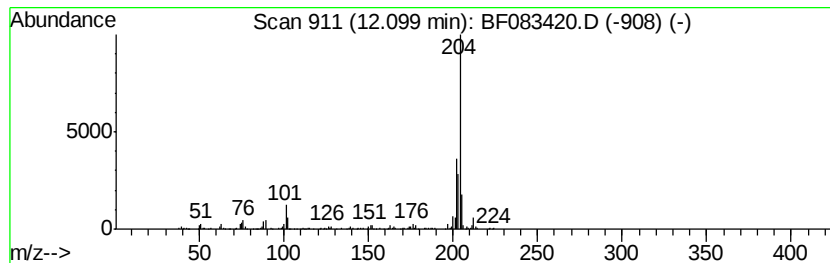
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Naphthalene, 2-phenyl- Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.10 | 6.11 ng | 531533 | Phenanthrene-d10 | 11.12 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2-phenyl-              | 204 | C16H12  | 000612-94-2 | 96   |
| 2    |    |   | 2-Phenylnaphthalene                 | 204 | C16H12  | 035465-71-5 | 94   |
| 3    |    |   | 5,16[1',2']8,13[1'',2'']-Dibenz...  | 408 | C32H24  | 005672-97-9 | 86   |
| 4    |    |   | Naphthalene, 1-phenyl-              | 204 | C16H12  | 000605-02-7 | 83   |
| 5    |    |   | Tricyclo[8.2.2.2(4,7)]hexadeca-2... | 204 | C16H12  | 006572-60-7 | 72   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120215\  
Data File : BF083420.D  
Acq On : 2 Dec 2015 16:12  
Operator : UM/IZ  
Sample : G4583-06  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

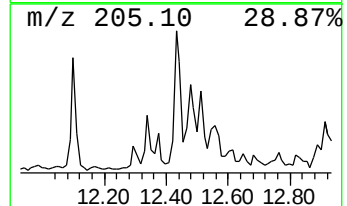
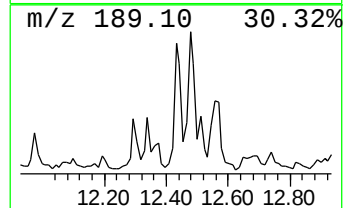
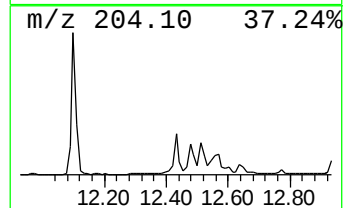
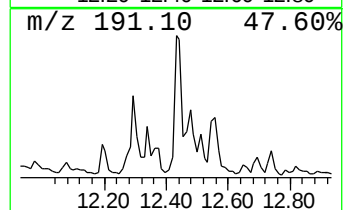
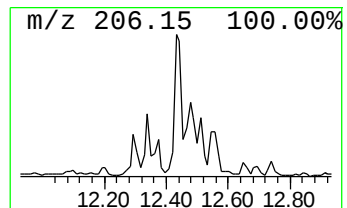
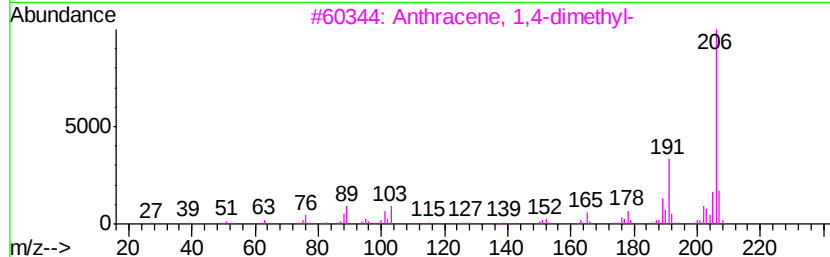
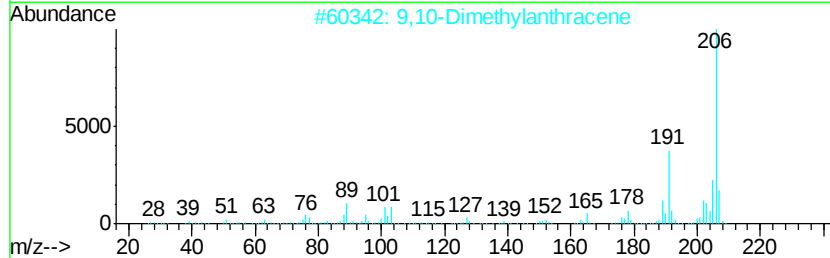
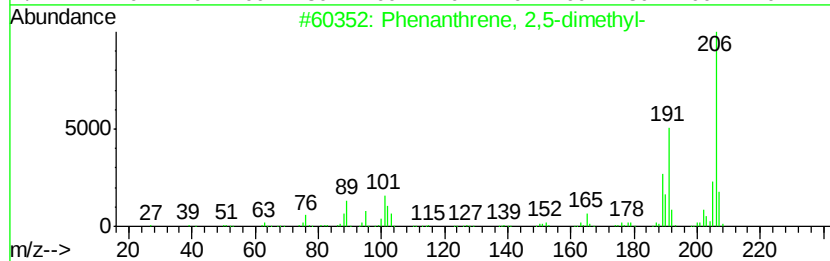
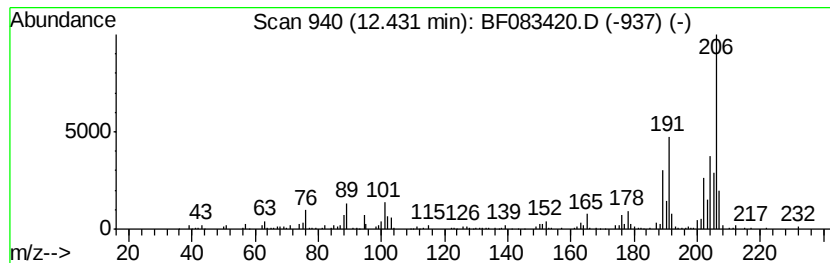
Instrument :  
BNA\_F  
ClientSampleId :  
SB-P3WC-06(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 18 Phenanthrene, 2,5-dimethyl- Concentration Rank 15

| R.T.      | EstConc                     | Area   | Relative to ISTD | R.T.        |      |
|-----------|-----------------------------|--------|------------------|-------------|------|
| 12.43     | 5.56 ng                     | 483776 | Phenanthrene-d10 | 11.12       |      |
| Hit# of 5 | Tentative ID                | MW     | MolForm          | CAS#        | Qual |
| 1         | Phenanthrene, 2,5-dimethyl- | 206    | C16H14           | 003674-66-6 | 94   |
| 2         | 9,10-Dimethylantracene      | 206    | C16H14           | 000781-43-1 | 93   |
| 3         | Anthracene, 1,4-dimethyl-   | 206    | C16H14           | 000781-92-0 | 93   |
| 4         | Phenanthrene, 1,7-dimethyl- | 206    | C16H14           | 000483-87-4 | 90   |
| 5         | Phenanthrene, 2,7-dimethyl- | 206    | C16H14           | 001576-69-8 | 90   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120215\  
Data File : BF083420.D  
Acq On : 2 Dec 2015 16:12  
Operator : UM/IZ  
Sample : G4583-06  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-P3WC-06(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

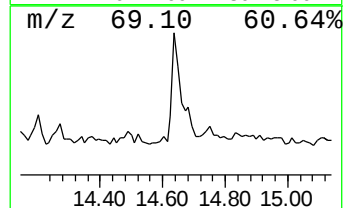
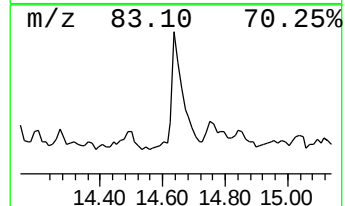
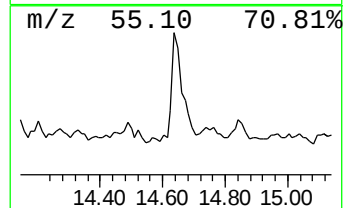
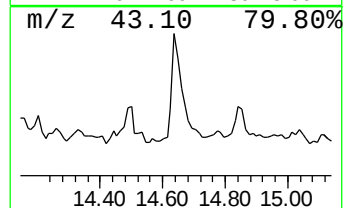
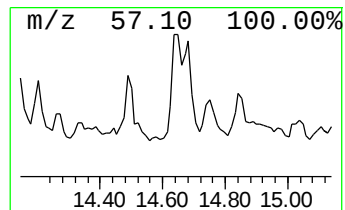
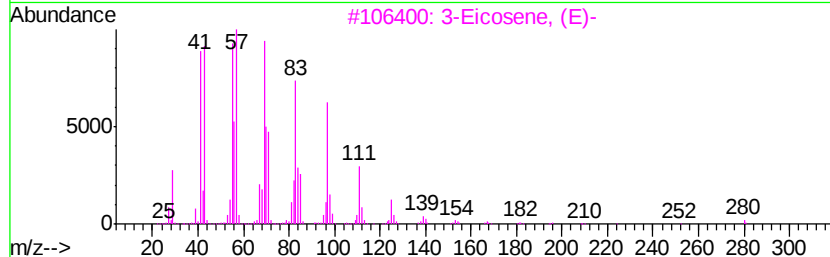
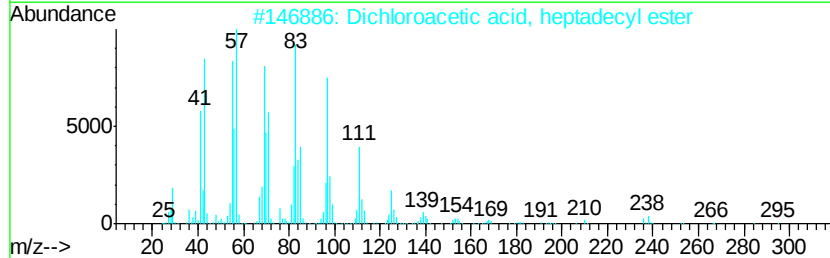
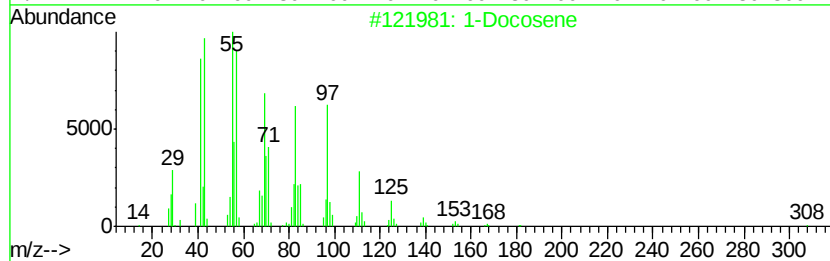
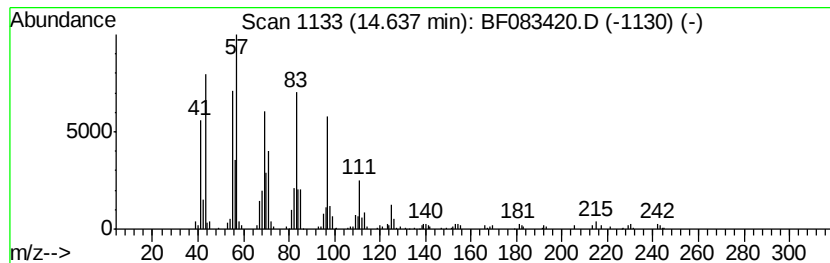
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 1-Docosene Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.64 | 4.47 ng | 428088 | Chrysene-d12     | 14.73 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | 1-Docosene                          | 308 | C22H44      | 001599-67-3  | 91   |
| 2    |    | Dichloroacetic acid, heptadecyl ... | 366 | C19H36Cl2O2 | 1000282-98-2 | 90   |
| 3    |    | 3-Eicosene, (E)-                    | 280 | C20H40      | 074685-33-9  | 87   |
| 4    |    | 1-Nonadecene                        | 266 | C19H38      | 018435-45-5  | 87   |
| 5    |    | Pentafluoropropionic acid, hepta... | 402 | C20H35F5O2  | 1000283-04-2 | 87   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120215\  
 Data File : BF083420.D  
 Acq On : 2 Dec 2015 16:12  
 Operator : UM/IZ  
 Sample : G4583-06  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-P3WC-06(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113015.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| 2-Pentanone, 4-hy... | 4.65  | 28.1    | ng    | 1312690  | 1                     | 6.54  | 934536  | 20.0 |
| Benzene, 1,2,3-tr... | 6.36  | 6.6     | ng    | 306423   | 1                     | 6.54  | 934536  | 20.0 |
| Naphthalene, 1-et... | 9.09  | 11.1    | ng    | 852686   | 3                     | 9.57  | 1538640 | 20.0 |
| Naphthalene, 1,6-... | 9.16  | 8.2     | ng    | 628747   | 3                     | 9.57  | 1538640 | 20.0 |
| Naphthalene, 2,3-... | 9.24  | 11.4    | ng    | 877228   | 3                     | 9.57  | 1538640 | 20.0 |
| Naphthalene, 2,3,... | 9.78  | 5.7     | ng    | 441229   | 3                     | 9.57  | 1538640 | 20.0 |
| Naphthalene, 1,4,... | 9.89  | 4.3     | ng    | 331096   | 3                     | 9.57  | 1538640 | 20.0 |
| Pentadecane, 2,6,... | 10.25 | 6.2     | ng    | 475878   | 3                     | 9.57  | 1538640 | 20.0 |
| Tridecane, 5-propyl- | 10.52 | 5.9     | ng    | 515449   | 4                     | 11.12 | 1740070 | 20.0 |
| Phenanthrene, 1-m... | 11.72 | 6.5     | ng    | 566699   | 4                     | 11.12 | 1740070 | 20.0 |
| Anthracene, 2-met... | 11.74 | 6.5     | ng    | 568747   | 4                     | 11.12 | 1740070 | 20.0 |
| Anthracene, 9-met... | 11.81 | 4.2     | ng    | 363164   | 4                     | 11.12 | 1740070 | 20.0 |
| 4H-Cyclopenta[def... | 11.85 | 10.5    | ng    | 916009   | 4                     | 11.12 | 1740070 | 20.0 |
| Naphthalene, 2-ph... | 12.10 | 6.1     | ng    | 531533   | 4                     | 11.12 | 1740070 | 20.0 |
| Phenanthrene, 2,5... | 12.43 | 5.6     | ng    | 483776   | 4                     | 11.12 | 1740070 | 20.0 |
| 1-Docosene           | 14.64 | 4.5     | ng    | 428088   | 5                     | 14.73 | 1915450 | 20.0 |