

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF103018\  
 Data File : BF110316.D  
 Acq On : 30 Oct 2018 21:12  
 Operator : JU/SJ  
 Sample : J5711-02 2X  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-2

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:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF102318.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 2.263       | 26            | 30          | 39           | rVB      | 118431         | 175653        | 19.30%          | 1.301%        |
| 2         | 3.751       | 273           | 283         | 288          | rBV4     | 6096           | 11352         | 1.25%           | 0.084%        |
| 3         | 3.875       | 297           | 304         | 310          | rBV      | 10216          | 17062         | 1.88%           | 0.126%        |
| 4         | 4.093       | 335           | 341         | 351          | rVB      | 307243         | 405619        | 44.58%          | 3.003%        |
| 5         | 4.322       | 376           | 380         | 385          | rBV2     | 11513          | 15887         | 1.75%           | 0.118%        |
| 6         | 4.722       | 444           | 448         | 451          | rBV      | 17690          | 20664         | 2.27%           | 0.153%        |
| 7         | 5.116       | 507           | 515         | 517          | rBV      | 19898          | 24092         | 2.65%           | 0.178%        |
| 8         | 5.193       | 524           | 528         | 535          | rBV      | 42291          | 52838         | 5.81%           | 0.391%        |
| 9         | 5.251       | 535           | 538         | 543          | rVB      | 9330           | 12012         | 1.32%           | 0.089%        |
| 10        | 5.381       | 556           | 560         | 564          | rBV3     | 14769          | 28147         | 3.09%           | 0.208%        |
| 11        | 5.422       | 564           | 567         | 570          | rVB2     | 16378          | 19358         | 2.13%           | 0.143%        |
| 12        | 5.463       | 570           | 574         | 576          | rBV      | 23719          | 22857         | 2.51%           | 0.169%        |
| 13        | 5.487       | 576           | 578         | 581          | rVB      | 12939          | 11246         | 1.24%           | 0.083%        |
| 14        | 5.581       | 589           | 594         | 601          | rBV      | 753663         | 801386        | 88.07%          | 5.934%        |
| 15        | 5.698       | 611           | 614         | 618          | rVB      | 12033          | 13655         | 1.50%           | 0.101%        |
| 16        | 5.740       | 618           | 621         | 627          | rVB5     | 15328          | 26687         | 2.93%           | 0.198%        |
| 17        | 5.787       | 627           | 629         | 632          | rVB      | 15942          | 11905         | 1.31%           | 0.088%        |
| 18        | 5.834       | 632           | 637         | 642          | rBV4     | 25501          | 40004         | 4.40%           | 0.296%        |
| 19        | 5.969       | 657           | 660         | 662          | rBV2     | 18161          | 20707         | 2.28%           | 0.153%        |
| 20        | 5.992       | 662           | 664         | 667          | rVB      | 28011          | 24630         | 2.71%           | 0.182%        |
| 21        | 6.028       | 667           | 670         | 673          | rBV2     | 32479          | 38275         | 4.21%           | 0.283%        |
| 22        | 6.110       | 681           | 684         | 686          | rVB      | 23765          | 18950         | 2.08%           | 0.140%        |
| 23        | 6.157       | 690           | 692         | 696          | rVB2     | 30623          | 29361         | 3.23%           | 0.217%        |
| 24        | 6.204       | 696           | 700         | 705          | rBV3     | 50874          | 78213         | 8.60%           | 0.579%        |
| 25        | 6.257       | 705           | 709         | 712          | rVV3     | 25068          | 35549         | 3.91%           | 0.263%        |
| 26        | 6.292       | 712           | 715         | 719          | rVV3     | 10186          | 13135         | 1.44%           | 0.097%        |
| 27        | 6.351       | 723           | 725         | 728          | rVB2     | 12144          | 12671         | 1.39%           | 0.094%        |
| 28        | 6.392       | 728           | 732         | 737          | rBV4     | 59775          | 90897         | 9.99%           | 0.673%        |
| 29        | 6.451       | 737           | 742         | 746          | rVB      | 123521         | 134435        | 14.77%          | 0.995%        |
| 30        | 6.487       | 746           | 748         | 750          | rBV2     | 64660          | 58749         | 6.46%           | 0.435%        |
| 31        | 6.504       | 750           | 751         | 755          | rVV4     | 30674          | 28712         | 3.16%           | 0.213%        |
| 32        | 6.545       | 755           | 758         | 761          | rVV      | 60722          | 51109         | 5.62%           | 0.378%        |
| 33        | 6.587       | 761           | 765         | 772          | rVV      | 898282         | 891398        | 97.96%          | 6.600%        |
| 34        | 6.640       | 772           | 774         | 780          | rVV2     | 48240          | 47655         | 5.24%           | 0.353%        |

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF103018\  
 Data File : BF110316.D  
 Acq On : 30 Oct 2018 21:12  
 Operator : JU/SJ  
 Sample : J5711-02 2X  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-2

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:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF102318.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |       |      |      |      |      |         |        |         |        |
|----|-------|------|------|------|------|---------|--------|---------|--------|
| 35 | 6.704 | 780  | 785  | 787  | rVV3 | 21272   | 32958  | 3.62%   | 0.244% |
| 36 | 6.734 | 787  | 790  | 797  | rVV  | 1046691 | 909956 | 100.00% | 6.737% |
| 37 | 6.792 | 797  | 800  | 804  | rVV3 | 49762   | 67240  | 7.39%   | 0.498% |
| 38 | 6.851 | 808  | 810  | 811  | rVV  | 18137   | 14734  | 1.62%   | 0.109% |
| 39 | 6.881 | 812  | 815  | 816  | rVV2 | 27051   | 31378  | 3.45%   | 0.232% |
| 40 | 6.916 | 819  | 821  | 823  | rVV3 | 42426   | 48865  | 5.37%   | 0.362% |
| 41 | 6.969 | 826  | 830  | 834  | rVV  | 763218  | 649049 | 71.33%  | 4.806% |
| 42 | 7.010 | 834  | 837  | 840  | rVV  | 49060   | 71586  | 7.87%   | 0.530% |
| 43 | 7.045 | 840  | 843  | 844  | rVV2 | 29757   | 35921  | 3.95%   | 0.266% |
| 44 | 7.081 | 844  | 849  | 851  | rVV2 | 85508   | 129711 | 14.25%  | 0.960% |
| 45 | 7.122 | 853  | 856  | 864  | rVV  | 737768  | 698013 | 76.71%  | 5.168% |
| 46 | 7.181 | 864  | 866  | 868  | rVV  | 36713   | 34170  | 3.76%   | 0.253% |
| 47 | 7.228 | 868  | 874  | 877  | rVV  | 134353  | 174800 | 19.21%  | 1.294% |
| 48 | 7.257 | 877  | 879  | 881  | rVV2 | 50278   | 44038  | 4.84%   | 0.326% |
| 49 | 7.292 | 881  | 885  | 889  | rVV2 | 96611   | 124696 | 13.70%  | 0.923% |
| 50 | 7.339 | 889  | 893  | 895  | rVV2 | 63710   | 73570  | 8.09%   | 0.545% |
| 51 | 7.387 | 898  | 901  | 906  | rVV2 | 43843   | 67992  | 7.47%   | 0.503% |
| 52 | 7.445 | 909  | 911  | 914  | rVV2 | 38593   | 40964  | 4.50%   | 0.303% |
| 53 | 7.498 | 914  | 920  | 922  | rVV  | 63460   | 85946  | 9.45%   | 0.636% |
| 54 | 7.528 | 922  | 925  | 933  | rVV  | 568269  | 590272 | 64.87%  | 4.370% |
| 55 | 7.610 | 933  | 939  | 942  | rVV3 | 54814   | 88226  | 9.70%   | 0.653% |
| 56 | 7.651 | 942  | 946  | 952  | rVV5 | 76025   | 116276 | 12.78%  | 0.861% |
| 57 | 7.704 | 952  | 955  | 957  | rVV2 | 37814   | 46083  | 5.06%   | 0.341% |
| 58 | 7.734 | 957  | 960  | 961  | rVV  | 52248   | 48418  | 5.32%   | 0.358% |
| 59 | 7.751 | 961  | 963  | 967  | rVV3 | 45397   | 72288  | 7.94%   | 0.535% |
| 60 | 7.787 | 967  | 969  | 971  | rVV  | 30580   | 28405  | 3.12%   | 0.210% |
| 61 | 7.845 | 974  | 979  | 981  | rVV2 | 41682   | 57607  | 6.33%   | 0.427% |
| 62 | 7.892 | 984  | 987  | 990  | rVV3 | 38582   | 59917  | 6.58%   | 0.444% |
| 63 | 7.922 | 990  | 992  | 995  | rVV4 | 34734   | 36534  | 4.01%   | 0.270% |
| 64 | 7.951 | 995  | 997  | 998  | rVV  | 22873   | 20852  | 2.29%   | 0.154% |
| 65 | 7.986 | 999  | 1003 | 1005 | rVV2 | 54411   | 70279  | 7.72%   | 0.520% |
| 66 | 8.010 | 1005 | 1007 | 1008 | rVV  | 30528   | 23964  | 2.63%   | 0.177% |
| 67 | 8.028 | 1008 | 1010 | 1012 | rVV  | 26959   | 21155  | 2.32%   | 0.157% |
| 68 | 8.057 | 1012 | 1015 | 1019 | rVV3 | 21831   | 28184  | 3.10%   | 0.209% |
| 69 | 8.110 | 1022 | 1024 | 1028 | rVB2 | 19532   | 18427  | 2.03%   | 0.136% |
| 70 | 8.192 | 1036 | 1038 | 1040 | rBV3 | 16270   | 14371  | 1.58%   | 0.106% |
| 71 | 8.251 | 1040 | 1048 | 1053 | rVV  | 794016  | 758460 | 83.35%  | 5.616% |

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF103018\  
 Data File : BF110316.D  
 Acq On : 30 Oct 2018 21:12  
 Operator : JU/SJ  
 Sample : J5711-02 2X  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-2

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\_F\METHODS\8270-BF102318.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 72 | 8.292  | 1053 | 1055 | 1059 | rVV2 | 53160  | 57492  | 6.32%  | 0.426% |
| 73 | 8.328  | 1059 | 1061 | 1064 | rVB2 | 19571  | 16985  | 1.87%  | 0.126% |
| 74 | 8.534  | 1090 | 1096 | 1101 | rVB6 | 15499  | 30564  | 3.36%  | 0.226% |
| 75 | 8.710  | 1124 | 1126 | 1131 | rVB2 | 12787  | 13802  | 1.52%  | 0.102% |
| 76 | 9.322  | 1227 | 1230 | 1241 | rBV  | 866177 | 822699 | 90.41% | 6.091% |
| 77 | 9.398  | 1241 | 1243 | 1247 | rVB3 | 10691  | 10397  | 1.14%  | 0.077% |
| 78 | 9.716  | 1294 | 1297 | 1304 | rBV  | 92634  | 89832  | 9.87%  | 0.665% |
| 79 | 10.010 | 1344 | 1347 | 1357 | rBV  | 698406 | 708587 | 77.87% | 5.246% |
| 80 | 10.416 | 1413 | 1416 | 1419 | rVB2 | 16756  | 13794  | 1.52%  | 0.102% |
| 81 | 10.804 | 1478 | 1482 | 1491 | rBV  | 391956 | 406032 | 44.62% | 3.006% |
| 82 | 10.910 | 1496 | 1500 | 1502 | rVB4 | 13758  | 13127  | 1.44%  | 0.097% |
| 83 | 11.374 | 1577 | 1579 | 1584 | rBV4 | 10648  | 10975  | 1.21%  | 0.081% |
| 84 | 11.504 | 1597 | 1601 | 1609 | rBV  | 647461 | 619137 | 68.04% | 4.584% |
| 85 | 11.727 | 1636 | 1639 | 1641 | rBV3 | 9403   | 12754  | 1.40%  | 0.094% |
| 86 | 12.016 | 1684 | 1688 | 1689 | rBV3 | 9345   | 11953  | 1.31%  | 0.089% |
| 87 | 12.957 | 1845 | 1848 | 1854 | rVB3 | 53842  | 65381  | 7.19%  | 0.484% |
| 88 | 13.057 | 1863 | 1865 | 1867 | rBV3 | 27255  | 24569  | 2.70%  | 0.182% |
| 89 | 13.086 | 1867 | 1870 | 1877 | rVV  | 641133 | 574400 | 63.12% | 4.253% |
| 90 | 13.274 | 1899 | 1902 | 1905 | rBV4 | 46223  | 56598  | 6.22%  | 0.419% |
| 91 | 14.157 | 2049 | 2052 | 2056 | rVB  | 528836 | 488890 | 53.73% | 3.620% |
| 92 | 15.698 | 2311 | 2314 | 2319 | rVB2 | 307275 | 411772 | 45.25% | 3.049% |
| 93 | 16.427 | 2435 | 2438 | 2445 | rVB  | 181932 | 326197 | 35.85% | 2.415% |

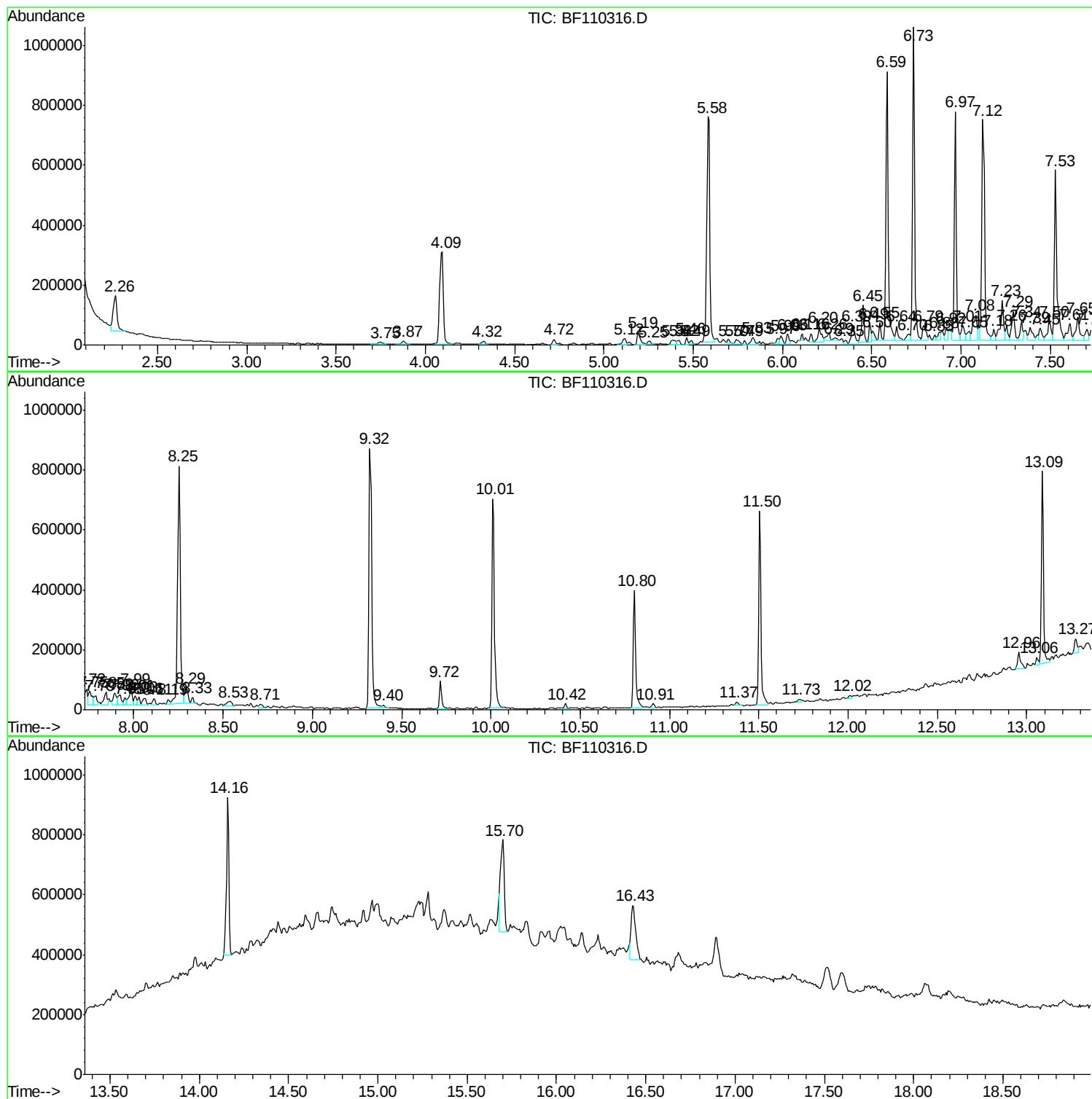
Sum of corrected areas: 13506112

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\  
Data File : BF110316.D  
Acq On : 30 Oct 2018 21:12  
Operator : JU/SJ  
Sample : J5711-02 2X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
SB-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\  
Data File : BF110316.D  
Acq On : 30 Oct 2018 21:12  
Operator : JU/SJ  
Sample : J5711-02 2X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-2

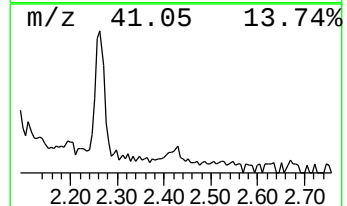
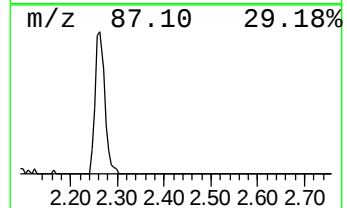
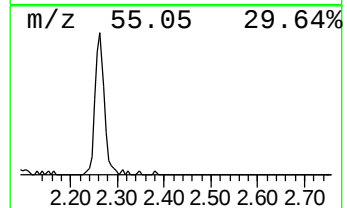
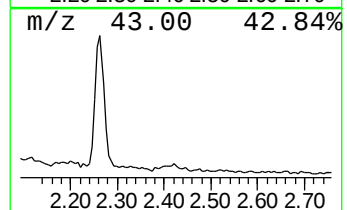
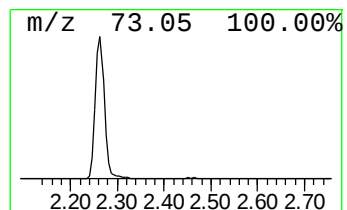
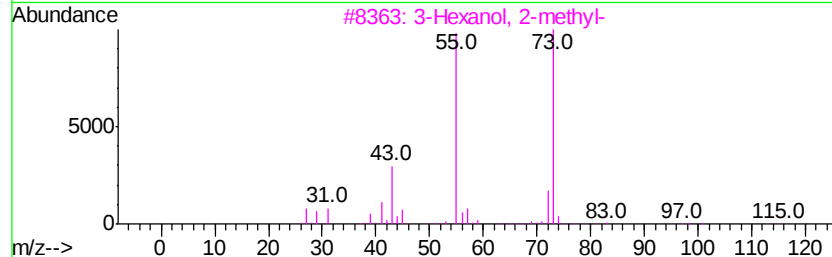
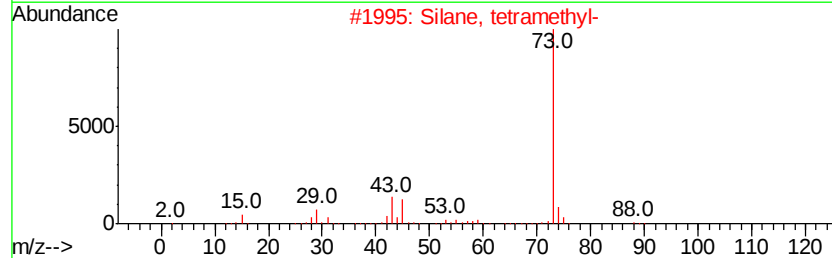
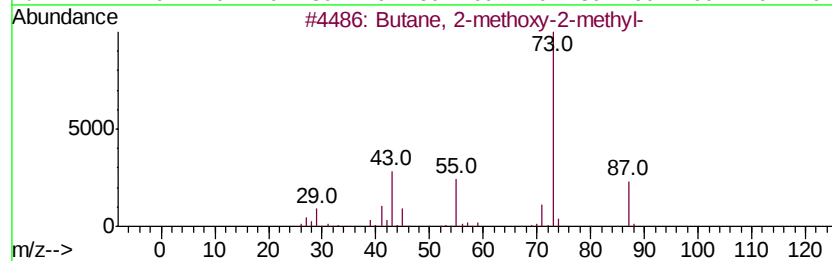
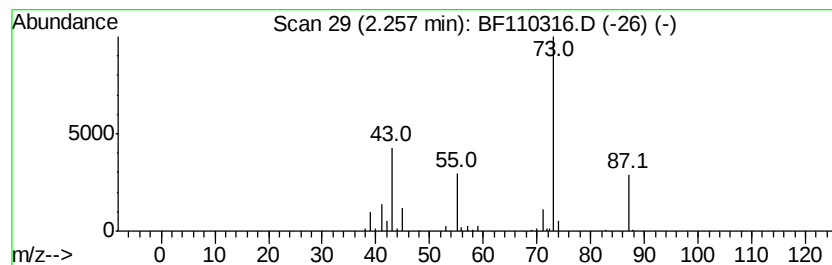
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 5

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 2.26 | 5.41 ng | 175653 | 1,4-Dichlorobenzene-d4 | 6.97 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 72   |
| 2    |    | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 35   |
| 3    |    | 3-Hexanol, 2-methyl-        | 116 | C7H16O  | 000617-29-8 | 25   |
| 4    |    | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 17   |
| 5    |    | N-Ethylformamide            | 73  | C3H7NO  | 000627-45-2 | 9    |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\  
Data File : BF110316.D  
Acq On : 30 Oct 2018 21:12  
Operator : JU/SJ  
Sample : J5711-02 2X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

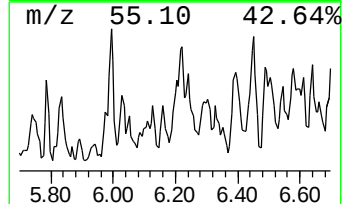
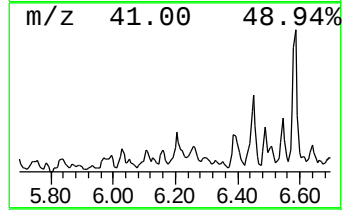
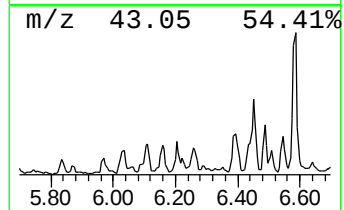
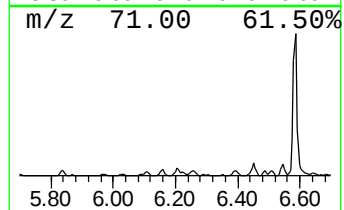
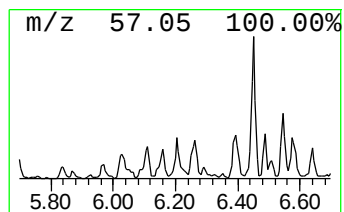
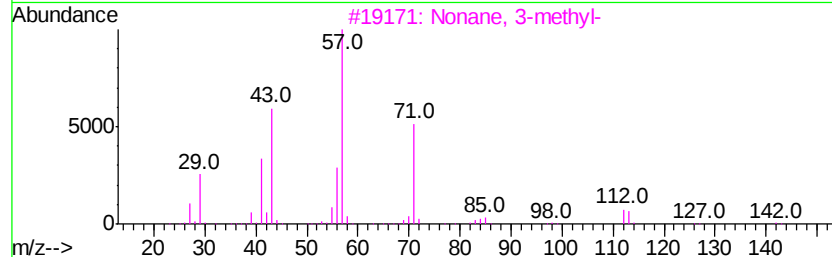
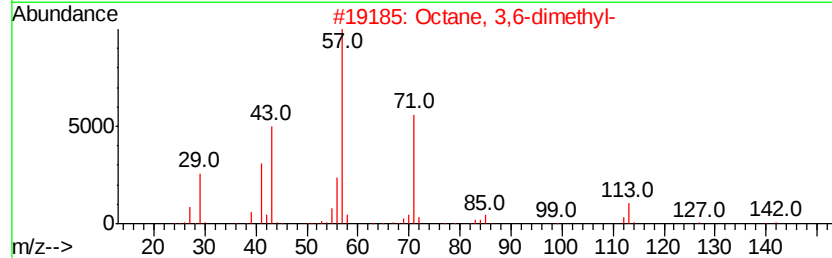
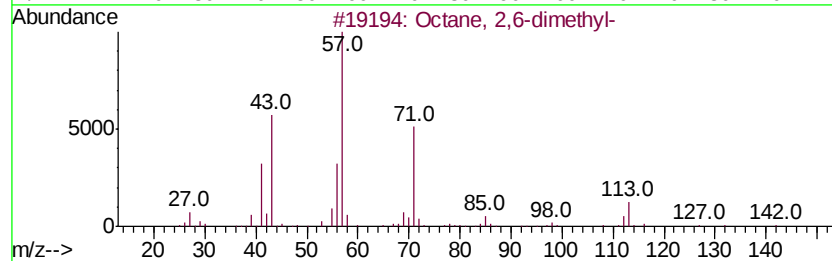
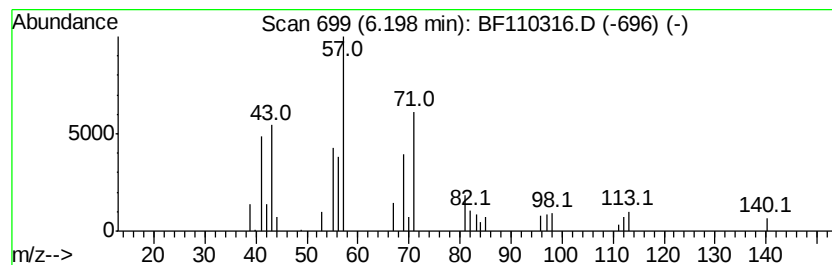
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Octane, 2,6-dimethyl- Concentration Rank 13

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 6.20 | 2.41 ng | 78213 | 1,4-Dichlorobenzene-d4 | 6.97 |

| Hit# | of | Tentative ID                      | MW  | MolForm   | CAS#         | Qual |
|------|----|-----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Octane, 2,6-dimethyl-             | 142 | C10H22    | 002051-30-1  | 72   |
| 2    |    | Octane, 3,6-dimethyl-             | 142 | C10H22    | 015869-94-0  | 59   |
| 3    |    | Nonane, 3-methyl-                 | 142 | C10H22    | 005911-04-6  | 59   |
| 4    |    | Undecane, 2,6-dimethyl-           | 184 | C13H28    | 017301-23-4  | 59   |
| 5    |    | Sulfurous acid, butyl octyl ester | 250 | C12H26O3S | 1000309-17-5 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\  
Data File : BF110316.D  
Acq On : 30 Oct 2018 21:12  
Operator : JU/SJ  
Sample : J5711-02 2X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-2

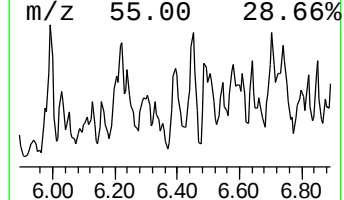
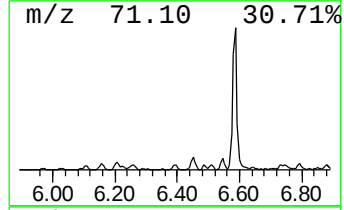
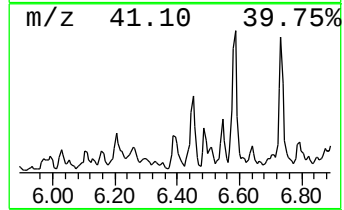
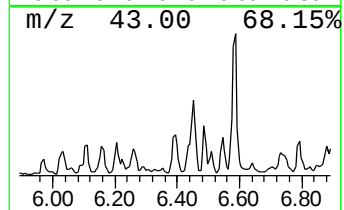
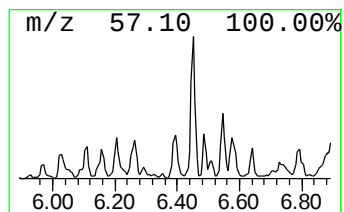
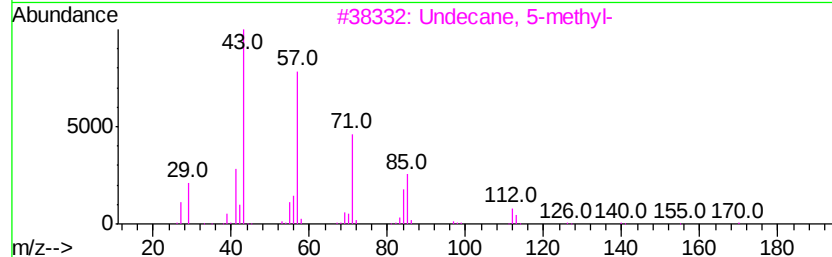
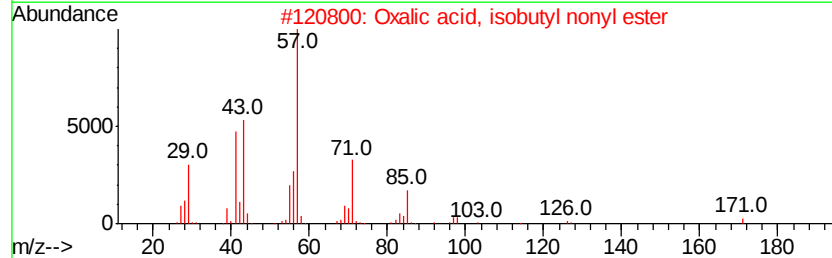
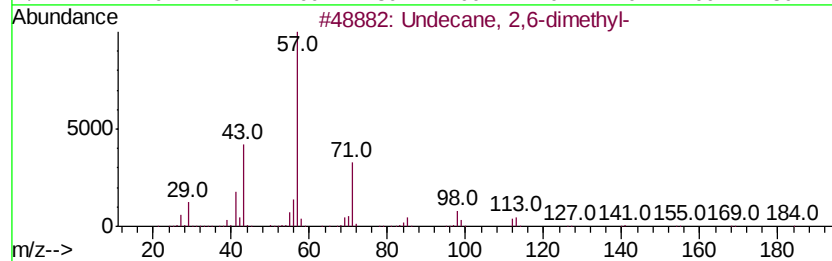
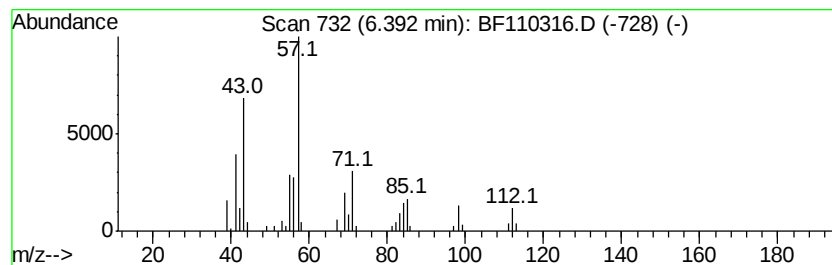
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 Undecane, 2,6-dimethyl- Concentration Rank 11

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 6.39 | 2.80 ng | 90897 | 1,4-Dichlorobenzene-d4 | 6.97 |

| Hit# | of | Tentative ID                      | MW  | MolForm  | CAS#         | Qual |
|------|----|-----------------------------------|-----|----------|--------------|------|
| 1    | 5  | Undecane, 2,6-dimethyl-           | 184 | C13H28   | 017301-23-4  | 53   |
| 2    |    | Oxalic acid, isobutyl nonyl ester | 272 | C15H28O4 | 1000309-37-4 | 52   |
| 3    |    | Undecane, 5-methyl-               | 170 | C12H26   | 001632-70-8  | 50   |
| 4    |    | Decane, 5-methyl-                 | 156 | C11H24   | 013151-35-4  | 47   |
| 5    |    | Octane, 2,4,6-trimethyl-          | 156 | C11H24   | 062016-37-9  | 47   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\  
Data File : BF110316.D  
Acq On : 30 Oct 2018 21:12  
Operator : JU/SJ  
Sample : J5711-02 2X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

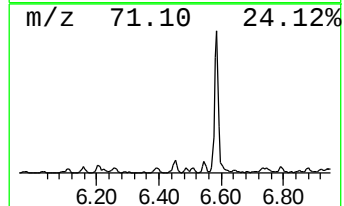
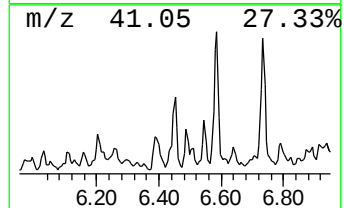
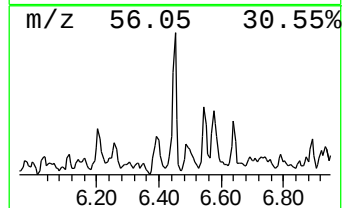
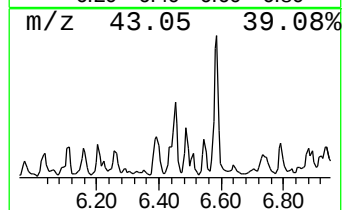
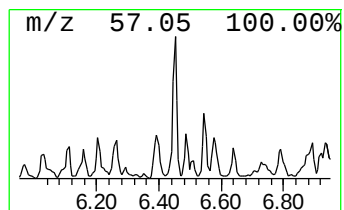
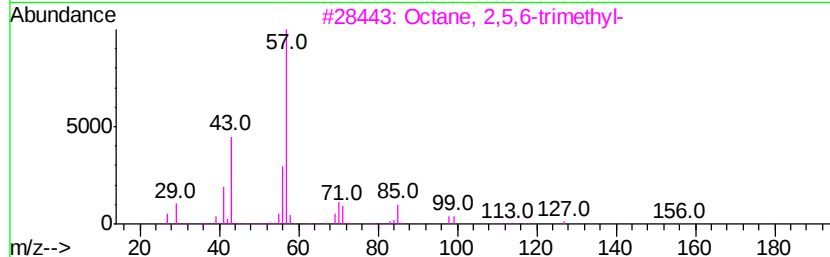
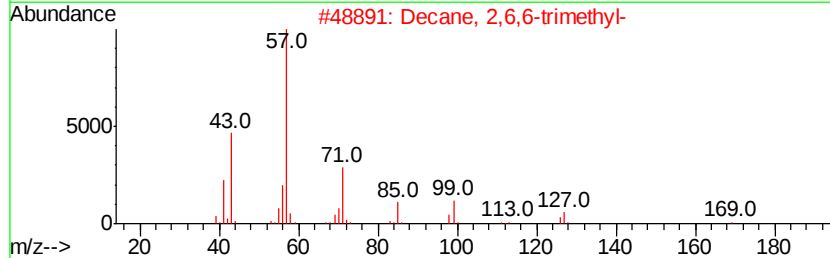
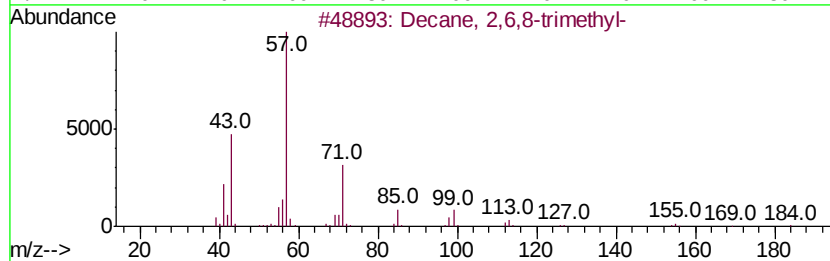
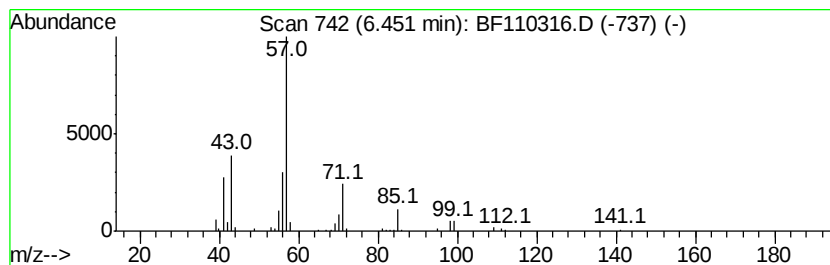
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Decane, 2,6,8-trimethyl- Concentration Rank 7

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.45 | 4.14 ng | 134435 | 1,4-Dichlorobenzene-d4 | 6.97 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 2,6,8-trimethyl- | 184 | C13H28  | 062108-26-3 | 59   |
| 2    |    |   | Decane, 2,6,6-trimethyl- | 184 | C13H28  | 062108-24-1 | 53   |
| 3    |    |   | Octane, 2,5,6-trimethyl- | 156 | C11H24  | 062016-14-2 | 53   |
| 4    |    |   | Octane, 4-ethyl-         | 142 | C10H22  | 015869-86-0 | 50   |
| 5    |    |   | Undecane, 2,6-dimethyl-  | 184 | C13H28  | 017301-23-4 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\  
Data File : BF110316.D  
Acq On : 30 Oct 2018 21:12  
Operator : JU/SJ  
Sample : J5711-02 2X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
SB-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

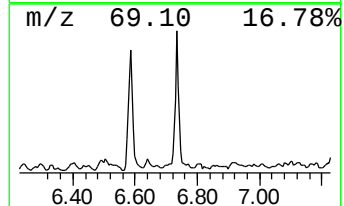
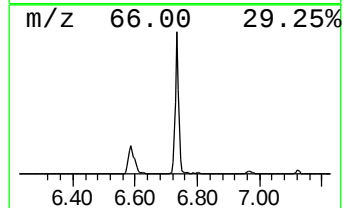
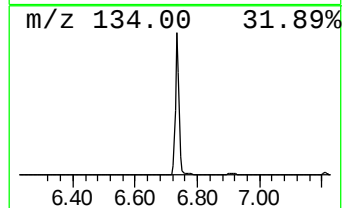
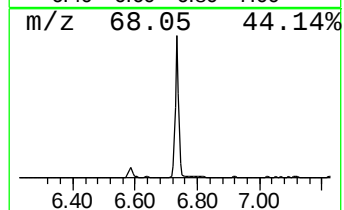
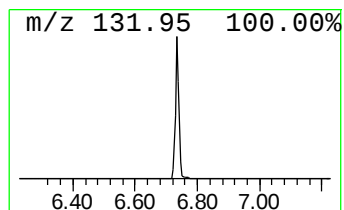
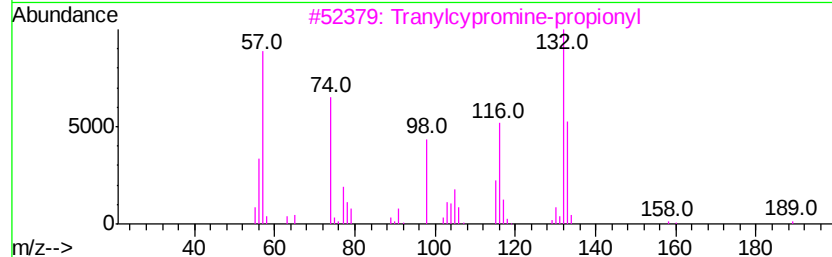
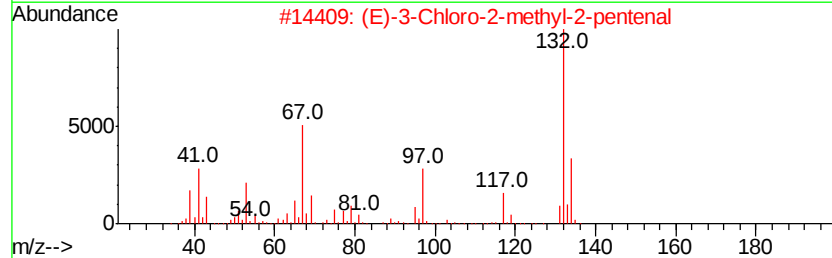
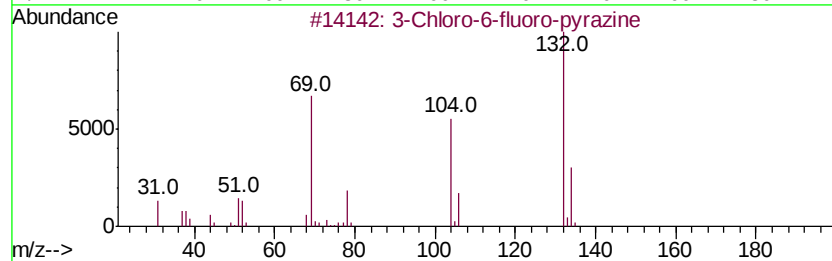
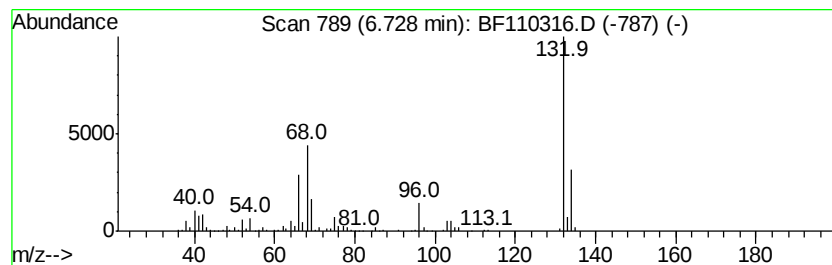
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 unknown6.73 Concentration Rank 1

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 6.73 | 28.04 ng | 909956 | 1,4-Dichlorobenzene-d4 | 6.97 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal    | 132 | C6H9ClO   | 031357-76-3  | 17   |
| 3    |    | Tranvlcypropine-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 10   |
| 4    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |
| 5    |    | 4-Chloro-2-fluoro-pyrimidine        | 132 | C4H2ClFN2 | 051422-00-5  | 9    |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\  
Data File : BF110316.D  
Acq On : 30 Oct 2018 21:12  
Operator : JU/SJ  
Sample : J5711-02 2X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-2

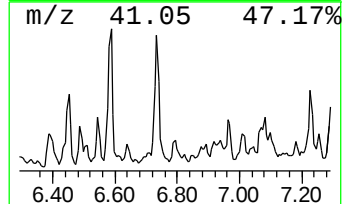
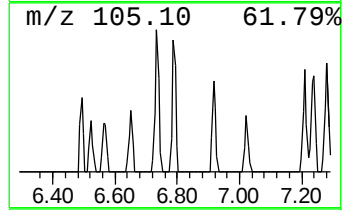
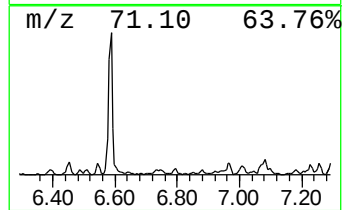
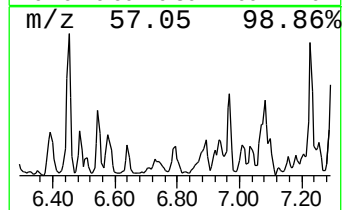
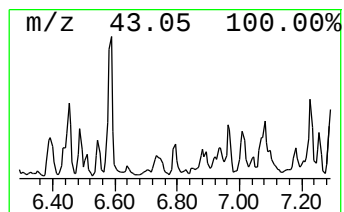
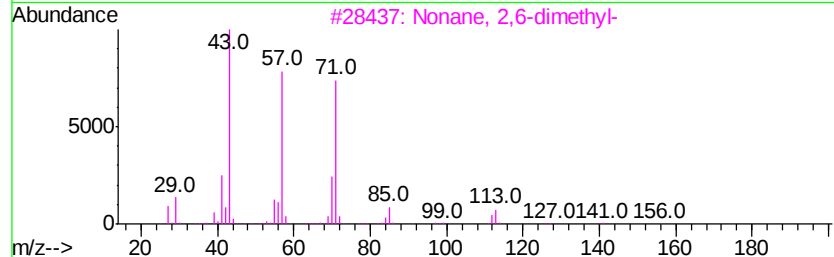
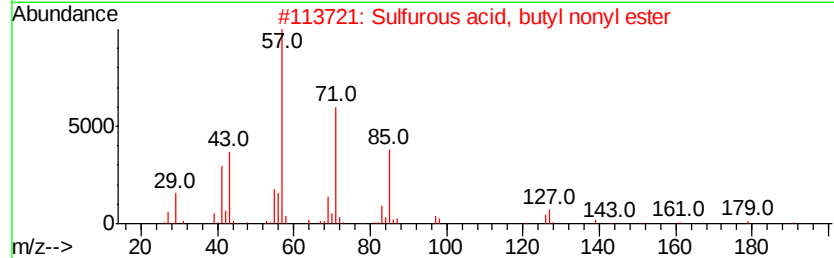
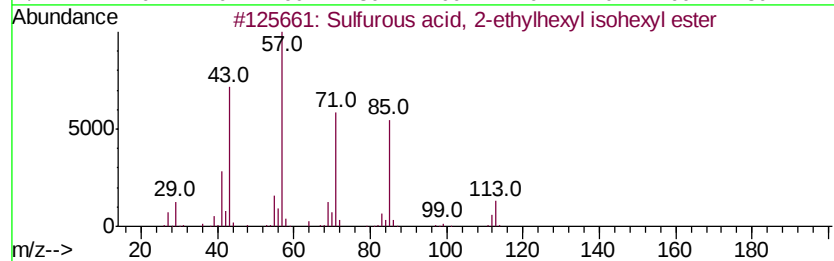
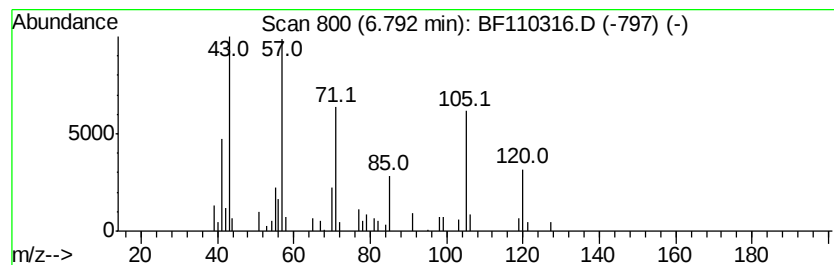
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 7 unknown6.79 Concentration Rank 18

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 6.79 | 2.07 ng | 67240 | 1,4-Dichlorobenzene-d4 | 6.97 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Sulfurous acid, 2-ethylhexyl iso... | 278 | C14H30O3S | 1000309-19-0 | 43   |
| 2    |    | Sulfurous acid, butyl nonyl ester   | 264 | C13H28O3S | 1000309-17-6 | 43   |
| 3    |    | Nonane, 2,6-dimethyl-               | 156 | C11H24    | 017302-28-2  | 43   |
| 4    |    | Heptane, 5-ethyl-2-methyl-          | 142 | C10H22    | 013475-78-0  | 43   |
| 5    |    | Decane, 4-methyl-                   | 156 | C11H24    | 002847-72-5  | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\  
Data File : BF110316.D  
Acq On : 30 Oct 2018 21:12  
Operator : JU/SJ  
Sample : J5711-02 2X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

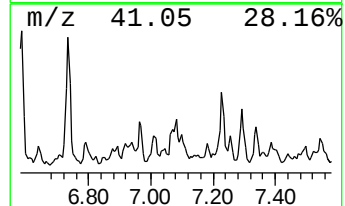
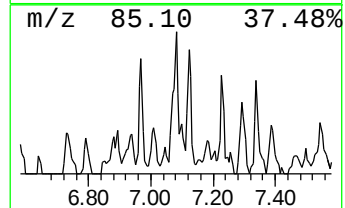
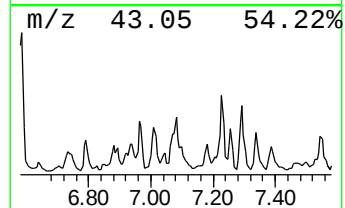
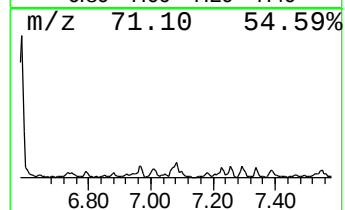
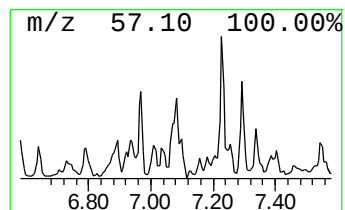
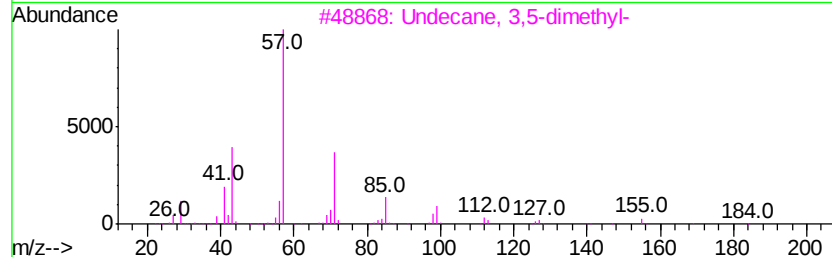
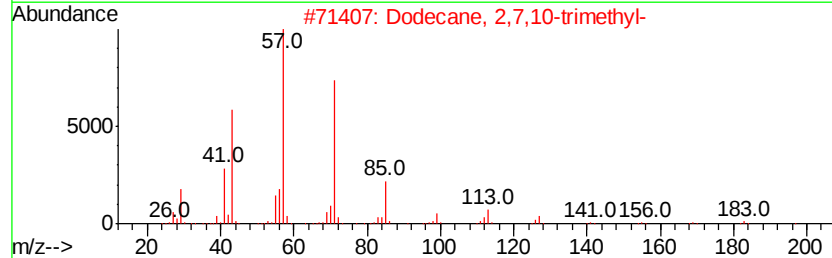
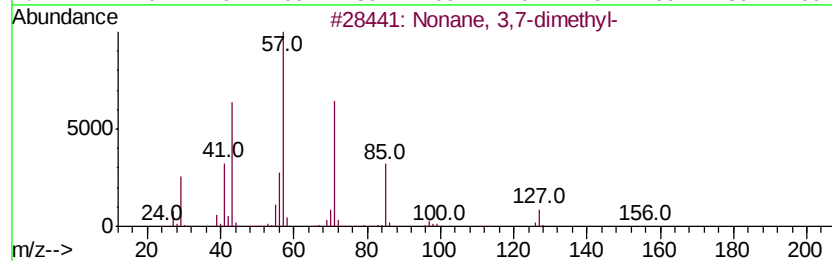
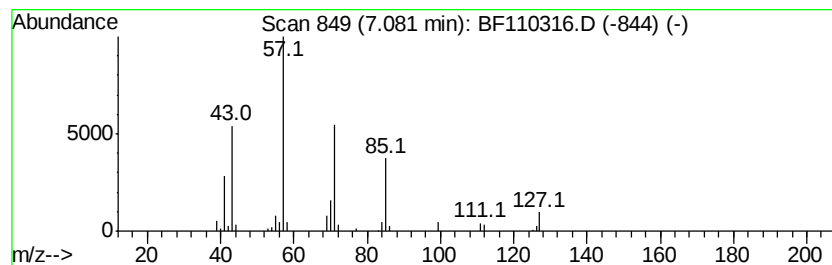
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 Nonane, 3,7-dimethyl- Concentration Rank 8

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 7.08 | 4.00 ng | 129711 | 1,4-Dichlorobenzene-d4 | 6.97 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Nonane, 3,7-dimethyl-       | 156 | C11H24  | 017302-32-8 | 72   |
| 2    |    | Dodecane, 2,7,10-trimethyl- | 212 | C15H32  | 074645-98-0 | 56   |
| 3    |    | Undecane, 3,5-dimethyl-     | 184 | C13H28  | 017312-81-1 | 56   |
| 4    |    | Undecane, 5-methyl-         | 170 | C12H26  | 001632-70-8 | 50   |
| 5    |    | Tetradecane                 | 198 | C14H30  | 000629-59-4 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\  
Data File : BF110316.D  
Acq On : 30 Oct 2018 21:12  
Operator : JU/SJ  
Sample : J5711-02 2X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

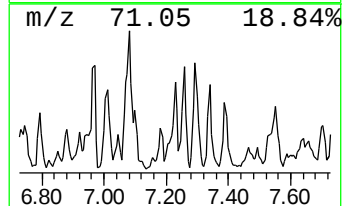
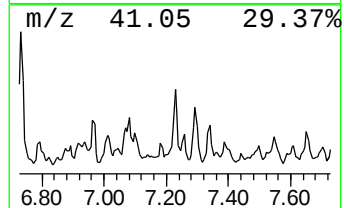
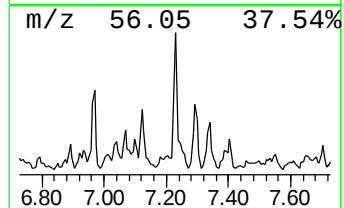
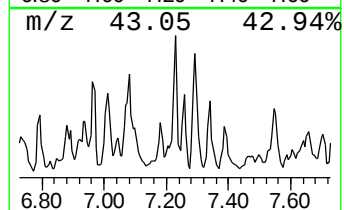
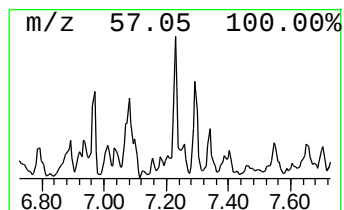
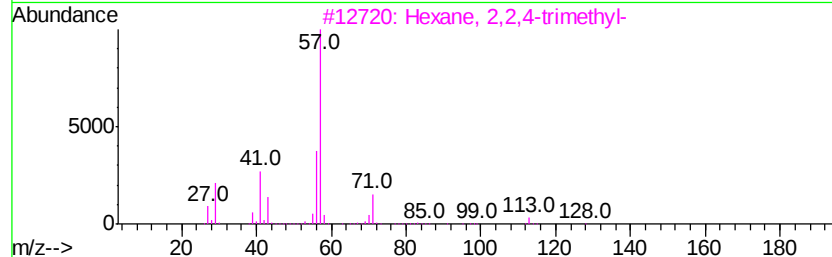
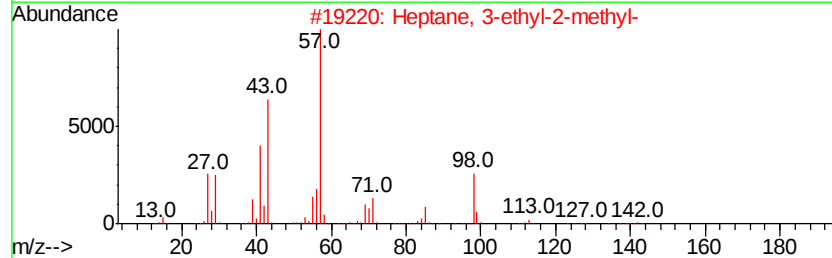
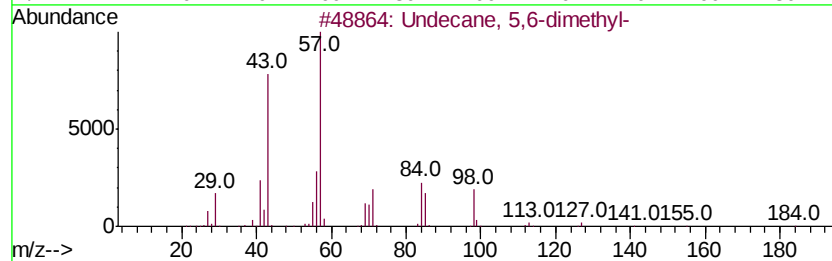
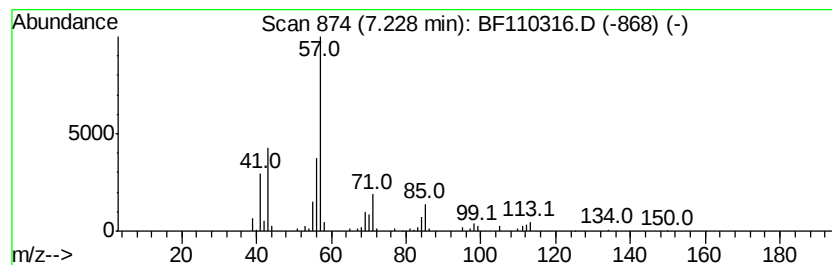
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Undecane, 5,6-dimethyl- Concentration Rank 6

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 7.23 | 5.39 ng | 174800 | 1,4-Dichlorobenzene-d4 | 6.97 |

| Hit# | of                         | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|----------------------------|-----|--------------|-------------|---------|------|------|
| 1    | Undecane, 5,6-dimethyl-    | 184 | C13H28       | 017615-91-7 | 64      |      |      |
| 2    | Heptane, 3-ethyl-2-methyl- | 142 | C10H22       | 014676-29-0 | 53      |      |      |
| 3    | Hexane, 2,2,4-trimethyl-   | 128 | C9H20        | 016747-26-5 | 47      |      |      |
| 4    | Hexane, 2,2,5-trimethyl-   | 128 | C9H20        | 003522-94-9 | 47      |      |      |
| 5    | Heptane, 2,2-dimethyl-     | 128 | C9H20        | 001071-26-7 | 47      |      |      |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\  
Data File : BF110316.D  
Acq On : 30 Oct 2018 21:12  
Operator : JU/SJ  
Sample : J5711-02 2X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

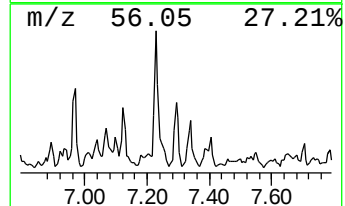
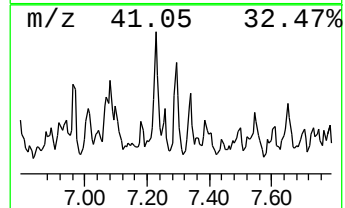
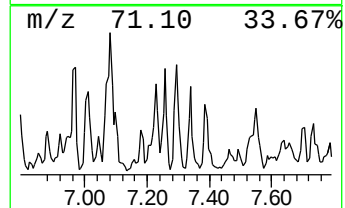
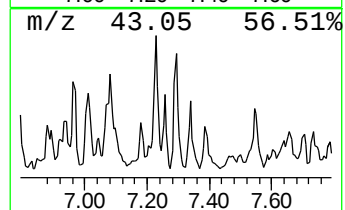
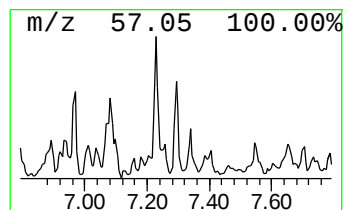
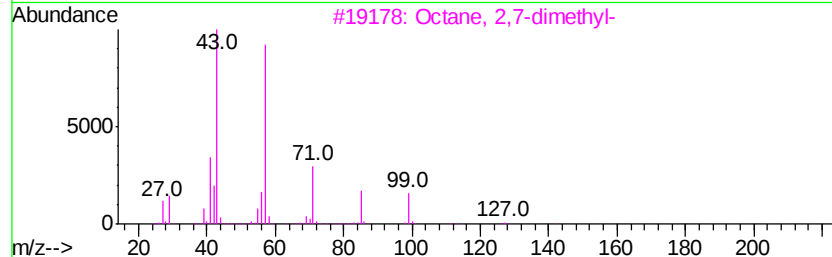
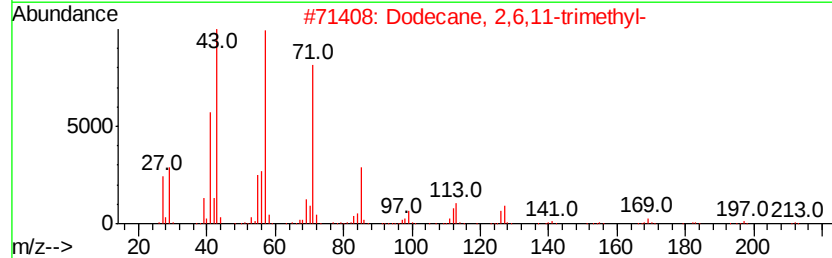
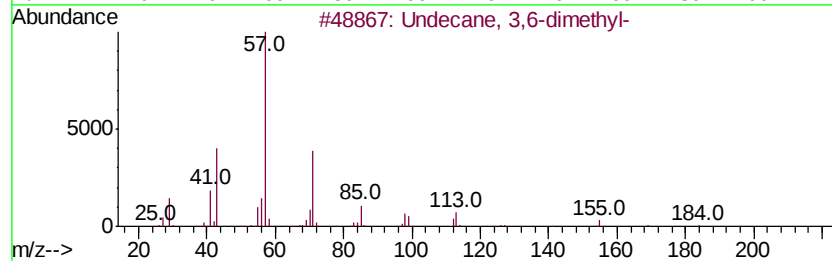
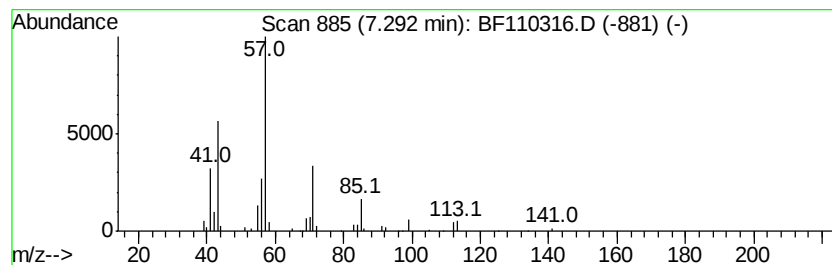
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Undecane, 3,6-dimethyl- Concentration Rank 9

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 7.29 | 3.84 ng | 124696 | 1,4-Dichlorobenzene-d4 | 6.97 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Undecane, 3,6-dimethyl-     | 184 | C13H28  | 017301-28-9 | 59   |
| 2    |    | Dodecane, 2,6,11-trimethyl- | 212 | C15H32  | 031295-56-4 | 59   |
| 3    |    | Octane, 2,7-dimethyl-       | 142 | C10H22  | 001072-16-8 | 59   |
| 4    |    | Decane, 2,2-dimethyl-       | 170 | C12H26  | 017302-37-3 | 53   |
| 5    |    | Octane, 2,6-dimethyl-       | 142 | C10H22  | 002051-30-1 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\  
Data File : BF110316.D  
Acq On : 30 Oct 2018 21:12  
Operator : JU/SJ  
Sample : J5711-02 2X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

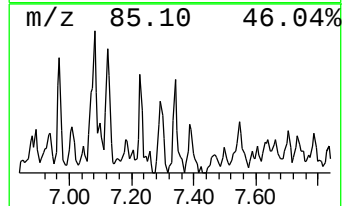
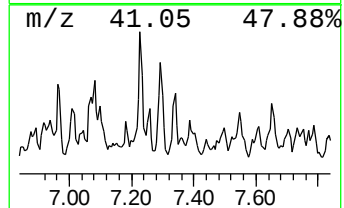
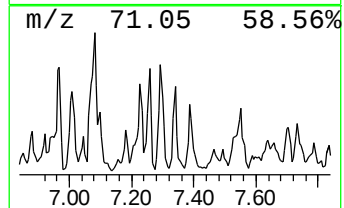
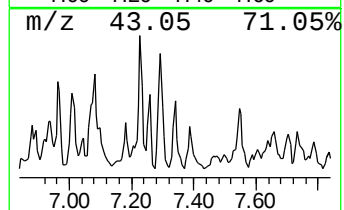
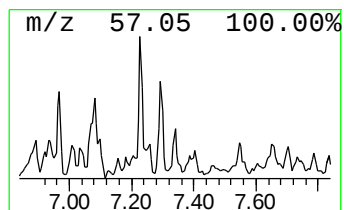
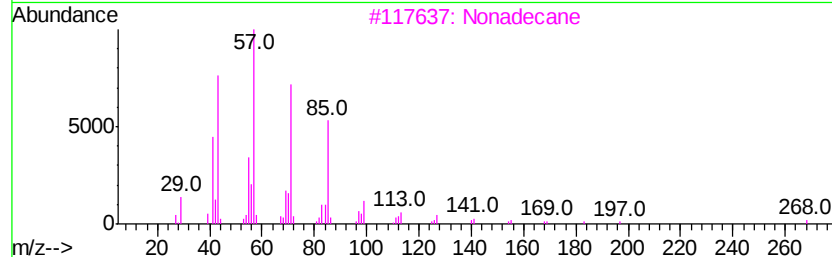
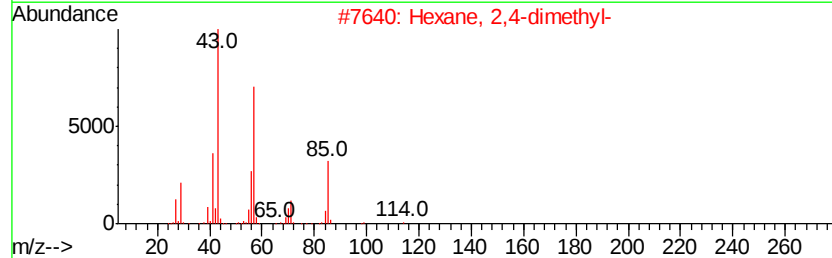
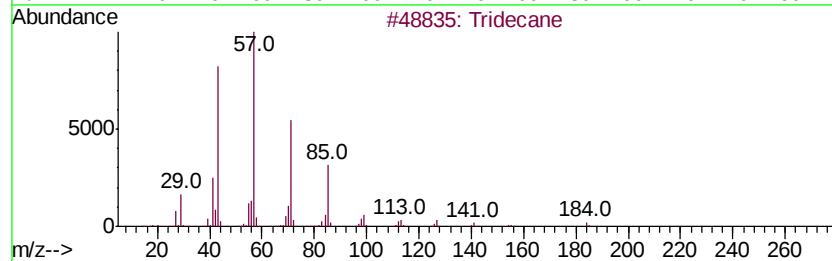
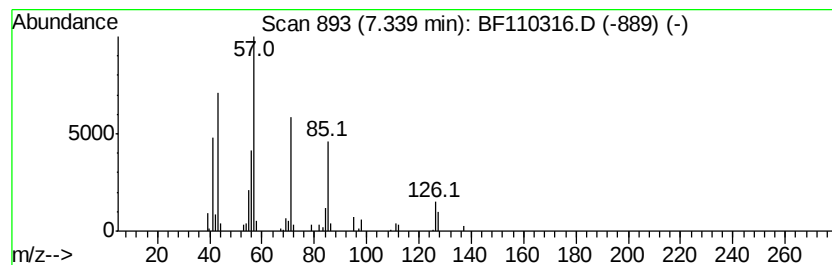
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Tridecane Concentration Rank 16

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 7.34 | 2.27 ng | 73570 | 1,4-Dichlorobenzene-d4 | 6.97 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Tridecane                           | 184 | C13H28    | 000629-50-5  | 53   |
| 2    |    | Hexane, 2,4-dimethyl-               | 114 | C8H18     | 000589-43-5  | 50   |
| 3    |    | Nonadecane                          | 268 | C19H40    | 000629-92-5  | 50   |
| 4    |    | Octane, 3-ethyl-2,7-dimethyl-       | 170 | C12H26    | 062183-55-5  | 50   |
| 5    |    | Sulfurous acid, 2-ethylhexyl iso... | 278 | C14H30O3S | 1000309-19-0 | 47   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\  
Data File : BF110316.D  
Acq On : 30 Oct 2018 21:12  
Operator : JU/SJ  
Sample : J5711-02 2X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
SB-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

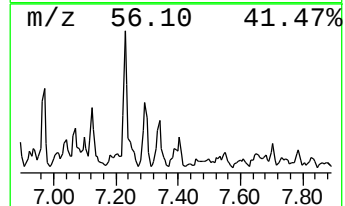
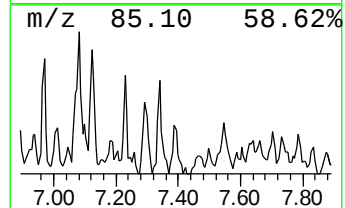
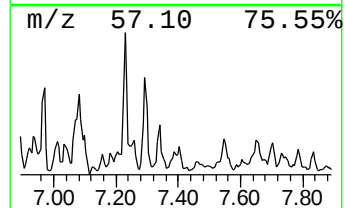
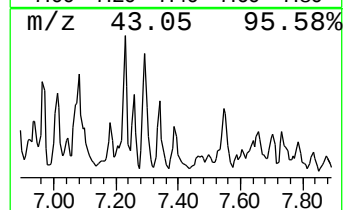
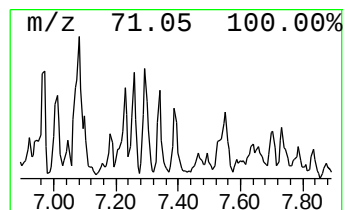
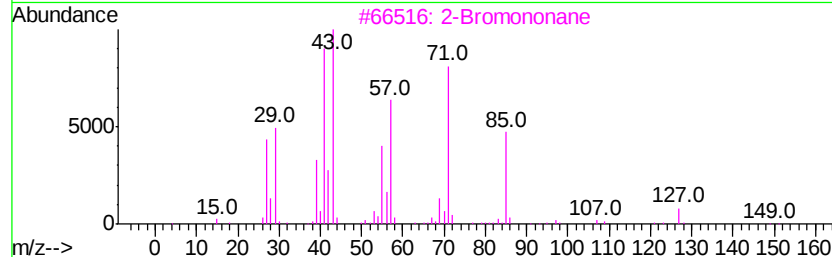
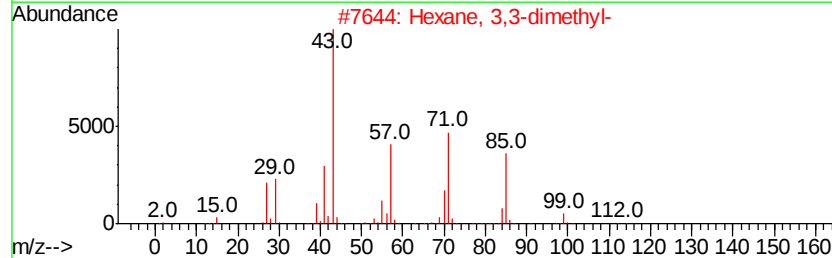
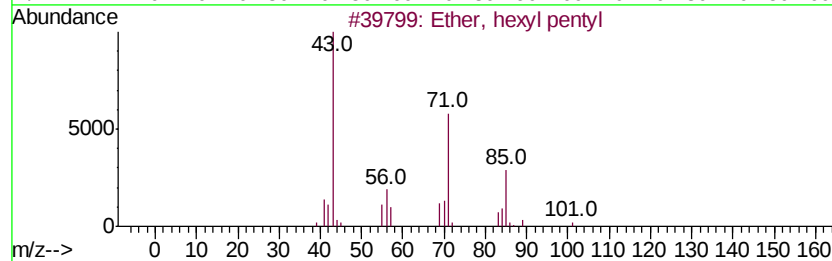
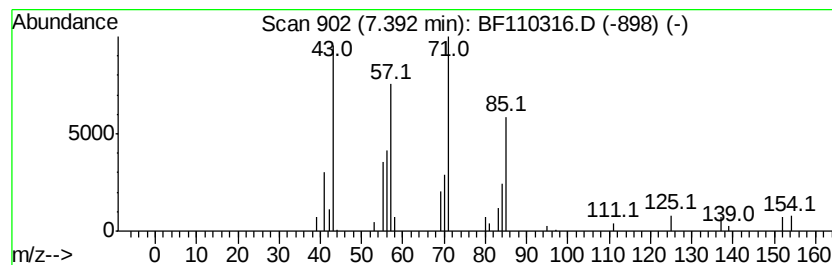
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Ether, hexyl pentyl Concentration Rank 17

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 7.39 | 2.10 ng | 67992 | 1,4-Dichlorobenzene-d4 | 6.97 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|-----|---------|-------------|------|
| 1    |    |   | Ether, hexyl pentyl          | 172 | C11H24O | 032357-83-8 | 80   |
| 2    |    |   | Hexane, 3,3-dimethyl-        | 114 | C8H18   | 000563-16-6 | 64   |
| 3    |    |   | 2-Bromononane                | 206 | C9H19Br | 002216-35-5 | 53   |
| 4    |    |   | Octane, 2,3,6,7-tetramethyl- | 170 | C12H26  | 052670-34-5 | 53   |
| 5    |    |   | Nonane, 2,6-dimethyl-        | 156 | C11H24  | 017302-28-2 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\  
Data File : BF110316.D  
Acq On : 30 Oct 2018 21:12  
Operator : JU/SJ  
Sample : J5711-02 2X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

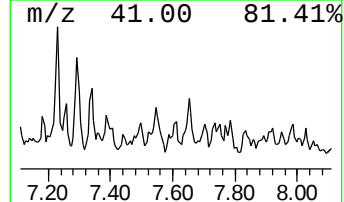
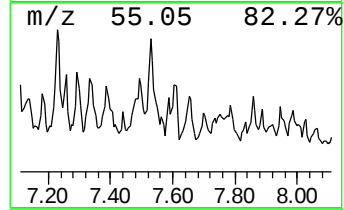
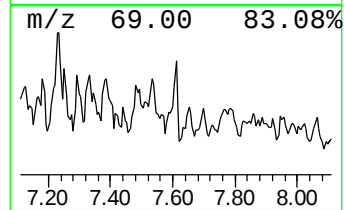
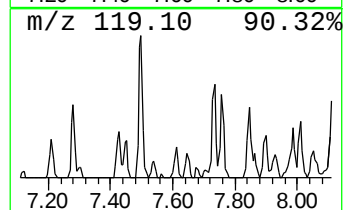
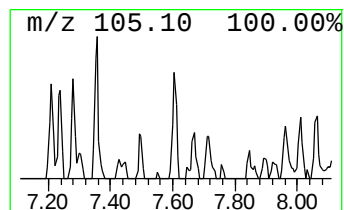
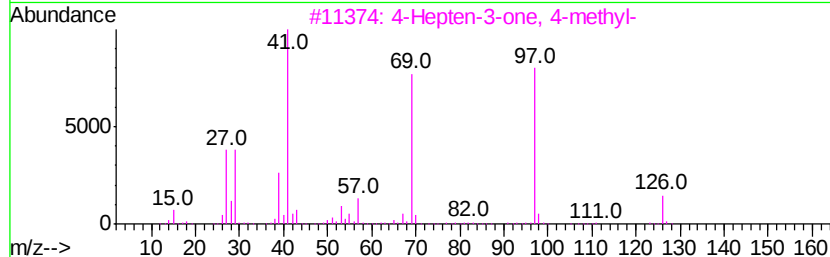
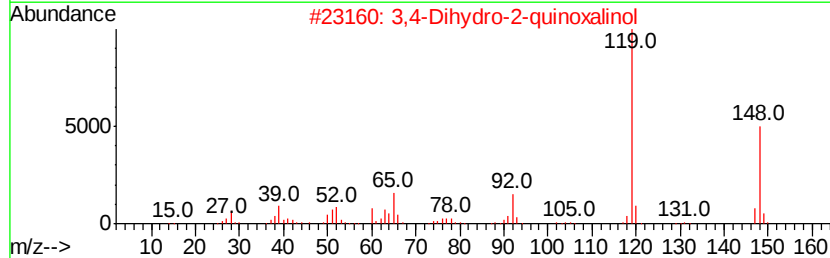
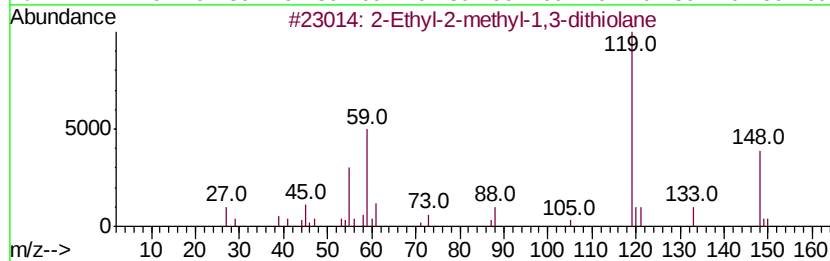
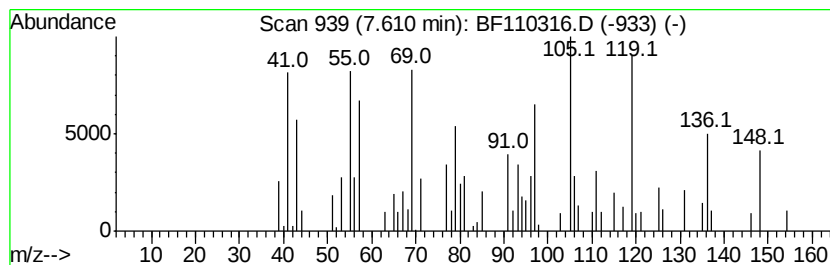
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 unknown7.61 Concentration Rank 14

| R.T. | EstConc | Area  | Relative to ISTD | R.T. |
|------|---------|-------|------------------|------|
| 7.61 | 2.33 ng | 88226 | Napthalene-d8    | 8.25 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm  | CAS#         | Qual |
|------|----|---|---------------------------------|-----|----------|--------------|------|
| 1    |    |   | 2-Ethyl-2-methyl-1,3-dithiolane | 148 | C6H12S2  | 006008-81-7  | 18   |
| 2    |    |   | 3,4-Dihydro-2-quinoxalinol      | 148 | C8H8N2O  | 1000332-36-8 | 14   |
| 3    |    |   | 4-Hepten-3-one, 4-methyl-       | 126 | C8H14O   | 022319-31-9  | 14   |
| 4    |    |   | 4-Methylphthalaldehyde          | 148 | C9H8O2   | 015158-36-8  | 14   |
| 5    |    |   | o-Toluic acid, 2-pentyl ester   | 206 | C13H18O2 | 212260-63-4  | 14   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\  
Data File : BF110316.D  
Acq On : 30 Oct 2018 21:12  
Operator : JU/SJ  
Sample : J5711-02 2X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

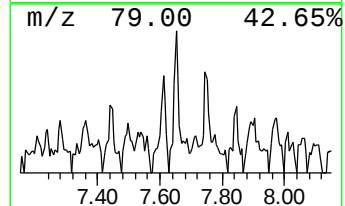
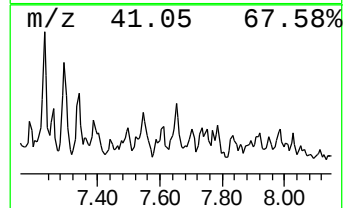
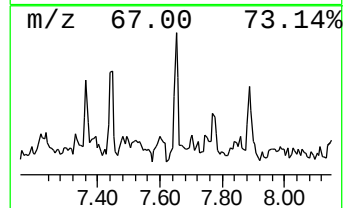
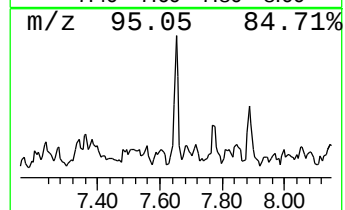
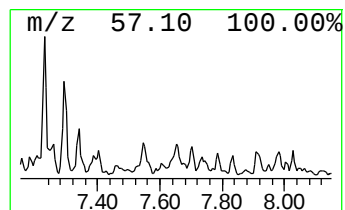
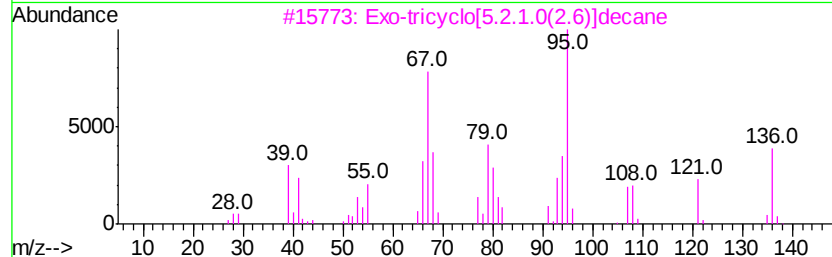
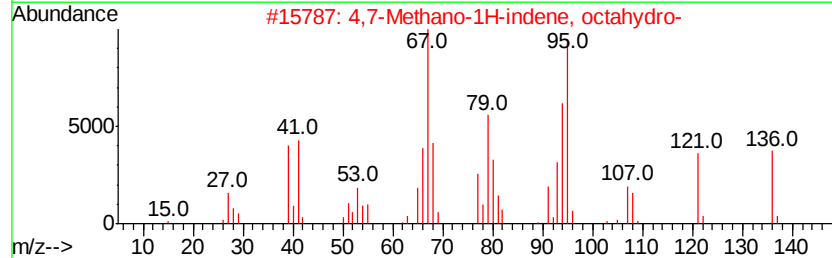
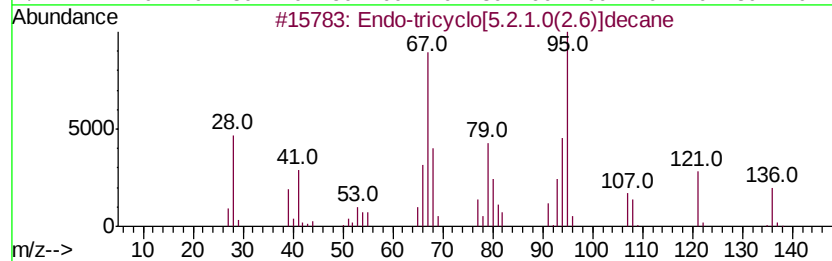
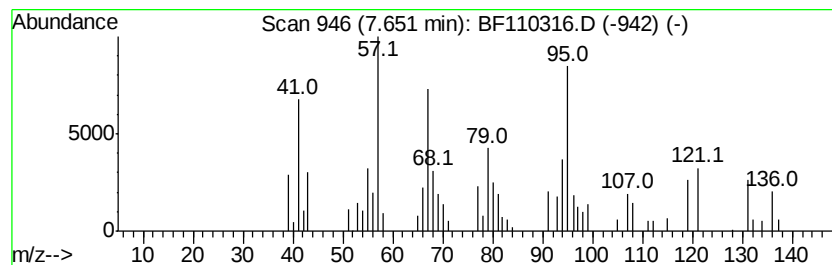
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Endo-tricyclo[5.2.1.0(2.6)]... Concentration Rank 10

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 7.65 | 3.07 ng | 116276 | Naphthalene-d8   | 8.25 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Endo-tricyclo[5.2.1.0(2.6)]decane   | 136 | C10H16  | 1000215-28-8 | 96   |
| 2    |    | 4,7-Methano-1H-indene, octahydro-   | 136 | C10H16  | 006004-38-2  | 95   |
| 3    |    | Exo-tricyclo[5.2.1.0(2.6)]decane    | 136 | C10H16  | 1000215-28-7 | 93   |
| 4    |    | 3-Cyclopentylcyclopentene           | 136 | C10H16  | 002690-17-7  | 72   |
| 5    |    | Bicyclo[2.2.1]heptane, 2-(1,1-di... | 164 | C12H20  | 069219-08-5  | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\  
Data File : BF110316.D  
Acq On : 30 Oct 2018 21:12  
Operator : JU/SJ  
Sample : J5711-02 2X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

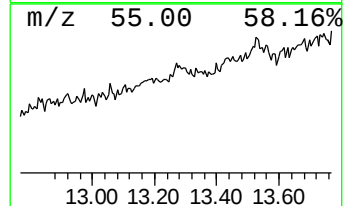
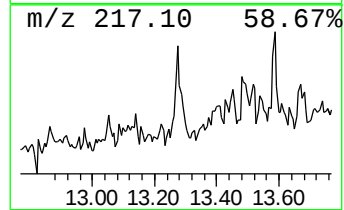
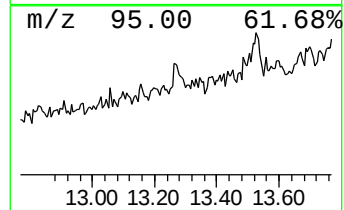
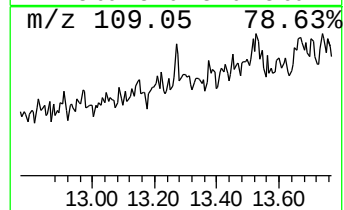
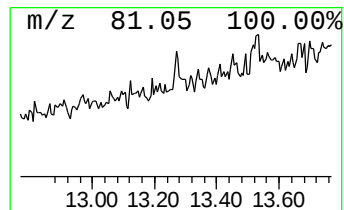
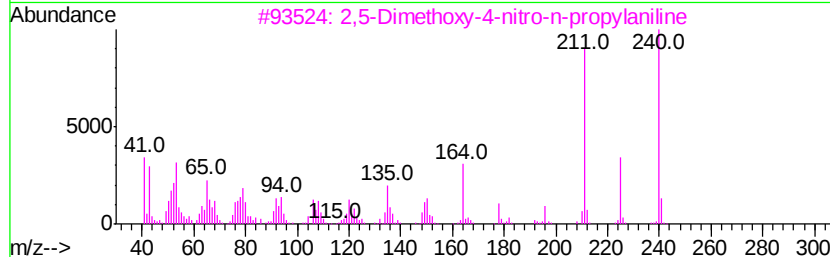
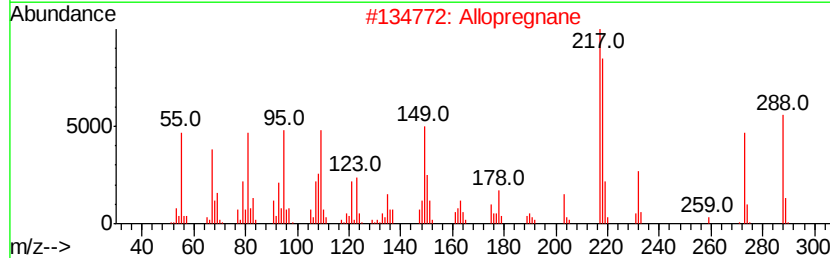
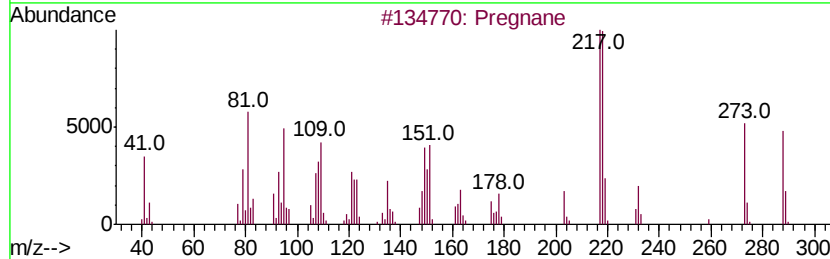
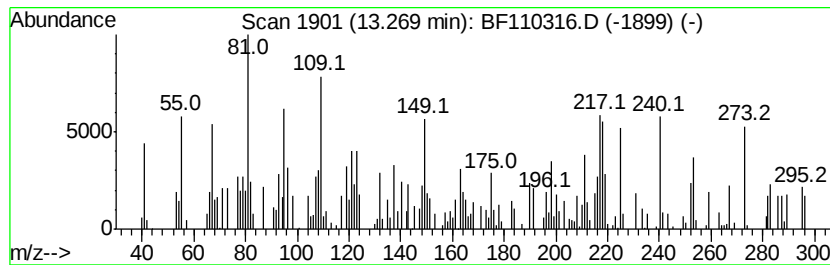
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 unknown13.27 Concentration Rank 15

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 13.27 | 2.32 ng | 56598 | Chrysene-d12     | 14.16 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Pregnane                            | 288 | C21H36     | 000481-26-5  | 25   |
| 2    |    | Allopregnane                        | 288 | C21H36     | 000641-85-0  | 25   |
| 3    |    | 2,5-Dimethoxy-4-nitro-n-propylan... | 240 | C11H16N2O4 | 1000213-86-5 | 14   |
| 4    |    | Cyclopentene-1-carbonitrile, 2-(... | 218 | C12H11ClN2 | 191980-01-5  | 14   |
| 5    |    | Cyclopropanecarbonitrile, 1-(p-c... | 253 | C16H12ClN  | 032589-55-2  | 12   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\  
Data File : BF110316.D  
Acq On : 30 Oct 2018 21:12  
Operator : JU/SJ  
Sample : J5711-02 2X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

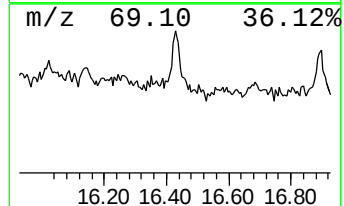
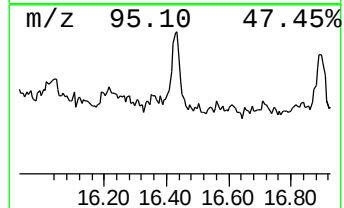
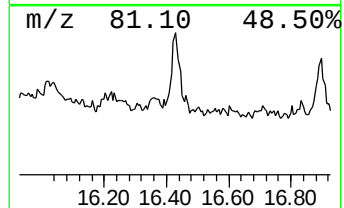
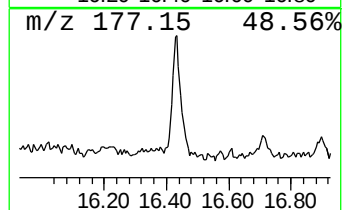
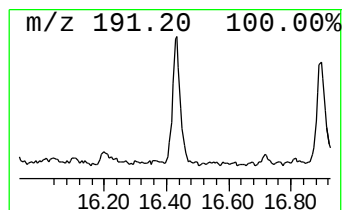
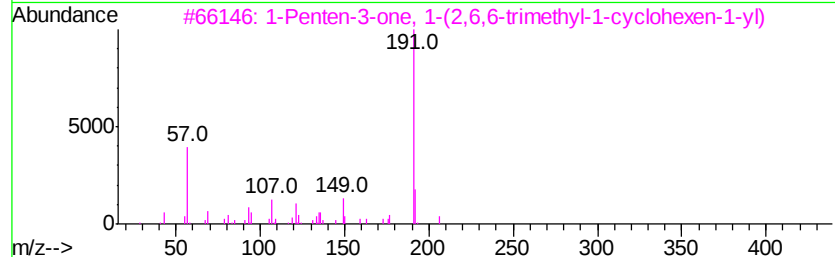
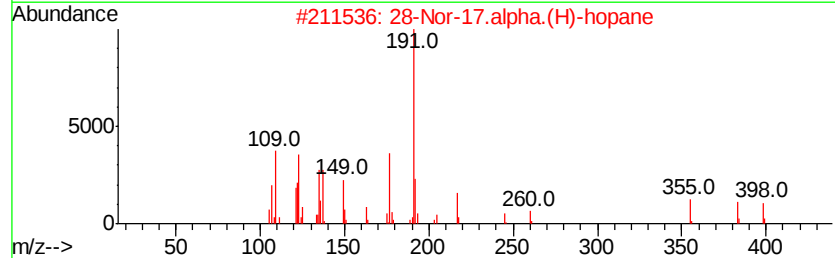
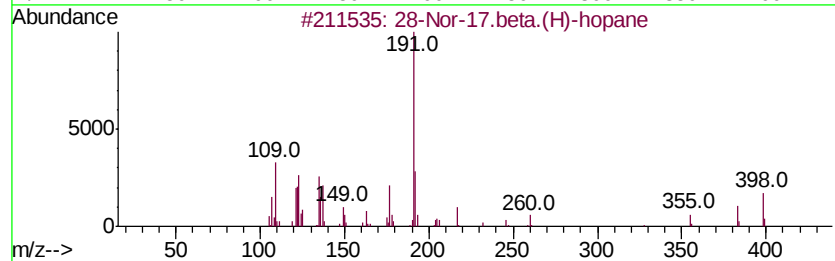
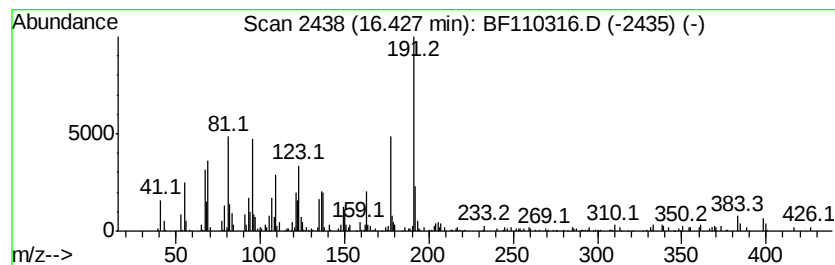
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 28-Nor-17.beta.(H)-hopane Concentration Rank 3

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 16.43 | 15.84 ng | 326197 | Perylene-d12     | 15.70 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 28-Nor-17.beta.(H)-hopane           | 398 | C29H50   | 036728-72-0 | 64   |
| 2    |    | 28-Nor-17.alpha.(H)-hopane          | 398 | C29H50   | 053584-60-4 | 46   |
| 3    |    | 1-Penten-3-one, 1-(2,6,6-trimeth... | 206 | C14H22O  | 000127-43-5 | 35   |
| 4    |    | 3,4-Dimethyl-2-(3-methyl-butyl)...  | 248 | C15H20O3 | 071940-29-9 | 22   |
| 5    |    | 5,5-Dimethyl-2-(2,2,2-trifluoro-... | 249 | C8H9F6NO | 020967-43-5 | 22   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF103018\  
 Data File : BF110316.D  
 Acq On : 30 Oct 2018 21:12  
 Operator : JU/SJ  
 Sample : J5711-02 2X  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF102318.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| Butane, 2-methoxy... | 2.26  | 5.4     | ng    | 175653   | 1                     | 6.97  | 649049 | 20.0 |
| Octane, 2,6-dimet... | 6.20  | 2.4     | ng    | 78213    | 1                     | 6.97  | 649049 | 20.0 |
| Undecane, 2,6-dim... | 6.39  | 2.8     | ng    | 90897    | 1                     | 6.97  | 649049 | 20.0 |
| Decane, 2,6,8-tri... | 6.45  | 4.1     | ng    | 134435   | 1                     | 6.97  | 649049 | 20.0 |
| unknown6.73          | 6.73  | 28.0    | ng    | 909956   | 1                     | 6.97  | 649049 | 20.0 |
| unknown6.79          | 6.79  | 2.1     | ng    | 67240    | 1                     | 6.97  | 649049 | 20.0 |
| Nonane, 3,7-dimet... | 7.08  | 4.0     | ng    | 129711   | 1                     | 6.97  | 649049 | 20.0 |
| Undecane, 5,6-dim... | 7.23  | 5.4     | ng    | 174800   | 1                     | 6.97  | 649049 | 20.0 |
| Undecane, 3,6-dim... | 7.29  | 3.8     | ng    | 124696   | 1                     | 6.97  | 649049 | 20.0 |
| Tridecane            | 7.34  | 2.3     | ng    | 73570    | 1                     | 6.97  | 649049 | 20.0 |
| Ether, hexyl pentyl  | 7.39  | 2.1     | ng    | 67992    | 1                     | 6.97  | 649049 | 20.0 |
| unknown7.61          | 7.61  | 2.3     | ng    | 88226    | 2                     | 8.25  | 758460 | 20.0 |
| Endo-tricyclo[5.2... | 7.65  | 3.1     | ng    | 116276   | 2                     | 8.25  | 758460 | 20.0 |
| unknown13.27         | 13.27 | 2.3     | ng    | 56598    | 5                     | 14.16 | 488890 | 20.0 |
| 28-Nor-17.beta.(H... | 16.43 | 15.8    | ng    | 326197   | 6                     | 15.70 | 411772 | 20.0 |