

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF011215\  
 Data File : BF076818.D  
 Acq On : 12 Jan 2015 20:40  
 Operator : IZ / TP  
 Sample : G1044-10  
 Misc : W  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SP-10

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-BF010915.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         | 1.407       | 26            | 28          | 34           | rVB      | 8655           | 17059         | 1.17%           | 0.160%        |
| 2         | 1.487       | 34            | 35          | 46           | rBV      | 94924          | 166435        | 11.39%          | 1.558%        |
| 3         | 1.944       | 72            | 75          | 80           | rVB      | 19789          | 34609         | 2.37%           | 0.324%        |
| 4         | 4.241       | 274           | 276         | 283          | rBV      | 41912          | 84091         | 5.76%           | 0.787%        |
| 5         | 5.167       | 354           | 357         | 369          | rVB      | 216529         | 415401        | 28.44%          | 3.888%        |
| 6         | 6.859       | 502           | 505         | 512          | rBB      | 23654          | 46488         | 3.18%           | 0.435%        |
| 7         | 7.476       | 557           | 559         | 565          | rBV      | 126803         | 227533        | 15.58%          | 2.129%        |
| 8         | 7.579       | 565           | 568         | 574          | rVB      | 588046         | 918754        | 62.89%          | 8.599%        |
| 9         | 7.979       | 601           | 603         | 606          | rBB      | 12499          | 16902         | 1.16%           | 0.158%        |
| 10        | 8.082       | 609           | 612         | 616          | rVB      | 139405         | 210192        | 14.39%          | 1.967%        |
| 11        | 8.402       | 637           | 640         | 644          | rBV      | 555230         | 874165        | 59.84%          | 8.181%        |
| 12        | 8.516       | 646           | 650         | 653          | rBB      | 45417          | 66328         | 4.54%           | 0.621%        |
| 13        | 9.373       | 722           | 725         | 728          | rBV      | 444546         | 713470        | 48.84%          | 6.677%        |
| 14        | 10.985      | 863           | 866         | 870          | rBV      | 199876         | 302506        | 20.71%          | 2.831%        |
| 15        | 13.214      | 1059          | 1061        | 1065         | rBB2     | 12883          | 23560         | 1.61%           | 0.220%        |
| 16        | 13.499      | 1084          | 1086        | 1090         | rVB      | 11620          | 15642         | 1.07%           | 0.146%        |
| 17        | 13.637      | 1095          | 1098        | 1102         | rVV      | 839379         | 1460876       | 100.00%         | 13.672%       |
| 18        | 13.762      | 1106          | 1109        | 1112         | rVB      | 25677          | 34789         | 2.38%           | 0.326%        |
| 19        | 14.162      | 1140          | 1144        | 1148         | rBV      | 59764          | 111799        | 7.65%           | 1.046%        |
| 20        | 14.700      | 1189          | 1191        | 1194         | rVB      | 30529          | 41855         | 2.87%           | 0.392%        |
| 21        | 14.882      | 1203          | 1207        | 1209         | rBV2     | 10098          | 15706         | 1.08%           | 0.147%        |
| 22        | 15.123      | 1224          | 1228        | 1232         | rBV      | 239467         | 357527        | 24.47%          | 3.346%        |
| 23        | 15.351      | 1243          | 1248        | 1250         | rBV2     | 12122          | 22435         | 1.54%           | 0.210%        |
| 24        | 15.454      | 1253          | 1257        | 1261         | rBV2     | 8826           | 15373         | 1.05%           | 0.144%        |
| 25        | 15.934      | 1294          | 1299        | 1302         | rBV3     | 6564           | 15851         | 1.09%           | 0.148%        |
| 26        | 16.185      | 1318          | 1321        | 1325         | rBV      | 7477           | 16120         | 1.10%           | 0.151%        |
| 27        | 16.288      | 1325          | 1330        | 1332         | rBV2     | 21286          | 33821         | 2.32%           | 0.317%        |
| 28        | 16.711      | 1365          | 1367        | 1369         | rBV2     | 9837           | 16517         | 1.13%           | 0.155%        |
| 29        | 16.768      | 1369          | 1372        | 1375         | rVB      | 704705         | 887149        | 60.73%          | 8.303%        |
| 30        | 16.940      | 1384          | 1387        | 1390         | rBV3     | 28190          | 56287         | 3.85%           | 0.527%        |
| 31        | 16.986      | 1390          | 1391        | 1394         | rVB      | 9569           | 14682         | 1.01%           | 0.137%        |
| 32        | 17.363      | 1419          | 1424        | 1427         | rBV3     | 5452           | 15500         | 1.06%           | 0.145%        |
| 33        | 17.614      | 1445          | 1446        | 1450         | rVV4     | 10603          | 23432         | 1.60%           | 0.219%        |
| 34        | 17.706      | 1452          | 1454        | 1457         | rVB      | 50989          | 68664         | 4.70%           | 0.643%        |

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 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SP-10

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

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

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF010915.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |         |
|----|--------|------|------|------|------|---------|---------|--------|---------|
| 35 | 17.854 | 1464 | 1467 | 1469 | rVB  | 296193  | 339828  | 23.26% | 3.180%  |
| 36 | 18.174 | 1492 | 1495 | 1497 | rBV  | 45105   | 56542   | 3.87%  | 0.529%  |
| 37 | 18.437 | 1514 | 1518 | 1521 | rBV  | 77798   | 90914   | 6.22%  | 0.851%  |
| 38 | 18.803 | 1548 | 1550 | 1554 | rBV  | 97890   | 127509  | 8.73%  | 1.193%  |
| 39 | 18.871 | 1554 | 1556 | 1558 | rVV  | 71850   | 95947   | 6.57%  | 0.898%  |
| 40 | 18.917 | 1558 | 1560 | 1562 | rVB  | 172235  | 224971  | 15.40% | 2.105%  |
| 41 | 19.009 | 1566 | 1568 | 1570 | rBV2 | 13194   | 17990   | 1.23%  | 0.168%  |
| 42 | 19.054 | 1570 | 1572 | 1575 | rBV  | 33191   | 42987   | 2.94%  | 0.402%  |
| 43 | 19.534 | 1612 | 1614 | 1617 | rVV  | 27893   | 40554   | 2.78%  | 0.380%  |
| 44 | 19.660 | 1623 | 1625 | 1627 | rBV  | 23157   | 22901   | 1.57%  | 0.214%  |
| 45 | 20.003 | 1653 | 1655 | 1657 | rBV  | 22546   | 26570   | 1.82%  | 0.249%  |
| 46 | 20.072 | 1658 | 1661 | 1663 | rVV  | 1223240 | 1303044 | 89.20% | 12.195% |
| 47 | 20.129 | 1663 | 1666 | 1668 | rVB  | 28228   | 38828   | 2.66%  | 0.363%  |
| 48 | 20.232 | 1673 | 1675 | 1676 | rBV  | 46562   | 48450   | 3.32%  | 0.453%  |
| 49 | 20.266 | 1676 | 1678 | 1680 | rVV  | 99655   | 106825  | 7.31%  | 1.000%  |
| 50 | 20.654 | 1710 | 1712 | 1717 | rVB2 | 11868   | 15228   | 1.04%  | 0.143%  |
| 51 | 21.306 | 1766 | 1769 | 1770 | rBV  | 19793   | 30261   | 2.07%  | 0.283%  |
| 52 | 21.340 | 1770 | 1772 | 1775 | rVB  | 294492  | 342037  | 23.41% | 3.201%  |
| 53 | 21.523 | 1785 | 1788 | 1790 | rVB  | 63047   | 80313   | 5.50%  | 0.752%  |
| 54 | 22.998 | 1914 | 1917 | 1920 | rBV  | 216946  | 291742  | 19.97% | 2.730%  |

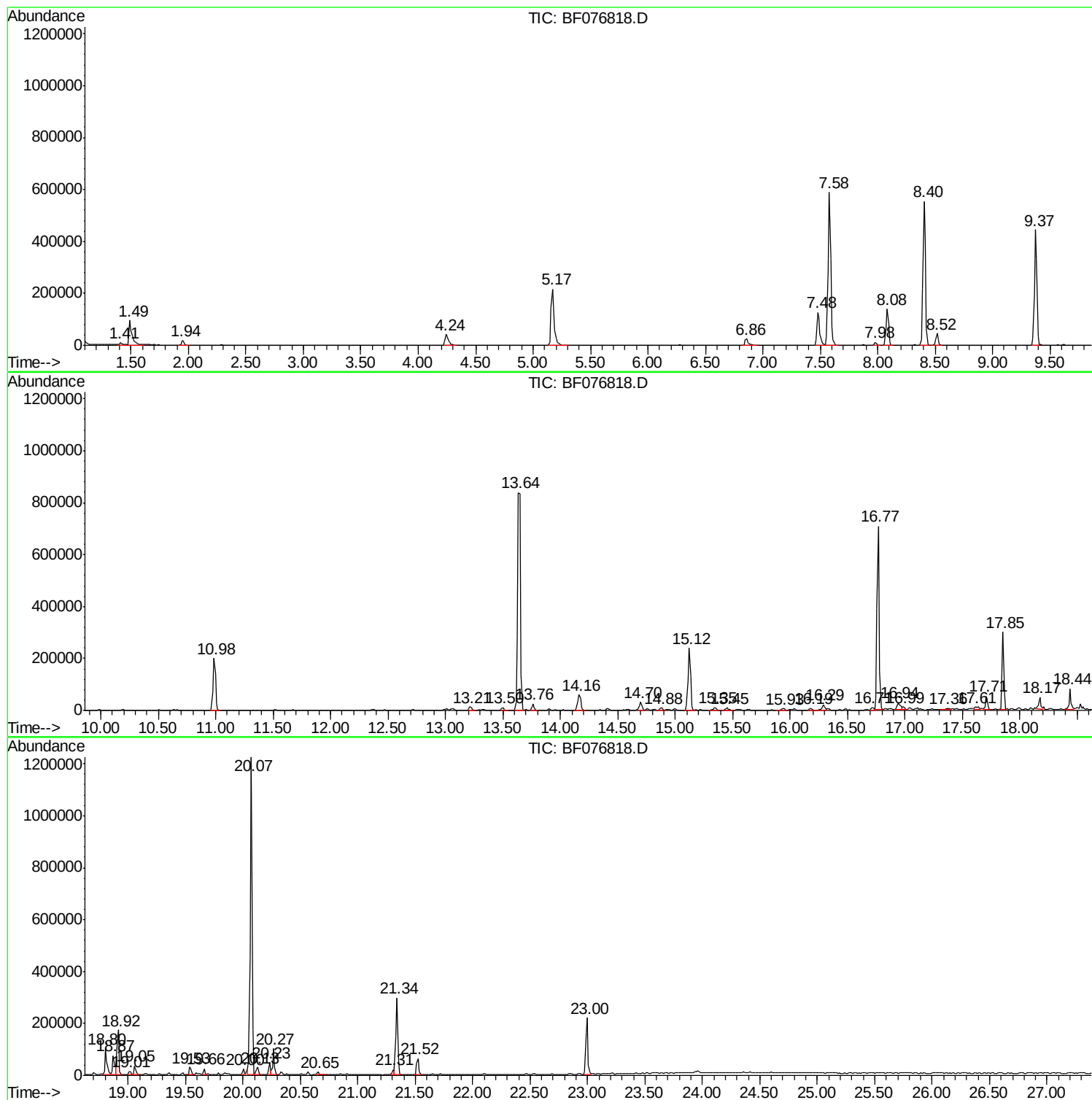
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Misc : W  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SP-10

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010915.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\BF011215\  
Data File : BF076818.D  
Acq On : 12 Jan 2015 20:40  
Operator : IZ / TP  
Sample : G1044-10  
Misc : W  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
SP-10

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF010915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

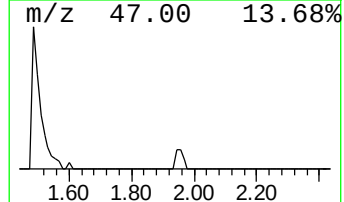
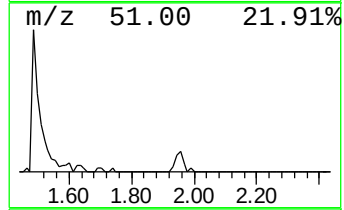
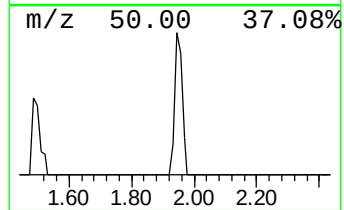
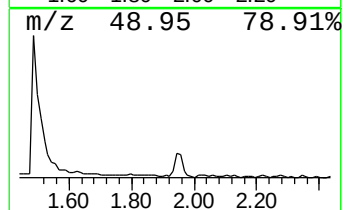
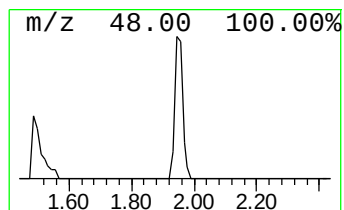
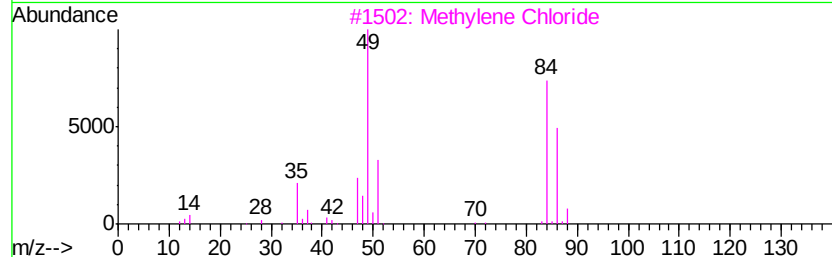
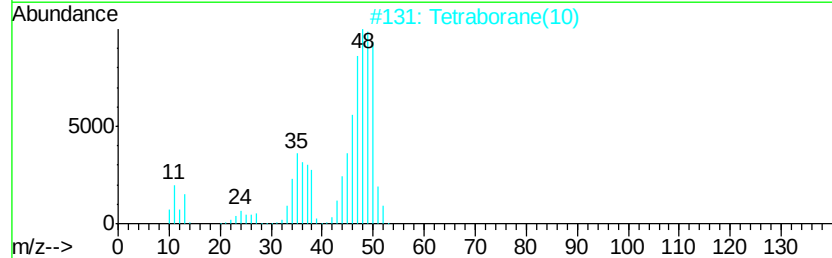
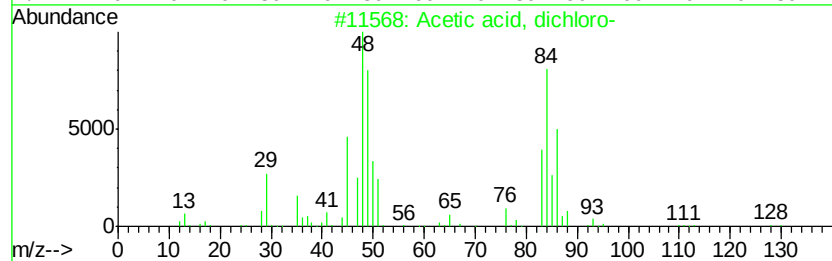
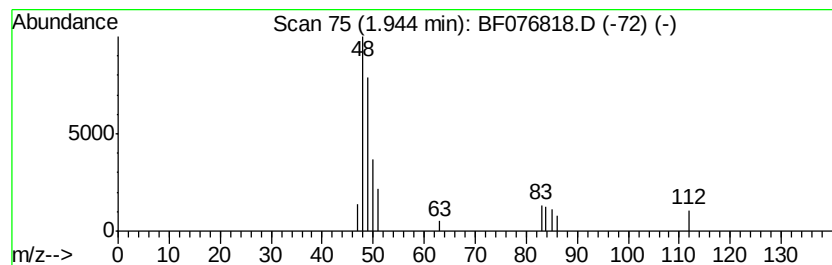
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 unknown1.94 Concentration Rank 16

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 1.94 | 3.29 ng | 34609 | 1,4-Dichlorobenzene-d4 | 8.08 |

| Hit# | of | Tentative ID           | MW  | MolForm   | CAS#        | Qual |
|------|----|------------------------|-----|-----------|-------------|------|
| 1    | 5  | Acetic acid, dichloro- | 128 | C2H2Cl2O2 | 000079-43-6 | 43   |
| 2    |    | Tetraborane(10)        | 54  | B4H10     | 018283-93-7 | 9    |
| 3    |    | Methylene Chloride     | 84  | CH2Cl2    | 000075-09-2 | 9    |
| 4    |    | Phosphine, methyl-     | 48  | CH5P      | 000593-54-4 | 5    |
| 5    |    | Ethyl Chloride         | 64  | C2H5Cl    | 000075-00-3 | 4    |



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Misc : W  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF010915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

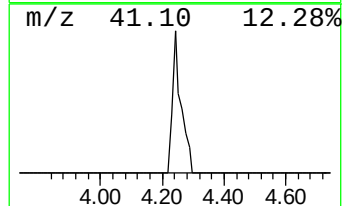
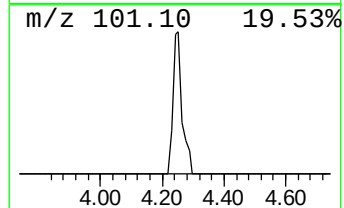
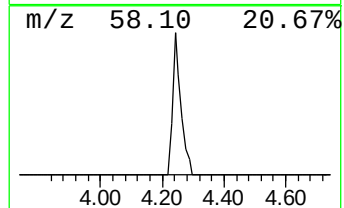
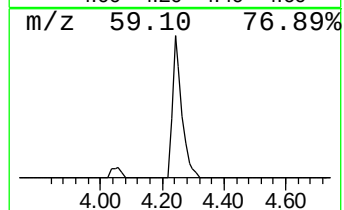
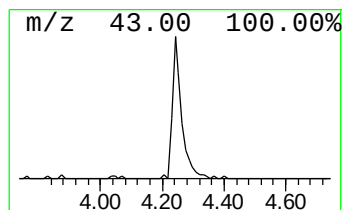
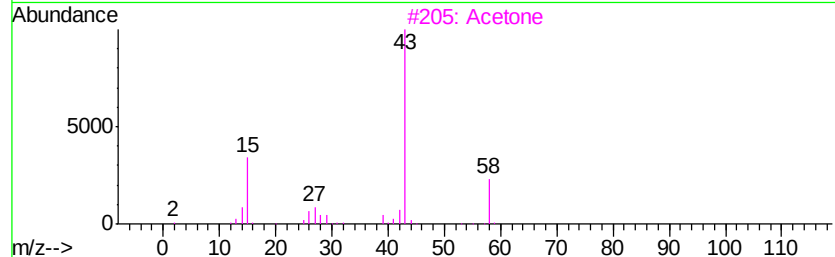
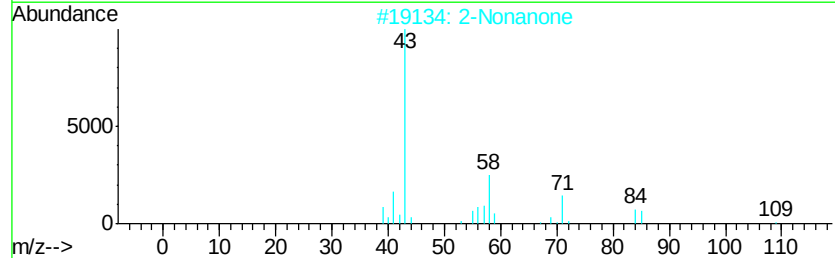
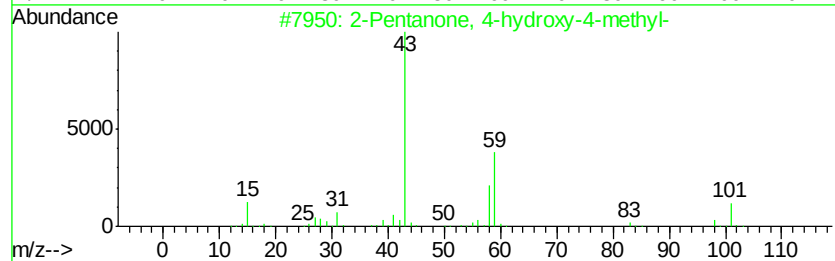
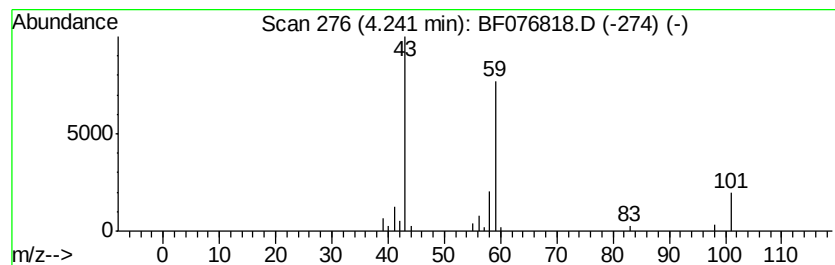
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 4.24 | 8.00 ng | 84091 | 1,4-Dichlorobenzene-d4 | 8.08 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 56   |
| 2    |    | 2-Nonanone                       | 142 | C9H18O  | 000821-55-6 | 10   |
| 3    |    | Acetone                          | 58  | C3H6O   | 000067-64-1 | 9    |
| 4    |    | Hexanamide                       | 115 | C6H13NO | 000628-02-4 | 9    |
| 5    |    | 3-Hexanol, 4-methyl-             | 116 | C7H16O  | 000615-29-2 | 9    |



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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF010915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

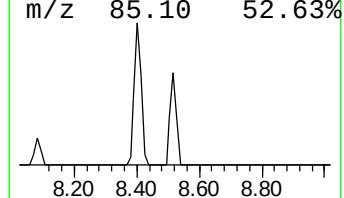
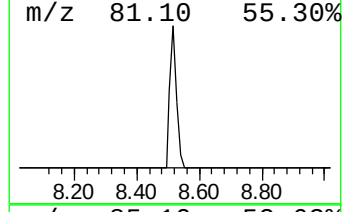
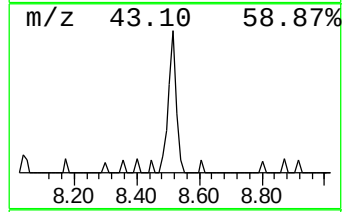
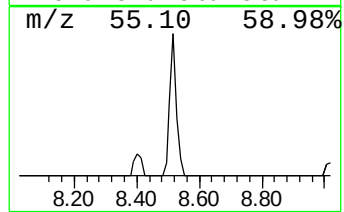
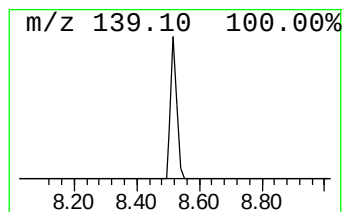
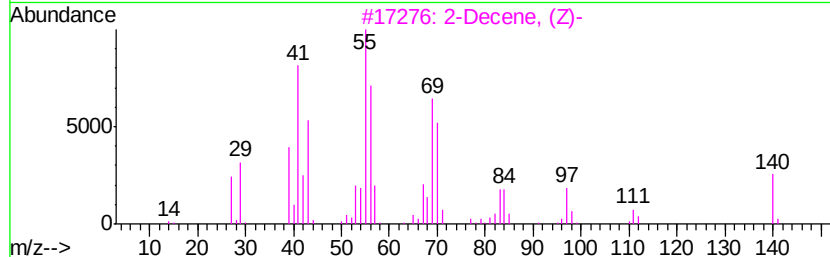
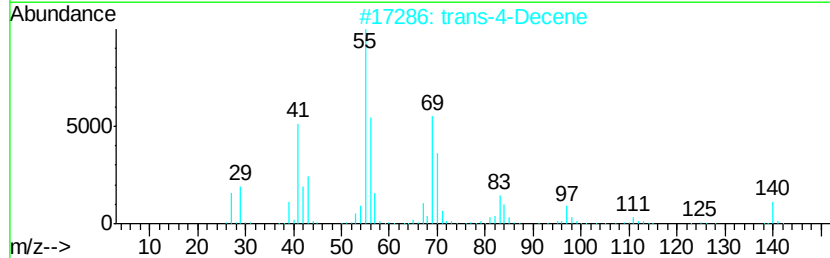
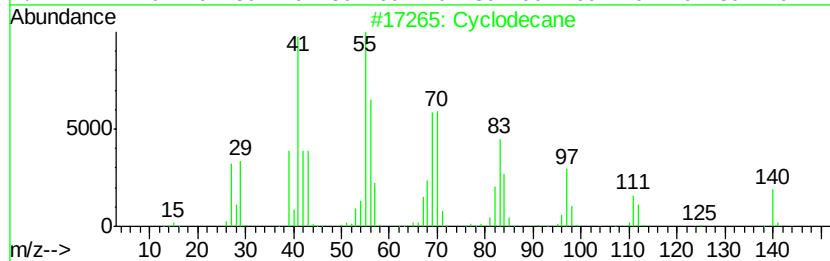
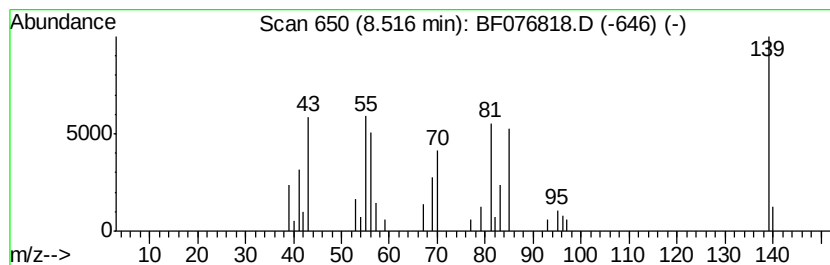
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 unknown8.52 Concentration Rank 6

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 8.52 | 6.31 ng | 66328 | 1,4-Dichlorobenzene-d4 | 8.08 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclodecane                         | 140 | C10H20  | 000293-96-9 | 30   |
| 2    |    | trans-4-Decene                      | 140 | C10H20  | 019398-89-1 | 27   |
| 3    |    | 2-Decene, (Z)-                      | 140 | C10H20  | 020348-51-0 | 22   |
| 4    |    | 1-Octene, 3,7-dimethyl-             | 140 | C10H20  | 004984-01-4 | 22   |
| 5    |    | 2-Propionyl-1,4,5,6-tetrahydropy... | 139 | C8H13NO | 080933-74-0 | 22   |



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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

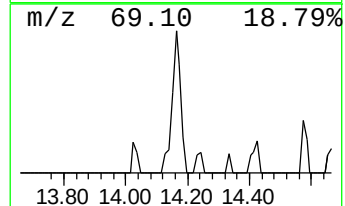
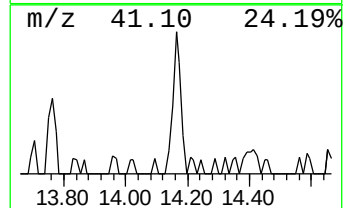
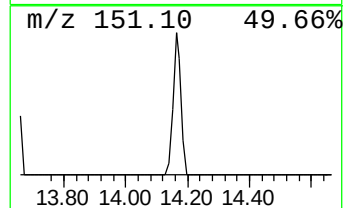
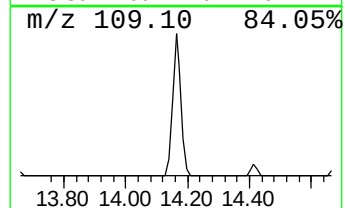
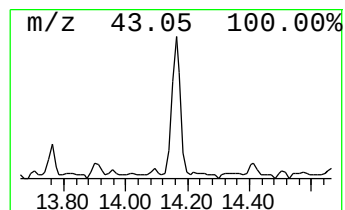
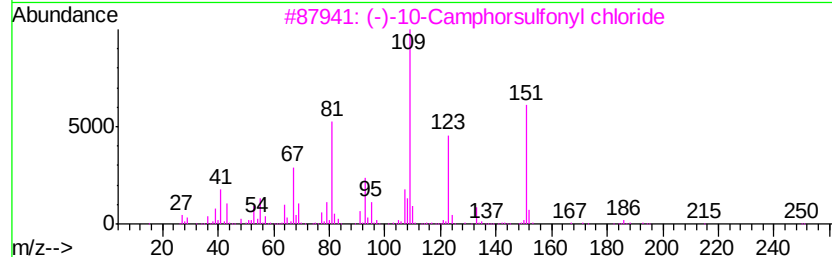
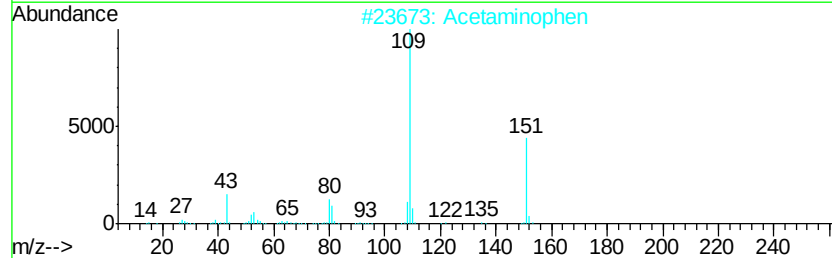
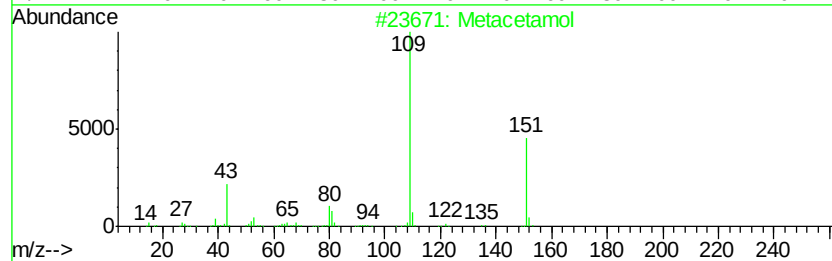
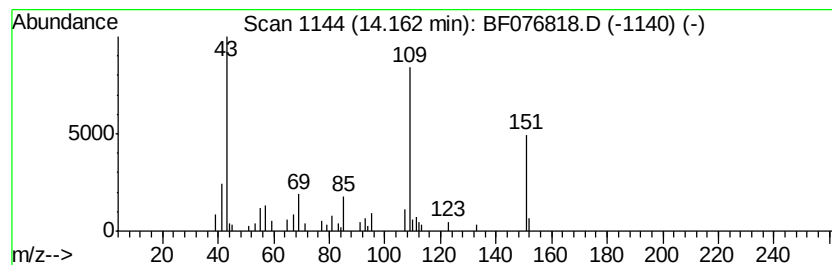
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Metacetamol Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.16 | 6.25 ng | 111799 | Acenaphthene-d10 | 15.12 |

| Hit# of | 5 | Tentative ID                    | MW  | MolForm     | CAS#        | Qual |
|---------|---|---------------------------------|-----|-------------|-------------|------|
| 1       |   | Metacetamol                     | 151 | C8H9NO2     | 000621-42-1 | 58   |
| 2       |   | Acetaminophen                   | 151 | C8H9NO2     | 000103-90-2 | 53   |
| 3       |   | (-)-10-Camphorsulfonyl chloride | 250 | C10H15ClO3S | 039262-22-1 | 50   |
| 4       |   | Pyridine, 2-butyl-, 1-oxide     | 151 | C9H13NO     | 031396-32-4 | 36   |
| 5       |   | Pyridine, 3-butyl-, 1-oxide     | 151 | C9H13NO     | 031396-33-5 | 36   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011215\  
Data File : BF076818.D  
Acq On : 12 Jan 2015 20:40  
Operator : IZ / TP  
Sample : G1044-10  
Misc : W  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SP-10

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

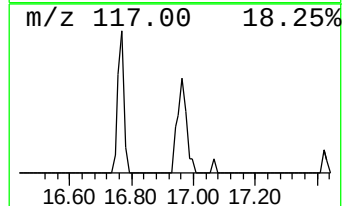
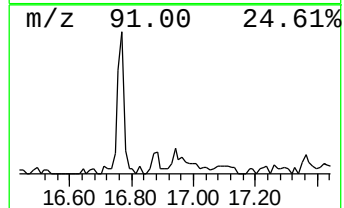
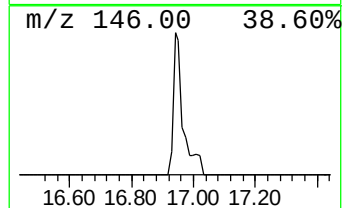
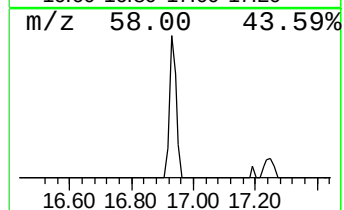
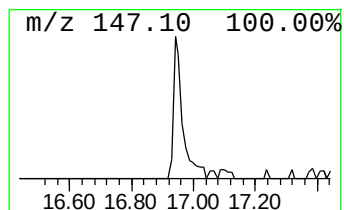
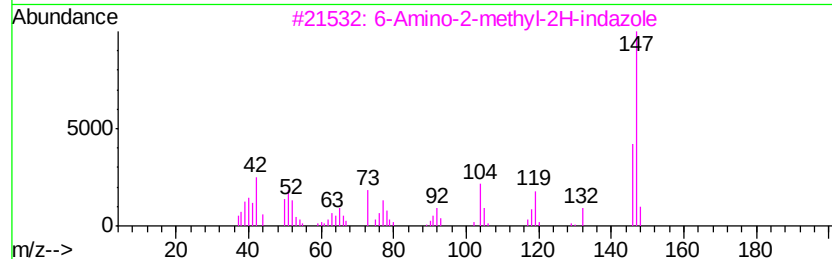
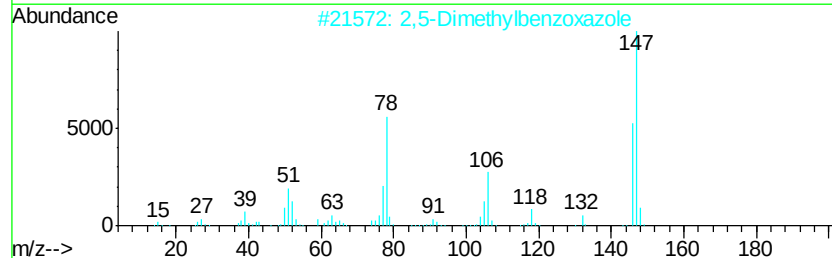
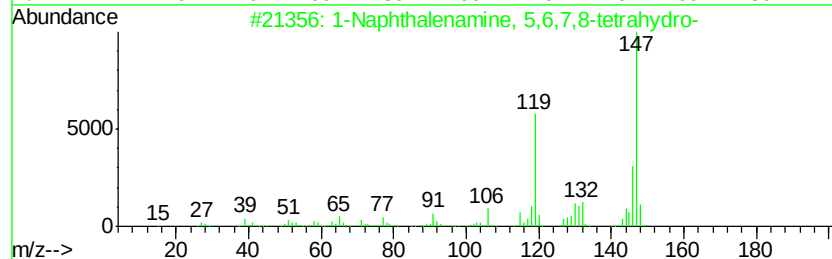
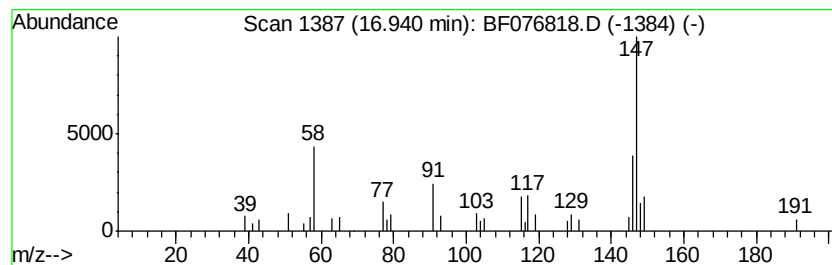
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 1-Naphthalenamine, 5,6,7,8-... Concentration Rank 15

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 16.94 | 3.31 ng | 56287 | Phenanthrene-d10 | 17.85 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 1-Naphthalenamine, 5,6,7,8-tetra... | 147 | C10H13N  | 002217-41-6  | 52   |
| 2    |    |   | 2,5-Dimethylbenzoxazole             | 147 | C9H9NO   | 005676-58-4  | 49   |
| 3    |    |   | 6-Amino-2-methyl-2H-indazole        | 147 | C8H9N3   | 050593-30-1  | 47   |
| 4    |    |   | 1,3,2-Benzoxazaborole, 2-ethyl-2... | 147 | C8H10BNO | 129363-62-8  | 40   |
| 5    |    |   | Indane, 2-methoxy-1-(2-methylallyl) | 202 | C14H18O  | 1000196-88-9 | 38   |





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Data File : BF076818.D  
Acq On : 12 Jan 2015 20:40  
Operator : IZ / TP  
Sample : G1044-10  
Misc : W  
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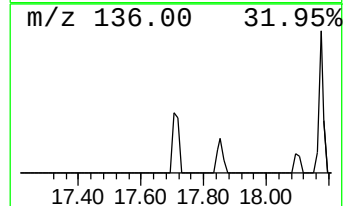
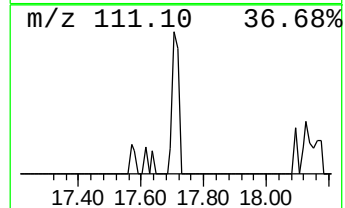
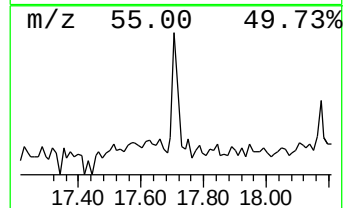
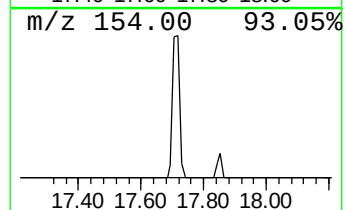
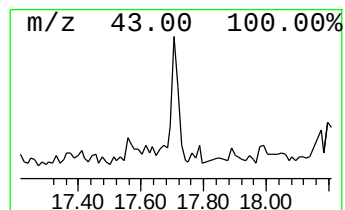
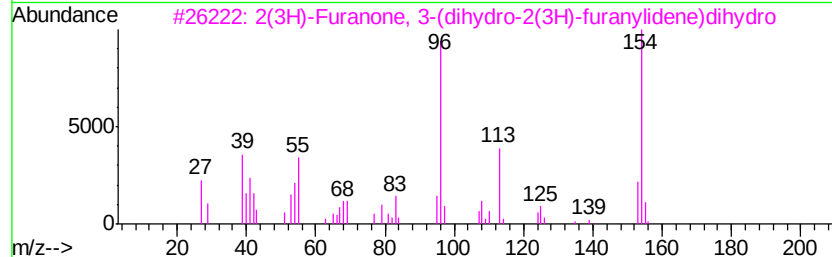
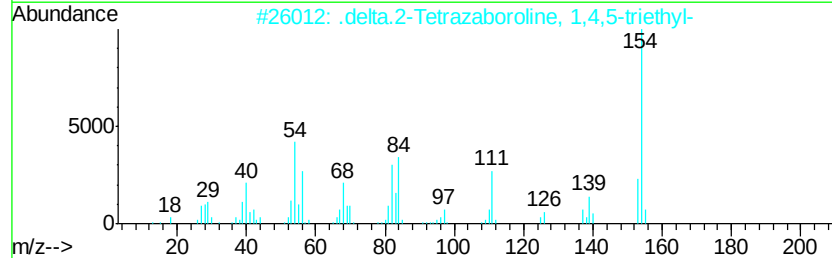
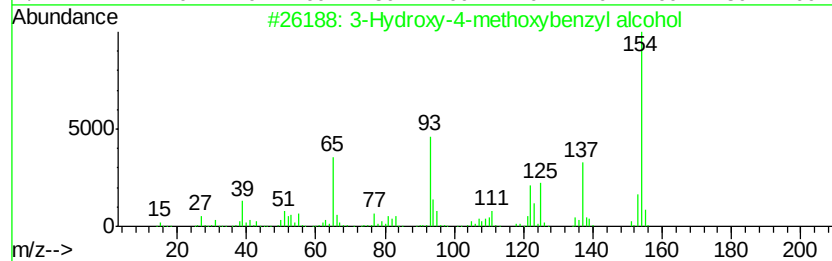
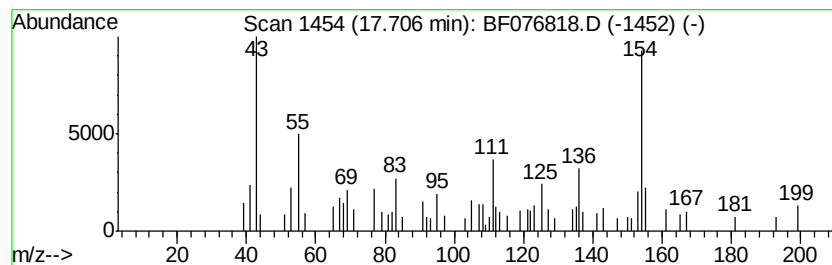
Instrument :  
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ClientSampleId :  
SP-10

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 unknown17.71 Concentration Rank 13

| R.T.      | EstConc                              | Area  | Relative to ISTD | R.T.        |      |
|-----------|--------------------------------------|-------|------------------|-------------|------|
| 17.71     | 4.04 ng                              | 68664 | Phenanthrene-d10 | 17.85       |      |
| Hit# of 5 | Tentative ID                         | MW    | MolForm          | CAS#        | Qual |
| 1         | 3-Hydroxy-4-methoxybenzyl alcohol    | 154   | C8H10O3          | 004383-06-6 | 43   |
| 2         | .delta.2-Tetrazaboroline, 1,4,5-...  | 154   | C6H15BN4         | 020534-04-7 | 38   |
| 3         | 2(3H)-Furanone, 3-(dihydro-2(3H))... | 154   | C8H10O3          | 055164-40-4 | 37   |
| 4         | 3-Hepten-2-one, 3-propyl-            | 154   | C10H18O          | 032064-69-0 | 32   |
| 5         | cis-1,2-Bis(acetamidomethyl)cycl...  | 226   | C12H22N2O2       | 070924-84-4 | 32   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011215\  
Data File : BF076818.D  
Acq On : 12 Jan 2015 20:40  
Operator : IZ / TP  
Sample : G1044-10  
Misc : W  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SP-10

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

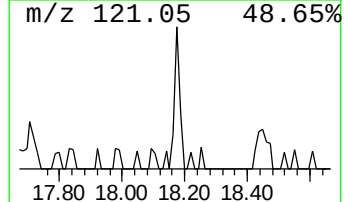
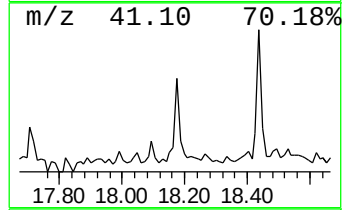
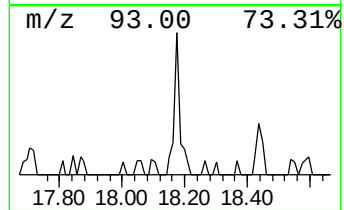
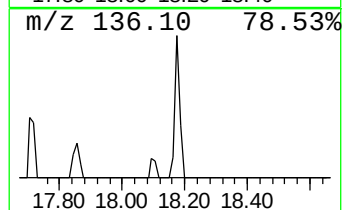
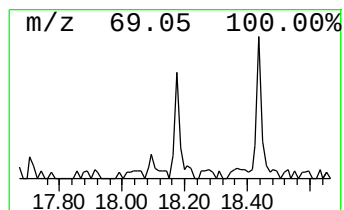
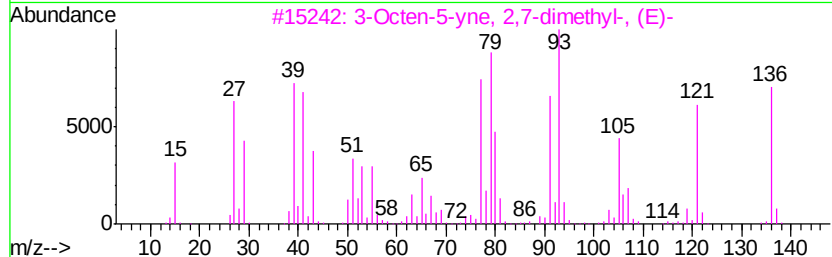
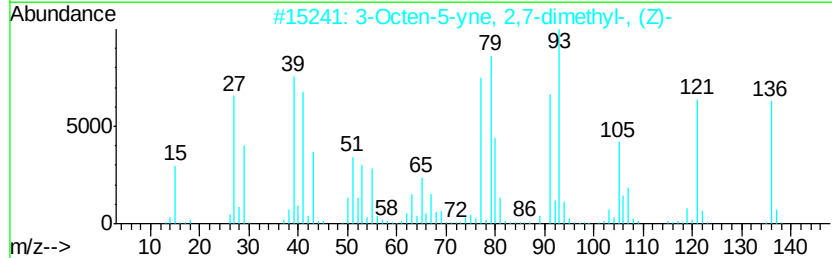
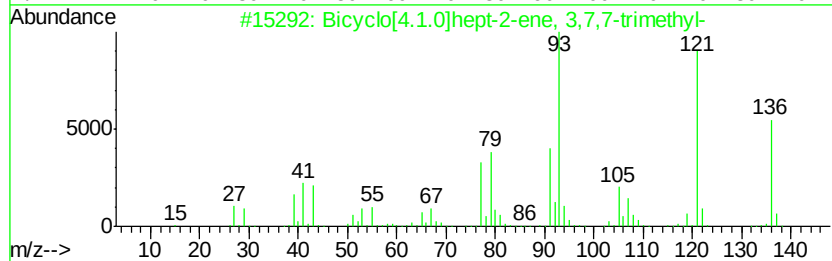
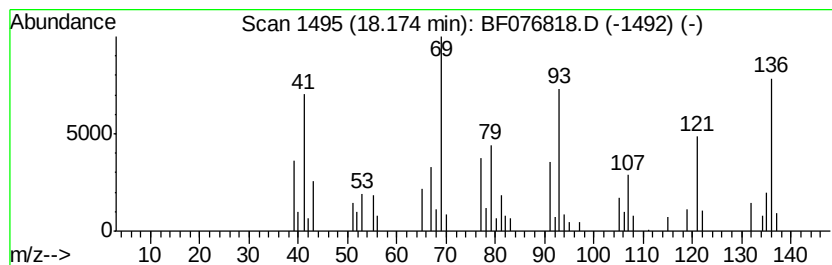
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TIC Integration Parameters: LSCINT.P

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Peak Number 10 Bicyclo[4.1.0]hept-2-ene, 3... Concentration Rank 14

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 18.17 | 3.33 ng | 56542 | Phenanthrene-d10 | 17.85 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Bicyclo[4.1.0]hept-2-ene, 3,7,7-... | 136 | C10H16  | 000554-61-0 | 91   |
| 2    |    | 3-Octen-5-yne, 2,7-dimethyl-, (Z)-  | 136 | C10H16  | 028935-76-4 | 60   |
| 3    |    | 3-Octen-5-yne, 2,7-dimethyl-, (E)-  | 136 | C10H16  | 055956-33-7 | 60   |
| 4    |    | (+)-4-Carene                        | 136 | C10H16  | 029050-33-7 | 60   |
| 5    |    | 1,4-Cyclohexadiene, 1-methyl-4-(... | 136 | C10H16  | 000099-85-4 | 55   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF011215\  
Data File : BF076818.D  
Acq On : 12 Jan 2015 20:40  
Operator : IZ / TP  
Sample : G1044-10  
Misc : W  
ALS Vial : 14 Sample Multiplier: 1

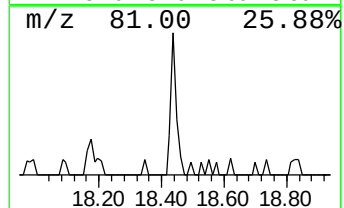
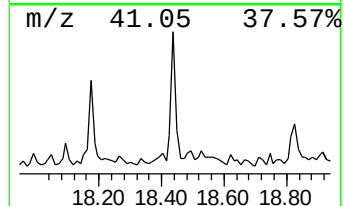
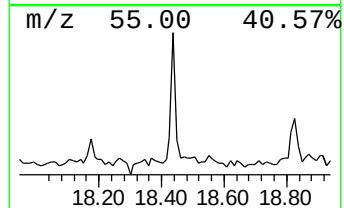
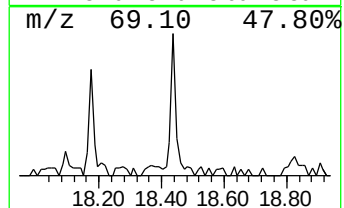
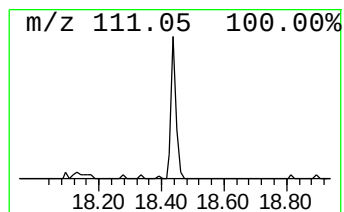
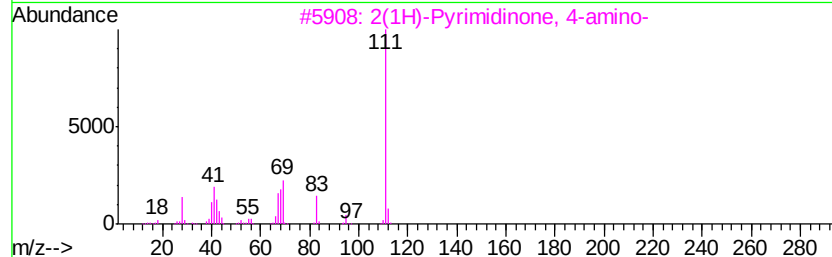
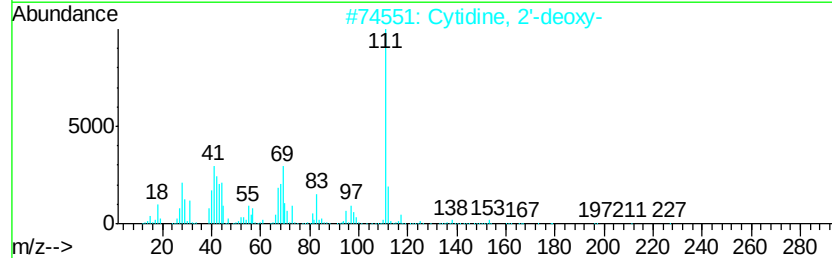
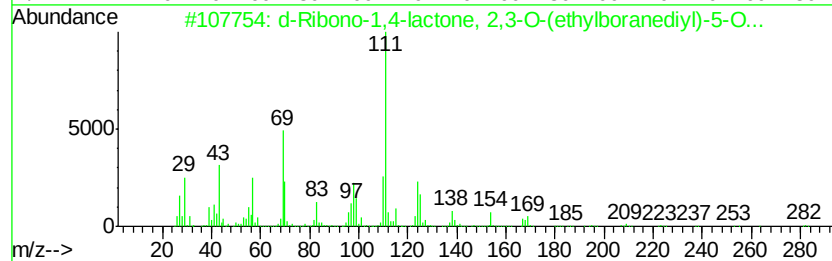
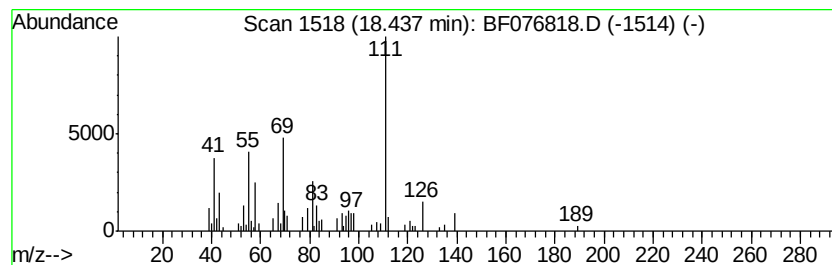
Instrument :  
BNA\_F  
ClientSampleId :  
SP-10

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF010915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 11 unknown18.44 Concentration Rank 10

| R.T.      | EstConc                             | Area  | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|-------|------------------|--------------|------|
| 18.44     | 5.35 ng                             | 90914 | Phenanthrene-d10 | 17.85        |      |
| Hit# of 5 | Tentative ID                        | MW    | MolForm          | CAS#         | Qual |
| 1         | d-Ribono-1,4-lactone, 2,3-O-(eth... | 282   | C9H10BF306       | 1000150-02-0 | 40   |
| 2         | Cytidine, 2'-deoxy-                 | 227   | C9H13N304        | 000951-77-9  | 38   |
| 3         | 2(1H)-Pvrimidinone, 4-amino-        | 111   | C4H5N30          | 000071-30-7  | 38   |
| 4         | Cyclohexane, 1,4-dimethyl-2-octa... | 364   | C26H52           | 055282-02-5  | 38   |
| 5         | 2-Thiophenecarboxylic acid, 2-tr... | 310   | C18H3002S        | 1000280-65-9 | 38   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011215\  
Data File : BF076818.D  
Acq On : 12 Jan 2015 20:40  
Operator : IZ / TP  
Sample : G1044-10  
Misc : W  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

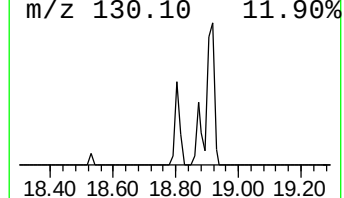
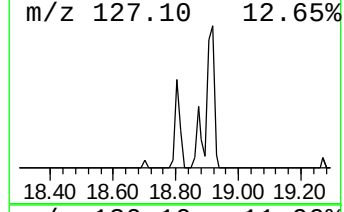
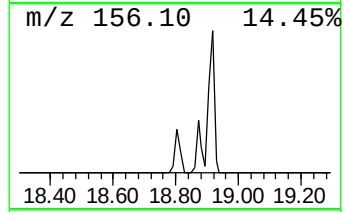
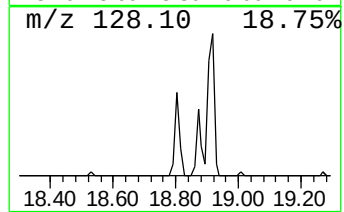
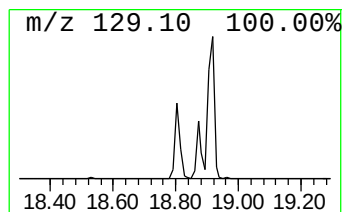
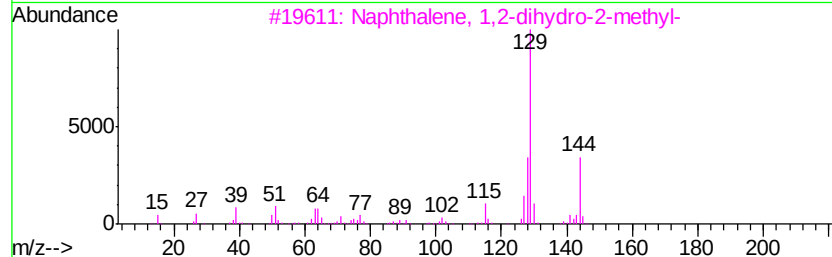
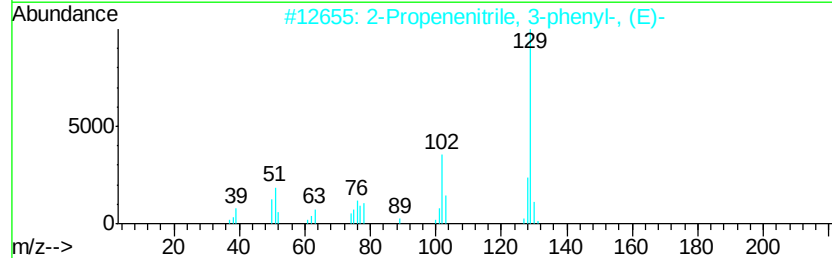
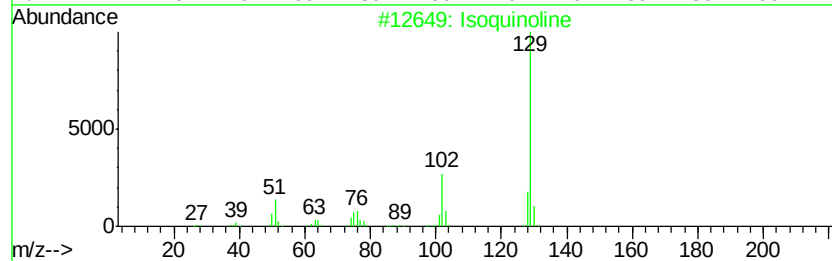
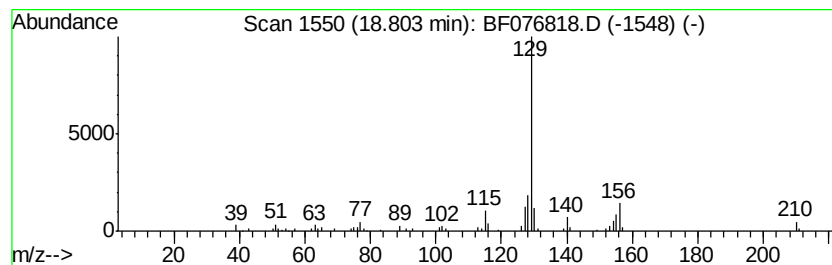
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Isoquinoline Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.80 | 7.50 ng | 127509 | Phenanthrene-d10 | 17.85 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Isoquinoline                       | 129 | C9H7N   | 000119-65-3 | 52   |
| 2    |    | 2-Propenenitrile, 3-phenyl-, (E)-  | 129 | C9H7N   | 001885-38-7 | 42   |
| 3    |    | Naphthalene, 1,2-dihydro-2-methyl- | 144 | C11H12  | 021564-79-4 | 42   |
| 4    |    | Benzene, 1,3-hexadienyl-           | 158 | C12H14  | 041635-77-2 | 42   |
| 5    |    | Quinoline                          | 129 | C9H7N   | 000091-22-5 | 40   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011215\  
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Sample : G1044-10  
Misc : W  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SP-10

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

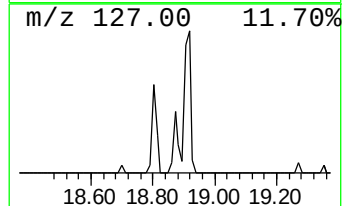
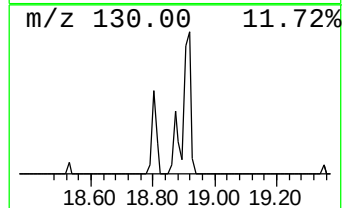
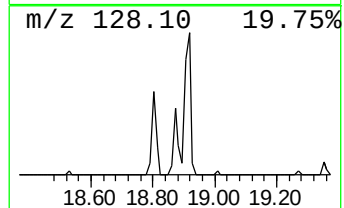
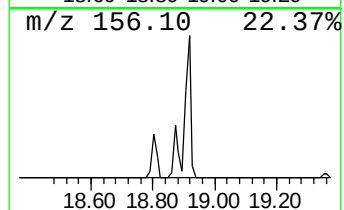
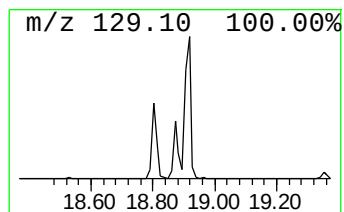
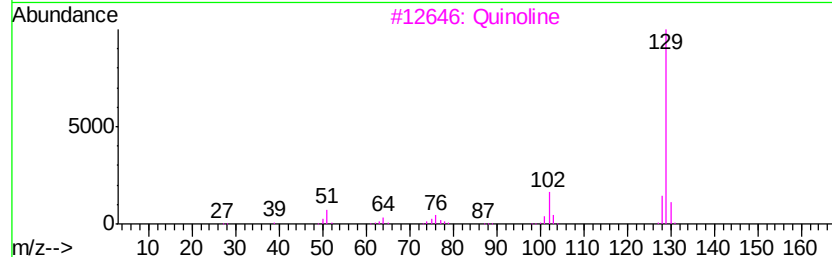
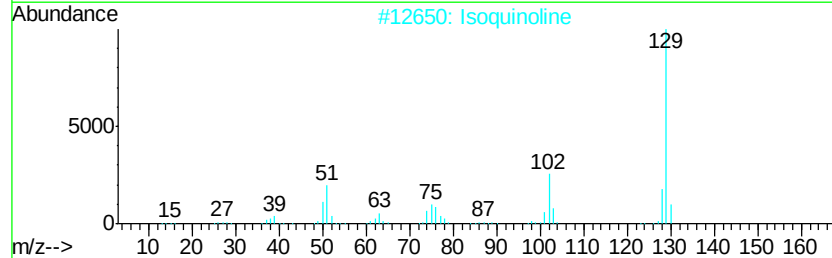
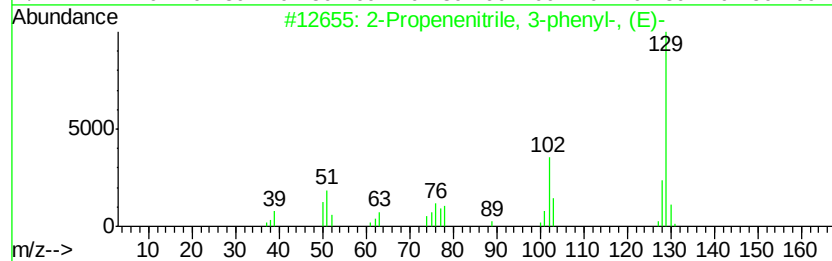
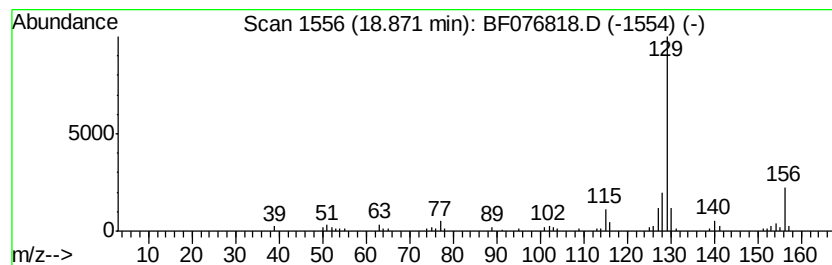
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 2-Propenenitrile, 3-phenyl-... Concentration Rank 9

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 18.87 | 5.65 ng | 95947 | Phenanthrene-d10 | 17.85 |

| Hit# of 5 | Tentative ID                      | MW  | MolForm | CAS#         | Qual |
|-----------|-----------------------------------|-----|---------|--------------|------|
| 1         | 2-Propenenitrile, 3-phenyl-, (E)- | 129 | C9H7N   | 001885-38-7  | 50   |
| 2         | Isoquinoline                      | 129 | C9H7N   | 000119-65-3  | 50   |
| 3         | Quinoline                         | 129 | C9H7N   | 000091-22-5  | 49   |
| 4         | 1,2-Benzenediacetonitrile         | 156 | C10H8N2 | 000613-73-0  | 39   |
| 5         | 4,5-Dimethylthiazole S-oxide      | 129 | C5H7NOS | 1000112-53-4 | 38   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011215\  
Data File : BF076818.D  
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Operator : IZ / TP  
Sample : G1044-10  
Misc : W  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SP-10

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

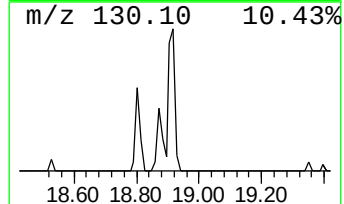
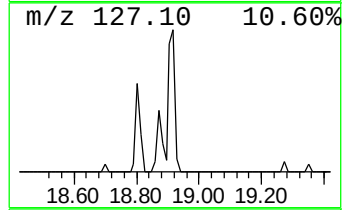
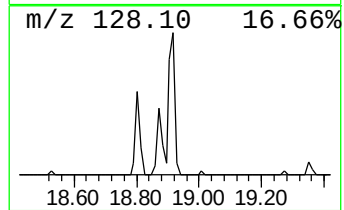
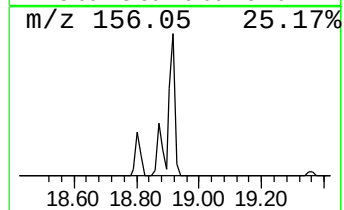
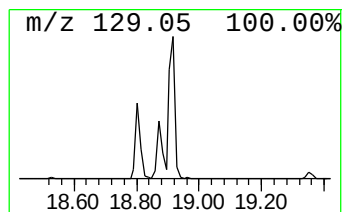
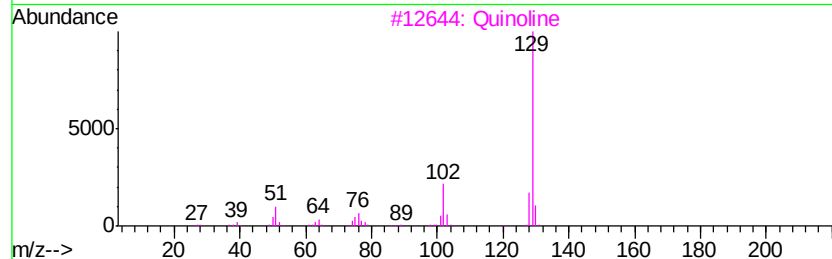
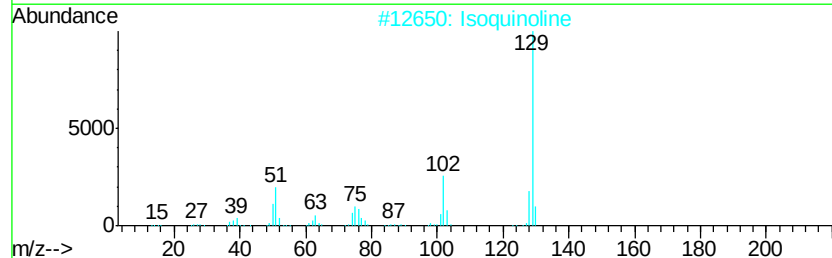
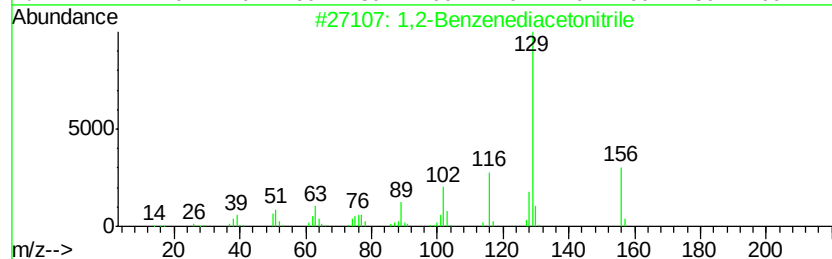
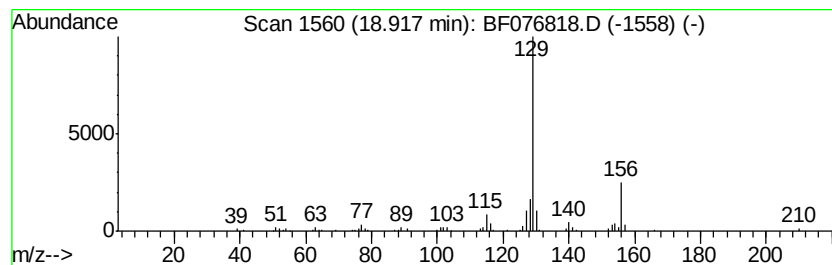
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 1,2-Benzenediacetonitrile Concentration Rank 3

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 18.92 | 13.24 ng | 224971 | Phenanthrene-d10 | 17.85 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | 1,2-Benzenediacetonitrile           | 156 | C10H8N2    | 000613-73-0 | 59   |
| 2    |    | Isoquinoline                        | 129 | C9H7N      | 000119-65-3 | 50   |
| 3    |    | Quinoline                           | 129 | C9H7N      | 000091-22-5 | 40   |
| 4    |    | 2-Propenenitrile, 3-phenyl-, (E)-   | 129 | C9H7N      | 001885-38-7 | 40   |
| 5    |    | Benzeneacetic acid, .alpha.-cyan... | 228 | C13H12N2O2 | 134882-04-5 | 39   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011215\  
Data File : BF076818.D  
Acq On : 12 Jan 2015 20:40  
Operator : IZ / TP  
Sample : G1044-10  
Misc : W  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SP-10

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

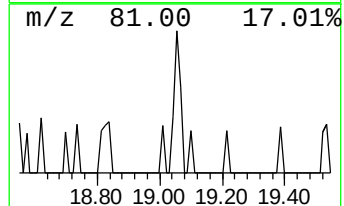
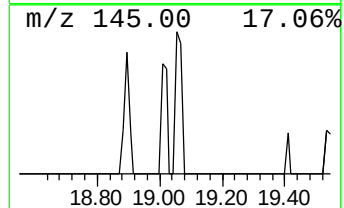
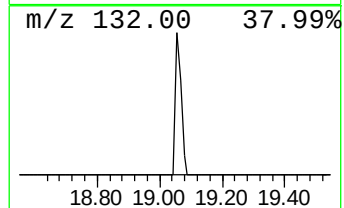
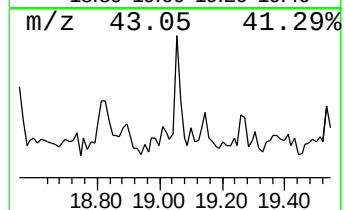
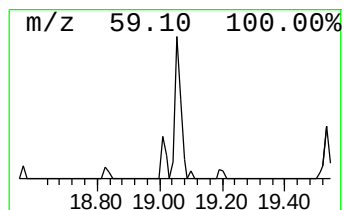
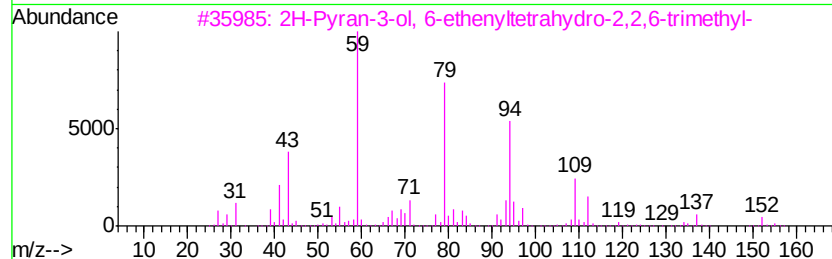
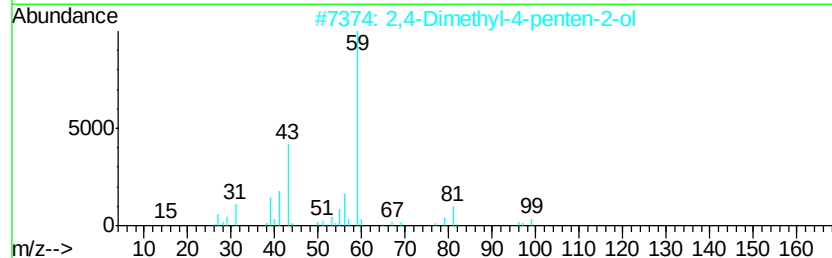
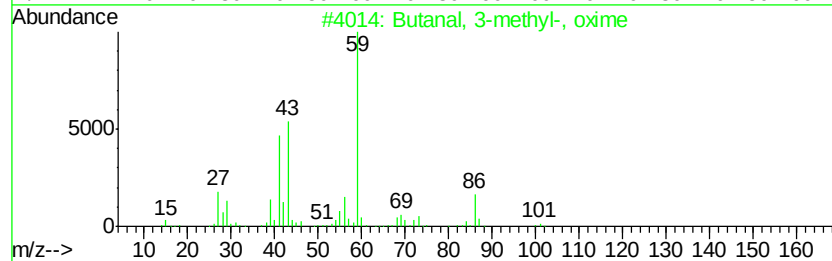
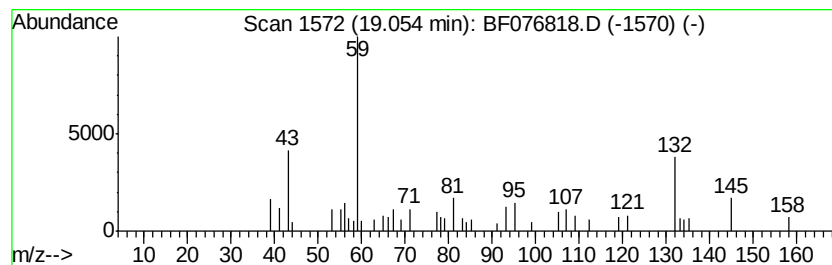
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 unknown19.05 Concentration Rank 18

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 19.05 | 2.53 ng | 42987 | Phenanthrene-d10 | 17.85 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|------------|-------------|------|
| 1    |    |   | Butanal, 3-methyl-, oxime           | 101 | C5H11NO    | 000626-90-4 | 32   |
| 2    |    |   | 2,4-Dimethyl-4-penten-2-ol          | 114 | C7H14O     | 019781-53-4 | 32   |
| 3    |    |   | 2H-Pyran-3-ol, 6-ethenvltetrahyd... | 170 | C10H18O2   | 014049-11-7 | 28   |
| 4    |    |   | 2-Octanol, 2-methyl-6-methylene-    | 156 | C10H20O    | 018479-59-9 | 25   |
| 5    |    |   | 2-(3,4-Dibromo-4-methylcyclohexy... | 312 | C10H18Br2O | 110202-13-6 | 25   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011215\  
Data File : BF076818.D  
Acq On : 12 Jan 2015 20:40  
Operator : IZ / TP  
Sample : G1044-10  
Misc : W  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SP-10

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

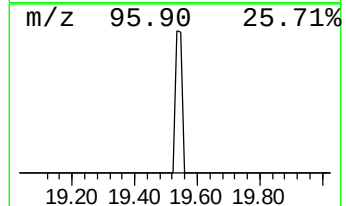
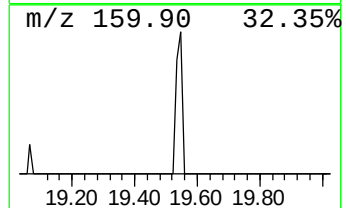
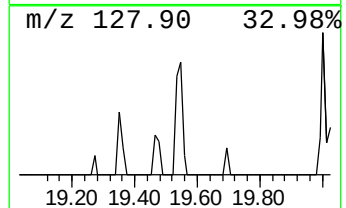
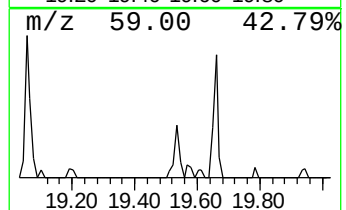
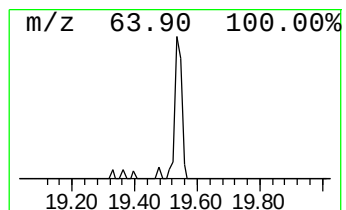
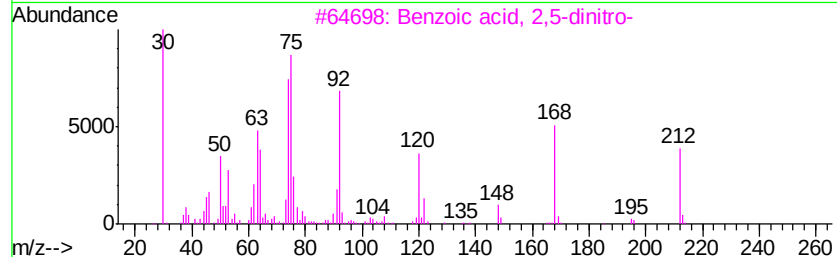
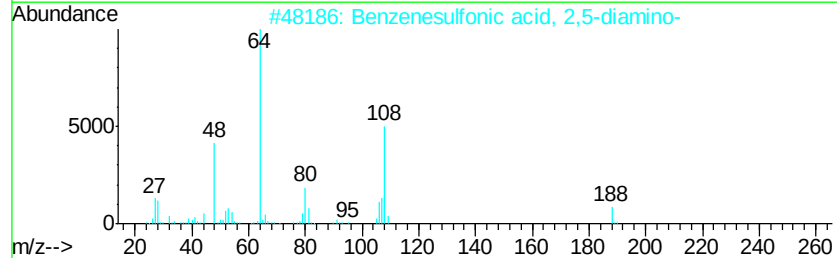
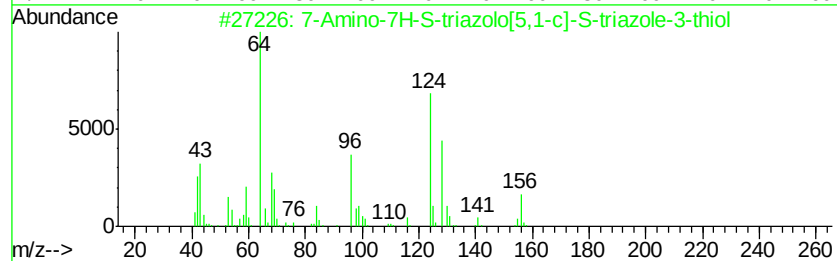
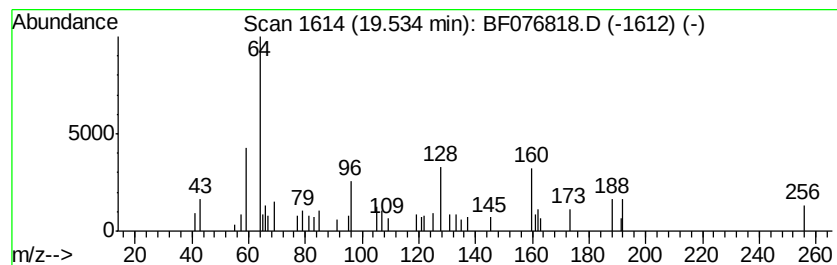
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 7-Amino-7H-S-triazolo[5,1-c... Concentration Rank 19

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 19.53 | 2.39 ng | 40554 | Phenanthrene-d10 | 17.85 |

| Hit# | of | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|------|----|-------------------------------------|-----|--------------|--------------|------|
| 1    | 5  | 7-Amino-7H-S-triazolo[5,1-c]-S-t... | 156 | C3H4N6S      | 013728-28-4  | 59   |
| 2    |    | Benzenesulfonic acid, 2,5-diamino-  | 188 | C6H8N2O3S    | 000088-45-9  | 53   |
| 3    |    | Benzoic acid, 2,5-dinitro-          | 212 | C7H4N2O6     | 000610-28-6  | 40   |
| 4    |    | Methanethiol, N-(3-phenoxypropyl... | 304 | C11H16N2O4S2 | 040283-89-4  | 36   |
| 5    |    | 1-[1-Bromo-2-(2,2,3,3-tetrafluor... | 328 | C12H10BrF5   | 1000223-18-3 | 36   |





Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF011215\  
Data File : BF076818.D  
Acq On : 12 Jan 2015 20:40  
Operator : IZ / TP  
Sample : G1044-10  
Misc : W  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SP-10

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF010915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

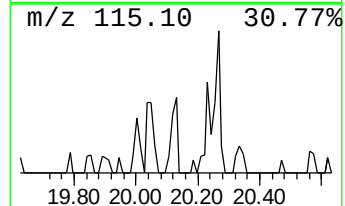
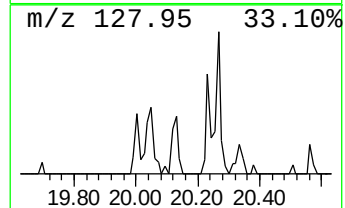
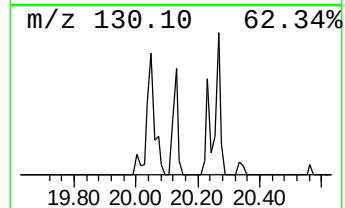
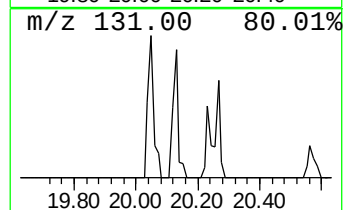
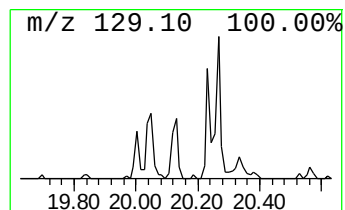
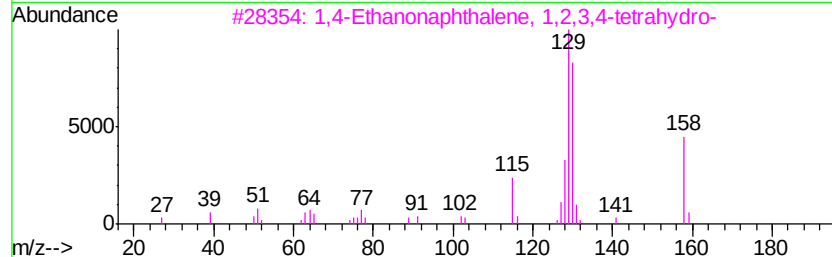
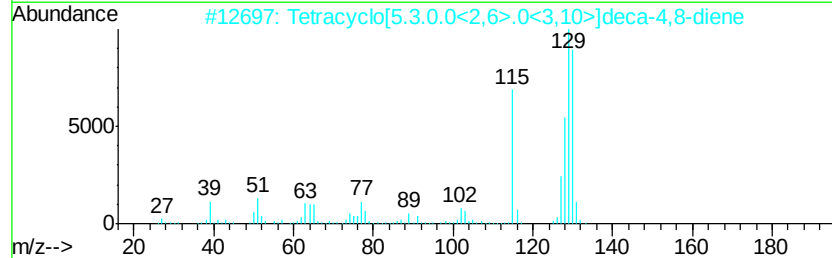
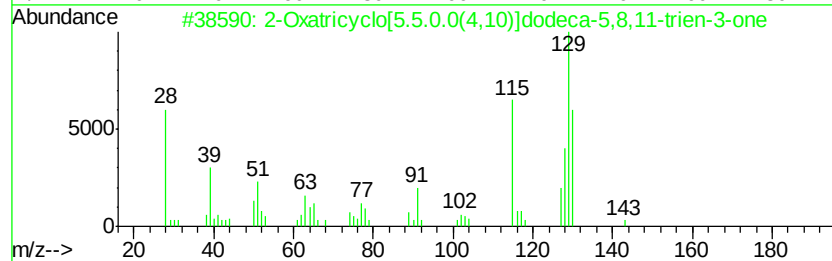
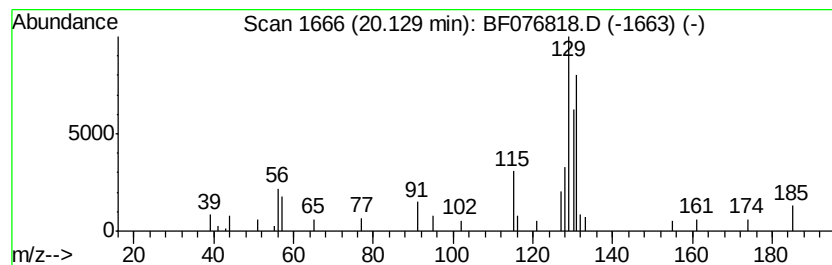
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 unknown20.13 Concentration Rank 20

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 20.13 | 2.27 ng | 38828 | Chrysene-d12     | 21.34 |

| Hit# | of | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Oxatricyclo[5.5.0.0(4,10)]dodec... | 174 | C11H10O2 | 056909-26-3 | 38   |
| 2    |    | Tetracyclo[5.3.0.0<2,6>.0<3,10>]...  | 130 | C10H10   | 034324-40-8 | 38   |
| 3    |    | 1,4-Ethanonaphthalene, 1,2,3,4-t...  | 158 | C12H14   | 004175-52-4 | 37   |
| 4    |    | 1H-Indole-3-ethanamine, N-methyl-    | 174 | C11H14N2 | 000061-49-4 | 35   |
| 5    |    | Benzene, (cyclopropylidenemethyl)-   | 130 | C10H10   | 007555-67-1 | 35   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF011215\  
Data File : BF076818.D  
Acq On : 12 Jan 2015 20:40  
Operator : IZ / TP  
Sample : G1044-10  
Misc : W  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SP-10

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF010915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

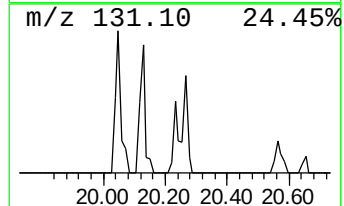
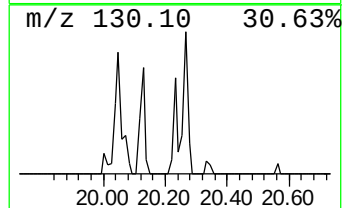
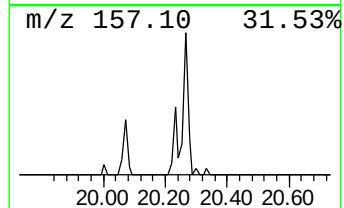
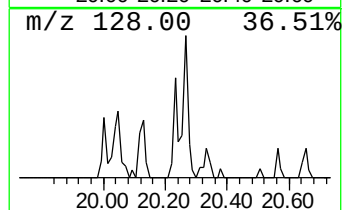
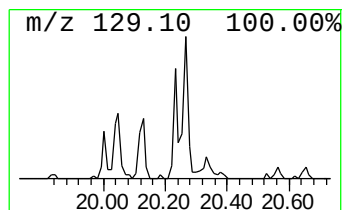
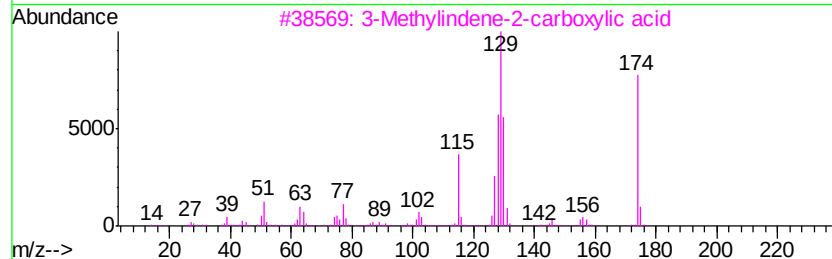
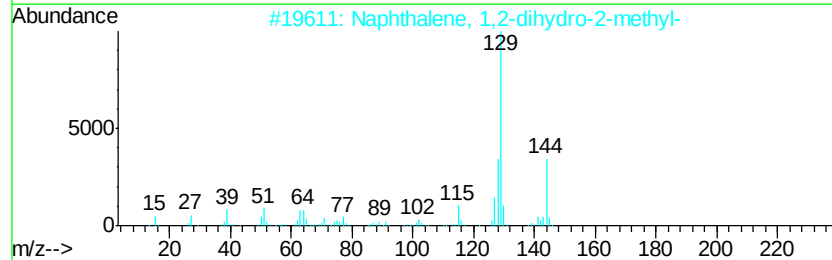
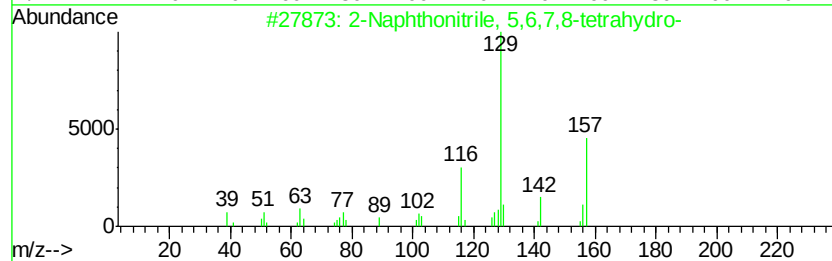
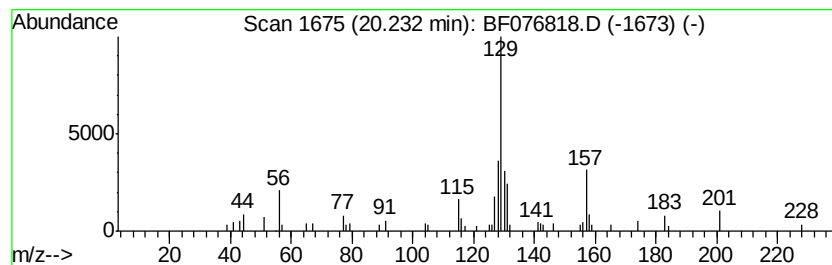
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 unknown20.23 Concentration Rank 17

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 20.23 | 2.83 ng | 48450 | Chrysene-d12     | 21.34 |

| Hit# | of                                                      | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|---------------------------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 2-Naphthonitrile, 5,6,7,8-tetrahydro-                   | 157 | C11H11N      | 017104-67-5  | 43      |      |      |
| 2    | Naphthalene, 1,2-dihydro-2-methyl-                      | 144 | C11H12       | 021564-79-4  | 43      |      |      |
| 3    | 3-Methylindene-2-carboxylic acid                        | 174 | C11H10O2     | 034225-81-5  | 38      |      |      |
| 4    | Hexa-2,4-dienylbenzene                                  | 158 | C12H14       | 079482-86-3  | 37      |      |      |
| 5    | Dimethyl 1,2-diethenyltricyclo[3.3.1.0 <sup>2,5</sup> ] | 248 | C14H16O4     | 1000222-75-3 | 37      |      |      |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF011215\  
Data File : BF076818.D  
Acq On : 12 Jan 2015 20:40  
Operator : IZ / TP  
Sample : G1044-10  
Misc : W  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SP-10

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF010915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

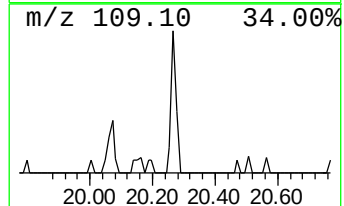
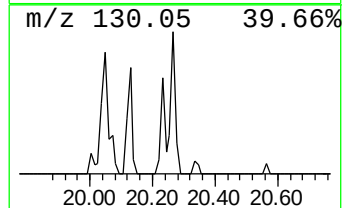
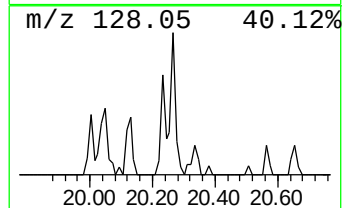
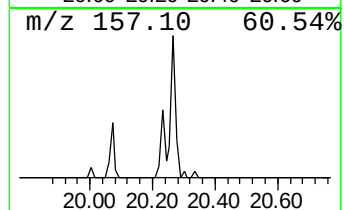
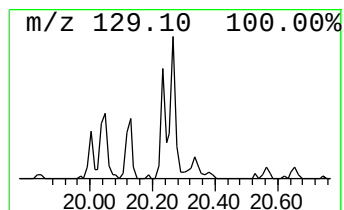
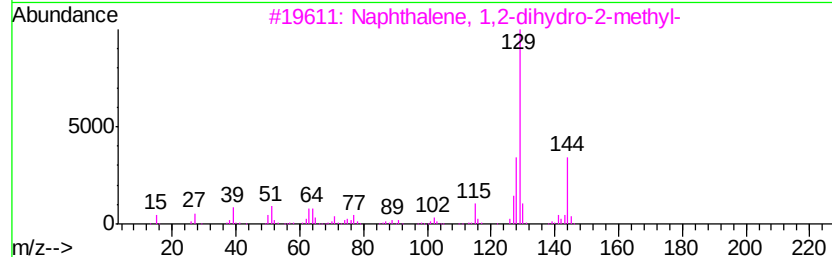
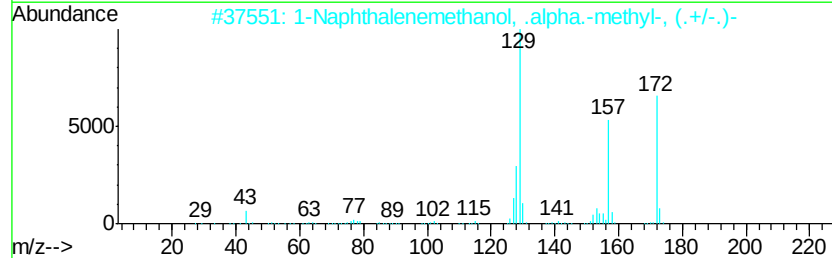
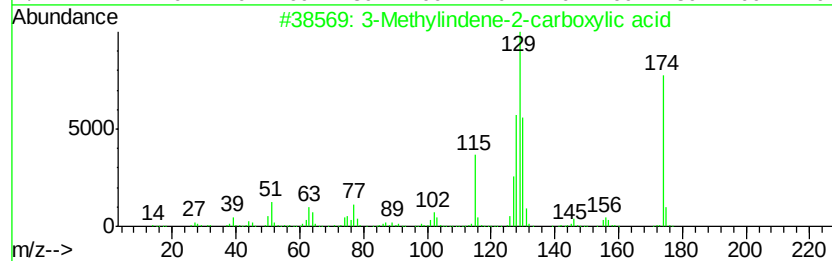
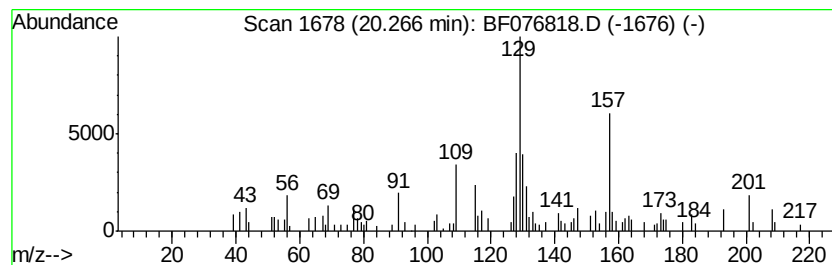
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 unknown20.27 Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 20.27 | 6.25 ng | 106825 | Chrysene-d12     | 21.34 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 3-Methylindene-2-carboxylic acid    | 174 | C11H10O2 | 034225-81-5 | 47   |
| 2    |    | 1-Naphthalenemethanol, .alpha.-m... | 172 | C12H12O  | 057605-95-5 | 47   |
| 3    |    | Naphthalene, 1,2-dihydro-2-methyl-  | 144 | C11H12   | 021564-79-4 | 35   |
| 4    |    | Benzene, 1,3-hexadienyl-            | 158 | C12H14   | 041635-77-2 | 35   |
| 5    |    | Hexa-2,4-dienylbenzene              | 158 | C12H14   | 079482-86-3 | 35   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011215\  
Data File : BF076818.D  
Acq On : 12 Jan 2015 20:40  
Operator : IZ / TP  
Sample : G1044-10  
Misc : W  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SP-10

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

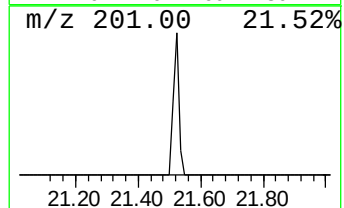
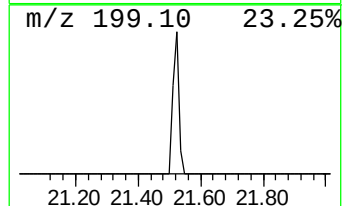
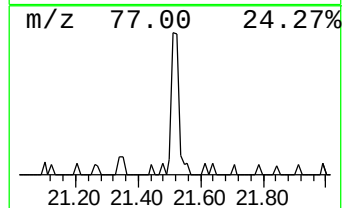
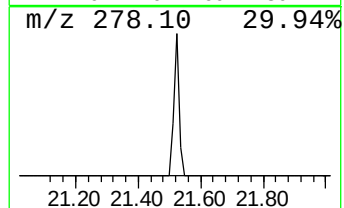
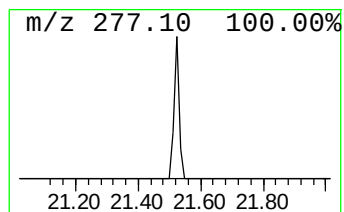
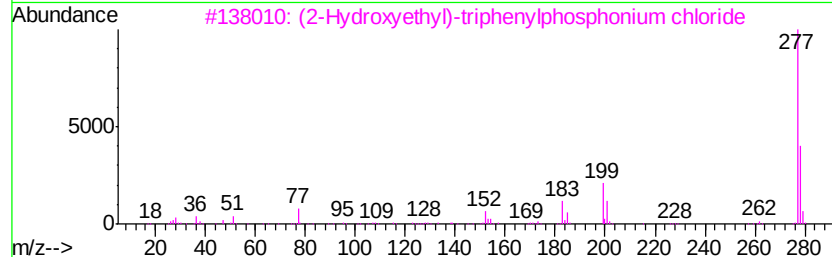
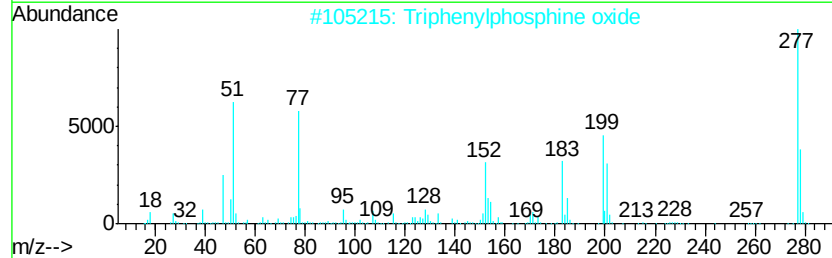
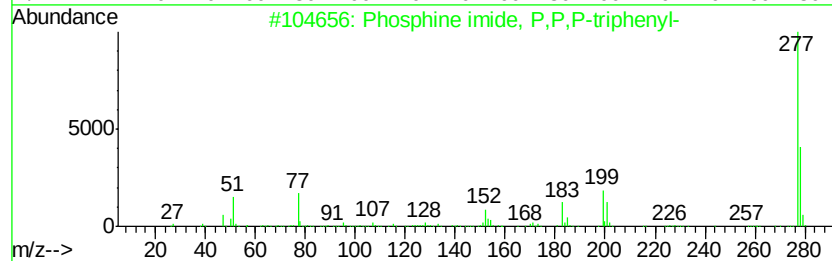
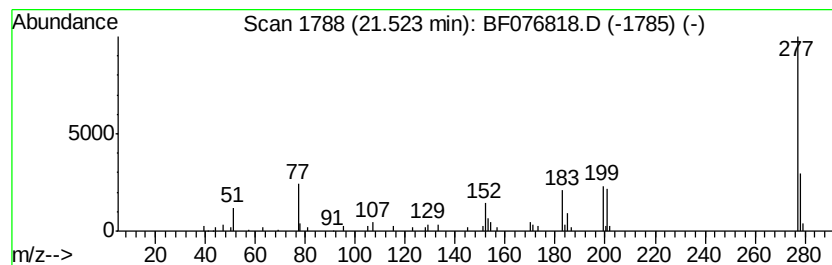
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 Phosphine imide, P,P,P-trip... Concentration Rank 11

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 21.52 | 4.70 ng | 80313 | Chrysene-d12     | 21.34 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm    | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|------------|-------------|------|
| 1    |    |   | Phosphine imide, P,P,P-triphenyl-    | 277 | C18H16NP   | 002240-47-3 | 68   |
| 2    |    |   | Triphenylphosphine oxide             | 278 | C18H15OP   | 000791-28-6 | 64   |
| 3    |    |   | (2-Hydroxyethyl)-triphenylphosph...  | 342 | C20H20ClOP | 023250-03-5 | 50   |
| 4    |    |   | Cobalt, (.eta.<5>-2,4-cyclopentad... | 277 | C16H15BCo  | 036682-12-9 | 43   |
| 5    |    |   | Formaldehyde, (triphenylphosphor...  | 304 | C19H17N2P  | 015990-54-2 | 42   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF011215\  
 Data File : BF076818.D  
 Acq On : 12 Jan 2015 20:40  
 Operator : IZ / TP  
 Sample : G1044-10  
 Misc : W  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SP-10

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF010915.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 |
| unknown1.94          | 1.94  | 3.3     | ng    | 34609    | 1                     | 8.08  | 210192 | 20.0 |
| 2-Pentanone, 4-hy... | 4.24  | 8.0     | ng    | 84091    | 1                     | 8.08  | 210192 | 20.0 |
| unknown8.52          | 8.52  | 6.3     | ng    | 66328    | 1                     | 8.08  | 210192 | 20.0 |
| Metacetamol          | 14.16 | 6.3     | ng    | 111799   | 3                     | 15.12 | 357527 | 20.0 |
| 1-Naphthalenamine... | 16.94 | 3.3     | ng    | 56287    | 4                     | 17.85 | 339828 | 20.0 |
| unknown17.71         | 17.71 | 4.0     | ng    | 68664    | 4                     | 17.85 | 339828 | 20.0 |
| Bicyclo[4.1.0]hep... | 18.17 | 3.3     | ng    | 56542    | 4                     | 17.85 | 339828 | 20.0 |
| unknown18.44         | 18.44 | 5.3     | ng    | 90914    | 4                     | 17.85 | 339828 | 20.0 |
| Isoquinoline         | 18.80 | 7.5     | ng    | 127509   | 4                     | 17.85 | 339828 | 20.0 |
| 2-Propenenitrile,... | 18.87 | 5.7     | ng    | 95947    | 4                     | 17.85 | 339828 | 20.0 |
| 1,2-Benzenediacet... | 18.92 | 13.2    | ng    | 224971   | 4                     | 17.85 | 339828 | 20.0 |
| unknown19.05         | 19.05 | 2.5     | ng    | 42987    | 4                     | 17.85 | 339828 | 20.0 |
| 7-Amino-7H-S-tria... | 19.53 | 2.4     | ng    | 40554    | 4                     | 17.85 | 339828 | 20.0 |
| unknown20.13         | 20.13 | 2.3     | ng    | 38828    | 5                     | 21.34 | 342037 | 20.0 |
| unknown20.23         | 20.23 | 2.8     | ng    | 48450    | 5                     | 21.34 | 342037 | 20.0 |
| unknown20.27         | 20.27 | 6.3     | ng    | 106825   | 5                     | 21.34 | 342037 | 20.0 |
| Phosphine imide, ... | 21.52 | 4.7     | ng    | 80313    | 5                     | 21.34 | 342037 | 20.0 |