

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020918\  
 Data File : BF102877.D  
 Acq On : 9 Feb 2018 10:02  
 Operator : SJ/JU  
 Sample : J1471-07  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 3N-18

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-BF020318.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         | 2.157       | 6             | 12          | 27           | rVB      | 525966         | 818033        | 16.95%          | 1.624%        |
| 2         | 5.116       | 512           | 515         | 525          | rVB      | 136823         | 184035        | 3.81%           | 0.365%        |
| 3         | 5.504       | 576           | 581         | 585          | rBV      | 4304601        | 4826885       | 100.00%         | 9.582%        |
| 4         | 6.516       | 747           | 753         | 756          | rBV      | 4261122        | 4708273       | 97.54%          | 9.347%        |
| 5         | 6.657       | 772           | 777         | 780          | rBV      | 4703369        | 4684668       | 97.05%          | 9.300%        |
| 6         | 6.887       | 812           | 816         | 819          | rBV      | 1297964        | 1120814       | 23.22%          | 2.225%        |
| 7         | 7.040       | 838           | 842         | 846          | rBV      | 3601599        | 3467076       | 71.83%          | 6.883%        |
| 8         | 7.445       | 906           | 911         | 915          | rBV      | 2941811        | 3208440       | 66.47%          | 6.369%        |
| 9         | 7.528       | 921           | 925         | 932          | rVB      | 52459          | 71502         | 1.48%           | 0.142%        |
| 10        | 7.622       | 938           | 941         | 944          | rVB      | 58366          | 51072         | 1.06%           | 0.101%        |
| 11        | 8.169       | 1030          | 1034        | 1037         | rBV      | 1655250        | 1493352       | 30.94%          | 2.965%        |
| 12        | 9.245       | 1212          | 1217        | 1220         | rBV      | 4396932        | 4467896       | 92.56%          | 8.870%        |
| 13        | 9.316       | 1226          | 1229        | 1231         | rVB      | 112089         | 80794         | 1.67%           | 0.160%        |
| 14        | 9.633       | 1279          | 1283        | 1285         | rBV      | 297275         | 272099        | 5.64%           | 0.540%        |
| 15        | 9.845       | 1315          | 1319        | 1322         | rVB      | 319145         | 254441        | 5.27%           | 0.505%        |
| 16        | 9.922       | 1328          | 1332        | 1336         | rBV2     | 1701945        | 1563741       | 32.40%          | 3.104%        |
| 17        | 10.075      | 1354          | 1358        | 1361         | rBV4     | 36538          | 49237         | 1.02%           | 0.098%        |
| 18        | 10.104      | 1361          | 1363        | 1366         | rBV2     | 67095          | 84443         | 1.75%           | 0.168%        |
| 19        | 10.328      | 1397          | 1401        | 1402         | rBV2     | 97841          | 118525        | 2.46%           | 0.235%        |
| 20        | 10.345      | 1402          | 1404        | 1407         | rVB      | 351042         | 259890        | 5.38%           | 0.516%        |
| 21        | 10.563      | 1437          | 1441        | 1444         | rBV2     | 104068         | 148720        | 3.08%           | 0.295%        |
| 22        | 10.669      | 1457          | 1459        | 1461         | rBV      | 60488          | 58555         | 1.21%           | 0.116%        |
| 23        | 10.716      | 1463          | 1467        | 1470         | rVV      | 2664706        | 2654905       | 55.00%          | 5.270%        |
| 24        | 10.816      | 1481          | 1484        | 1486         | rBV      | 241583         | 240680        | 4.99%           | 0.478%        |
| 25        | 10.833      | 1486          | 1487        | 1490         | rVV      | 285130         | 224871        | 4.66%           | 0.446%        |
| 26        | 10.886      | 1493          | 1496        | 1499         | rBV      | 584940         | 483866        | 10.02%          | 0.961%        |
| 27        | 10.922      | 1499          | 1502        | 1505         | rVV2     | 529291         | 611741        | 12.67%          | 1.214%        |
| 28        | 10.951      | 1505          | 1507        | 1510         | rVV2     | 564353         | 535161        | 11.09%          | 1.062%        |
| 29        | 10.980      | 1510          | 1512        | 1514         | rVV      | 350639         | 277956        | 5.76%           | 0.552%        |
| 30        | 11.022      | 1514          | 1519        | 1520         | rVV2     | 254831         | 334876        | 6.94%           | 0.665%        |
| 31        | 11.069      | 1523          | 1527        | 1530         | rVV3     | 553979         | 683361        | 14.16%          | 1.357%        |
| 32        | 11.104      | 1530          | 1533        | 1535         | rVV      | 680658         | 541517        | 11.22%          | 1.075%        |
| 33        | 11.122      | 1535          | 1536        | 1538         | rVV      | 171167         | 152113        | 3.15%           | 0.302%        |
| 34        | 11.145      | 1538          | 1540        | 1544         | rVB2     | 382676         | 304163        | 6.30%           | 0.604%        |

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ALS Vial : 3 Sample Multiplier: 1

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BNA\_F  
ClientSampleId :  
3N-18

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-BF020318.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 11.269 | 1558 | 1561 | 1563 | rBV  | 180164  | 161062  | 3.34%  | 0.320% |
| 36 | 11.292 | 1563 | 1565 | 1568 | rVB3 | 63923   | 57839   | 1.20%  | 0.115% |
| 37 | 11.410 | 1581 | 1585 | 1588 | rBV  | 1608654 | 1475469 | 30.57% | 2.929% |
| 38 | 11.433 | 1588 | 1589 | 1593 | rVB  | 119497  | 85257   | 1.77%  | 0.169% |
| 39 | 11.692 | 1630 | 1633 | 1638 | rBV  | 62961   | 60868   | 1.26%  | 0.121% |
| 40 | 11.927 | 1668 | 1673 | 1681 | rBV  | 669609  | 815131  | 16.89% | 1.618% |
| 41 | 12.080 | 1696 | 1699 | 1701 | rBV3 | 86230   | 120824  | 2.50%  | 0.240% |
| 42 | 12.104 | 1701 | 1703 | 1711 | rVB2 | 179694  | 258783  | 5.36%  | 0.514% |
| 43 | 12.621 | 1788 | 1791 | 1797 | rBV  | 241613  | 246084  | 5.10%  | 0.489% |
| 44 | 12.716 | 1803 | 1807 | 1815 | rBV  | 255273  | 296577  | 6.14%  | 0.589% |
| 45 | 12.851 | 1825 | 1830 | 1832 | rBV  | 192803  | 242181  | 5.02%  | 0.481% |
| 46 | 12.874 | 1832 | 1834 | 1838 | rVB  | 305985  | 276567  | 5.73%  | 0.549% |
| 47 | 12.998 | 1850 | 1855 | 1858 | rBV  | 3476682 | 3424064 | 70.94% | 6.797% |
| 48 | 13.127 | 1871 | 1877 | 1883 | rBV  | 180398  | 335748  | 6.96%  | 0.667% |
| 49 | 13.180 | 1884 | 1886 | 1895 | rVB6 | 56789   | 87087   | 1.80%  | 0.173% |
| 50 | 13.469 | 1931 | 1935 | 1938 | rBV  | 921156  | 800860  | 16.59% | 1.590% |
| 51 | 13.880 | 2002 | 2005 | 2009 | rBV  | 337441  | 328451  | 6.80%  | 0.652% |
| 52 | 14.051 | 2030 | 2034 | 2037 | rBV  | 1138578 | 1001593 | 20.75% | 1.988% |
| 53 | 15.163 | 2220 | 2223 | 2233 | rVB3 | 118105  | 227351  | 4.71%  | 0.451% |
| 54 | 15.539 | 2282 | 2287 | 2294 | rVB  | 788126  | 1033793 | 21.42% | 2.052% |

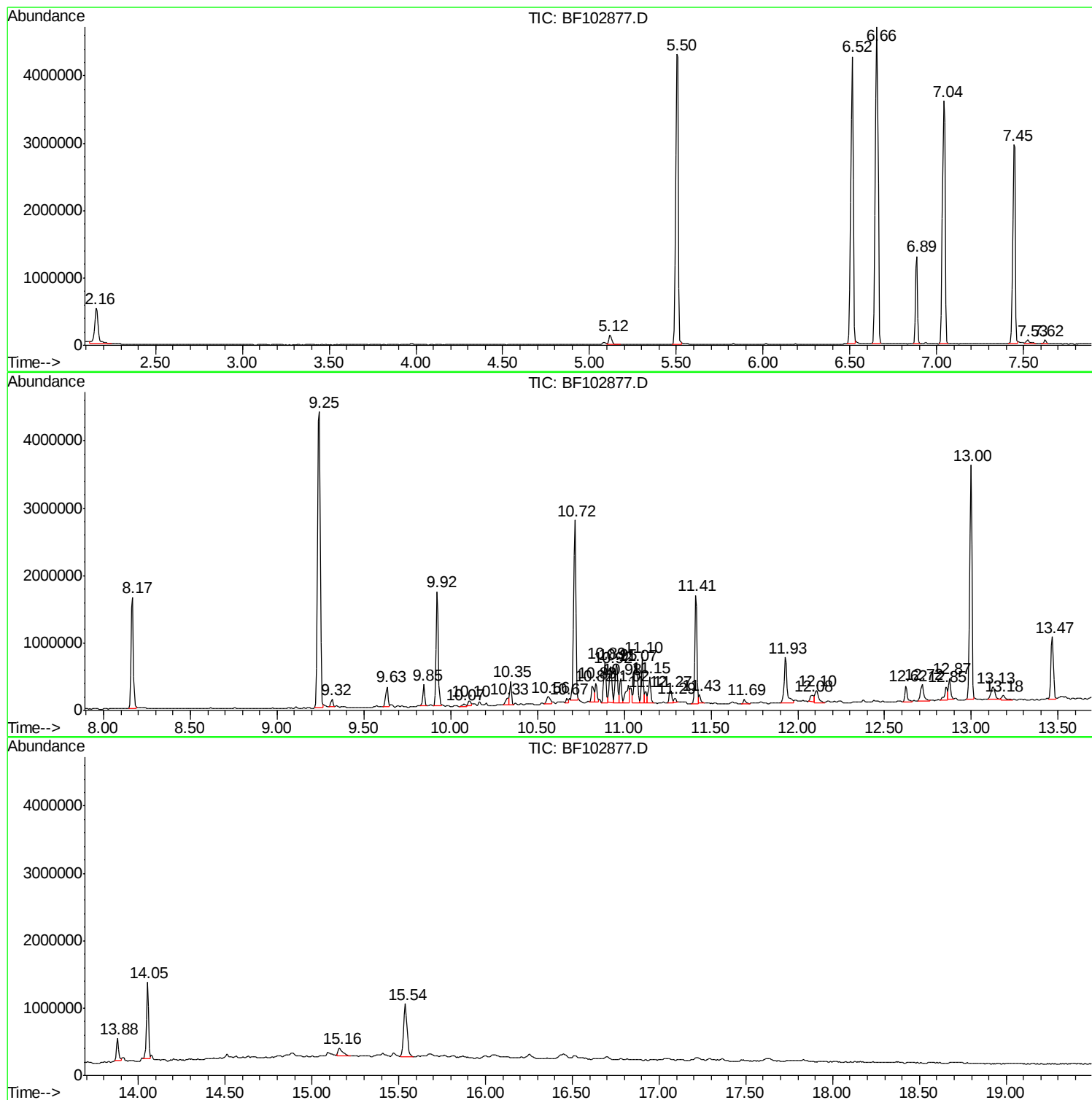
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Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
3N-18

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020918\  
Data File : BF102877.D  
Acq On : 9 Feb 2018 10:02  
Operator : SJ/JU  
Sample : J1471-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
3N-18

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

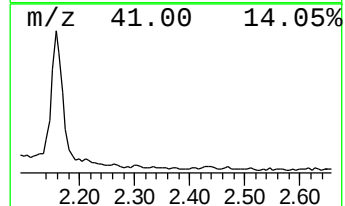
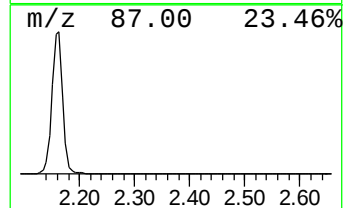
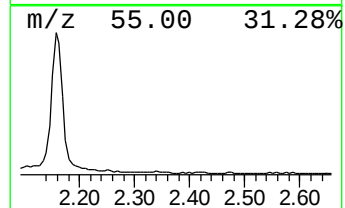
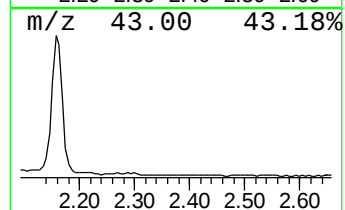
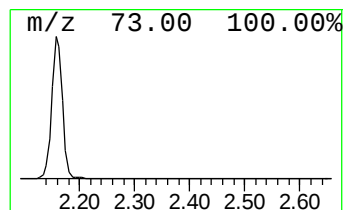
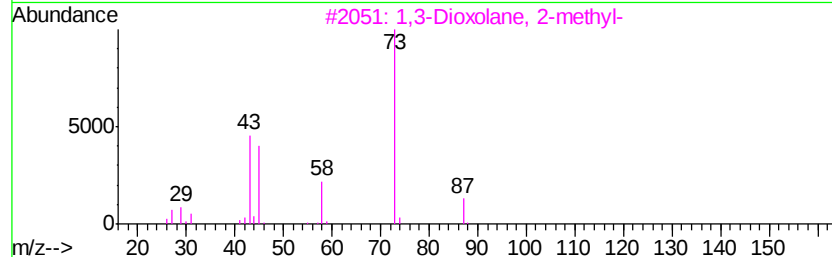
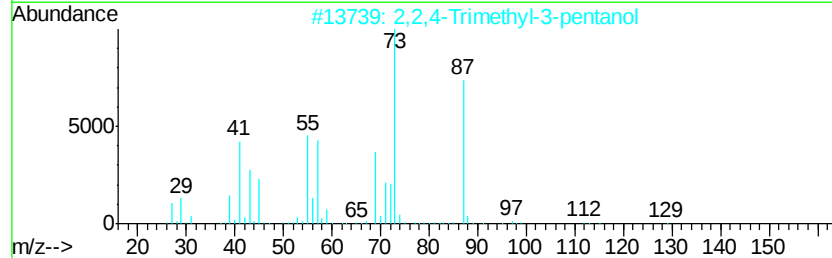
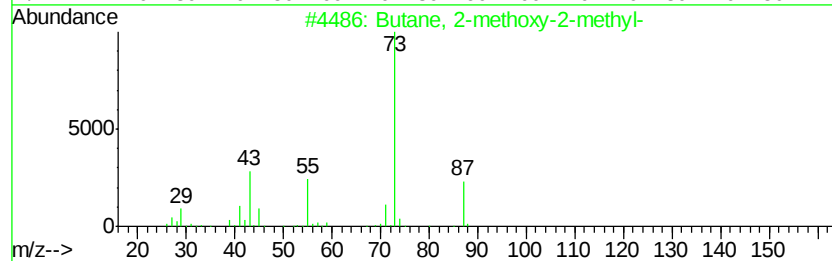
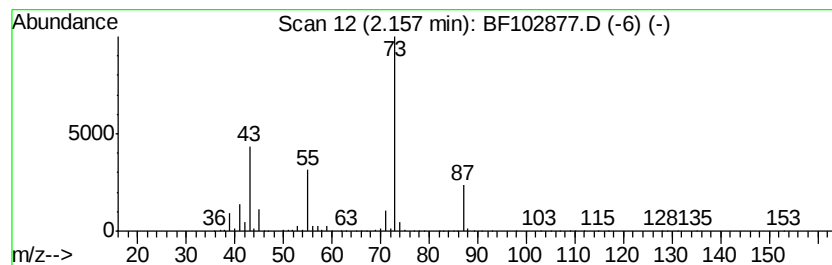
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 2.16 | 14.60 ng | 818033 | 1,4-Dichlorobenzene-d4 | 6.89 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl-        | 102 | C6H14O  | 000994-05-8 | 83   |
| 2    |    | 2,2,4-Trimethyl-3-pentanol         | 130 | C8H18O  | 005162-48-1 | 40   |
| 3    |    | 1,3-Dioxolane, 2-methyl-           | 88  | C4H8O2  | 000497-26-7 | 23   |
| 4    |    | Butanoic acid, 2-hydroxy-2-methyl- | 132 | C6H12O3 | 032793-34-3 | 23   |
| 5    |    | Acetamide, N-ethyl-                | 87  | C4H9NO  | 000625-50-3 | 9    |



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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

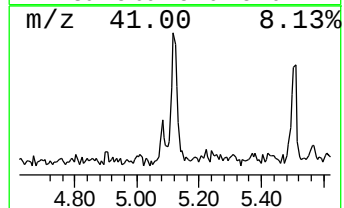
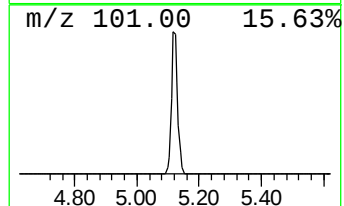
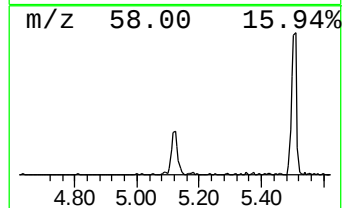
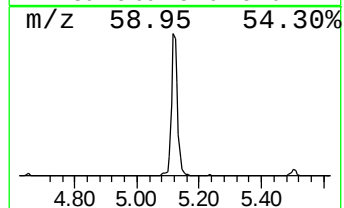
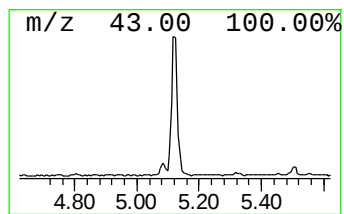
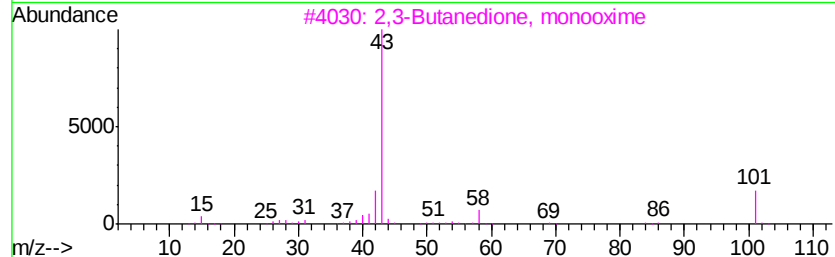
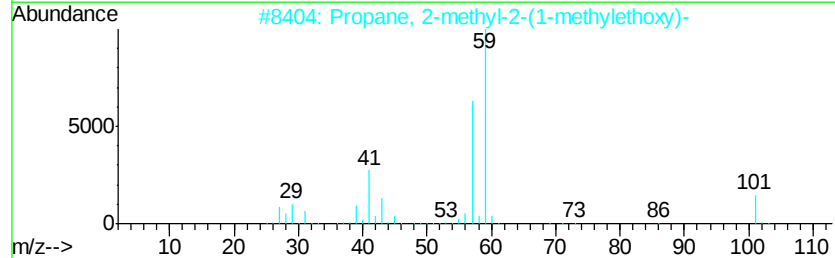
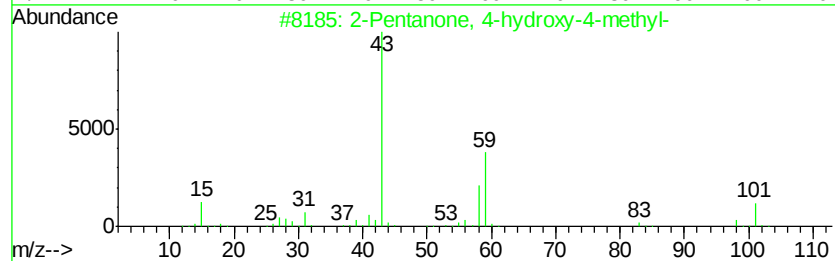
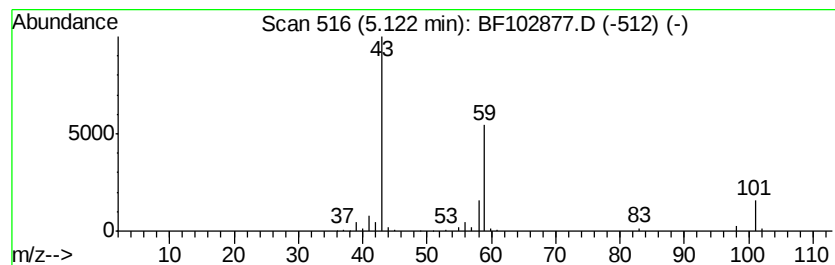
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 19

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 5.12 | 3.28 ng | 184035 | 1,4-Dichlorobenzene-d4 | 6.89 |

| Hit# of | 5                                   | Tentative ID | MW      | MolForm     | CAS# | Qual |
|---------|-------------------------------------|--------------|---------|-------------|------|------|
| 1       | 2-Pentanone, 4-hydroxy-4-methyl-    | 116          | C6H12O2 | 000123-42-2 | 59   |      |
| 2       | Propane, 2-methyl-2-(1-methyleth... | 116          | C7H16O  | 017348-59-3 | 42   |      |
| 3       | 2,3-Butanedione, monooxime          | 101          | C4H7NO2 | 000057-71-6 | 16   |      |
| 4       | 1-Propen-2-ol, acetate              | 100          | C5H8O2  | 000108-22-5 | 10   |      |
| 5       | Butane, 1-ethoxy-                   | 102          | C6H14O  | 000628-81-9 | 9    |      |



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ClientSampleId :  
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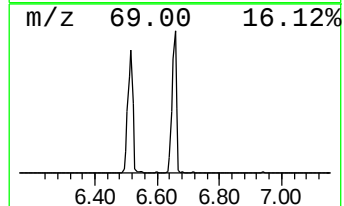
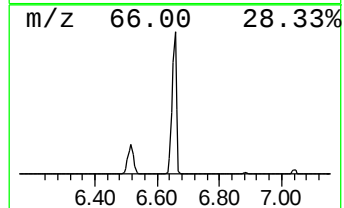
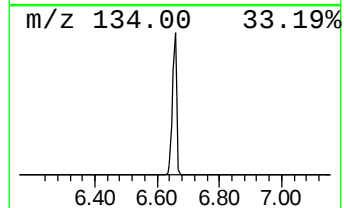
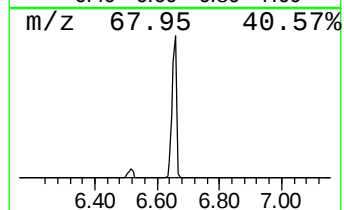
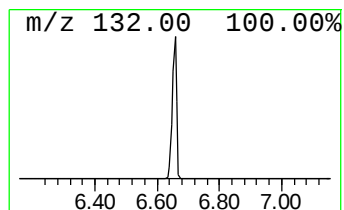
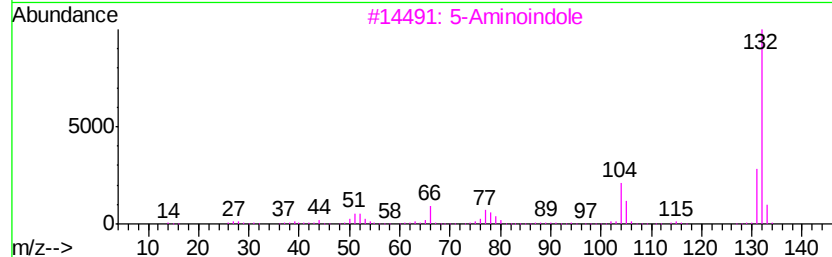
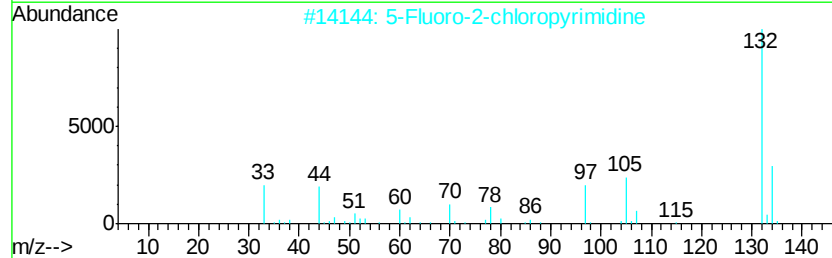
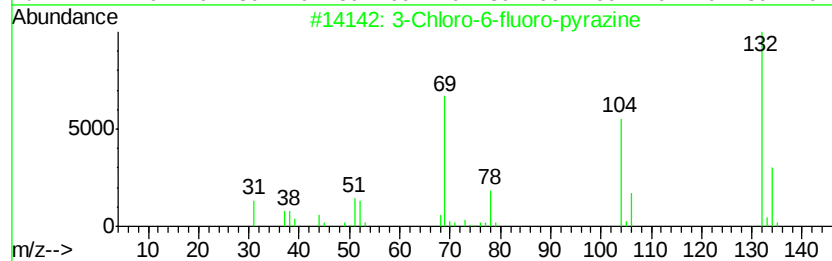
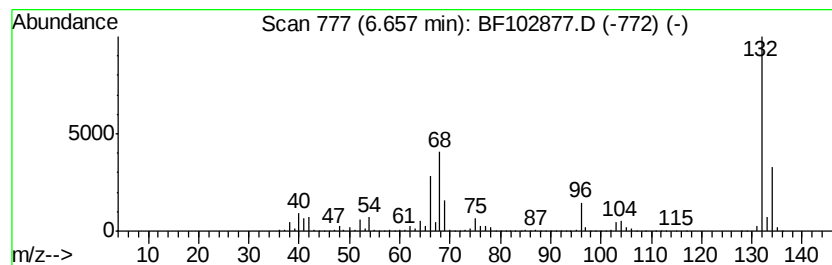
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 3 unknown6.66 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.66 | 83.59 ng | 4684670 | 1,4-Dichlorobenzene-d4 | 6.89 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | 5-Fluoro-2-chloropyrimidine         | 132 | C4H2ClFN2 | 062802-42-0  | 16   |
| 3    |    | 5-Aminoindole                       | 132 | C8H8N2    | 005192-03-0  | 12   |
| 4    |    | Xenon                               | 132 | Xe        | 007440-63-3  | 9    |
| 5    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |



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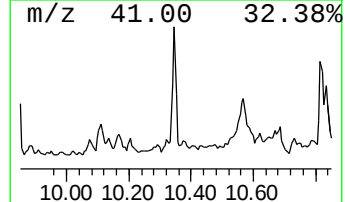
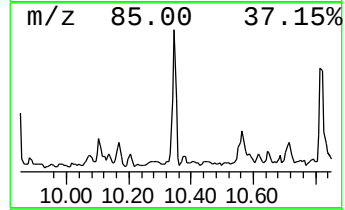
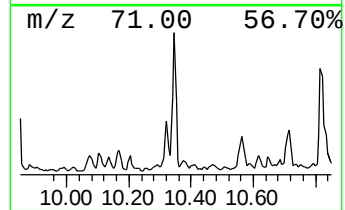
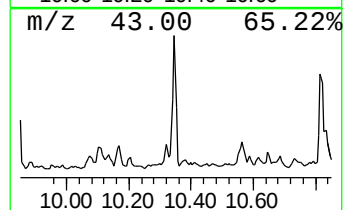
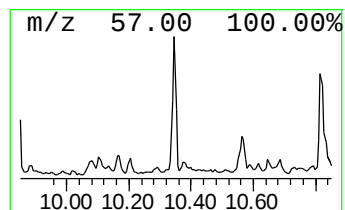
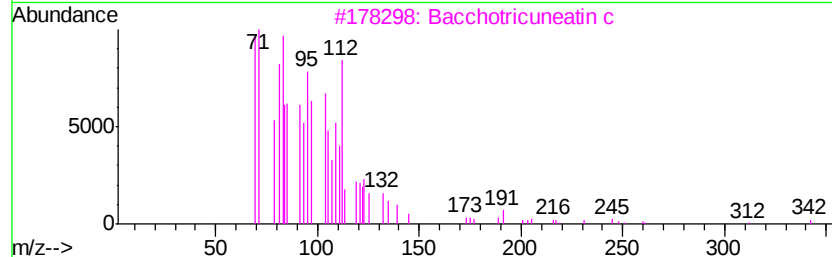
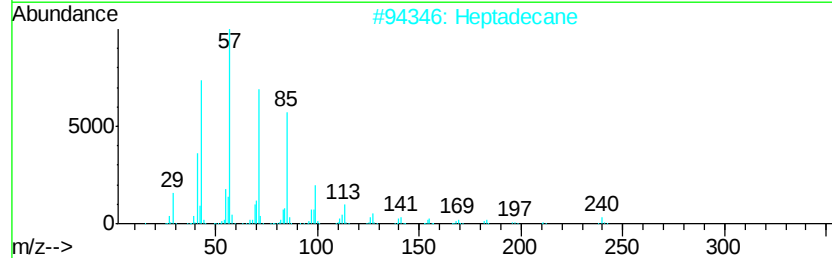
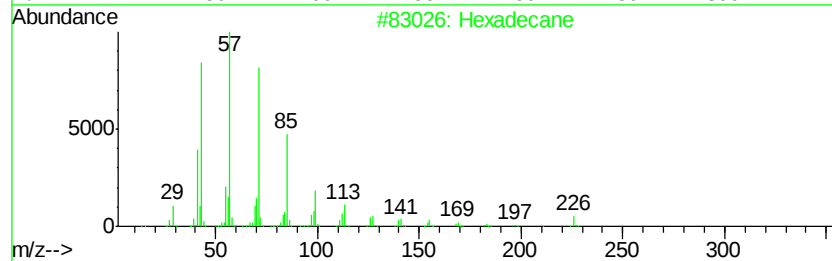
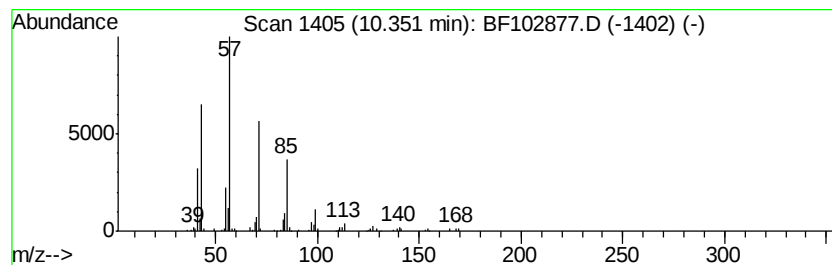
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BNA\_F  
ClientSampleId :  
3N-18

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 Hexadecane Concentration Rank 18

| R.T.    | EstConc                      | Area         | Relative to ISTD | R.T.      |
|---------|------------------------------|--------------|------------------|-----------|
| 10.35   | 3.32 ng                      | 259890       | Acenaphthene-d10 | 9.92      |
| Hit# of | 5                            | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Hexadecane                   | 226 C16H34   | 000544-76-3      | 93        |
| 2       | Heptadecane                  | 240 C17H36   | 000629-78-7      | 83        |
| 3       | Bacchotricuneatin c          | 342 C20H22O5 | 066563-30-2      | 76        |
| 4       | Dodecane, 2-methyl-6-propyl- | 226 C16H34   | 055045-08-4      | 72        |
| 5       | 9-methylheptadecane          | 254 C18H38   | 026741-18-4      | 72        |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020918\  
Data File : BF102877.D  
Acq On : 9 Feb 2018 10:02  
Operator : SJ/JU  
Sample : J1471-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
3N-18

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

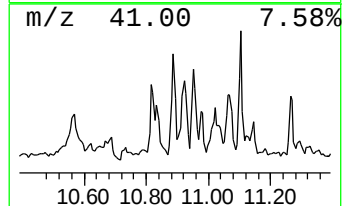
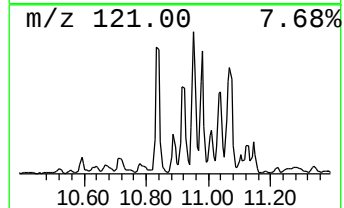
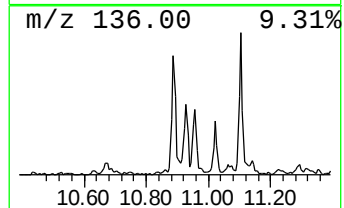
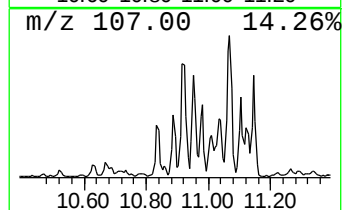
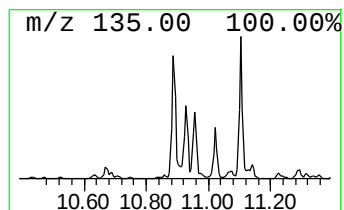
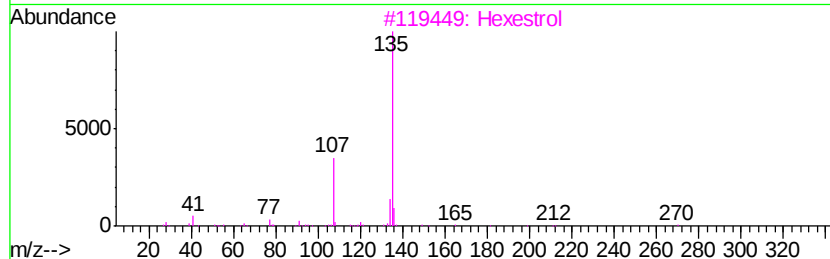
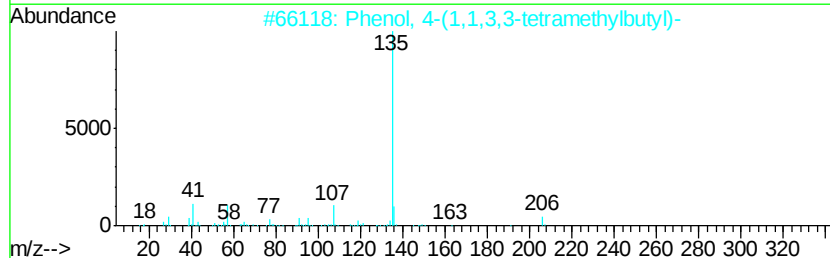
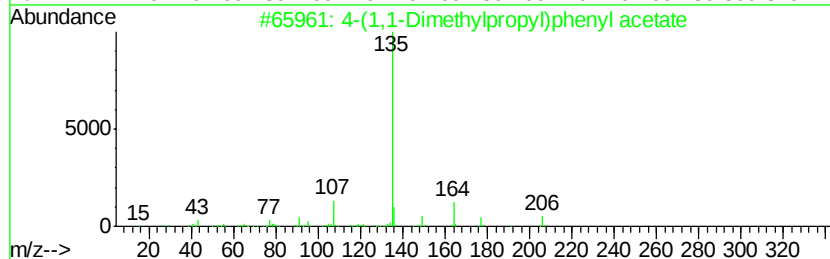
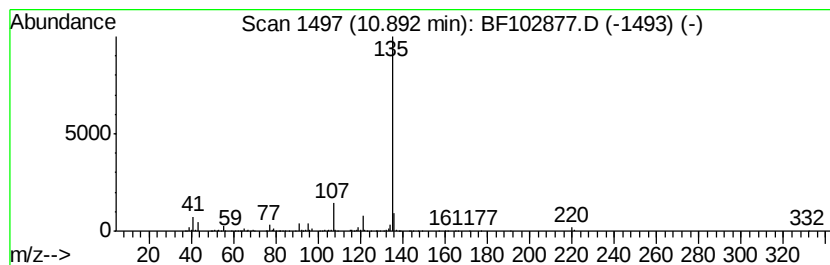
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 4-(1,1-Dimethylpropyl)pheny... Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 10.89 | 6.56 ng | 483866 | Phenanthrene-d10 | 11.41 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 4-(1,1-Dimethylpropyl)phenyl ace... | 206 | C13H18O2  | 1000373-45-3 | 78   |
| 2    |    |   | Phenol, 4-(1,1,3,3-tetramethylbu... | 206 | C14H22O   | 000140-66-9  | 78   |
| 3    |    |   | Hexestrol                           | 270 | C18H22O2  | 000084-16-2  | 72   |
| 4    |    |   | m-Anisic acid, 4-nitrophenyl ester  | 273 | C14H11NO5 | 1000307-80-4 | 64   |
| 5    |    |   | Phenol, 4-(1,1-dimethylpropyl)-     | 164 | C11H16O   | 000080-46-6  | 64   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020918\  
Data File : BF102877.D  
Acq On : 9 Feb 2018 10:02  
Operator : SJ/JU  
Sample : J1471-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

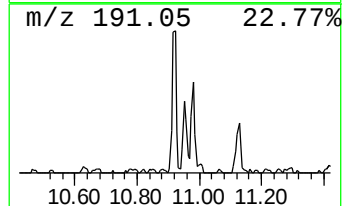
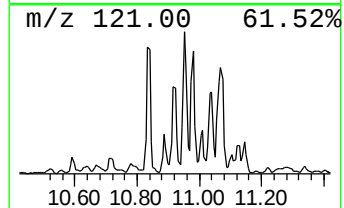
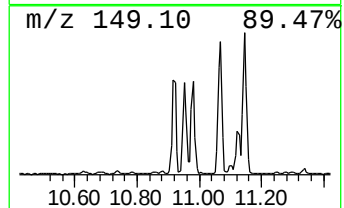
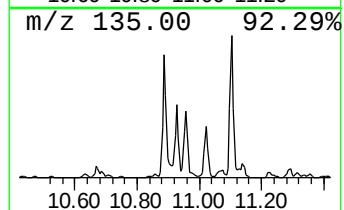
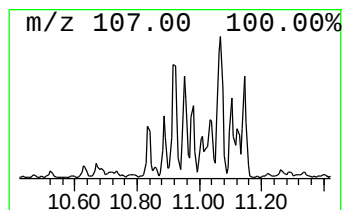
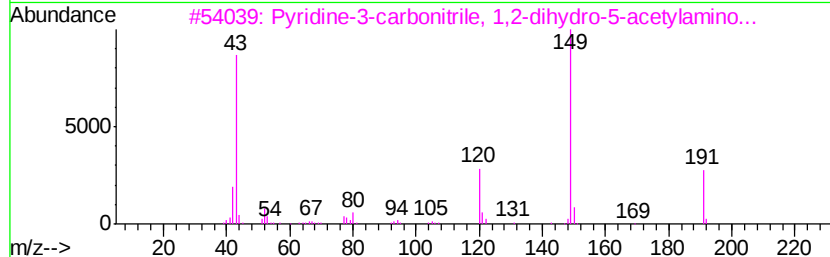
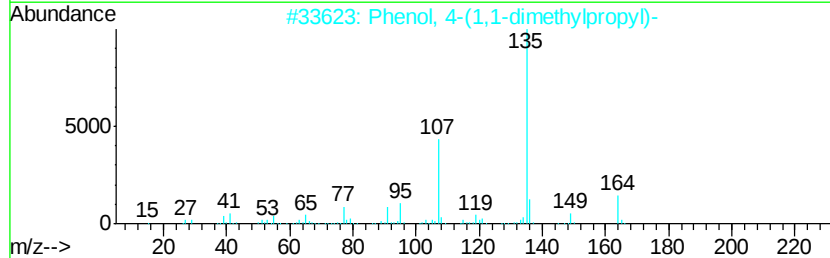
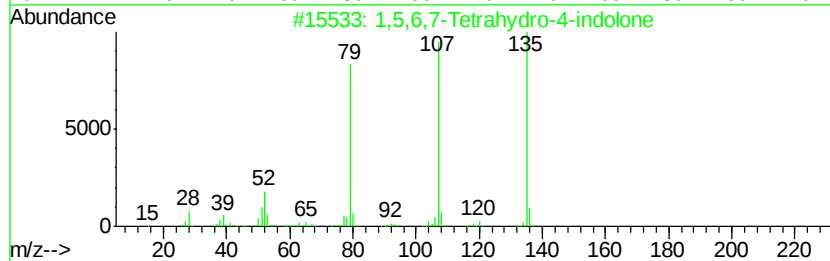
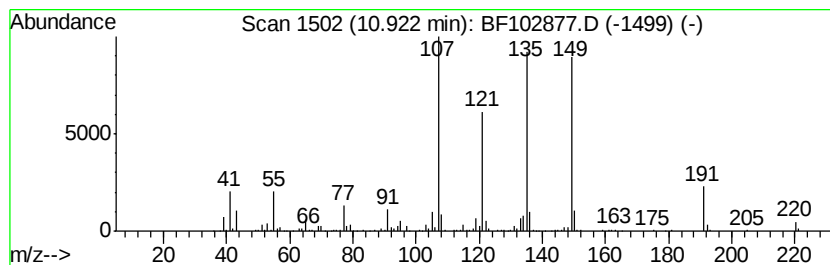
Instrument :  
BNA\_F  
ClientSampleId :  
3N-18

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 8 unknown10.92 Concentration Rank 6

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 10.92     | 8.29 ng                             | 611741 | Phenanthrene-d10 | 11.41        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 1,5,6,7-Tetrahydro-4-indolone       | 135    | C8H9NO           | 013754-86-4  | 46   |
| 2         | Phenol, 4-(1,1-dimethylpropyl)-     | 164    | C11H16O          | 000080-46-6  | 43   |
| 3         | Pvridine-3-carbonitrile, 1,2-dih... | 191    | C9H9N3O2         | 220098-92-0  | 25   |
| 4         | 5H-Pyrrolo(3,2-d)pvrimidine-2,4-... | 149    | C6H7N5           | 1000244-21-4 | 22   |
| 5         | Acetamide, N-(4-methylphenyl)-      | 149    | C9H11NO          | 000103-89-9  | 22   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020918\  
Data File : BF102877.D  
Acq On : 9 Feb 2018 10:02  
Operator : SJ/JU  
Sample : J1471-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

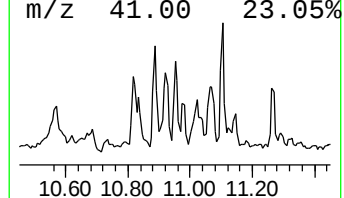
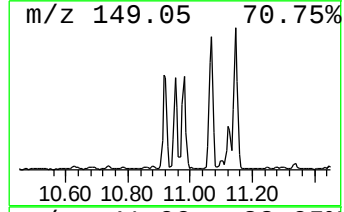
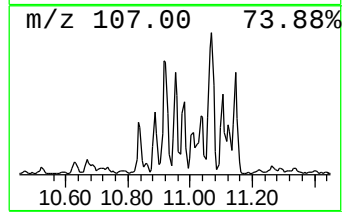
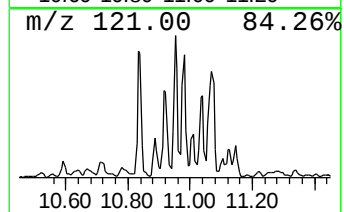
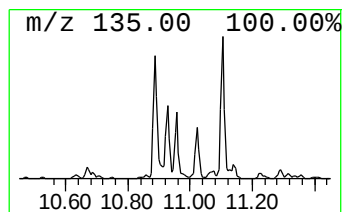
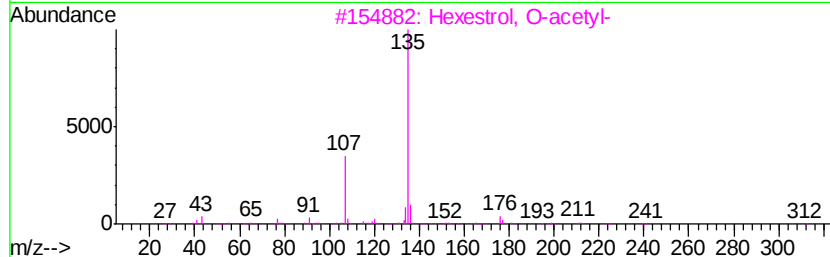
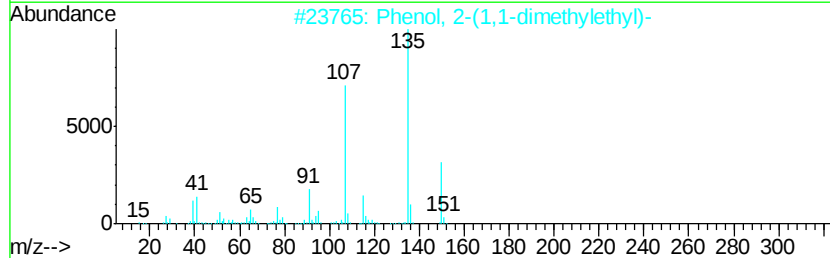
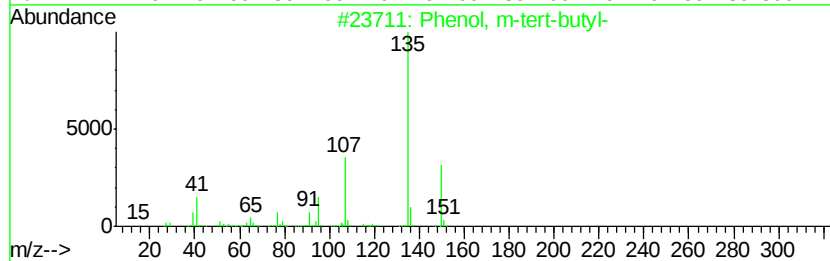
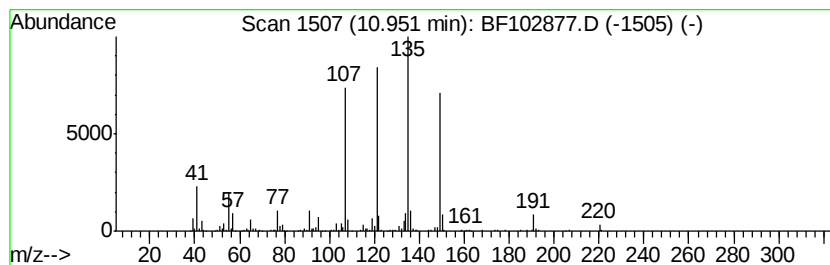
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ClientSampleId :  
3N-18

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 9 Phenol, m-tert-butyl- Concentration Rank 8

| R.T.    | EstConc | Area                           | Relative to ISTD | R.T.     |              |      |
|---------|---------|--------------------------------|------------------|----------|--------------|------|
| 10.95   | 7.25 ng | 535161                         | Phenanthrene-d10 | 11.41    |              |      |
| Hit# of | 5       | Tentative ID                   | MW               | MolForm  | CAS#         | Qual |
| 1       |         | Phenol, m-tert-butyl-          | 150              | C10H14O  | 000585-34-2  | 55   |
| 2       |         | Phenol, 2-(1,1-dimethylethyl)- | 150              | C10H14O  | 000088-18-6  | 49   |
| 3       |         | Hexestrol, O-acetyl-           | 312              | C20H24O3 | 1000374-87-8 | 43   |
| 4       |         | Neodihydrocarveol              | 154              | C10H18O  | 018675-34-8  | 38   |
| 5       |         | Acetamide, N-(3-methylphenyl)- | 149              | C9H11NO  | 000537-92-8  | 30   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020918\  
Data File : BF102877.D  
Acq On : 9 Feb 2018 10:02  
Operator : SJ/JU  
Sample : J1471-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

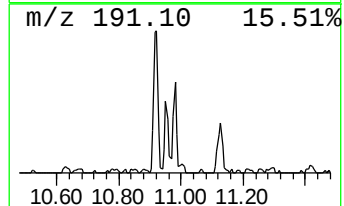
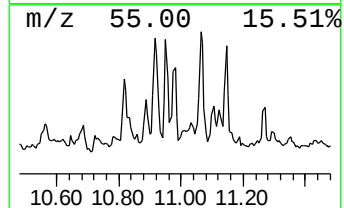
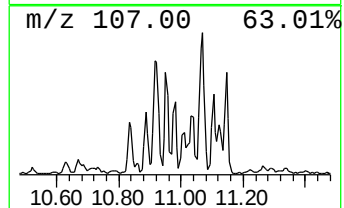
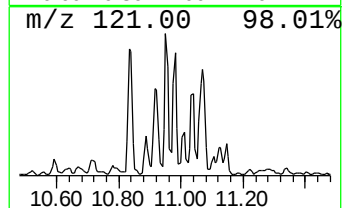
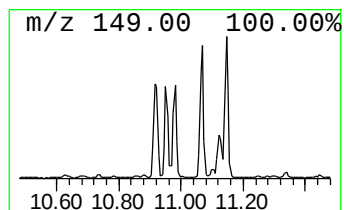
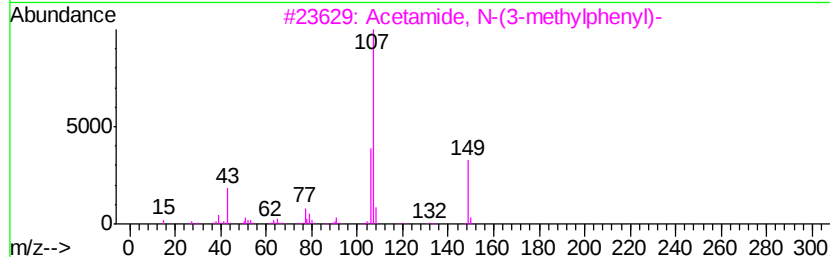
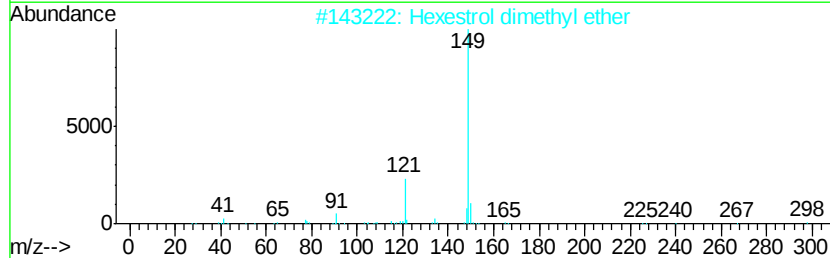
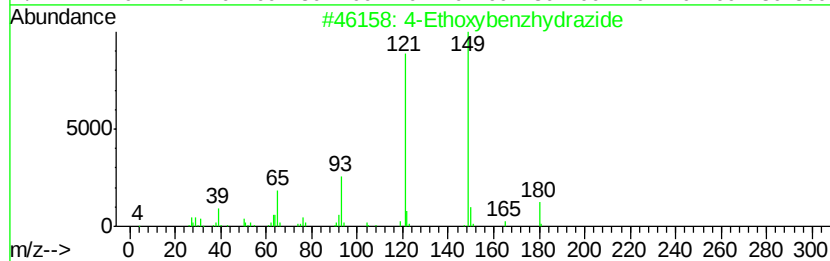
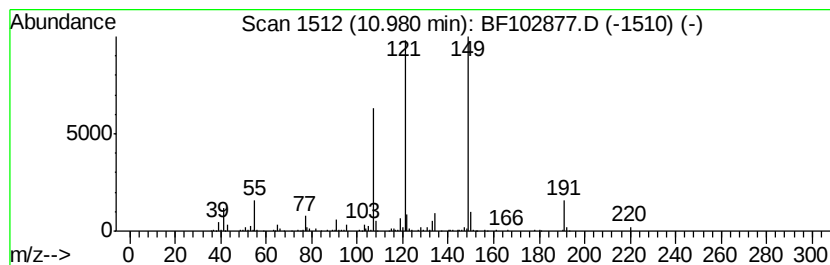
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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 10 4-Ethoxybenzhydrazide Concentration Rank 16

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 10.98     | 3.77 ng                             | 277956 | Phenanthrene-d10 | 11.41        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 4-Ethoxybenzhydrazide               | 180    | C9H12N2O2        | 058586-81-5  | 53   |
| 2         | Hexestrol dimethyl ether            | 298    | C20H26O2         | 000130-78-9  | 45   |
| 3         | Acetamide, N-(3-methylphenyl)-      | 149    | C9H11NO          | 000537-92-8  | 43   |
| 4         | Phenol, 2-methyl-4-(1,1,3,3-tetr... | 220    | C15H24O          | 002219-84-3  | 38   |
| 5         | Isoxazole, 4,5-dihydro-5-[(4-met... | 191    | C11H13NO2        | 1000362-14-0 | 38   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020918\  
Data File : BF102877.D  
Acq On : 9 Feb 2018 10:02  
Operator : SJ/JU  
Sample : J1471-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
3N-18

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

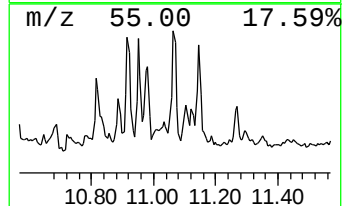
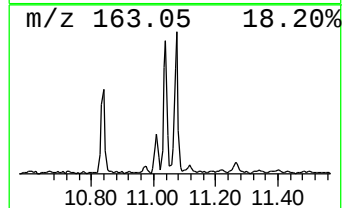
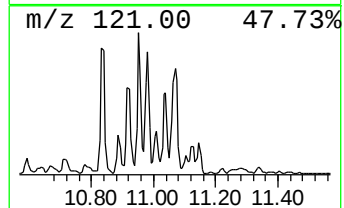
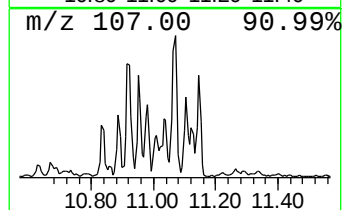
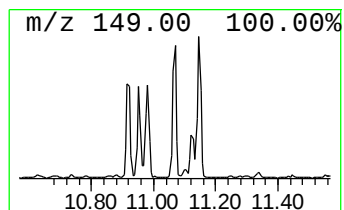
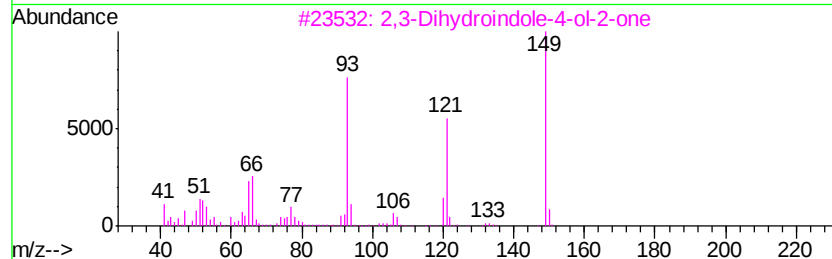
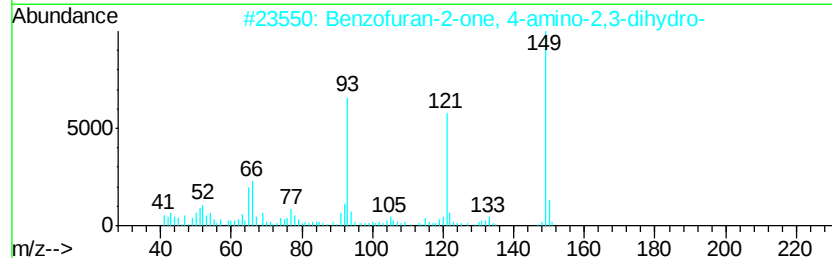
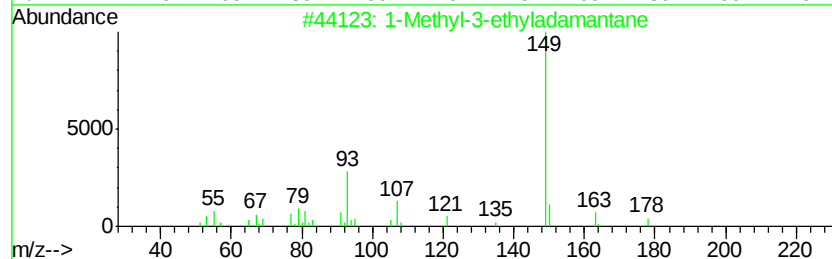
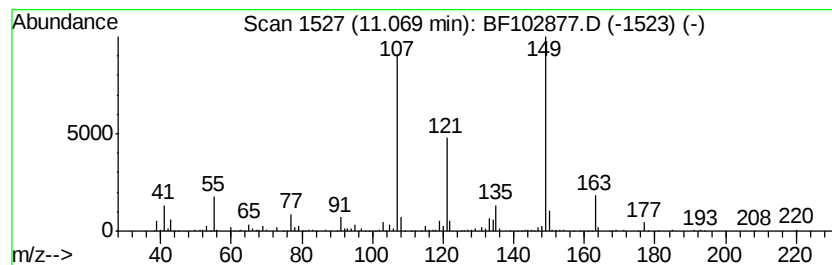
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 unknown11.07 Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.07 | 9.26 ng | 683361 | Phenanthrene-d10 | 11.41 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 1-Methyl-3-ethyladamantane          | 178 | C13H22   | 001687-34-9  | 43   |
| 2    |    | Benzofuran-2-one, 4-amino-2,3-di... | 149 | C8H7NO2  | 075990-99-7  | 38   |
| 3    |    | 2,3-Dihydroindole-4-ol-2-one        | 149 | C8H7NO2  | 013402-55-6  | 38   |
| 4    |    | Phthalic acid, butyl 3-methoxybe... | 342 | C20H22O5 | 1000377-97-7 | 38   |
| 5    |    | Phenol, 4-(1,1-dimethylethyl)-2-... | 164 | C11H16O  | 000098-27-1  | 38   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020918\  
Data File : BF102877.D  
Acq On : 9 Feb 2018 10:02  
Operator : SJ/JU  
Sample : J1471-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
3N-18

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

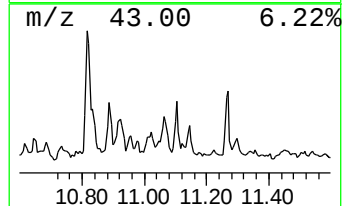
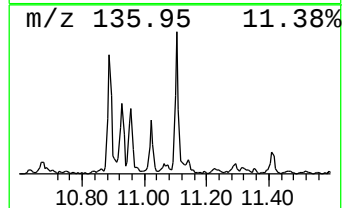
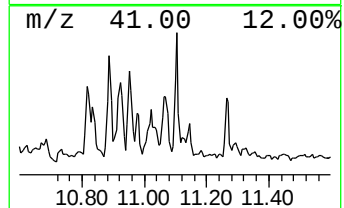
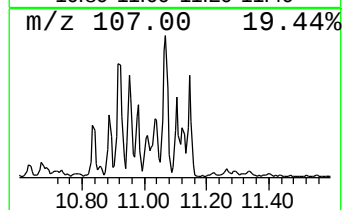
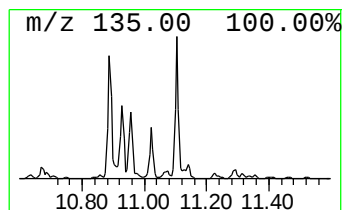
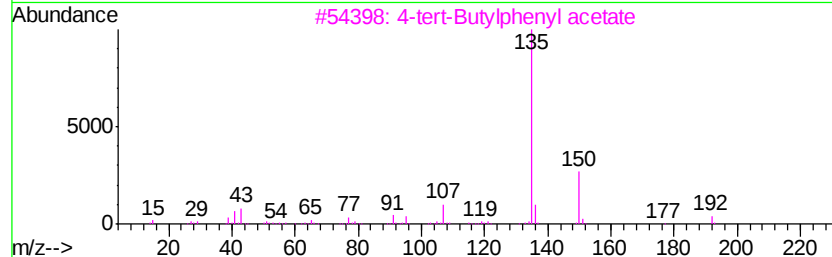
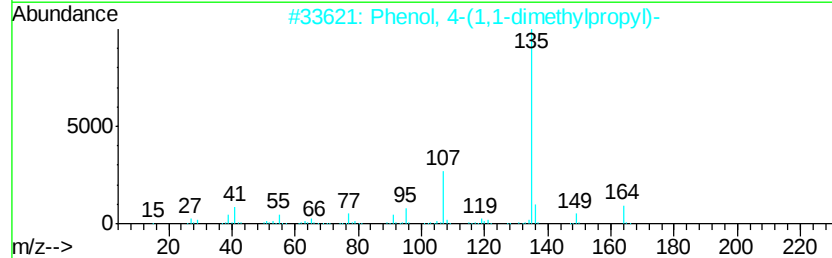
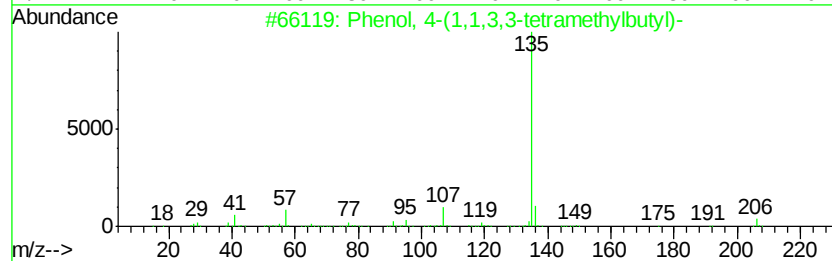
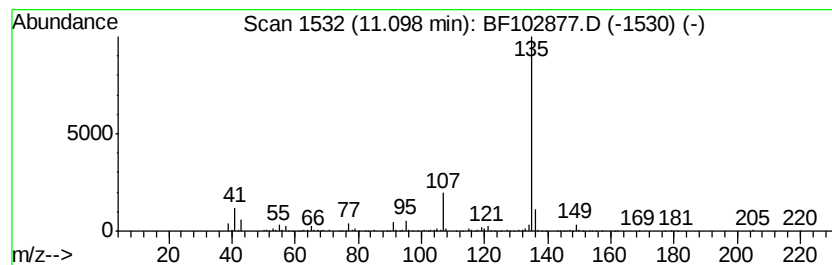
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.10 | 7.34 ng | 541517 | Phenanthrene-d10 | 11.41 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Phenol, 4-(1,1,3,3-tetramethylbu... | 206 | C14H22O     | 000140-66-9  | 78   |
| 2    |    |   | Phenol, 4-(1,1-dimethylpropyl)-     | 164 | C11H16O     | 000080-46-6  | 78   |
| 3    |    |   | 4-tert-Butylphenyl acetate          | 192 | C12H16O2    | 003056-64-2  | 64   |
| 4    |    |   | 1-Adamantanecarboxylic acid, 3,4... | 324 | C17H18Cl2O2 | 1000307-66-4 | 64   |
| 5    |    |   | Hexestrol                           | 270 | C18H22O2    | 005635-50-7  | 64   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020918\  
Data File : BF102877.D  
Acq On : 9 Feb 2018 10:02  
Operator : SJ/JU  
Sample : J1471-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
3N-18

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

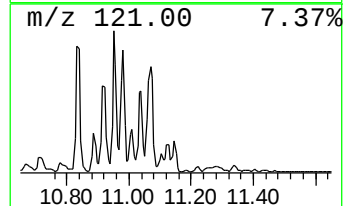
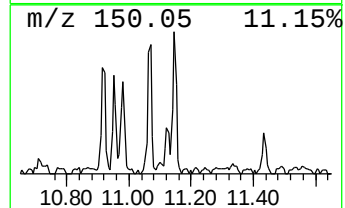
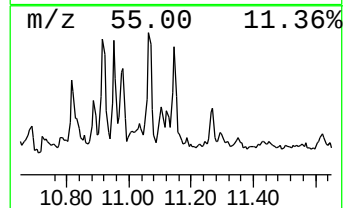
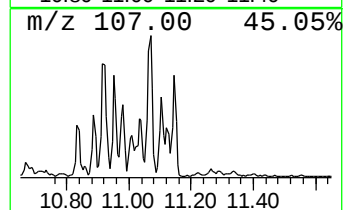
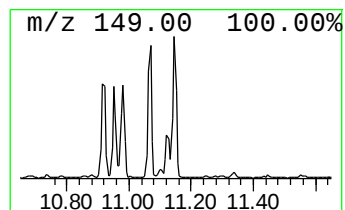
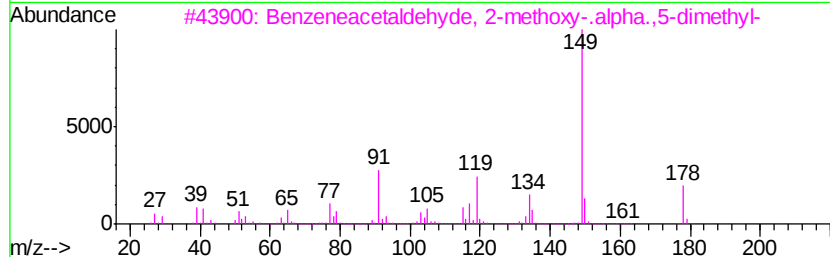
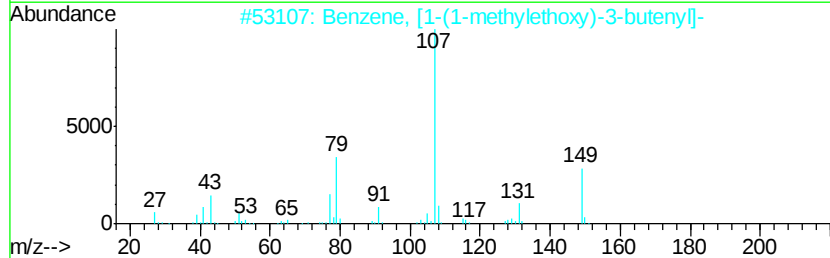
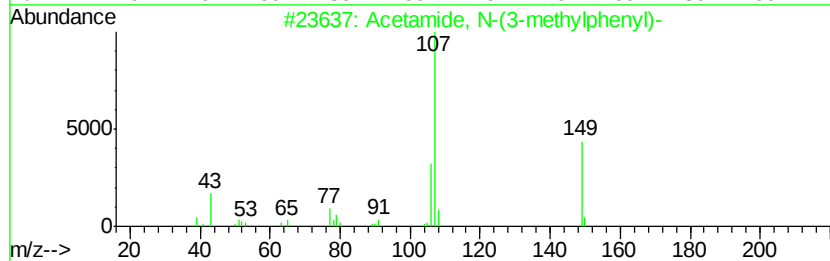
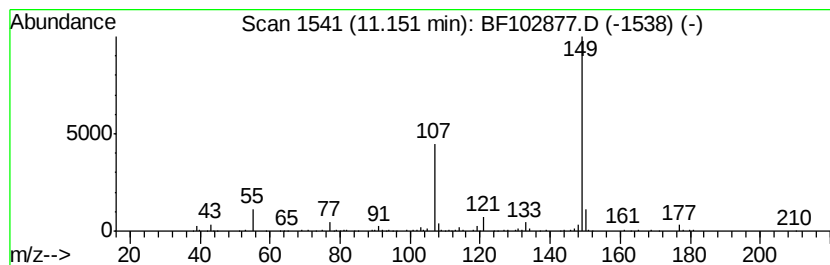
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Acetamide, N-(3-methylphenyl)- Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.15 | 4.12 ng | 304163 | Phenanthrene-d10 | 11.41 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Acetamide, N-(3-methylphenyl)-      | 149 | C9H11NO  | 000537-92-8  | 64   |
| 2    |    |   | Benzene, [1-(1-methylethoxy)-3-b... | 190 | C13H18O  | 098088-47-2  | 50   |
| 3    |    |   | Benzeneacetaldehyde, 2-methoxy-.... | 178 | C11H14O2 | 053155-90-1  | 47   |
| 4    |    |   | Phthalic acid, isobutyl propyl e... | 264 | C15H20O4 | 1000308-97-1 | 43   |
| 5    |    |   | Phthalic acid, cyclobutyl dodecy... | 388 | C24H36O4 | 1000314-90-7 | 43   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020918\  
Data File : BF102877.D  
Acq On : 9 Feb 2018 10:02  
Operator : SJ/JU  
Sample : J1471-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
3N-18

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

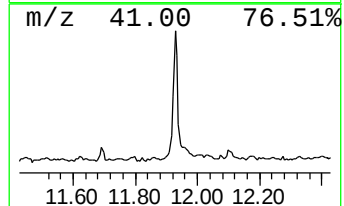
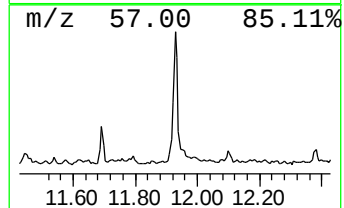
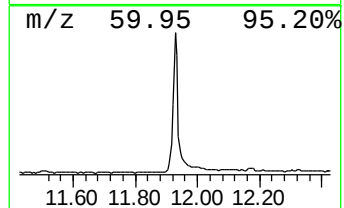
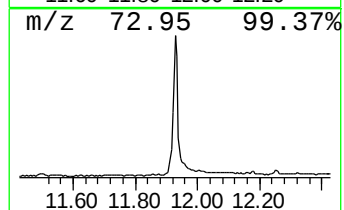
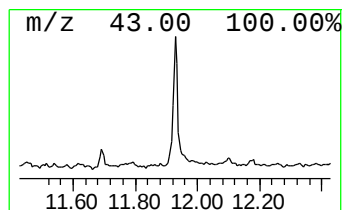
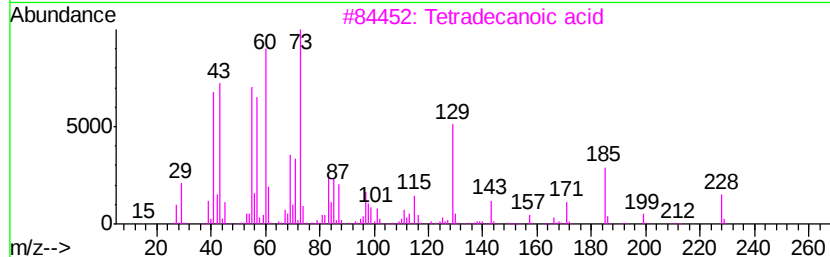
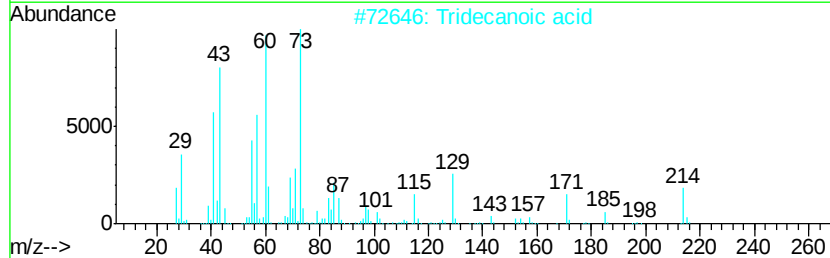
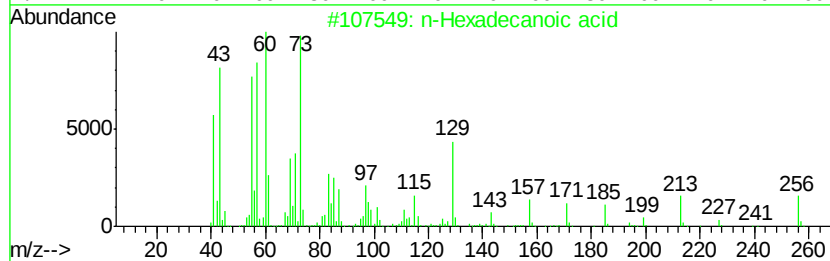
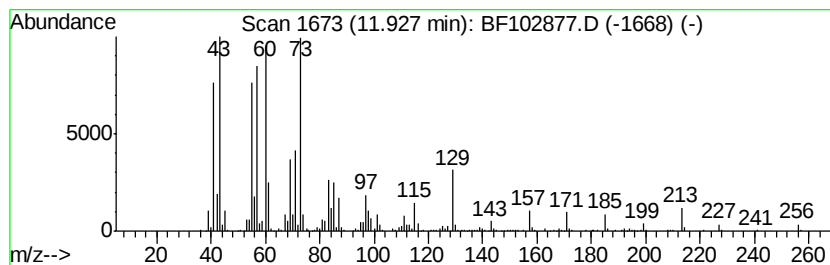
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TIC Integration Parameters: LSCINT.P

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Peak Number 15 n-Hexadecanoic acid Concentration Rank 4

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 11.93 | 11.05 ng | 815131 | Phenanthrene-d10 | 11.41 |

| Hit# | of | Tentative ID                       | MW  | MolForm  | CAS#         | Qual |
|------|----|------------------------------------|-----|----------|--------------|------|
| 1    | 5  | n-Hexadecanoic acid                | 256 | C16H32O2 | 000057-10-3  | 99   |
| 2    |    | Tridecanoic acid                   | 214 | C13H26O2 | 000638-53-9  | 93   |
| 3    |    | Tetradecanoic acid                 | 228 | C14H28O2 | 000544-63-8  | 93   |
| 4    |    | Pentadecanoic acid                 | 242 | C15H30O2 | 001002-84-2  | 90   |
| 5    |    | Oxalic acid, propyl tridecyl ester | 314 | C18H34O4 | 1000309-26-6 | 18   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020918\  
Data File : BF102877.D  
Acq On : 9 Feb 2018 10:02  
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Sample : J1471-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
3N-18

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

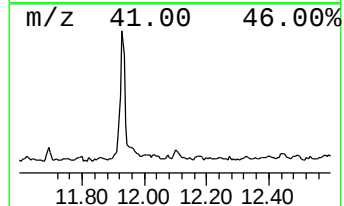
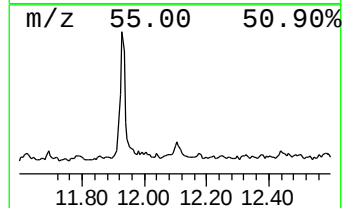
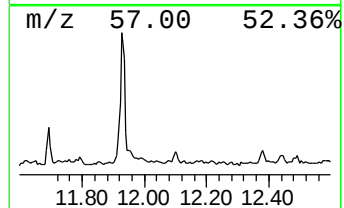
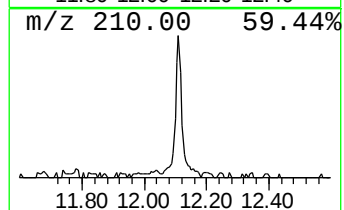
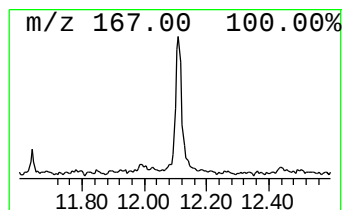
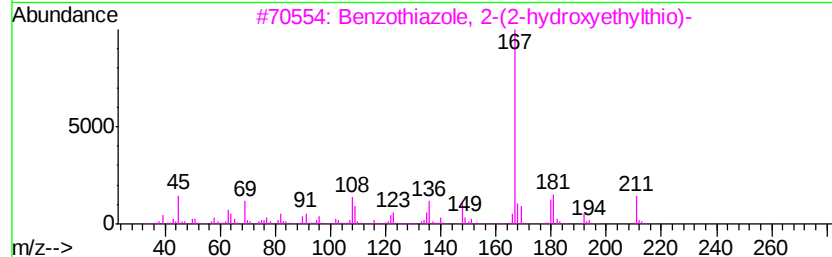
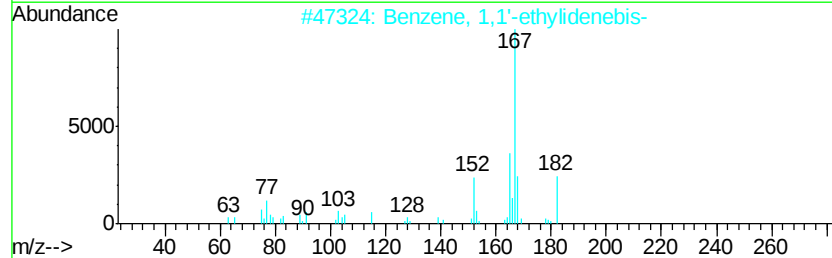
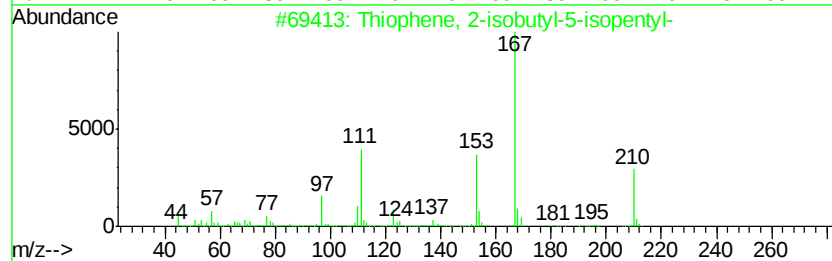
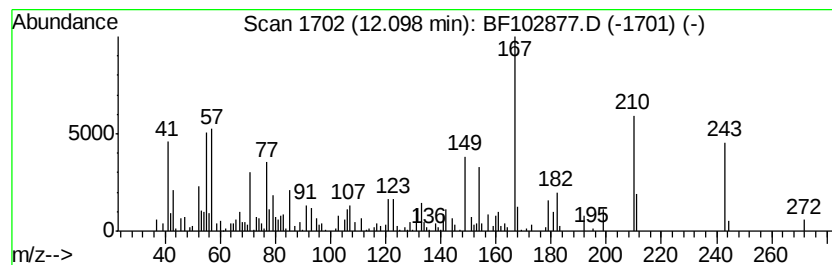
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 unknown12.10 Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.10 | 3.51 ng | 258783 | Phenanthrene-d10 | 11.41 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Thiophene, 2-isobutyl-5-isopentyl-  | 210 | C13H22S   | 004806-10-4 | 43   |
| 2    |    | Benzene, 1,1'-ethylidenebis-        | 182 | C14H14    | 000612-00-0 | 35   |
| 3    |    | Benzothiazole, 2-(2-hydroxyethyl)-  | 211 | C9H9NOS2  | 004665-63-8 | 35   |
| 4    |    | 9-Methyl-1-phenazinol               | 210 | C13H10N2O | 013860-43-0 | 30   |
| 5    |    | 3,5-Dimethoxy-4-hydroxyphenylace... | 212 | C10H12O5  | 004385-56-2 | 30   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020918\  
Data File : BF102877.D  
Acq On : 9 Feb 2018 10:02  
Operator : SJ/JU  
Sample : J1471-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
3N-18

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

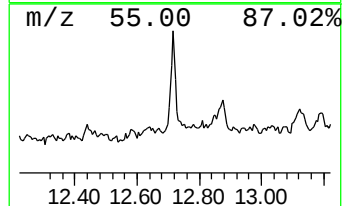
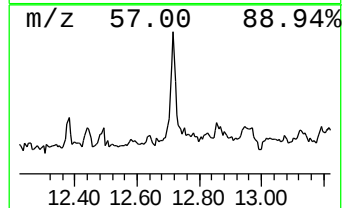
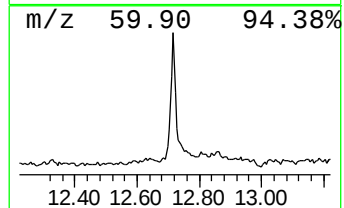
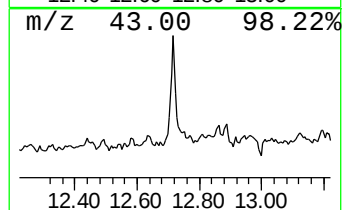
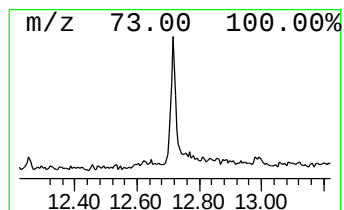
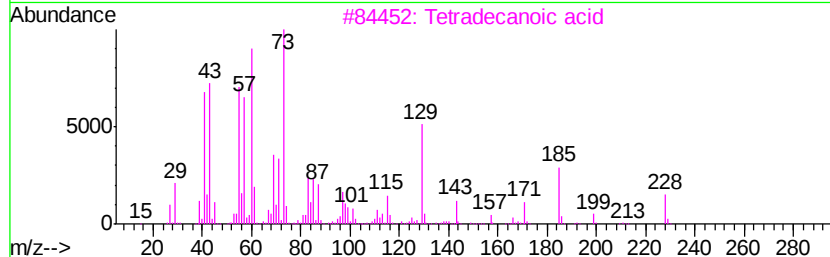
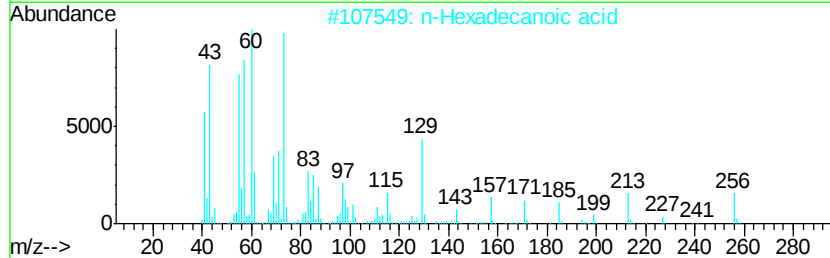
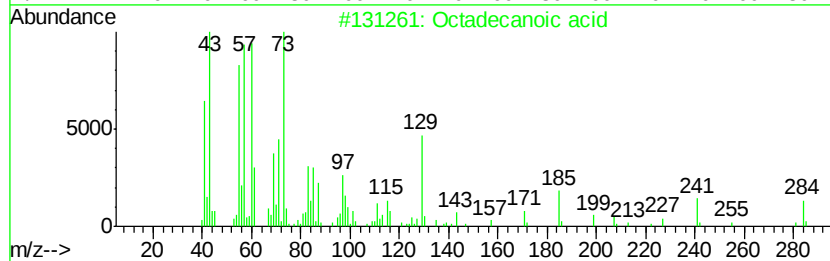
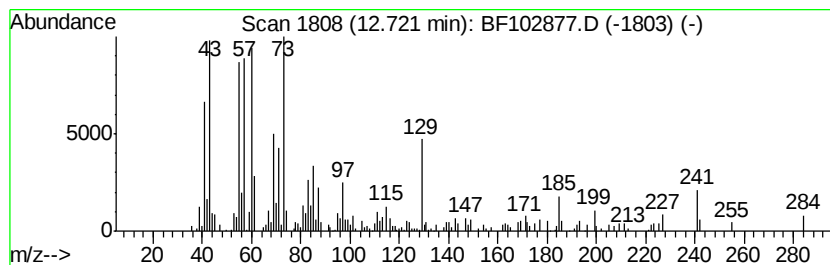
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TIC Integration Parameters: LSCINT.P

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Peak Number 17 Octadecanoic acid Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.72 | 4.02 ng | 296577 | Phenanthrene-d10 | 11.41 |

| Hit# | of | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------|-----|----------|-------------|------|
| 1    | 5  | Octadecanoic acid   | 284 | C18H36O2 | 000057-11-4 | 98   |
| 2    |    | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 87   |
| 3    |    | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 87   |
| 4    |    | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 78   |
| 5    |    | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 72   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020918\  
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Sample : J1471-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
3N-18

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

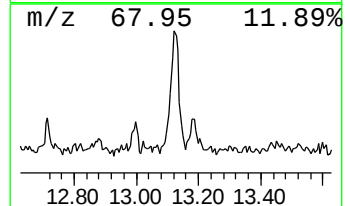
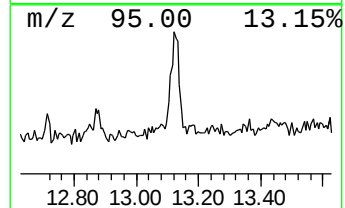
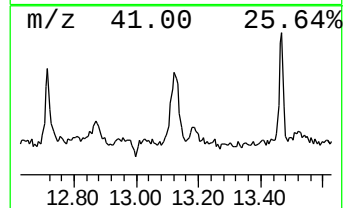
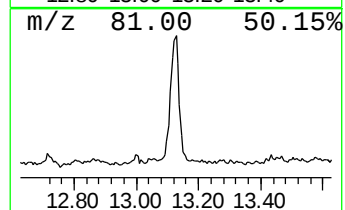
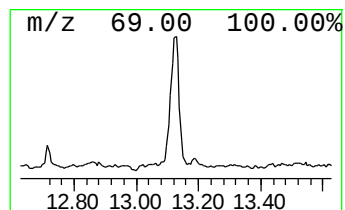
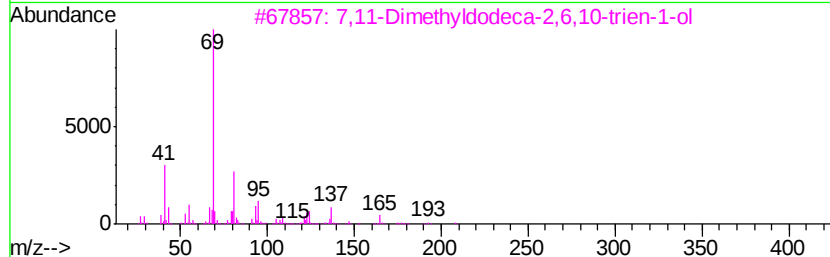
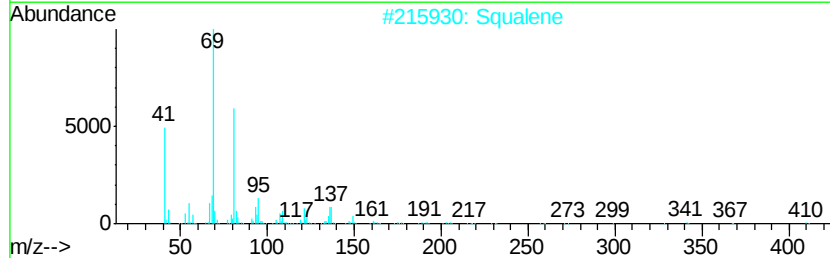
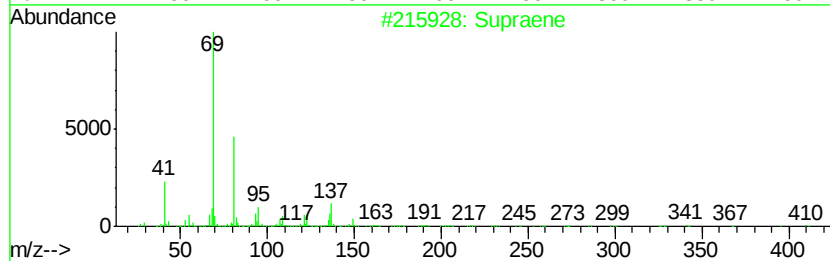
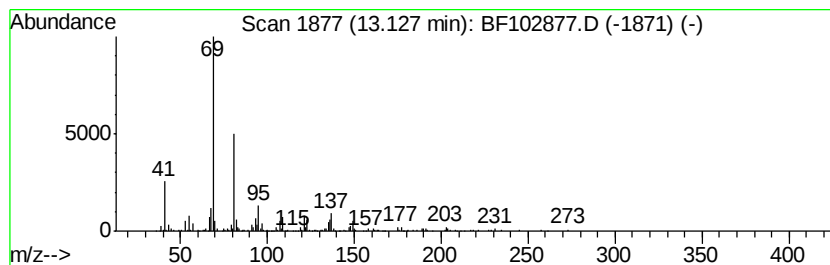
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Supraene Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.13 | 6.70 ng | 335748 | Chrysene-d12     | 14.05 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Supraene                            | 410 | C30H50     | 007683-64-9  | 91   |
| 2    |    | Squalene                            | 410 | C30H50     | 000111-02-4  | 91   |
| 3    |    | 7,11-Dimethyldodeca-2,6,10-trien... | 208 | C14H24O    | 125529-10-4  | 80   |
| 4    |    | 6,10-Dodecadien-1-yn-3-ol, 3,7,1... | 220 | C15H24O    | 002387-68-0  | 72   |
| 5    |    | (.+/-)-Lavandulol, pentafluorop...  | 300 | C13H17F5O2 | 1000352-63-1 | 64   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020918\  
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Acq On : 9 Feb 2018 10:02  
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Sample : J1471-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
3N-18

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

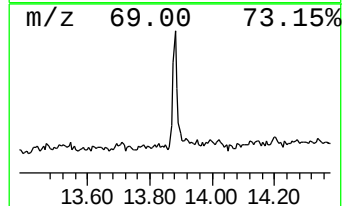
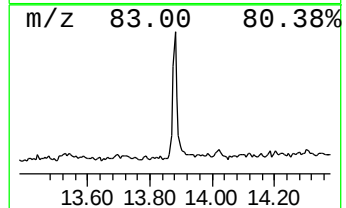
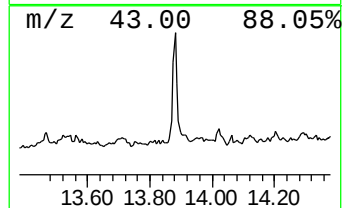
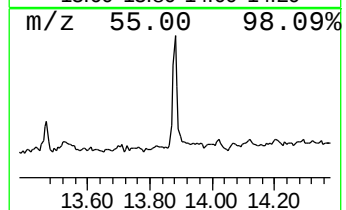
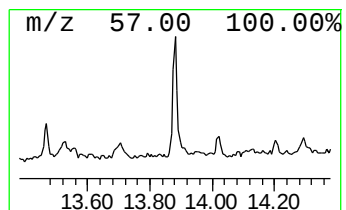
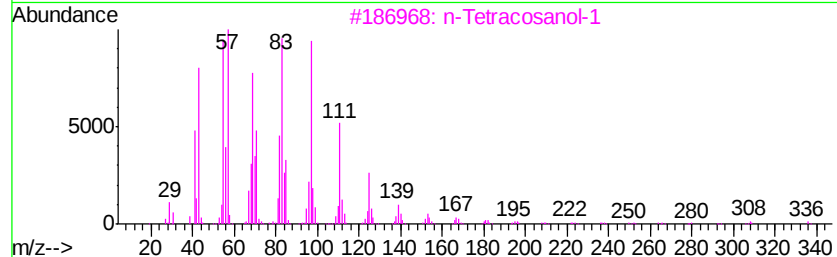
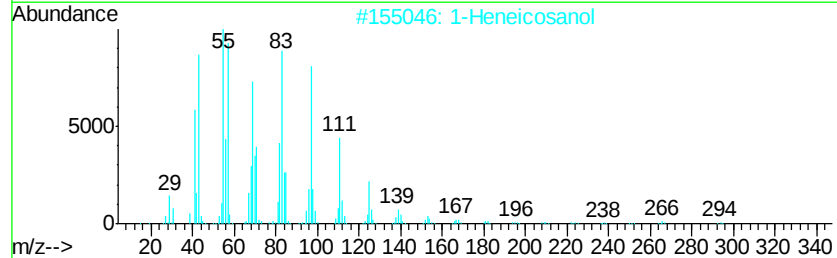
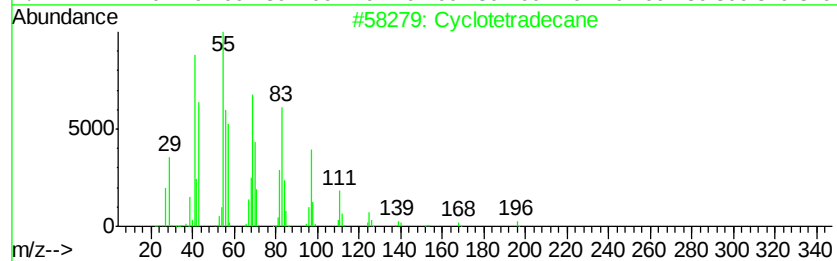
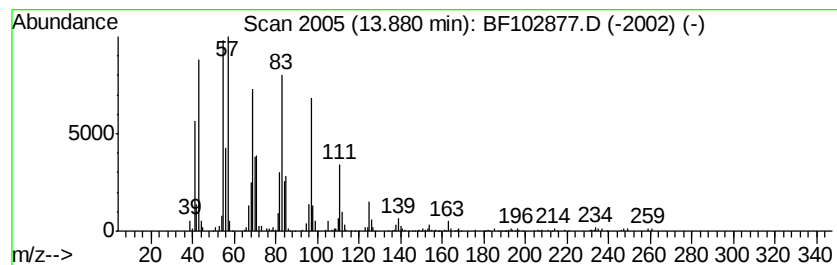
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Cyclotetradecane Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.88 | 6.56 ng | 328451 | Chrysene-d12     | 14.05 |

| Hit# | of | 5 | Tentative ID     | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclotetradecane | 196 | C14H28  | 000295-17-0 | 95   |
| 2    |    |   | 1-Heneicosanol   | 312 | C21H44O | 015594-90-8 | 94   |
| 3    |    |   | n-Tetracosanol-1 | 354 | C24H50O | 000506-51-4 | 91   |
| 4    |    |   | n-Nonadecanol-1  | 284 | C19H40O | 001454-84-8 | 91   |
| 5    |    |   | 1-Octadecanol    | 270 | C18H38O | 000112-92-5 | 91   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020918\  
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Acq On : 9 Feb 2018 10:02  
Operator : SJ/JU  
Sample : J1471-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
3N-18

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020318.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

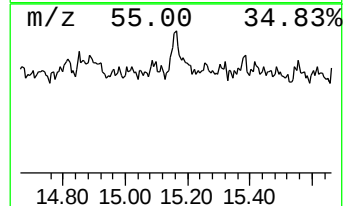
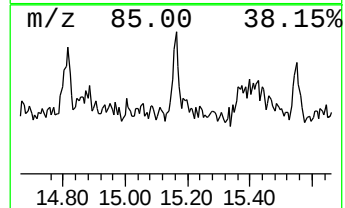
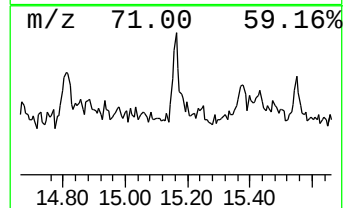
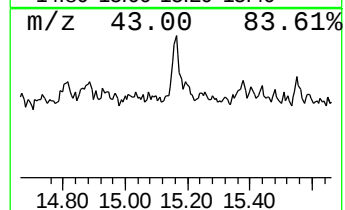
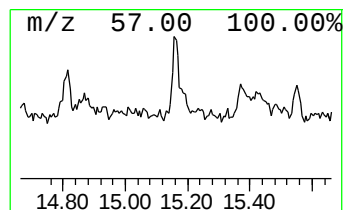
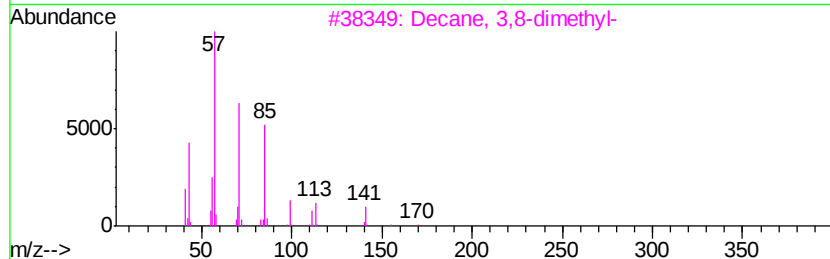
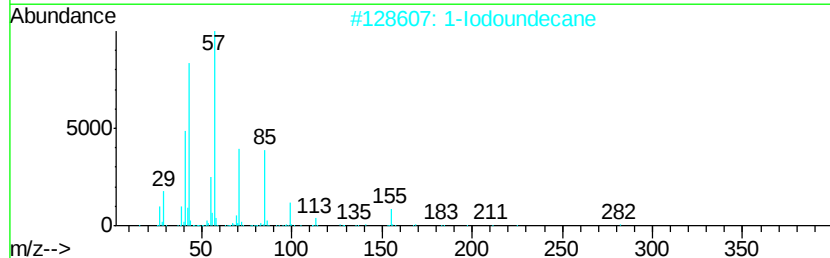
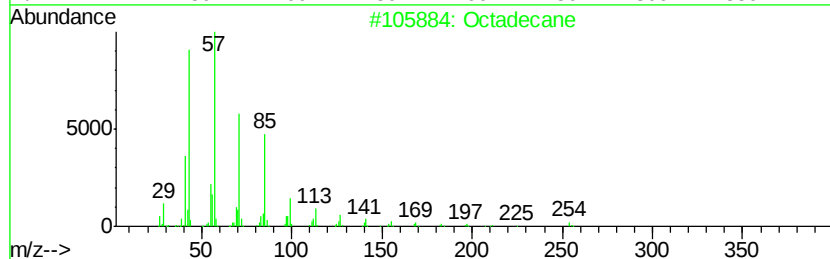
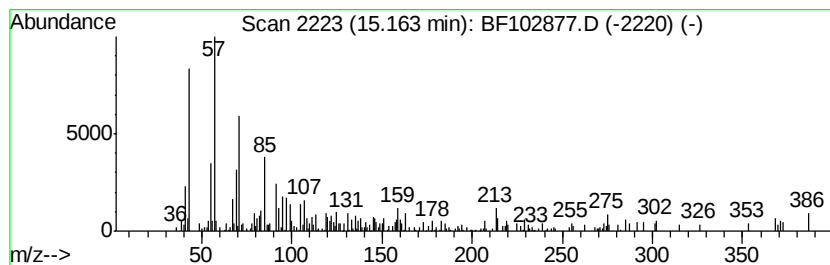
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Octadecane Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.16 | 4.40 ng | 227351 | Perylene-d12     | 15.54 |

| Hit# | of | Tentative ID           | MW  | MolForm  | CAS#        | Qual |
|------|----|------------------------|-----|----------|-------------|------|
| 1    | 5  | Octadecane             | 254 | C18H38   | 000593-45-3 | 60   |
| 2    |    | 1-Iodoundecane         | 282 | C11H23I  | 004282-44-4 | 60   |
| 3    |    | Decane, 3,8-dimethyl-  | 170 | C12H26   | 017312-55-9 | 53   |
| 4    |    | Tetradecane, 1-chloro- | 232 | C14H29Cl | 002425-54-9 | 53   |
| 5    |    | Nonadecane, 1-chloro-  | 302 | C19H39Cl | 062016-76-6 | 53   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020918\  
 Data File : BF102877.D  
 Acq On : 9 Feb 2018 10:02  
 Operator : SJ/JU  
 Sample : J1471-07  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 3N-18

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020318.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... | 2.16  | 14.6    | ng    | 818033   | 1                     | 6.89  | 1120810 | 20.0 |
| 2-Pentanone, 4-hy... | 5.12  | 3.3     | ng    | 184035   | 1                     | 6.89  | 1120810 | 20.0 |
| unknown6.66          | 6.66  | 83.6    | ng    | 4684670  | 1                     | 6.89  | 1120810 | 20.0 |
| Hexadecane           | 10.35 | 3.3     | ng    | 259890   | 3                     | 9.92  | 1563740 | 20.0 |
| 4-(1,1-Dimethylpr... | 10.89 | 6.6     | ng    | 483866   | 4                     | 11.41 | 1475470 | 20.0 |
| unknown10.92         | 10.92 | 8.3     | ng    | 611741   | 4                     | 11.41 | 1475470 | 20.0 |
| Phenol, m-tert-bu... | 10.95 | 7.3     | ng    | 535161   | 4                     | 11.41 | 1475470 | 20.0 |
| 4-Ethoxybenzhydra... | 10.98 | 3.8     | ng    | 277956   | 4                     | 11.41 | 1475470 | 20.0 |
| unknown11.07         | 11.07 | 9.3     | ng    | 683361   | 4                     | 11.41 | 1475470 | 20.0 |
| Phenol, 4-(1,1,3,... | 11.10 | 7.3     | ng    | 541517   | 4                     | 11.41 | 1475470 | 20.0 |
| Acetamide, N-(3-m... | 11.15 | 4.1     | ng    | 304163   | 4                     | 11.41 | 1475470 | 20.0 |
| n-Hexadecanoic acid  | 11.93 | 11.1    | ng    | 815131   | 4                     | 11.41 | 1475470 | 20.0 |
| unknown12.10         | 12.10 | 3.5     | ng    | 258783   | 4                     | 11.41 | 1475470 | 20.0 |
| Octadecanoic acid    | 12.72 | 4.0     | ng    | 296577   | 4                     | 11.41 | 1475470 | 20.0 |
| Supraene             | 13.13 | 6.7     | ng    | 335748   | 5                     | 14.05 | 1001590 | 20.0 |
| Cyclotetradecane     | 13.88 | 6.6     | ng    | 328451   | 5                     | 14.05 | 1001590 | 20.0 |
| Octadecane           | 15.16 | 4.4     | ng    | 227351   | 6                     | 15.54 | 1033790 | 20.0 |