

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA\_F\DATA\BF022818\  
 Data File : BF103317.D  
 Acq On : 1 Mar 2018 2:04  
 Operator : SJ/JU  
 Sample : J1705-03  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 100MVAR

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-BF022218.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         | 2.152       | 6             | 11          | 20           | rVB      | 228280         | 296875        | 8.71%           | 1.103%        |
| 2         | 5.110       | 510           | 514         | 533          | rVB      | 98451          | 138423        | 4.06%           | 0.515%        |
| 3         | 5.498       | 576           | 580         | 599          | rBV      | 3005444        | 3407053       | 100.00%         | 12.664%       |
| 4         | 6.510       | 747           | 752         | 771          | rBV      | 3102681        | 3353339       | 98.42%          | 12.464%       |
| 5         | 6.645       | 771           | 775         | 779          | rBV      | 3325087        | 3177436       | 93.26%          | 11.810%       |
| 6         | 6.875       | 810           | 814         | 821          | rBV      | 1056378        | 907942        | 26.65%          | 3.375%        |
| 7         | 7.028       | 836           | 840         | 844          | rBV      | 2502750        | 2428650       | 71.28%          | 9.027%        |
| 8         | 7.439       | 905           | 910         | 925          | rBV      | 2204646        | 2149017       | 63.08%          | 7.988%        |
| 9         | 8.157       | 1028          | 1032        | 1049         | rBV      | 1213830        | 1154301       | 33.88%          | 4.290%        |
| 10        | 9.227       | 1210          | 1214        | 1224         | rBV      | 2744438        | 2772558       | 81.38%          | 10.305%       |
| 11        | 9.298       | 1224          | 1226        | 1230         | rVB      | 56147          | 52442         | 1.54%           | 0.195%        |
| 12        | 9.622       | 1276          | 1281        | 1289         | rBV      | 192591         | 227413        | 6.67%           | 0.845%        |
| 13        | 9.827       | 1313          | 1316        | 1321         | rBV      | 119733         | 114227        | 3.35%           | 0.425%        |
| 14        | 9.916       | 1327          | 1331        | 1339         | rBV      | 1036873        | 986230        | 28.95%          | 3.666%        |
| 15        | 10.327      | 1398          | 1401        | 1404         | rVB      | 136582         | 107217        | 3.15%           | 0.399%        |
| 16        | 10.551      | 1434          | 1439        | 1441         | rBV      | 35446          | 40104         | 1.18%           | 0.149%        |
| 17        | 10.710      | 1462          | 1466        | 1474         | rBV      | 1198435        | 1309455       | 38.43%          | 4.867%        |
| 18        | 10.804      | 1479          | 1482        | 1492         | rVB      | 84541          | 101670        | 2.98%           | 0.378%        |
| 19        | 11.404      | 1581          | 1584        | 1588         | rBV      | 757189         | 773506        | 22.70%          | 2.875%        |
| 20        | 12.621      | 1788          | 1791        | 1796         | rBV      | 119860         | 119014        | 3.49%           | 0.442%        |
| 21        | 12.857      | 1826          | 1831        | 1838         | rVB      | 108974         | 127018        | 3.73%           | 0.472%        |
| 22        | 12.992      | 1850          | 1854        | 1857         | rBV      | 1740588        | 1744434       | 51.20%          | 6.484%        |
| 23        | 13.874      | 2000          | 2004        | 2008         | rBV2     | 95818          | 126053        | 3.70%           | 0.469%        |
| 24        | 14.057      | 2031          | 2035        | 2038         | rBV      | 725318         | 743437        | 21.82%          | 2.763%        |
| 25        | 15.557      | 2284          | 2290        | 2299         | rVB      | 355761         | 545948        | 16.02%          | 2.029%        |

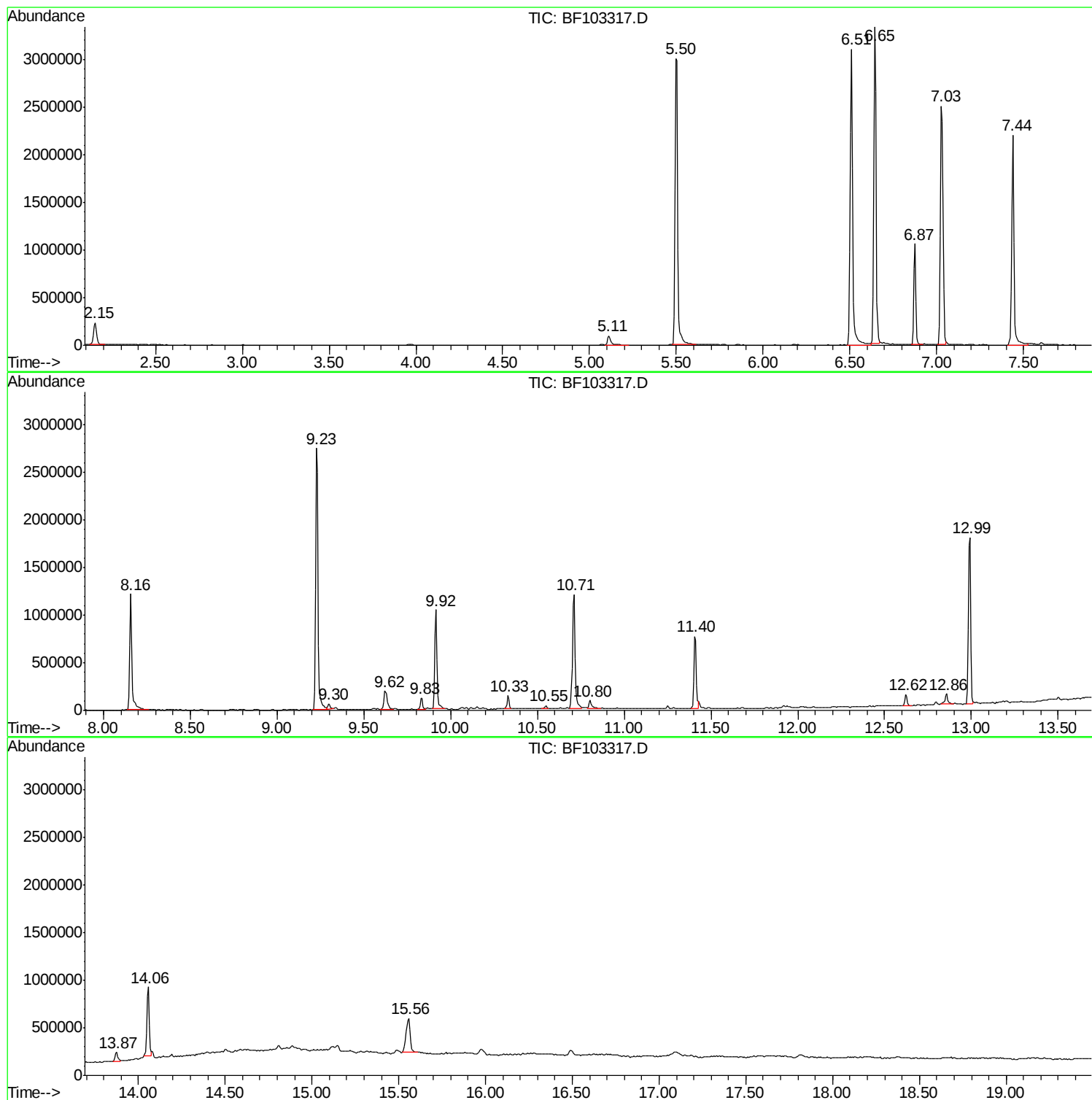
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF022218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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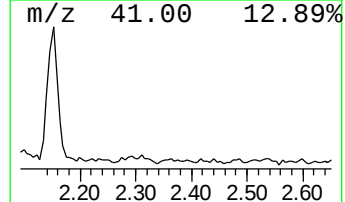
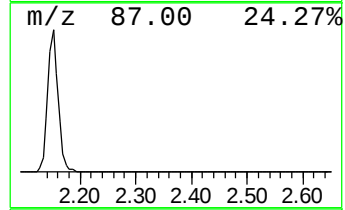
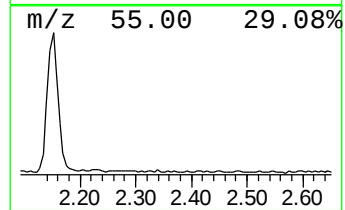
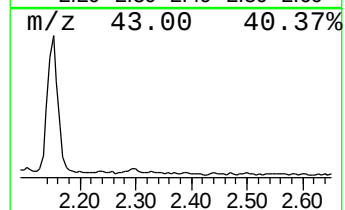
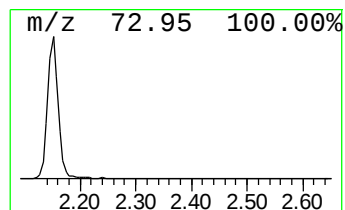
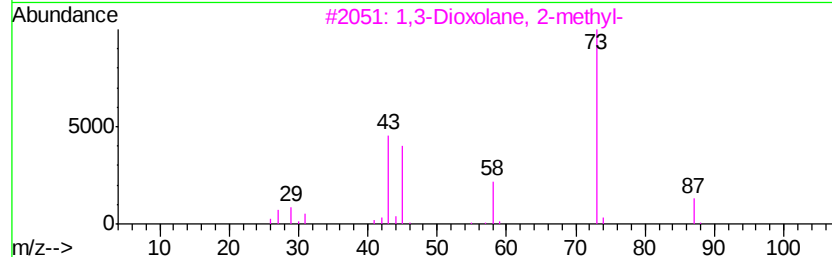
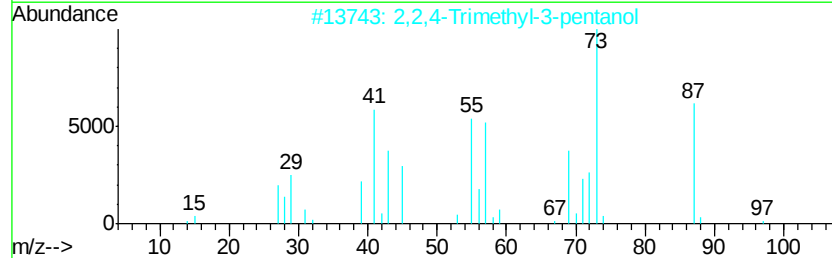
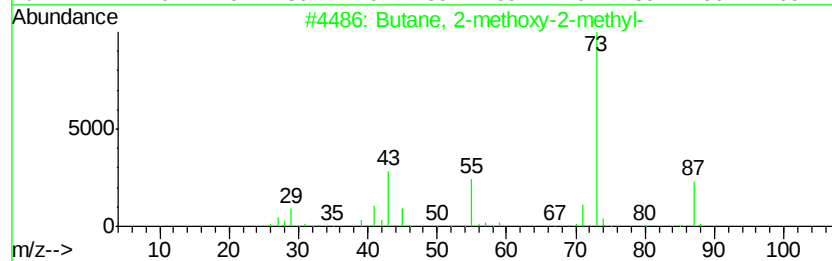
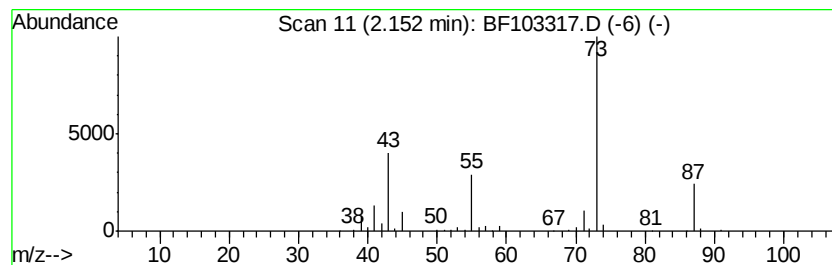
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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. |
|------|---------|--------|------------------------|------|
| 2.15 | 6.54 ng | 296875 | 1,4-Dichlorobenzene-d4 | 6.87 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Butane, 2-methoxy-2-methyl-         | 102 | C6H14O   | 000994-05-8 | 83   |
| 2    |    |   | 2,2,4-Trimethyl-3-pentanol          | 130 | C8H18O   | 005162-48-1 | 39   |
| 3    |    |   | 1,3-Dioxolane, 2-methyl-            | 88  | C4H8O2   | 000497-26-7 | 37   |
| 4    |    |   | Boronic acid, ethyl-, dimethyl e... | 102 | C4H11BO2 | 007318-82-3 | 17   |
| 5    |    |   | Silane, tetramethyl-                | 88  | C4H12Si  | 000075-76-3 | 12   |



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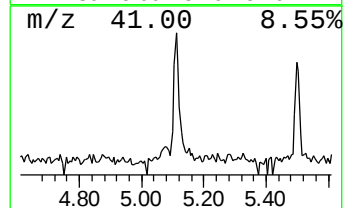
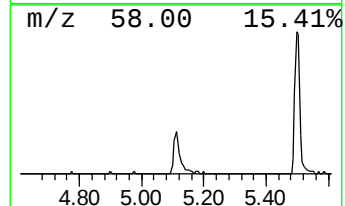
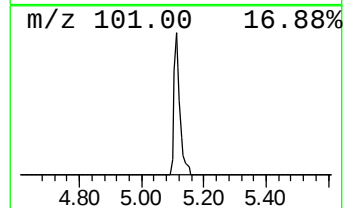
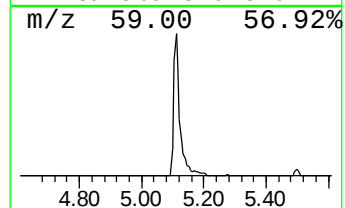
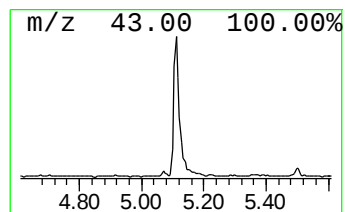
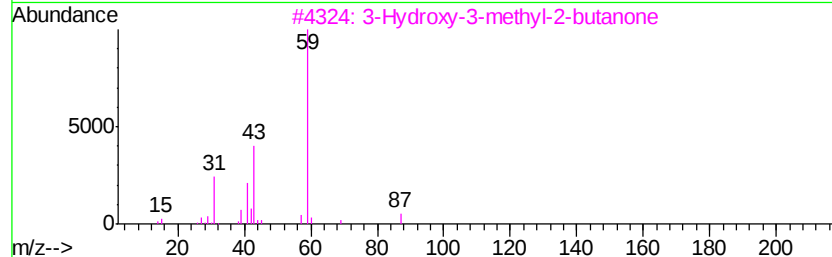
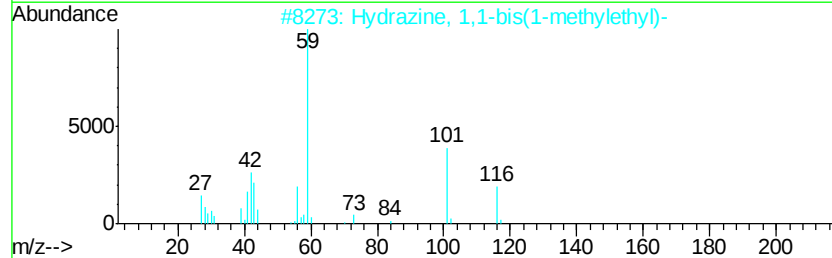
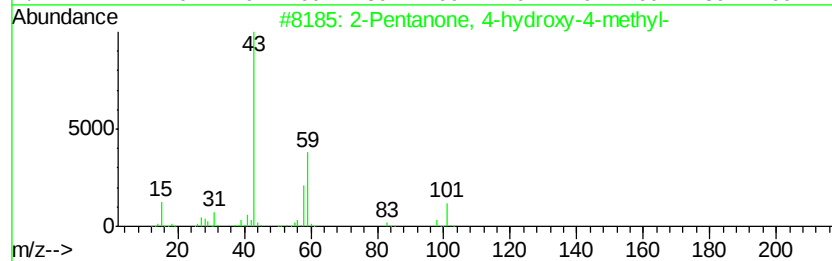
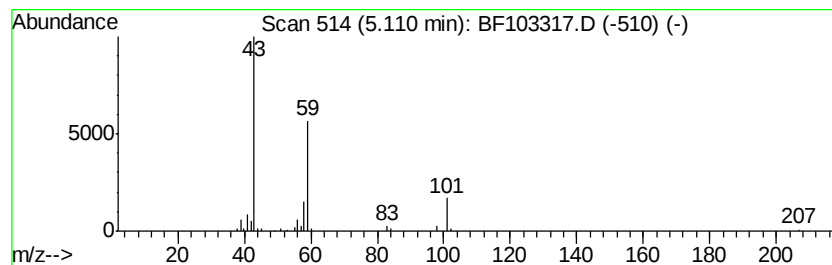
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 5.11 | 3.05 ng | 138423 | 1,4-Dichlorobenzene-d4 | 6.87 |

| Hit# of | 5                                  | Tentative ID | MW      | MolForm     | CAS# | Qual |
|---------|------------------------------------|--------------|---------|-------------|------|------|
| 1       | 2-Pentanone, 4-hydroxy-4-methyl-   | 116          | C6H12O2 | 000123-42-2 | 50   |      |
| 2       | Hydrazine, 1,1-bis(1-methylethyl)- | 116          | C6H16N2 | 000921-14-2 | 25   |      |
| 3       | 3-Hydroxy-3-methyl-2-butanone      | 102          | C5H10O2 | 000115-22-0 | 23   |      |
| 4       | Morpholine, 4-methyl-              | 101          | C5H11NO | 000109-02-4 | 9    |      |
| 5       | 2,3-Butanedione, monooxime         | 101          | C4H7NO2 | 000057-71-6 | 9    |      |



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100MVAR

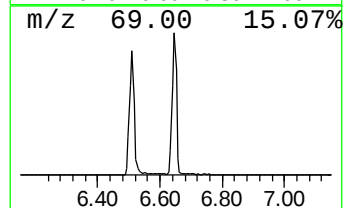
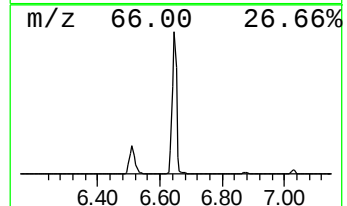
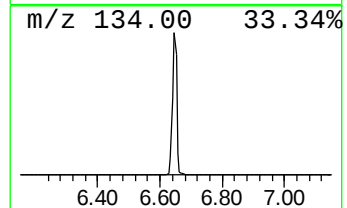
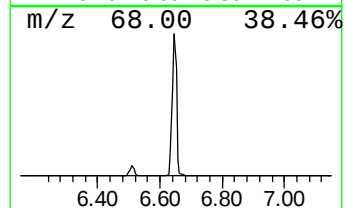
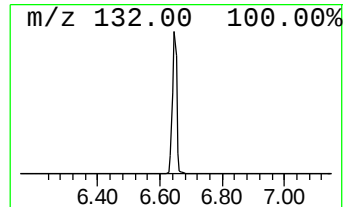
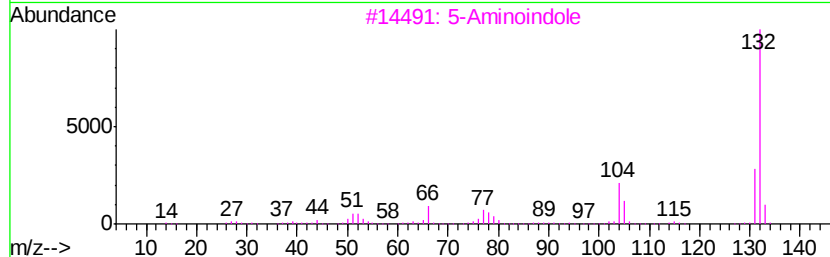
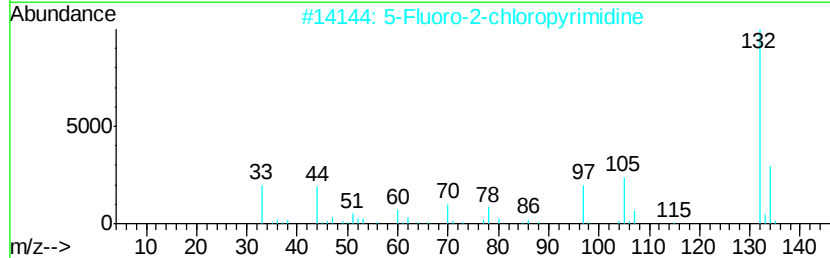
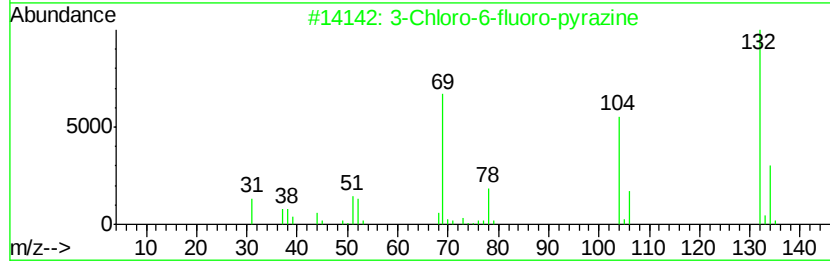
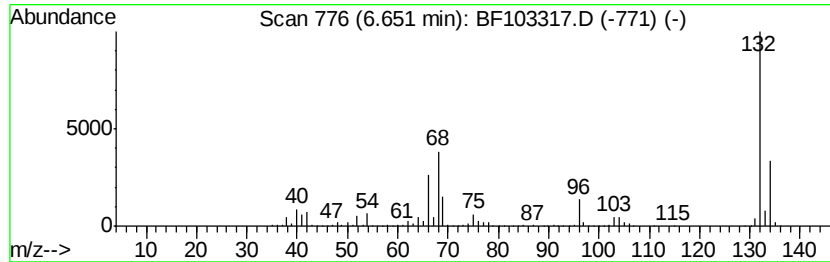
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TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 unknown6.65 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.65 | 69.99 ng | 3177440 | 1,4-Dichlorobenzene-d4 | 6.87 |

| 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    |    | Tranylcypromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 10   |
| 5    |    | Benzene, 1-ethenyl-3,5-dimethyl- | 132 | C10H12    | 005379-20-4  | 9    |



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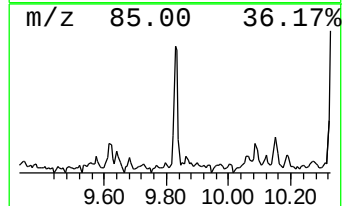
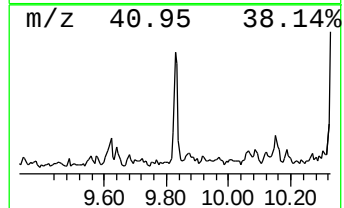
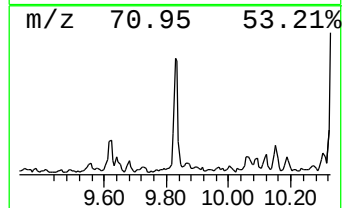
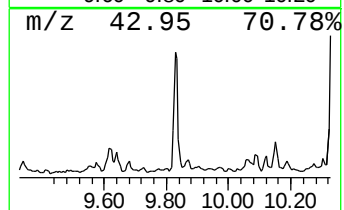
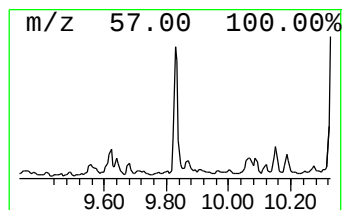
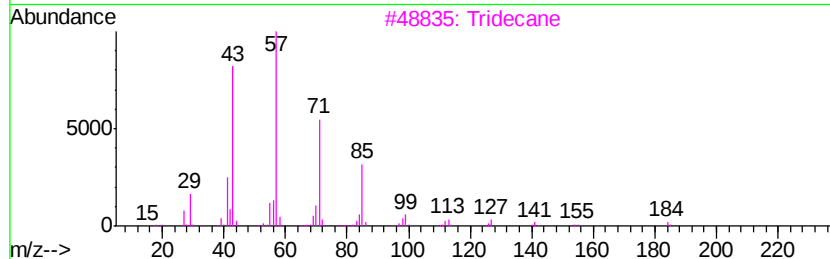
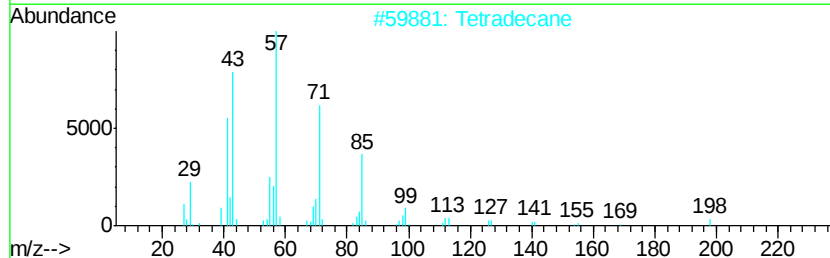
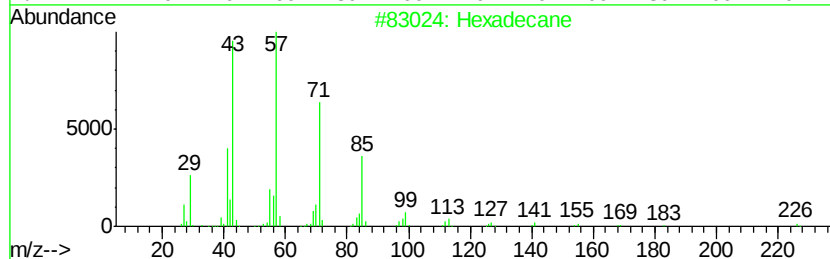
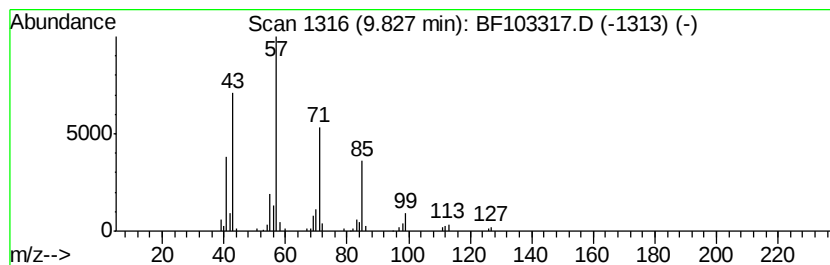
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Peak Number 5 Hexadecane Concentration Rank 7

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.83 | 2.32 ng | 114227 | Acenaphthene-d10 | 9.92 |

| Hit# of | 5                                   | Tentative ID | MW  | MolForm   | CAS#         | Qual |
|---------|-------------------------------------|--------------|-----|-----------|--------------|------|
| 1       | Hexadecane                          |              | 226 | C16H34    | 000544-76-3  | 90   |
| 2       | Tetradecane                         |              | 198 | C14H30    | 000629-59-4  | 86   |
| 3       | Tridecane                           |              | 184 | C13H28    | 000629-50-5  | 72   |
| 4       | Sulfurous acid, 2-ethylhexyl hex... |              | 278 | C14H30O3S | 1000309-20-2 | 59   |
| 5       | Heptadecane                         |              | 240 | C17H36    | 000629-78-7  | 53   |



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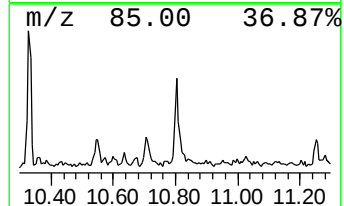
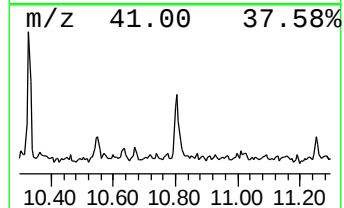
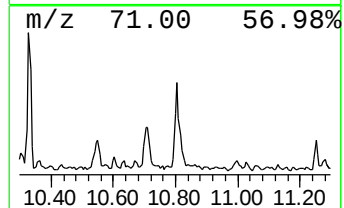
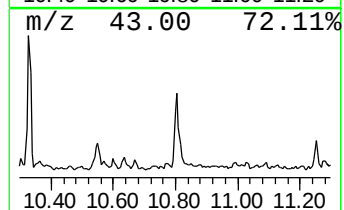
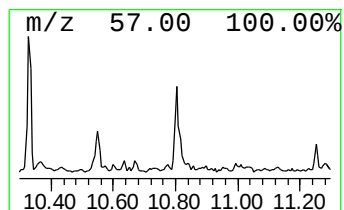
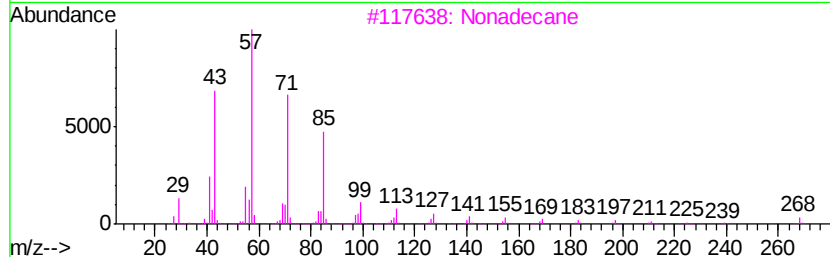
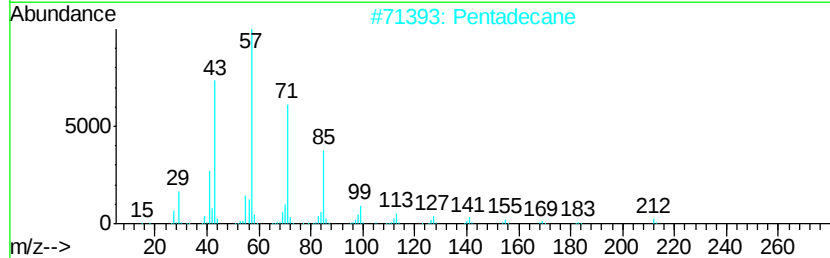
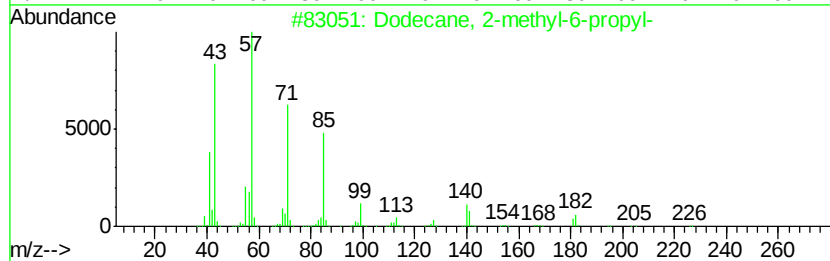
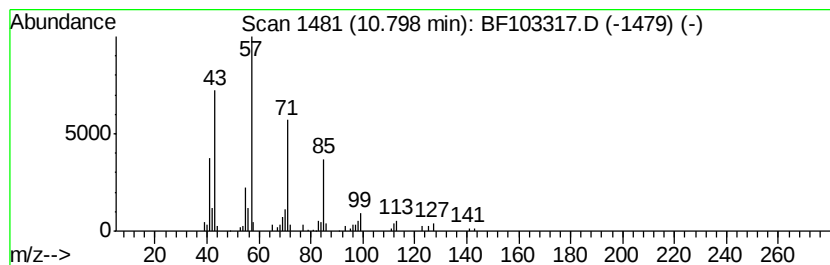
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ClientSampleId :  
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Peak Number 7 Dodecane, 2-methyl-6-propyl- Concentration Rank 6

| R.T.      | EstConc                      | Area   | Relative to ISTD | R.T.        |      |
|-----------|------------------------------|--------|------------------|-------------|------|
| 10.80     | 2.63 ng                      | 101670 | Phenanthrene-d10 | 11.40       |      |
| Hit# of 5 | Tentative ID                 | MW     | MolForm          | CAS#        | Qual |
| 1         | Dodecane, 2-methyl-6-propyl- | 226    | C16H34           | 055045-08-4 | 86   |
| 2         | Pentadecane                  | 212    | C15H32           | 000629-62-9 | 86   |
| 3         | Nonadecane                   | 268    | C19H40           | 000629-92-5 | 86   |
| 4         | Heneicosane                  | 296    | C21H44           | 000629-94-7 | 86   |
| 5         | Dodecane, 1-iodo-            | 296    | C12H25I          | 004292-19-7 | 86   |



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF022818\  
Data File : BF103317.D  
Acq On : 1 Mar 2018 2:04  
Operator : SJ/JU  
Sample : J1705-03  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

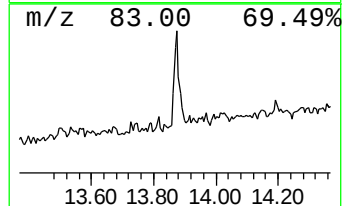
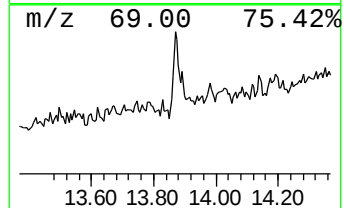
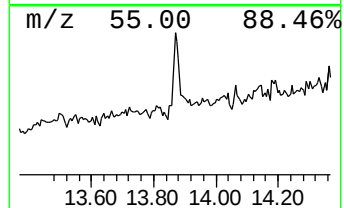
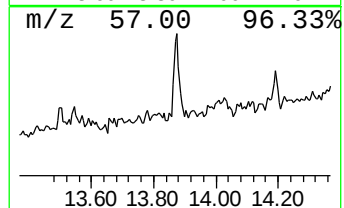
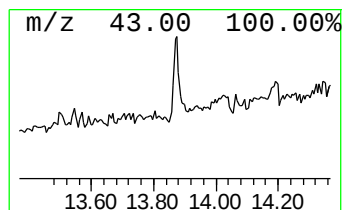
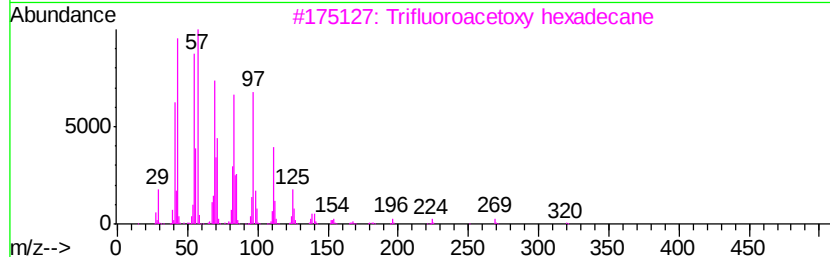
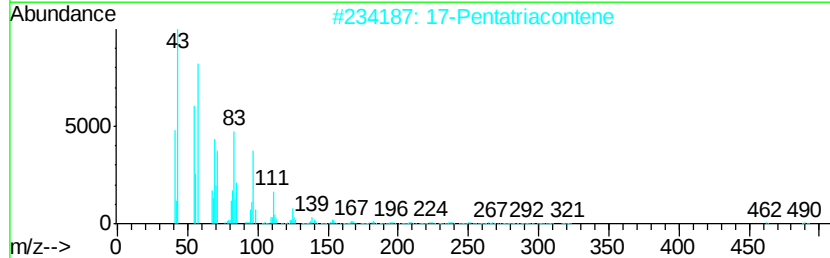
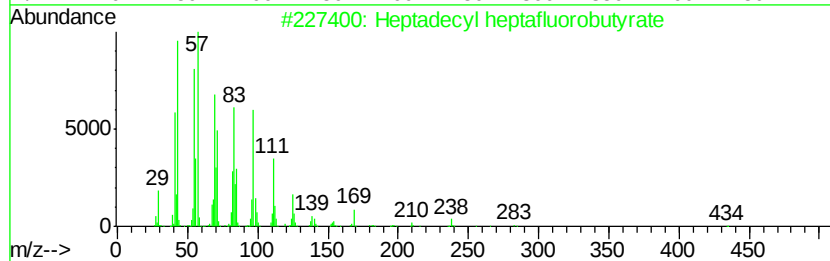
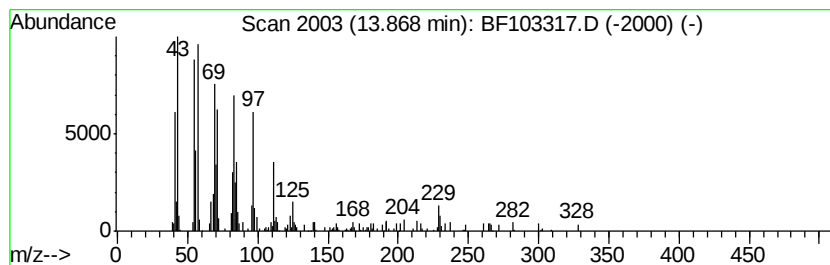
Instrument :  
BNA\_F  
ClientSampleId :  
100MVAR

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF022218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 8 Heptadecyl heptafluorobutyrate Concentration Rank 4

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 13.87     | 3.39 ng                             | 126053 | Chrysene-d12     | 14.06       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Heptadecyl heptafluorobutvrate      | 452    | C21H35F7O2       | 959085-66-6 | 91   |
| 2         | 17-Pentatriacontene                 | 491    | C35H70           | 006971-40-0 | 74   |
| 3         | Trifluoroacetoxv hexadecane         | 338    | C18H33F3O2       | 006222-03-3 | 70   |
| 4         | Octadecyl trifluoroacetate          | 366    | C20H37F3O2       | 079392-43-1 | 68   |
| 5         | Pentafluoropropionic acid, octad... | 416    | C21H37F5O2       | 959261-25-7 | 68   |



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA\_F\DATA\BF022818\  
Data File : BF103317.D  
Acq On : 1 Mar 2018 2:04  
Operator : SJ/JU  
Sample : J1705-03  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
100MVAR

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF022218.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... | 2.15  | 6.5     | ng    | 296875   | 1                       | 6.87  | 907942 | 20.0 |
| 2-Pentanone, 4-hy... | 5.11  | 3.0     | ng    | 138423   | 1                       | 6.87  | 907942 | 20.0 |
| unknown6.65          | 6.65  | 70.0    | ng    | 3177440  | 1                       | 6.87  | 907942 | 20.0 |
| Hexadecane           | 9.83  | 2.3     | ng    | 114227   | 3                       | 9.92  | 986230 | 20.0 |
| Dodecane, 2-methy... | 10.80 | 2.6     | ng    | 101670   | 4                       | 11.40 | 773506 | 20.0 |
| Heptadecyl heptaf... | 13.87 | 3.4     | ng    | 126053   | 5                       | 14.06 | 743437 | 20.0 |