

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF080918\  
 Data File : BF108073.D  
 Acq On : 10 Aug 2018 1:35  
 Operator : JU/SJ  
 Sample : J4378-02  
 Misc :  
 ALS Vial : 33 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 080618-TIPC-E-U-M

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF080818.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         | 2.366       | 3             | 5           | 13           | rVB      | 2073722        | 2250948       | 24.11%          | 3.957%        |
| 2         | 5.248       | 492           | 495         | 502          | rVB      | 207534         | 217043        | 2.32%           | 0.382%        |
| 3         | 5.637       | 556           | 561         | 564          | rBV      | 3491993        | 3201551       | 34.29%          | 5.628%        |
| 4         | 6.625       | 724           | 729         | 732          | rBV      | 2534690        | 2234214       | 23.93%          | 3.928%        |
| 5         | 6.784       | 751           | 756         | 759          | rBV      | 5773244        | 5241102       | 56.13%          | 9.214%        |
| 6         | 7.013       | 792           | 795         | 799          | rVB      | 1907023        | 1686140       | 18.06%          | 2.964%        |
| 7         | 7.172       | 818           | 822         | 826          | rVB      | 6802867        | 6894805       | 73.84%          | 12.121%       |
| 8         | 7.578       | 885           | 891         | 894          | rBV      | 5449405        | 5847211       | 62.62%          | 10.279%       |
| 9         | 8.301       | 1009          | 1014        | 1017         | rBV      | 2139919        | 2015411       | 21.58%          | 3.543%        |
| 10        | 9.378       | 1191          | 1197        | 1200         | rBV      | 9559194        | 9337890       | 100.00%         | 16.415%       |
| 11        | 9.760       | 1259          | 1262        | 1266         | rBV      | 335397         | 298622        | 3.20%           | 0.525%        |
| 12        | 10.060      | 1310          | 1313        | 1317         | rVB2     | 2190500        | 1926066       | 20.63%          | 3.386%        |
| 13        | 10.724      | 1422          | 1426        | 1429         | rBV      | 726461         | 556265        | 5.96%           | 0.978%        |
| 14        | 10.854      | 1443          | 1448        | 1457         | rBV      | 3346189        | 3018015       | 32.32%          | 5.306%        |
| 15        | 11.560      | 1564          | 1568        | 1571         | rBV      | 1866474        | 1668991       | 17.87%          | 2.934%        |
| 16        | 13.148      | 1833          | 1838        | 1841         | rBV      | 7563879        | 6955398       | 74.49%          | 12.227%       |
| 17        | 14.042      | 1986          | 1990        | 2001         | rBV3     | 98957          | 254017        | 2.72%           | 0.447%        |
| 18        | 14.213      | 2015          | 2019        | 2023         | rBV      | 2074510        | 1821885       | 19.51%          | 3.203%        |
| 19        | 15.771      | 2279          | 2284        | 2291         | rBV      | 995864         | 1358396       | 14.55%          | 2.388%        |
| 20        | 16.259      | 2363          | 2367        | 2372         | rBV      | 62874          | 100647        | 1.08%           | 0.177%        |

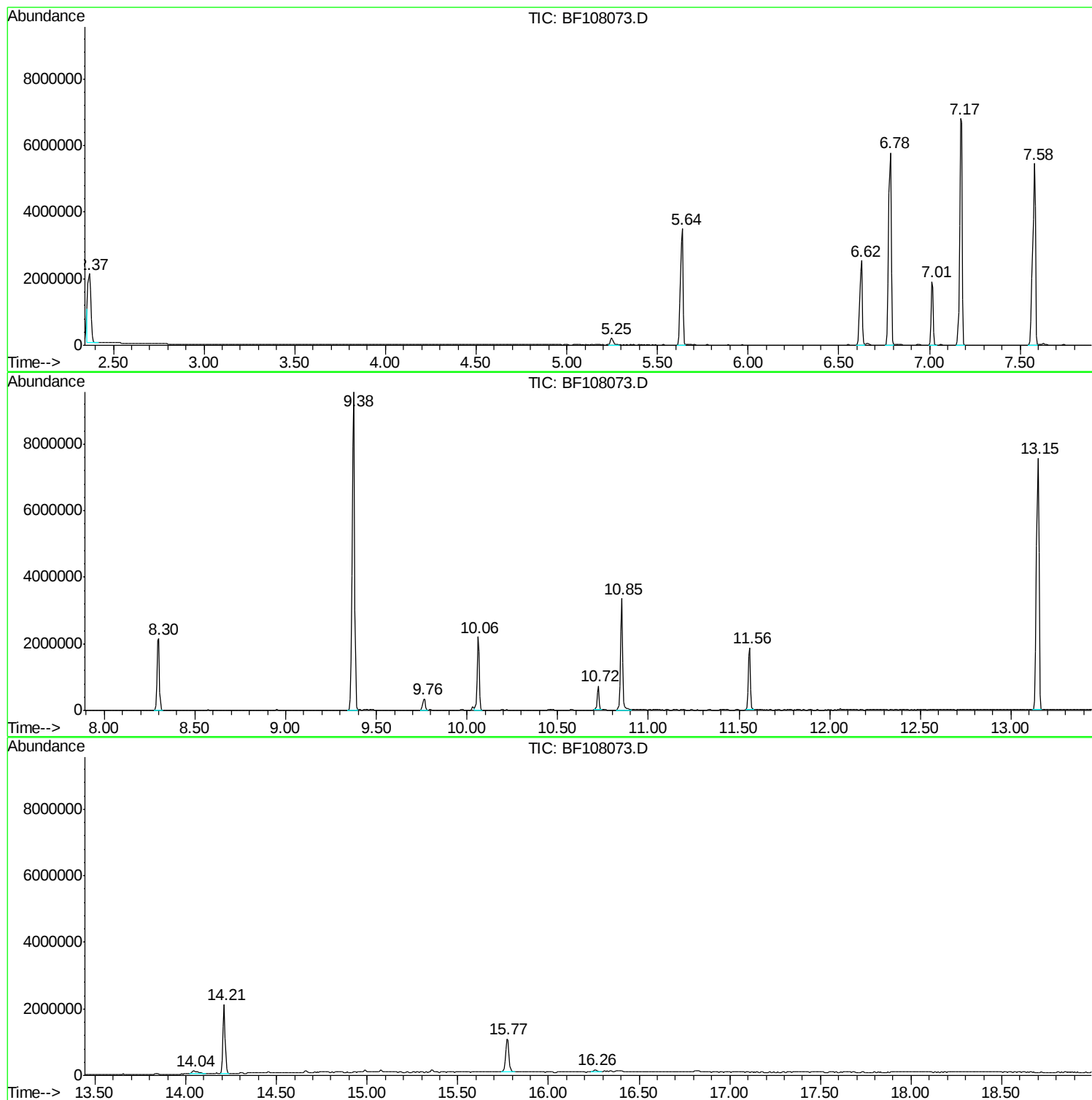
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080618-TIPC-E-U-M

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.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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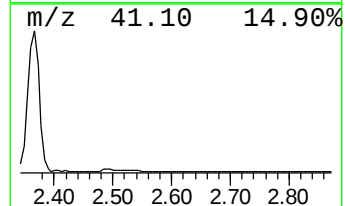
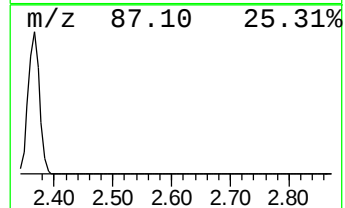
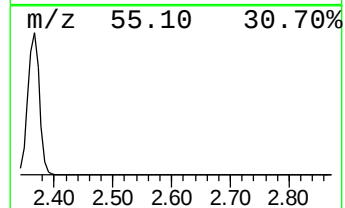
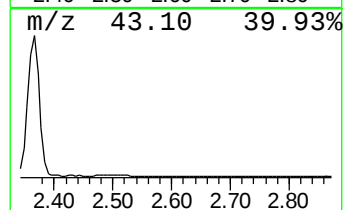
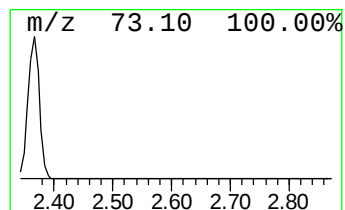
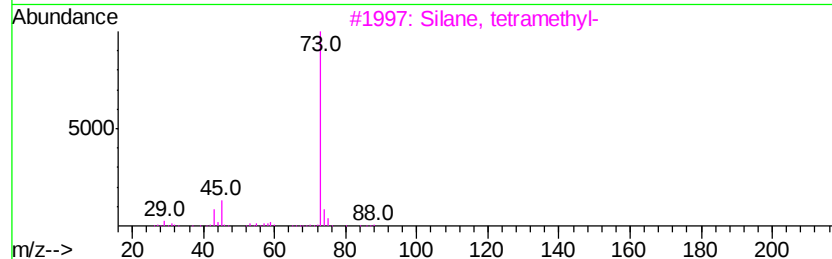
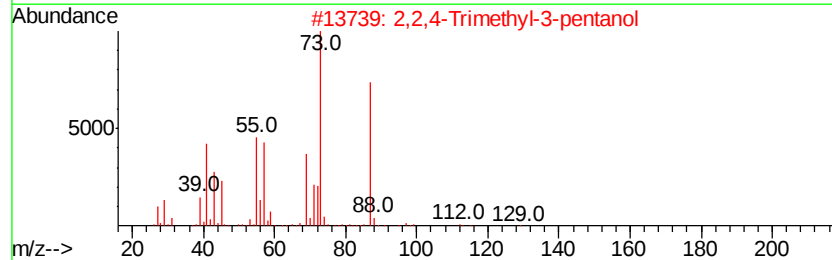
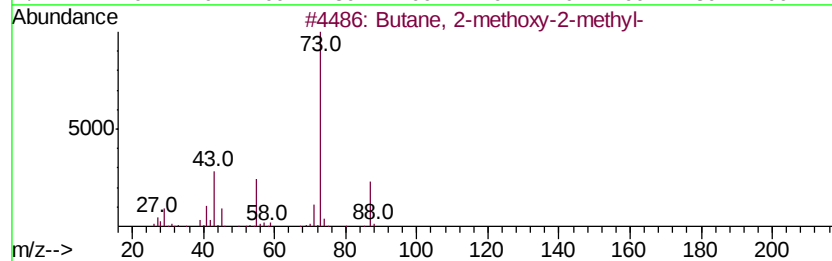
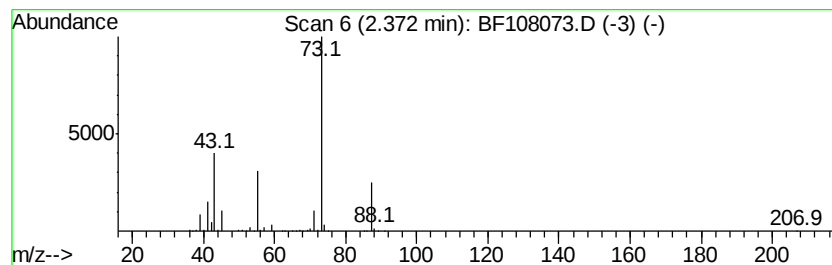
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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. |
|------|----------|---------|------------------------|------|
| 2.37 | 26.70 ng | 2250950 | 1,4-Dichlorobenzene-d4 | 7.01 |

| 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    |    |   | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 38   |
| 4    |    |   | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 25   |
| 5    |    |   | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 12   |



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

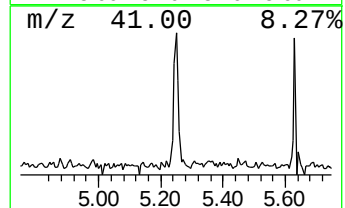
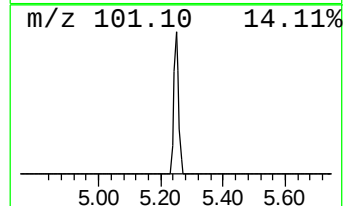
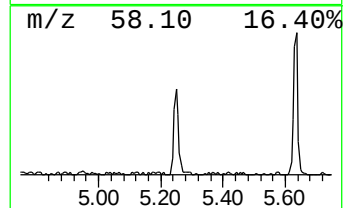
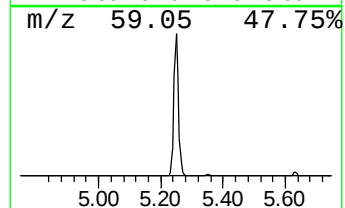
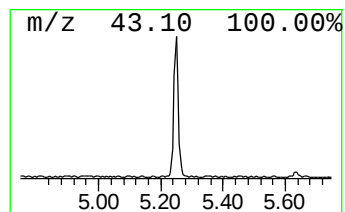
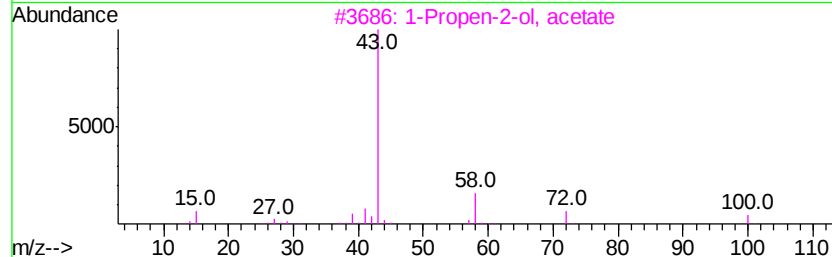
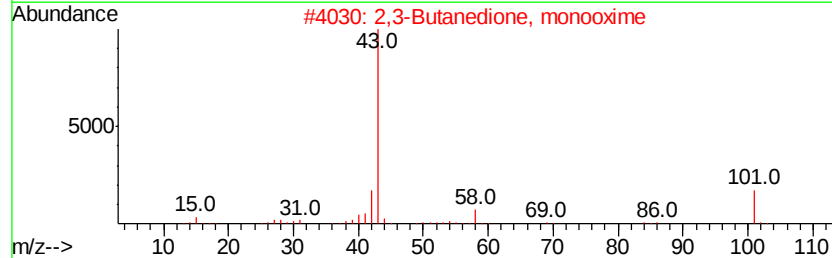
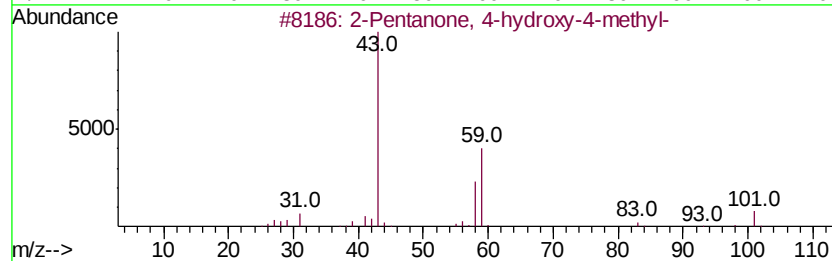
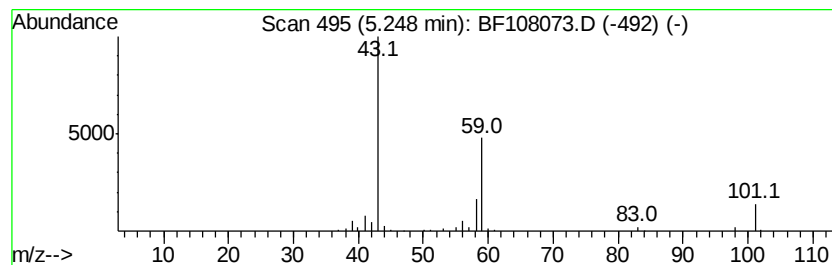
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 5.25 | 2.57 ng | 217043 | 1,4-Dichlorobenzene-d4 | 7.01 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 50   |
| 2    |    | 2,3-Butanedione, monooxime       | 101 | C4H7NO2 | 000057-71-6 | 10   |
| 3    |    | 1-Propen-2-ol, acetate           | 100 | C5H8O2  | 000108-22-5 | 10   |
| 4    |    | Diacetamide                      | 101 | C4H7NO2 | 000625-77-4 | 9    |
| 5    |    | Allyl acetate                    | 100 | C5H8O2  | 000591-87-7 | 9    |



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

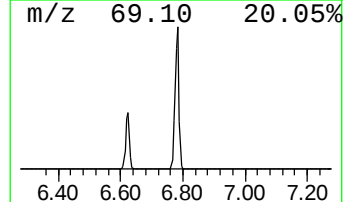
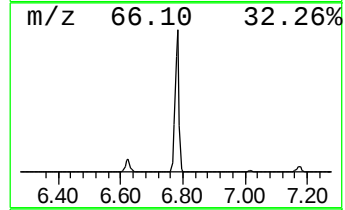
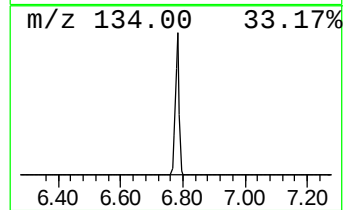
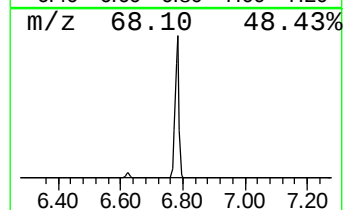
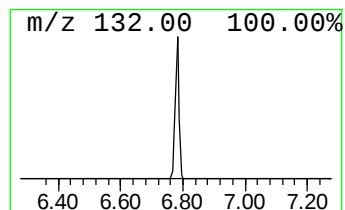
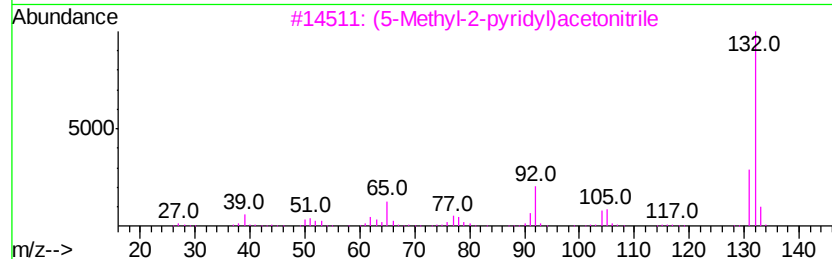
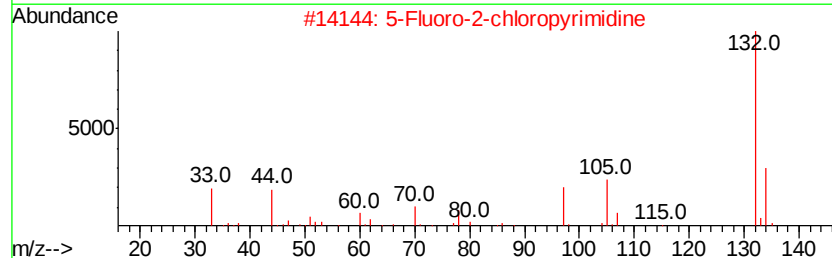
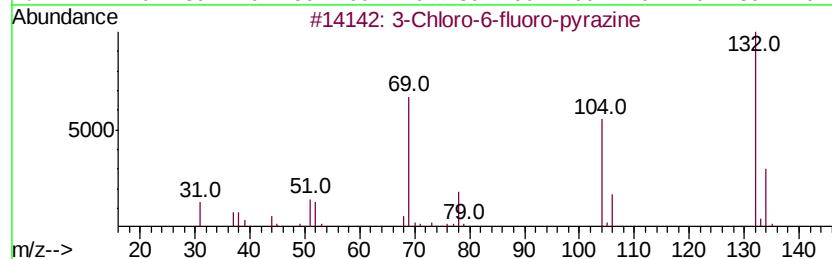
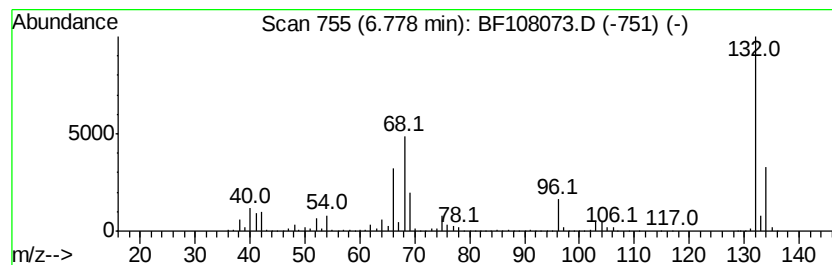
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 unknown6.78 Concentration Rank 2

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.78 | 62.17 ng | 5241100 | 1,4-Dichlorobenzene-d4 | 7.01 |

| 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-Methyl-2-pyridyl)acetonitrile | 132 | C8H8N2    | 1000241-93-9 | 10   |
| 4    |    | 1,3,5-Trifluorobenzene           | 132 | C6H3F3    | 000372-38-3  | 9    |
| 5    |    | Xenon                            | 132 | Xe        | 007440-63-3  | 9    |



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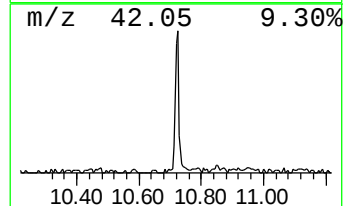
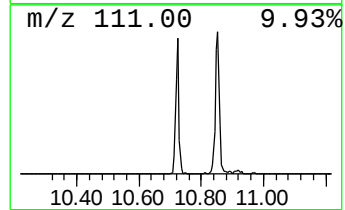
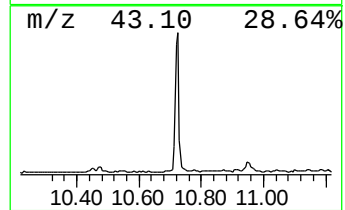
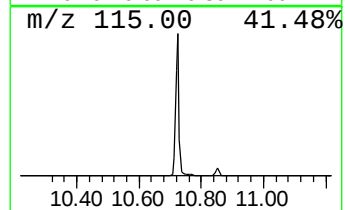
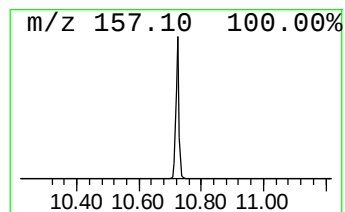
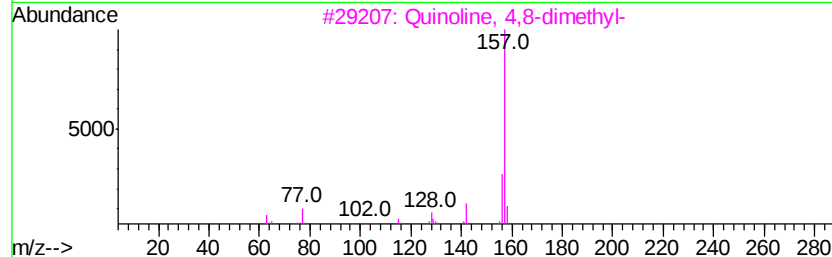
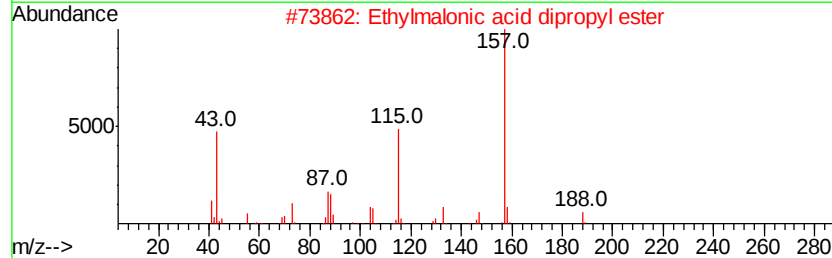
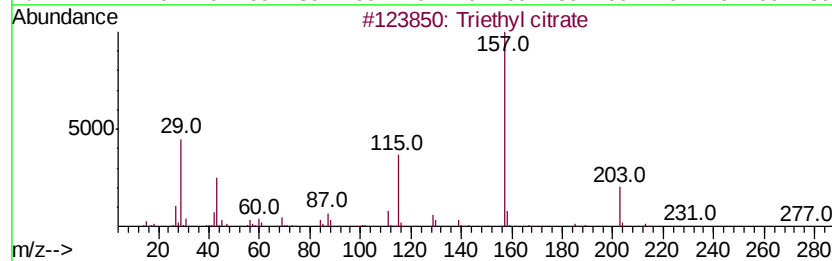
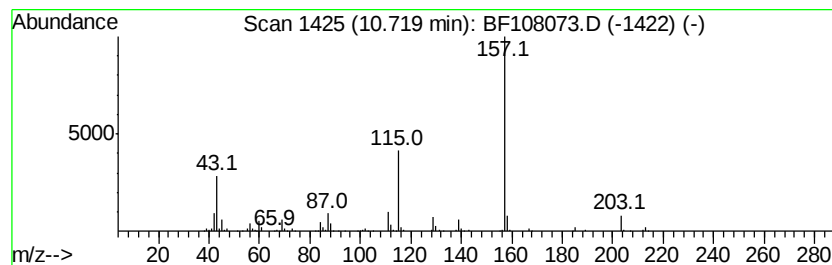
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TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Triethyl citrate Concentration Rank 4

| R.T.    | EstConc | Area                                | Relative to ISTD |               | R.T.         |      |
|---------|---------|-------------------------------------|------------------|---------------|--------------|------|
| 10.72   | 5.78 ng | 556265                              | Acenaphthene-d10 |               | 10.06        |      |
| Hit# of | 5       | Tentative ID                        | MW               | MolForm       | CAS#         | Qual |
| 1       |         | Triethyl citrate                    | 276              | C12H20O7      | 000077-93-0  | 86   |
| 2       |         | Ethylmalonic acid dipropyl ester    | 216              | C11H20O4      | 001113-91-3  | 64   |
| 3       |         | Quinoline, 4,8-dimethyl-            | 157              | C11H11N       | 013362-80-6  | 40   |
| 4       |         | Piperazine-1-carboxylic acid, 4-... | 362              | C14H19ClN2O5S | 1000303-79-9 | 40   |
| 5       |         | 3-Chloro2-fluorobenzoic acid, 4-... | 356              | C20H30ClF02   | 1000338-64-4 | 40   |



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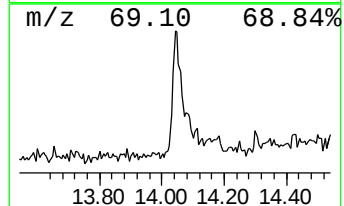
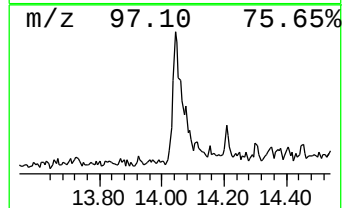
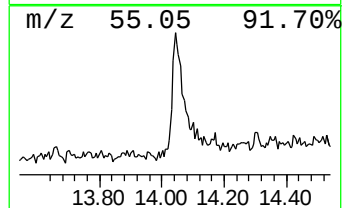
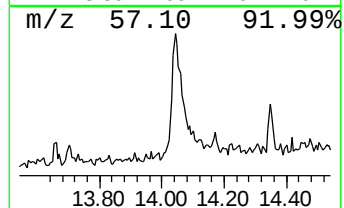
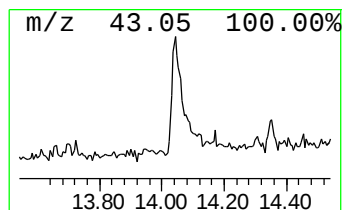
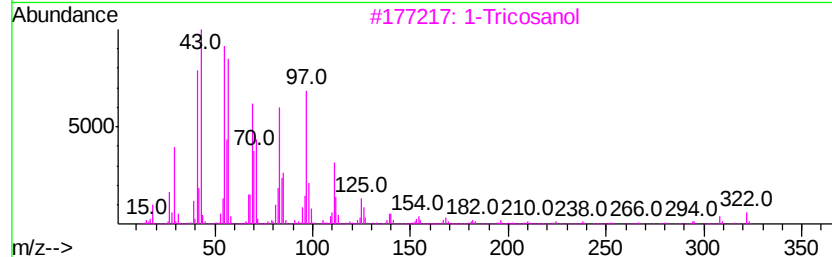
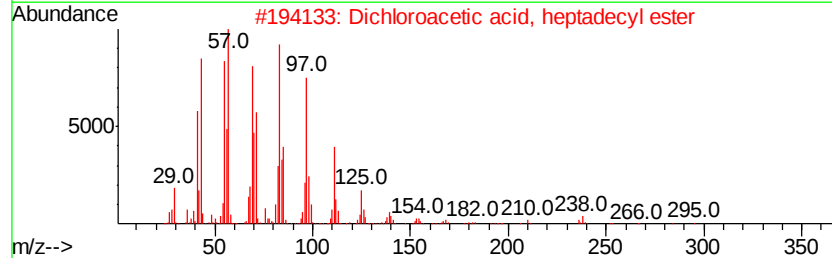
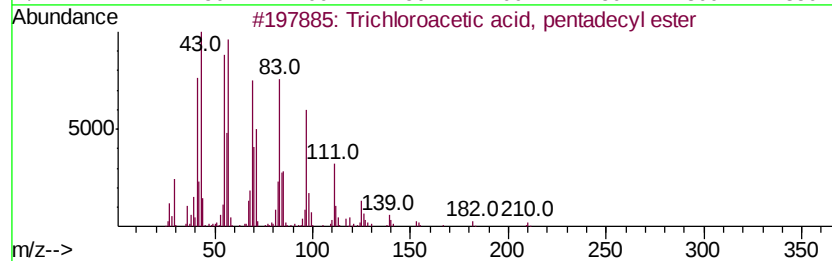
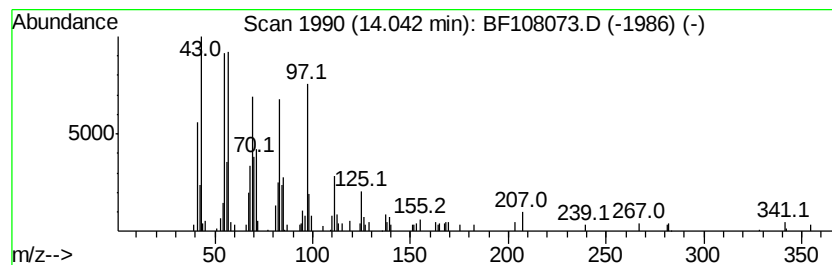
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Trichloroacetic acid, penta... Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.04 | 2.79 ng | 254017 | Chrysene-d12     | 14.21 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|---------|---|-------------------------------------|-----|-------------|--------------|------|
| 1       |   | Trichloroacetic acid, pentadecyl... | 372 | C17H31Cl3O2 | 074339-53-0  | 93   |
| 2       |   | Dichloroacetic acid, heptadecyl ... | 366 | C19H36Cl2O2 | 1000282-98-2 | 90   |
| 3       |   | 1-Tricosanol                        | 340 | C23H48O     | 003133-01-5  | 87   |
| 4       |   | 1-Eicosanol                         | 298 | C20H42O     | 000629-96-9  | 87   |
| 5       |   | 1-Docosene                          | 308 | C22H44      | 001599-67-3  | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF080918\  
Data File : BF108073.D  
Acq On : 10 Aug 2018 1:35  
Operator : JU/SJ  
Sample : J4378-02  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
080618-TIPC-E-U-M

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF080818.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... | 2.37  | 26.7    | ng    | 2250950  | 1                     | 7.01  | 1686140 | 20.0 |
| 2-Pentanone, 4-hy... | 5.25  | 2.6     | ng    | 217043   | 1                     | 7.01  | 1686140 | 20.0 |
| unknown6.78          | 6.78  | 62.2    | ng    | 5241100  | 1                     | 7.01  | 1686140 | 20.0 |
| Triethyl citrate     | 10.72 | 5.8     | ng    | 556265   | 3                     | 10.06 | 1926070 | 20.0 |
| Trichloroacetic a... | 14.04 | 2.8     | ng    | 254017   | 5                     | 14.21 | 1821890 | 20.0 |