

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF071024\  
 Data File : BF138507.D  
 Acq On : 10 Jul 2024 18:14  
 Operator : CG\JU  
 Sample : P3170-01  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SOIL-DRUMS-1

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-BF070924.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF138507.D\data.ms

| 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         | 2.175       | 7             | 15          | 20           | rBV      | 1688984        | 2219511       | 95.98%          | 9.337%        |
| 2         | 3.393       | 220           | 222         | 235          | rBV      | 30970          | 74040         | 3.20%           | 0.311%        |
| 3         | 5.087       | 506           | 510         | 516          | rVB      | 146756         | 169473        | 7.33%           | 0.713%        |
| 4         | 5.492       | 575           | 579         | 582          | rBV      | 2362922        | 2106856       | 91.11%          | 8.863%        |
| 5         | 6.498       | 745           | 750         | 753          | rBV      | 2330336        | 2155357       | 93.20%          | 9.067%        |
| 6         | 6.863       | 808           | 812         | 816          | rBV      | 681364         | 532119        | 23.01%          | 2.239%        |
| 7         | 7.428       | 903           | 908         | 911          | rBV      | 1530804        | 1431913       | 61.92%          | 6.024%        |
| 8         | 8.145       | 1025          | 1030        | 1033         | rBV      | 810653         | 663151        | 28.68%          | 2.790%        |
| 9         | 8.451       | 1078          | 1082        | 1086         | rBV      | 38776          | 43100         | 1.86%           | 0.181%        |
| 10        | 9.028       | 1176          | 1180        | 1182         | rBV3     | 41368          | 59287         | 2.56%           | 0.249%        |
| 11        | 9.098       | 1190          | 1192        | 1194         | rBV3     | 40030          | 47597         | 2.06%           | 0.200%        |
| 12        | 9.139       | 1196          | 1199        | 1200         | rVV3     | 73151          | 66109         | 2.86%           | 0.278%        |
| 13        | 9.163       | 1200          | 1203        | 1207         | rVB4     | 47112          | 52627         | 2.28%           | 0.221%        |
| 14        | 9.222       | 1207          | 1213        | 1216         | rBV      | 2684200        | 2312505       | 100.00%         | 9.728%        |
| 15        | 9.263       | 1216          | 1220        | 1222         | rBV2     | 40872          | 50712         | 2.19%           | 0.213%        |
| 16        | 9.298       | 1222          | 1226        | 1229         | rBV      | 215492         | 268181        | 11.60%          | 1.128%        |
| 17        | 9.492       | 1257          | 1259        | 1261         | rBV2     | 90701          | 97795         | 4.23%           | 0.411%        |
| 18        | 9.604       | 1276          | 1278        | 1283         | rBV2     | 648403         | 783365        | 33.88%          | 3.295%        |
| 19        | 9.663       | 1286          | 1288        | 1292         | rVB2     | 264250         | 305453        | 13.21%          | 1.285%        |
| 20        | 9.763       | 1302          | 1305        | 1308         | rBV2     | 709857         | 804111        | 34.77%          | 3.383%        |
| 21        | 9.810       | 1310          | 1313        | 1316         | rVB      | 1427951        | 1329660       | 57.50%          | 5.594%        |
| 22        | 9.851       | 1317          | 1320        | 1322         | rBV2     | 389450         | 405118        | 17.52%          | 1.704%        |
| 23        | 9.880       | 1323          | 1325        | 1326         | rVV2     | 410588         | 372906        | 16.13%          | 1.569%        |
| 24        | 9.898       | 1326          | 1328        | 1332         | rVB      | 843169         | 808491        | 34.96%          | 3.401%        |
| 25        | 9.945       | 1332          | 1336        | 1338         | rBV2     | 430665         | 654233        | 28.29%          | 2.752%        |
| 26        | 10.157      | 1370          | 1372        | 1375         | rBV3     | 487771         | 543703        | 23.51%          | 2.287%        |
| 27        | 10.239      | 1384          | 1386        | 1389         | rVB3     | 419054         | 282697        | 12.22%          | 1.189%        |
| 28        | 10.310      | 1396          | 1398        | 1402         | rBV      | 924417         | 784943        | 33.94%          | 3.302%        |
| 29        | 10.692      | 1461          | 1463        | 1466         | rBV      | 714508         | 652126        | 28.20%          | 2.743%        |
| 30        | 10.822      | 1483          | 1485        | 1488         | rVB      | 735160         | 665566        | 28.78%          | 2.800%        |
| 31        | 11.398      | 1581          | 1583        | 1587         | rBV3     | 634805         | 596566        | 25.80%          | 2.510%        |
| 32        | 12.974      | 1848          | 1851        | 1855         | rVB      | 1503910        | 1386000       | 59.94%          | 5.831%        |
| 33        | 14.027      | 2027          | 2030        | 2037         | rVB      | 501176         | 547124        | 23.66%          | 2.302%        |
| 34        | 15.486      | 2273          | 2278        | 2289         | rVB      | 392737         | 498707        | 21.57%          | 2.098%        |

Sum of corrected areas: 23771102

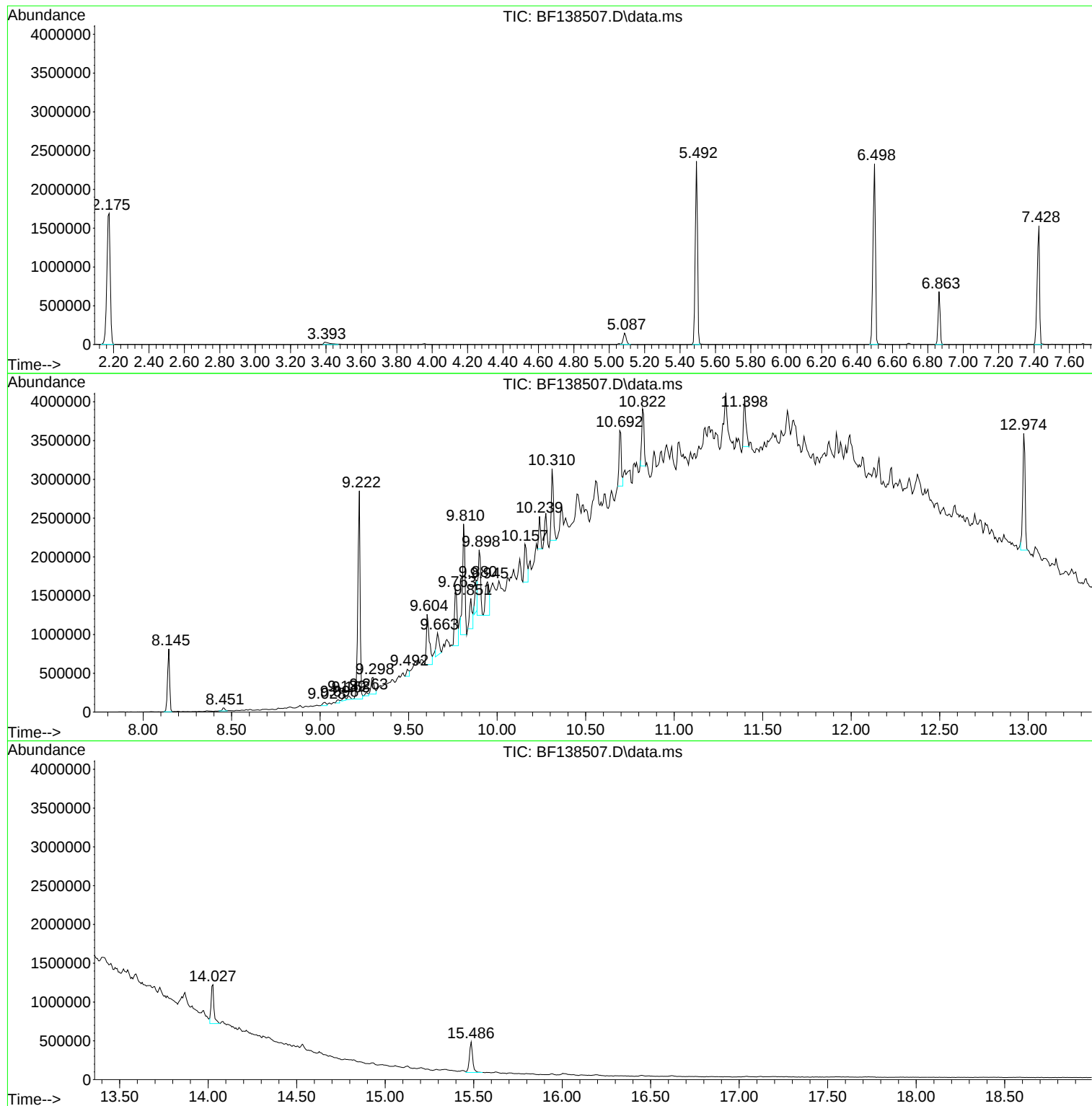
**Instrument :**  
BNA\_F  
**ClientSampleId :**  
SOIL-DRUMS-1

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Misc :  
ALS Vial : 16 Sample Multiplier: 1

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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF070924.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF071024\  
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SOIL-DRUMS-1

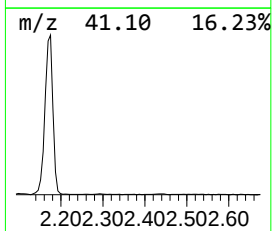
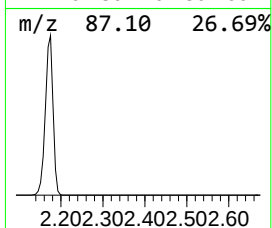
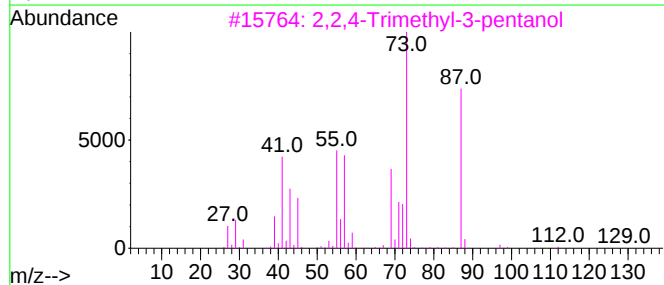
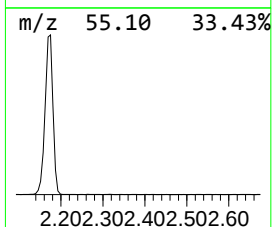
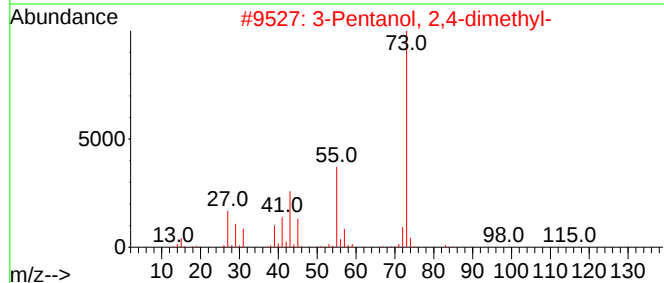
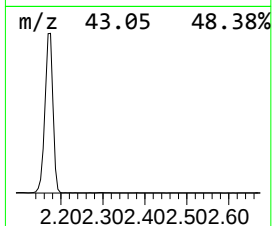
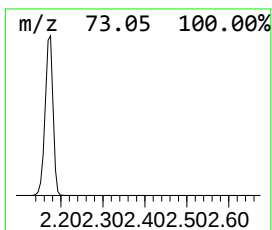
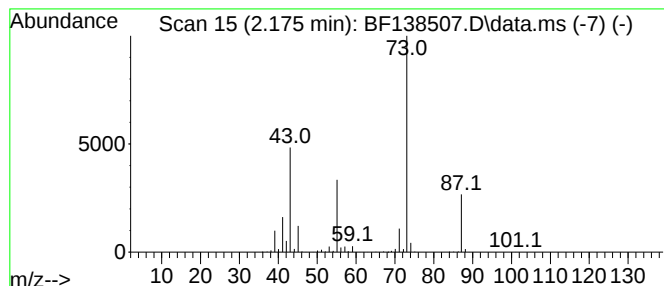
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TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 2.175 | 83.42 ng | 2219510 | 1,4-Dichlorobenzene-d4 | 6.863 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 83   |
| 2    |    |   | 3-Pentanol, 2,4-dimethyl-   | 116 | C7H16O  | 000600-36-2 | 43   |
| 3    |    |   | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 40   |
| 4    |    |   | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 38   |
| 5    |    |   | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 23   |



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SOIL-DRUMS-1

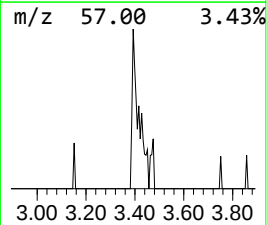
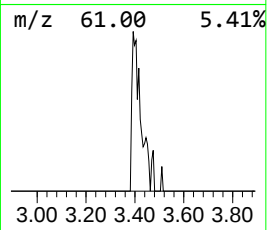
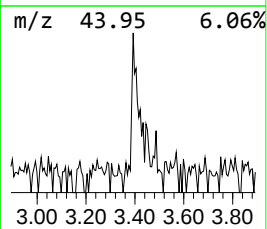
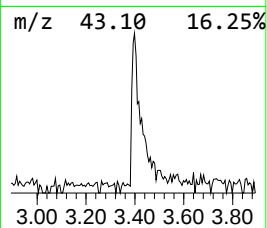
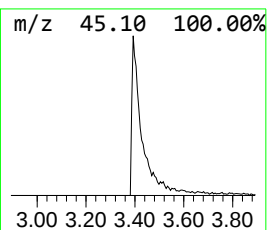
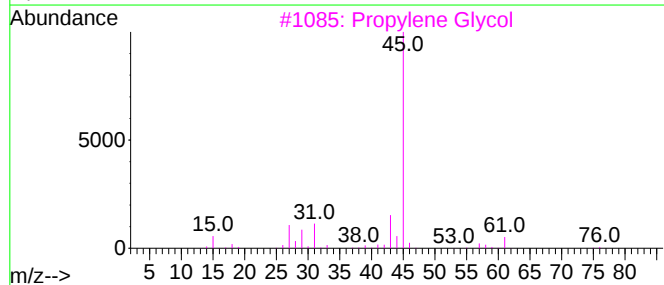
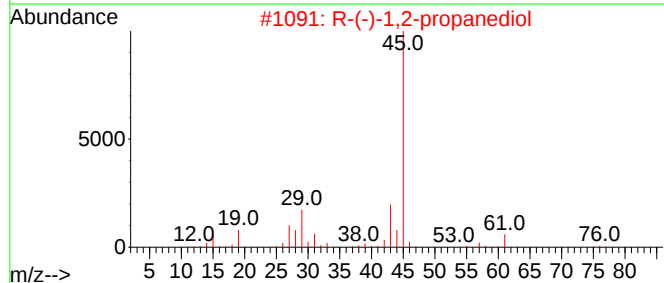
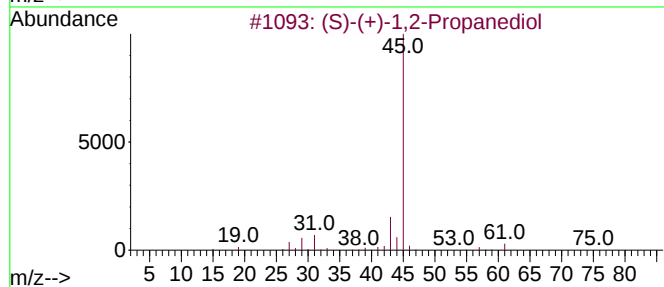
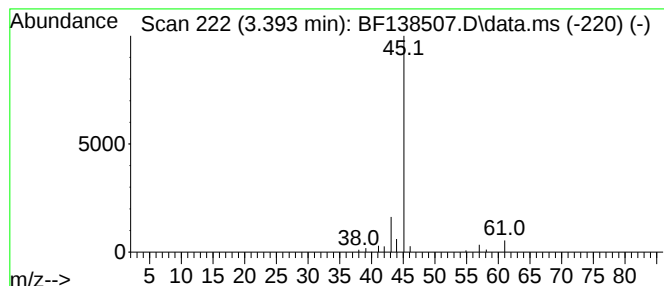
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TIC Library : C:\Database\NIST20.L  
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Peak Number 2 (S)-(+)-1,2-Propanediol Concentration Rank 13

| R.T.  | EstConc | Area  | Relative to ISTD       | R.T.  |
|-------|---------|-------|------------------------|-------|
| 3.393 | 2.78 ng | 74040 | 1,4-Dichlorobenzene-d4 | 6.863 |

| Hit# | of                      | 5 | Tentative ID | MW | MolForm | CAS#        | Qual |
|------|-------------------------|---|--------------|----|---------|-------------|------|
| 1    | (S)-(+)-1,2-Propanediol |   |              | 76 | C3H8O2  | 004254-15-3 | 64   |
| 2    | R-(-)-1,2-propanediol   |   |              | 76 | C3H8O2  | 004254-14-2 | 64   |
| 3    | Propylene Glycol        |   |              | 76 | C3H8O2  | 000057-55-6 | 64   |
| 4    | 2-Butanol, (R)-         |   |              | 74 | C4H10O  | 014898-79-4 | 9    |
| 5    | Formamide               |   |              | 45 | CH3NO   | 000075-12-7 | 7    |



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Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOIL-DRUMS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF070924.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

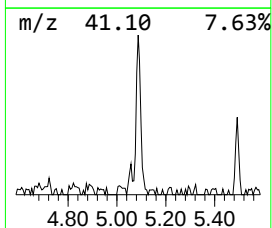
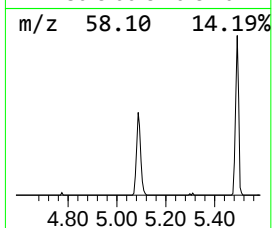
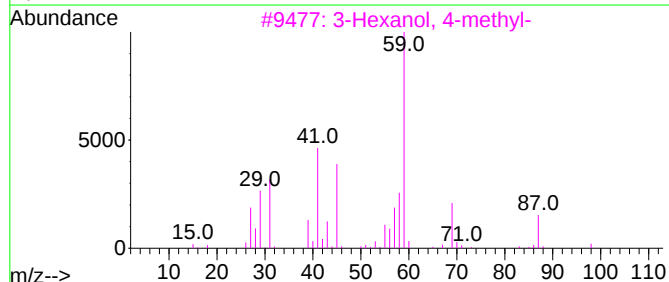
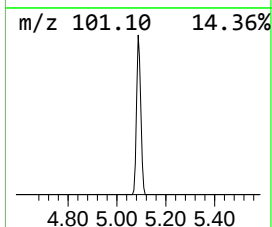
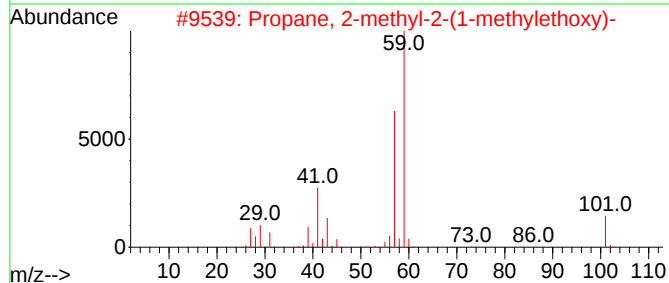
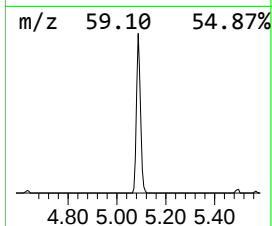
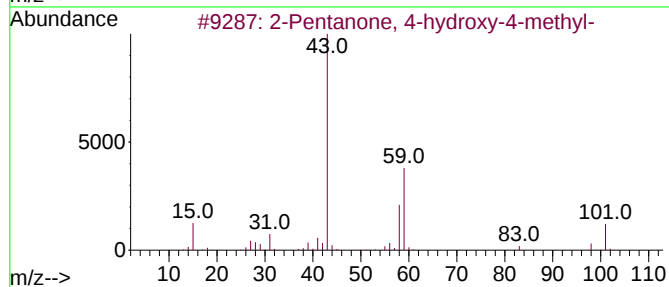
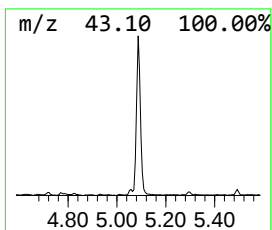
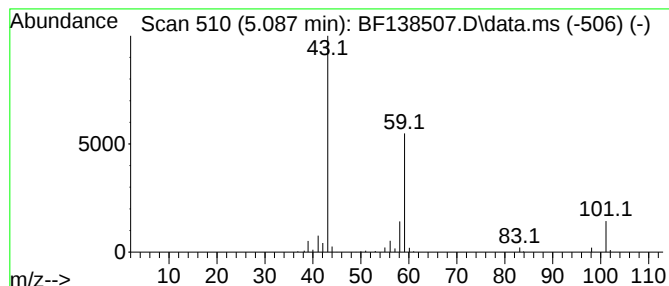
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 5.087 | 6.37 ng | 169473 | 1,4-Dichlorobenzene-d4 | 6.863 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2      | 000123-42-2 | 64      |      |      |
| 2    | Propane, 2-methyl-2-(1-methyleth... | 116 | C7H16O       | 017348-59-3 | 53      |      |      |
| 3    | 3-Hexanol, 4-methyl-                | 116 | C7H16O       | 000615-29-2 | 33      |      |      |
| 4    | 2-Hexanol, 2-methyl-                | 116 | C7H16O       | 000625-23-0 | 28      |      |      |
| 5    | 1-Propen-2-ol, acetate              | 100 | C5H8O2       | 000108-22-5 | 10      |      |      |



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SOIL-DRUMS-1

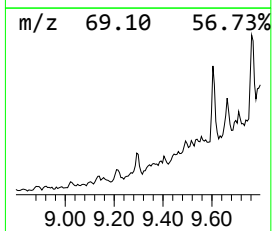
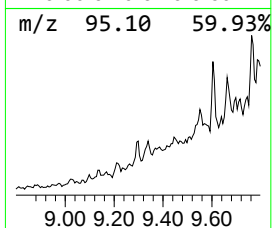
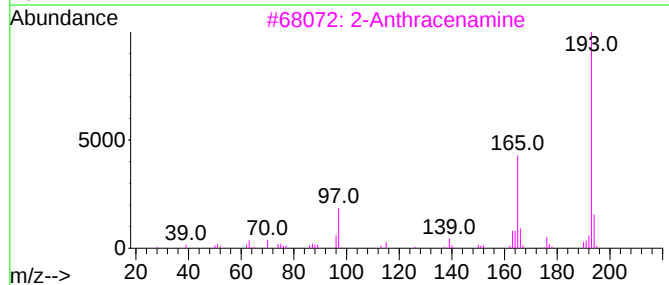
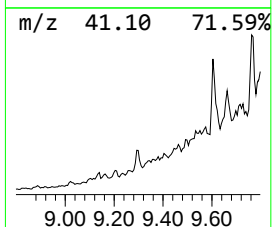
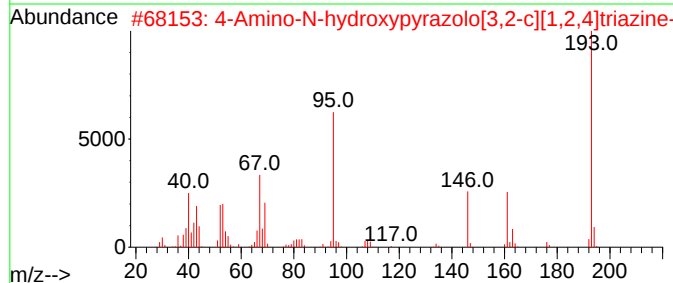
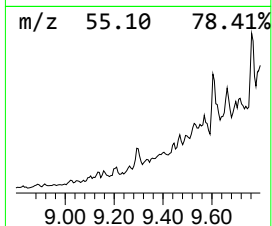
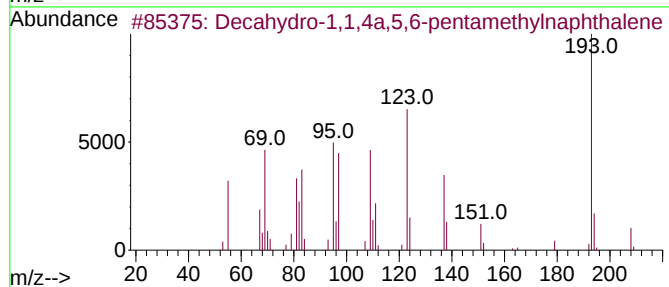
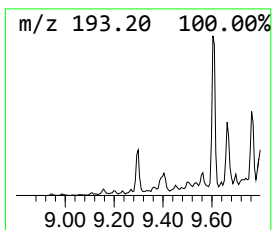
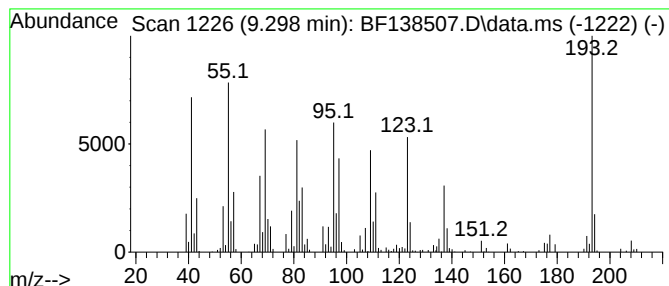
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF070924.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Decahydro-1,1,4a,5,6-pentam... Concentration Rank 11

| R.T.      | EstConc                             | Area   | Relative to ISTD |         | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|---------|--------------|------|
| 9.298     | 6.63 ng                             | 268181 | Acenaphthene-d10 |         | 9.904        |      |
| Hit# of 5 | Tentative ID                        |        | MW               | MolForm | CAS#         | Qual |
| 1         | Decahydro-1,1,4a,5,6-pentamethyl... |        | 208              | C15H28  | 080655-44-3  | 90   |
| 2         | 4-Amino-N-hydroxypyrazolo[3,2-c]... |        | 193              | C6H7N7O | 1000443-50-2 | 30   |
| 3         | 2-Anthracenamine                    |        | 193              | C14H11N | 000613-13-8  | 27   |
| 4         | 9-Phenanthrenamine                  |        | 193              | C14H11N | 000947-73-9  | 22   |
| 5         | Benzo[f]quinoline, 2-methyl-        |        | 193              | C14H11N | 039258-30-5  | 18   |



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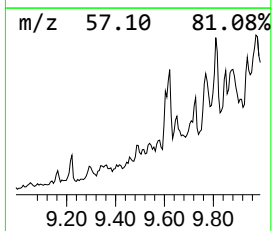
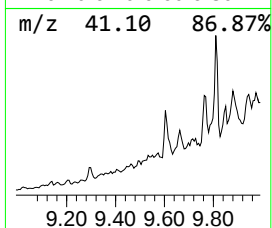
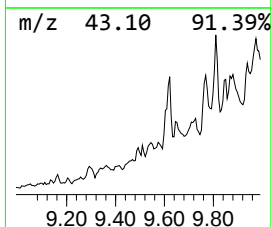
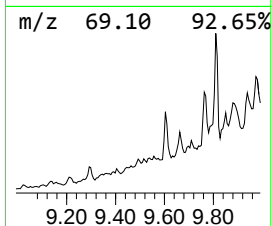
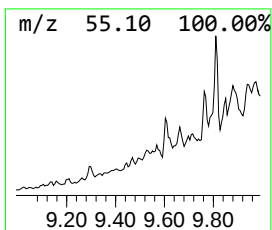
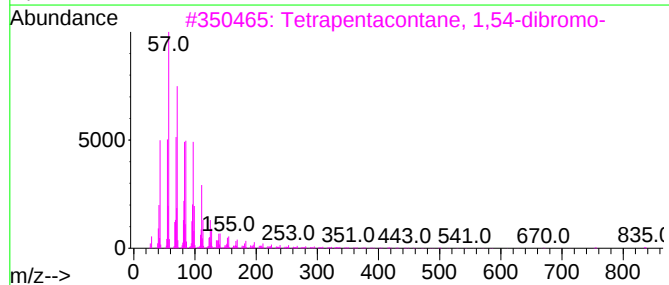
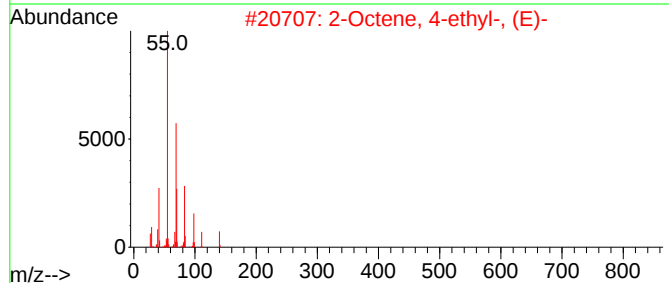
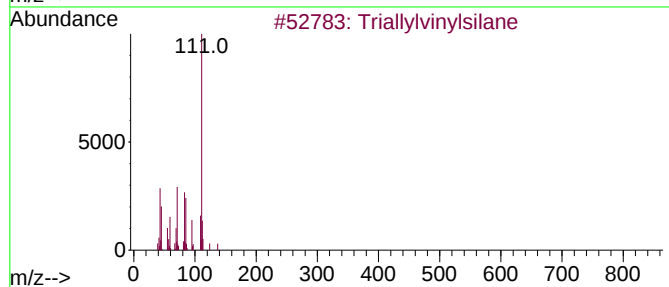
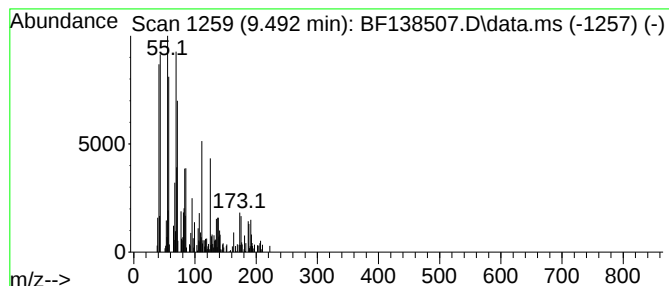
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 unknown9.492 Concentration Rank 14

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 9.492 | 2.42 ng | 97795 | Acenaphthene-d10 | 9.904 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm    | CAS#         | Qual |
|------|----|---|----------------------------------|-----|------------|--------------|------|
| 1    |    |   | Triallylvinylsilane              | 178 | C11H18Si   | 017988-47-5  | 35   |
| 2    |    |   | 2-Octene, 4-ethyl-, (E)-         | 140 | C10H20     | 074630-09-4  | 35   |
| 3    |    |   | Tetrapentacontane, 1,54-dibromo- | 915 | C54H108Br2 | 1000156-09-4 | 30   |
| 4    |    |   | Dodecane, 4-methyl-              | 184 | C13H28     | 006117-97-1  | 25   |
| 5    |    |   | Nonane, 1,9-dibromo-             | 284 | C9H18Br2   | 004549-33-1  | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF071024\  
Data File : BF138507.D  
Acq On : 10 Jul 2024 18:14  
Operator : CG\JU  
Sample : P3170-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

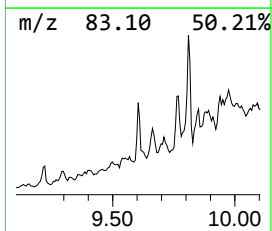
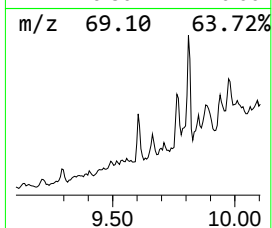
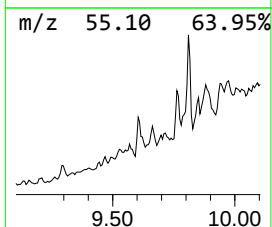
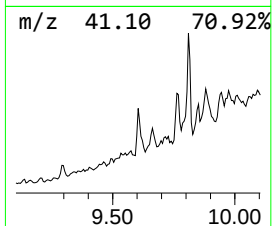
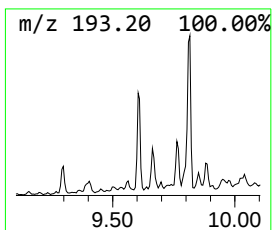
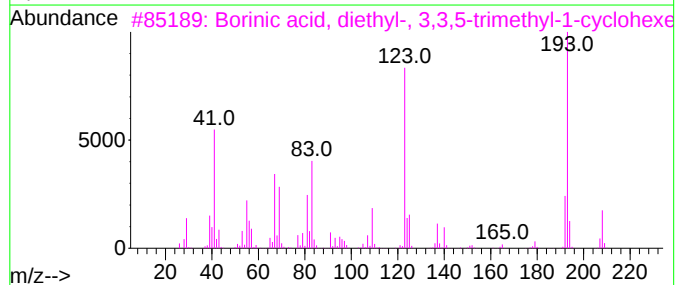
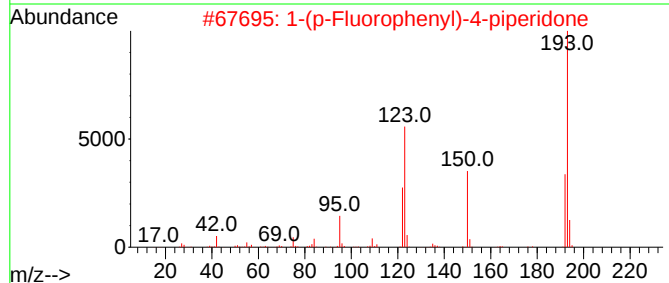
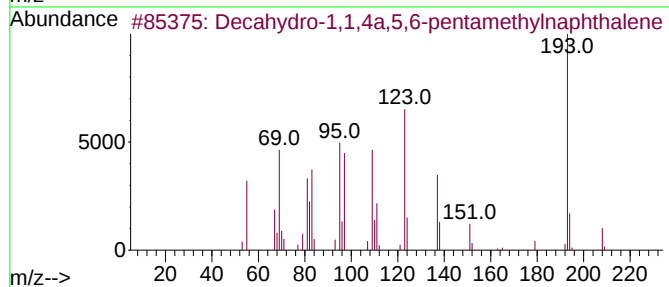
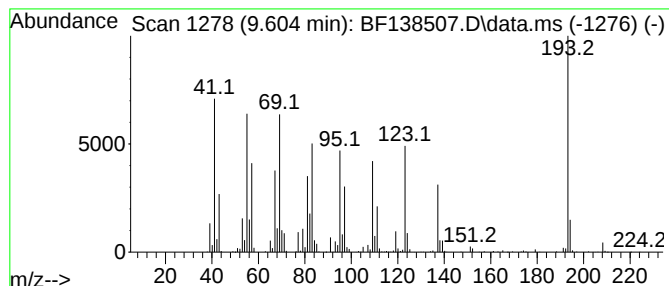
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BNA\_F  
ClientSampleId :  
SOIL-DRUMS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF070924.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 unknown9.604 Concentration Rank 6

| R.T.      | EstConc                             | Area   | Relative to ISTD |            |              | R.T.  |
|-----------|-------------------------------------|--------|------------------|------------|--------------|-------|
| 9.604     | 19.38 ng                            | 783365 | Acenaphthene-d10 |            |              | 9.904 |
| Hit# of 5 | Tentative ID                        |        | MW               | MolForm    | CAS#         | Qual  |
| 1         | Decahydro-1,1,4a,5,6-pentamethyl... |        | 208              | C15H28     | 080655-44-3  | 38    |
| 2         | 1-(p-Fluorophenyl)-4-piperidone     |        | 193              | C11H12FNO  | 1000238-56-7 | 38    |
| 3         | Borinic acid, diethyl-, 3,3,5-tr... |        | 208              | C13H25BO   | 057387-76-5  | 32    |
| 4         | 4H-1,3,2-Dioxaborin, 4-ethenyl-2... |        | 208              | C12H21BO2  | 074630-04-9  | 25    |
| 5         | Butan-2-one, 1-[3-(3,3-dimethyl-... |        | 278              | C16H26N2O2 | 1000303-34-2 | 22    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF071024\  
Data File : BF138507.D  
Acq On : 10 Jul 2024 18:14  
Operator : CG\JU  
Sample : P3170-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

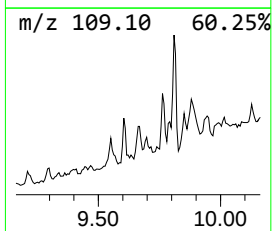
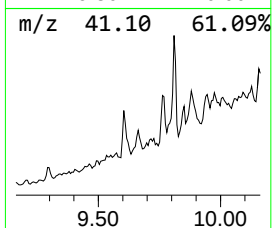
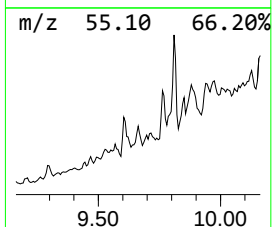
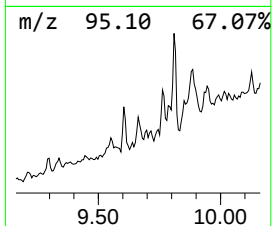
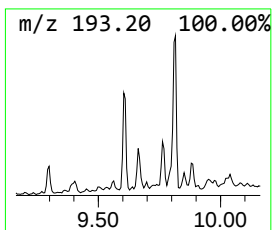
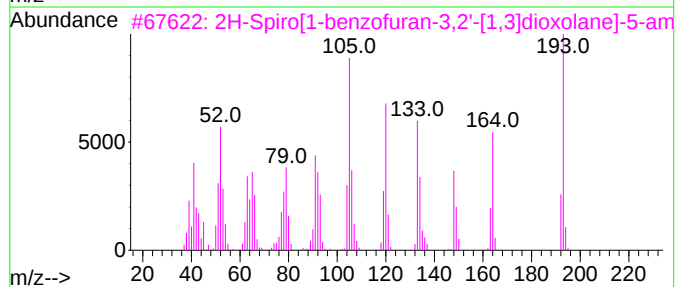
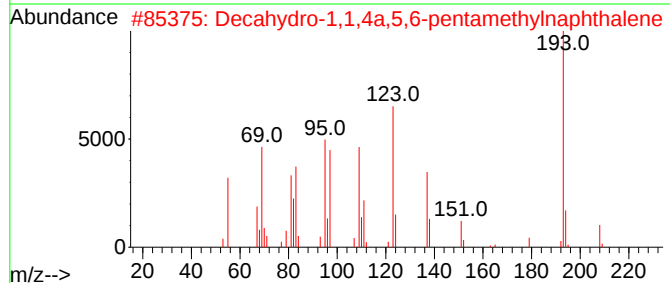
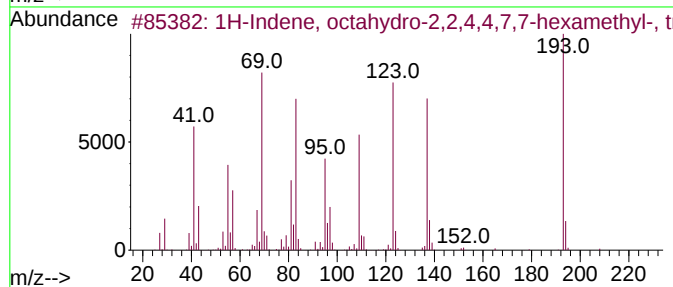
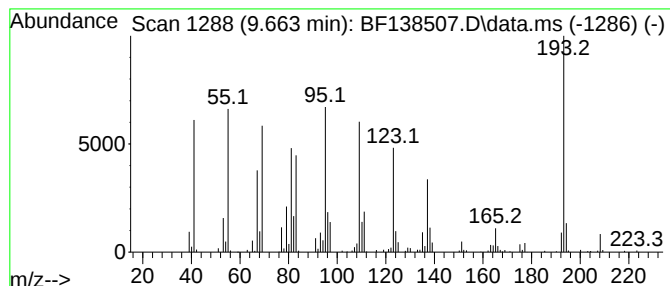
Instrument :  
BNA\_F  
ClientSampleId :  
SOIL-DRUMS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF070924.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 1H-Indene, octahydro-2,2,4,... Concentration Rank 9

| R.T.      | EstConc                             | Area   | Relative to ISTD |           | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|-----------|--------------|------|
| 9.663     | 7.56 ng                             | 305453 | Acenaphthene-d10 |           | 9.904        |      |
| Hit# of 5 | Tentative ID                        |        | MW               | MolForm   | CAS#         | Qual |
| 1         | 1H-Indene, octahydro-2,2,4,4,7,7... |        | 208              | C15H28    | 054832-83-6  | 50   |
| 2         | Decahydro-1,1,4a,5,6-pentamethyl... |        | 208              | C15H28    | 080655-44-3  | 49   |
| 3         | 2H-Spiro[1-benzofuran-3,2'-[1,3]... |        | 193              | C10H11NO3 | 1000387-33-4 | 35   |
| 4         | Acetylene, 1-(2-aminophenyl)-2-p... |        | 193              | C14H11N   | 1000380-73-4 | 27   |
| 5         | Naphthalene, decahydro-1,8a-dime... |        | 208              | C15H28    | 015404-63-4  | 27   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF071024\  
Data File : BF138507.D  
Acq On : 10 Jul 2024 18:14  
Operator : CG\JU  
Sample : P3170-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOIL-DRUMS-1

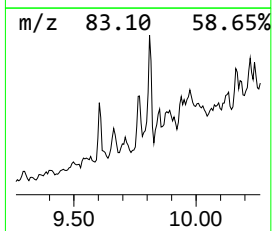
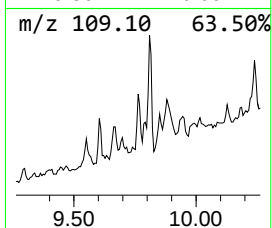
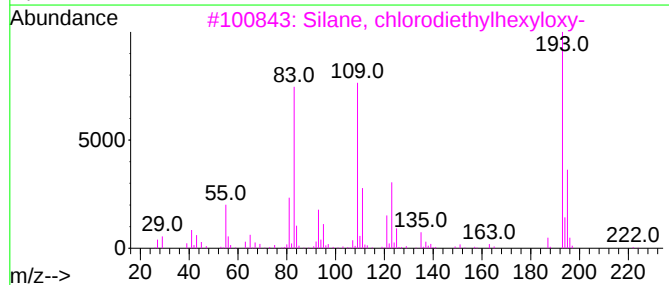
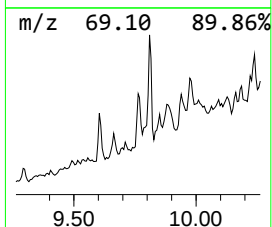
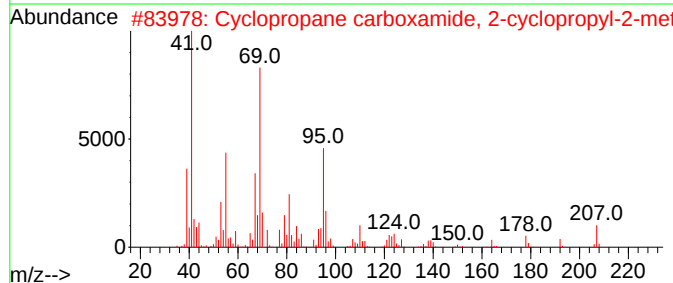
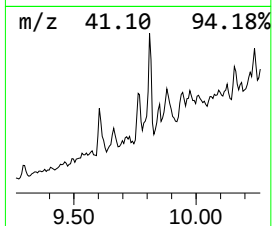
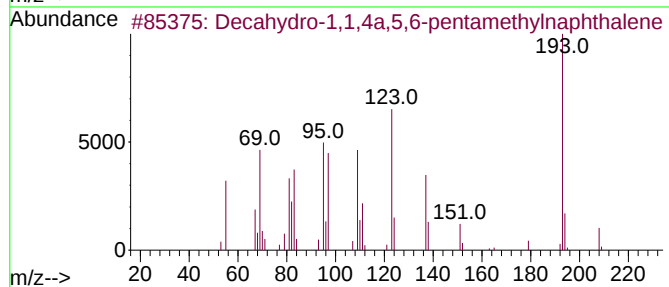
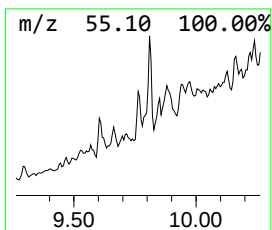
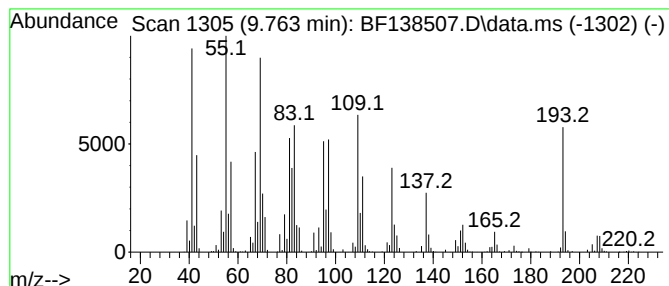
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 unknown9.763 Concentration Rank 4

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 9.763     | 19.89 ng                            | 804111 | Acenaphthene-d10 | 9.904        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | Decahydro-1,1,4a,5,6-pentamethyl... | 208    | C15H28           | 080655-44-3  | 60   |
| 2         | Cyclopropane carboxamide, 2-cycl... | 207    | C13H21NO         | 331416-19-4  | 38   |
| 3         | Silane, chlorodiethylhexyloxy-      | 222    | C10H23ClOSi      | 1000363-74-4 | 35   |
| 4         | 2-Pentanone, 4-cyclohexylidene-3... | 222    | C15H26O          | 313253-65-5  | 30   |
| 5         | 1,7-Dimethyl-4-(1-methylethyl)cy... | 210    | C15H30           | 000645-10-3  | 27   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF071024\  
Data File : BF138507.D  
Acq On : 10 Jul 2024 18:14  
Operator : CG\JU  
Sample : P3170-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOIL-DRUMS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF070924.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

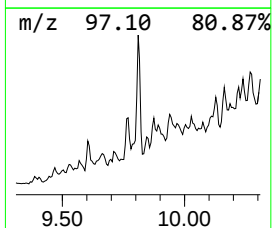
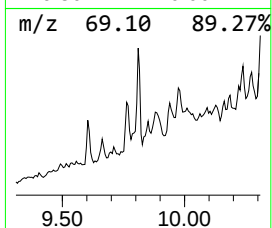
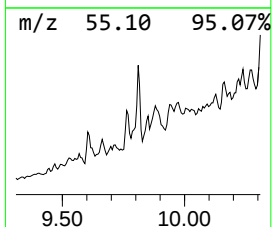
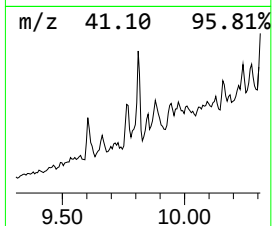
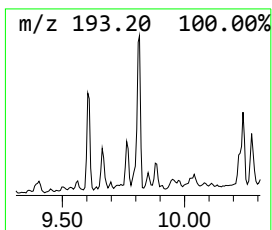
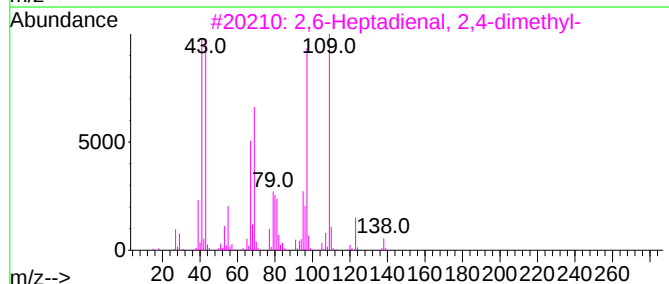
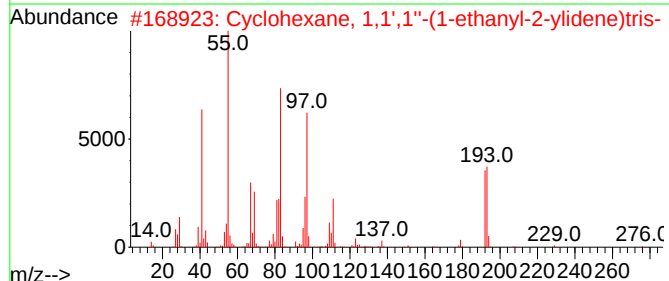
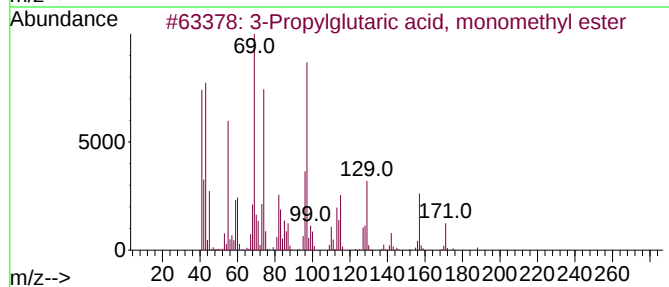
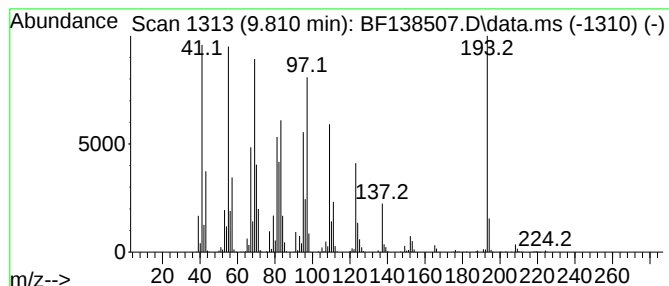
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Peak Number 9 unknown9.810 Concentration Rank 2

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 9.810 | 32.89 ng | 1329660 | Acenaphthene-d10 | 9.904 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|------|-------------------------------------|-----|----------|--------------|------|
| 1    |      | 3-Propylglutaric acid, monomethy... | 188 | C9H16O4  | 1000254-25-9 | 27   |
| 2    |      | Cyclohexane, 1,1',1''-(1-ethanyl... | 276 | C20H36   | 055682-86-5  | 27   |
| 3    |      | 2,6-Heptadienal, 2,4-dimethyl-      | 138 | C9H14O   | 085136-08-9  | 25   |
| 4    |      | Cyclopentanecarboxylic acid, 3-p... | 324 | C21H40O2 | 1000280-59-1 | 22   |
| 5    |      | 3-Heptene, 4-propyl-                | 140 | C10H20   | 004485-13-6  | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF071024\  
Data File : BF138507.D  
Acq On : 10 Jul 2024 18:14  
Operator : CG\JU  
Sample : P3170-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOIL-DRUMS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF070924.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

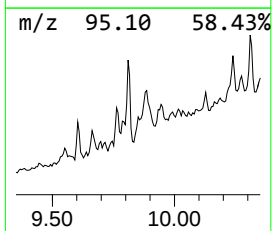
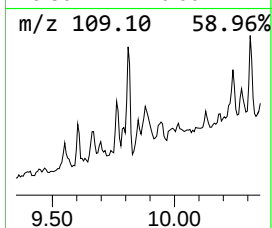
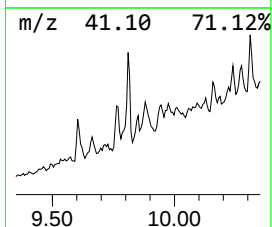
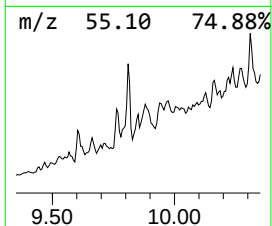
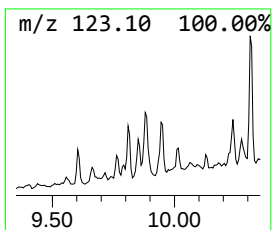
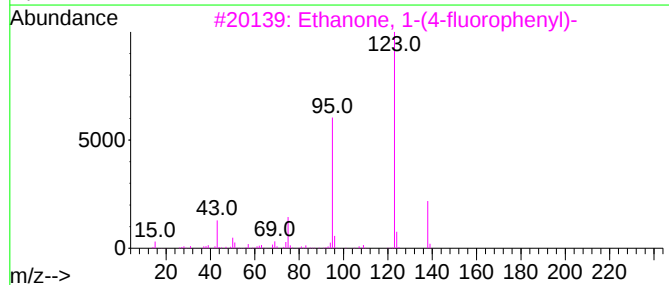
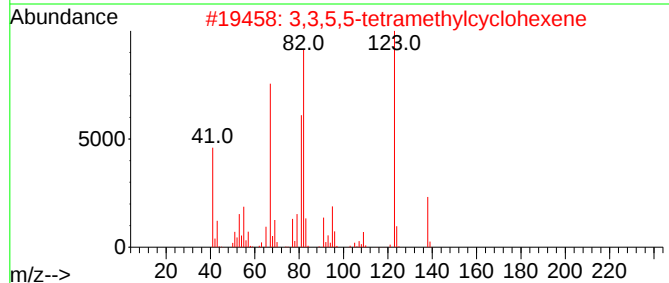
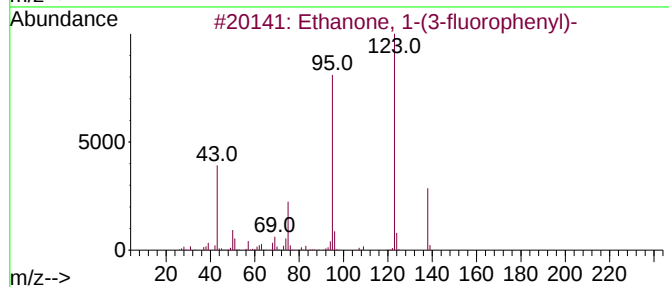
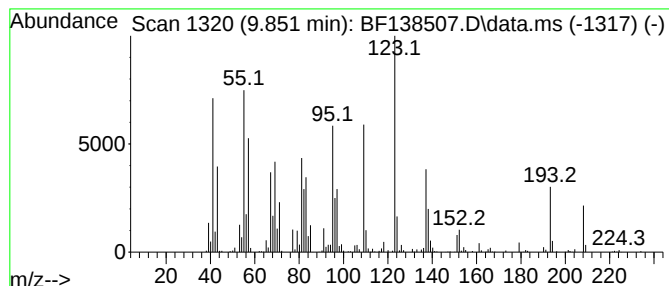
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TIC Integration Parameters: LSCINT.P

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Peak Number 10 unknown9.851 Concentration Rank 8

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 9.851 | 10.02 ng | 405118 | Acenaphthene-d10 | 9.904 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|------|-------------------------------------|-----|---------|--------------|------|
| 1    |      | Ethanone, 1-(3-fluorophenyl)-       | 138 | C8H7FO  | 000455-36-7  | 38   |
| 2    |      | 3,3,5,5-tetramethylcyclohexene      | 138 | C10H18  | 1000510-60-7 | 38   |
| 3    |      | Ethanone, 1-(4-fluorophenyl)-       | 138 | C8H7FO  | 000403-42-9  | 35   |
| 4    |      | Bicyclo[2.2.1]heptane, 2,2,3-tri... | 138 | C10H18  | 000473-19-8  | 27   |
| 5    |      | 1-Propanone, 1-(4-fluorophenyl)-    | 152 | C9H9FO  | 000456-03-1  | 27   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF071024\  
Data File : BF138507.D  
Acq On : 10 Jul 2024 18:14  
Operator : CG\JU  
Sample : P3170-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

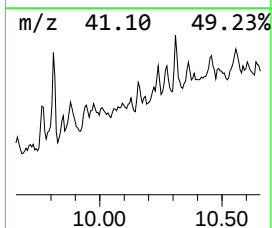
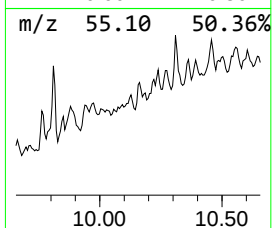
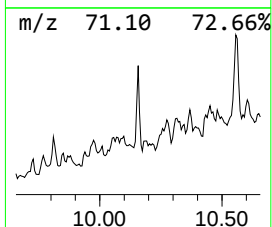
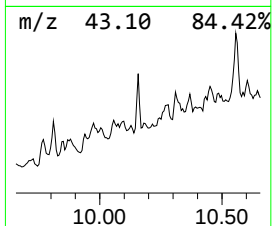
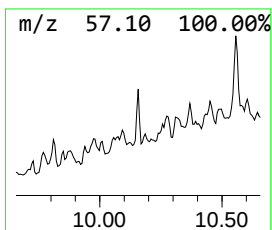
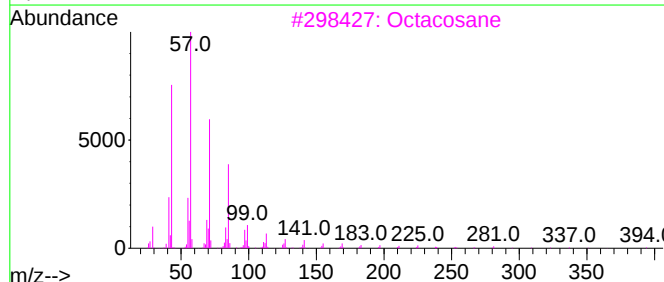
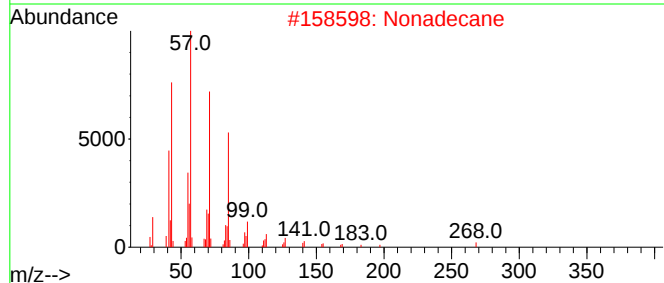
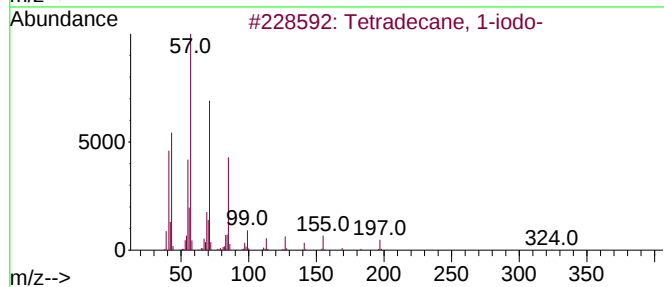
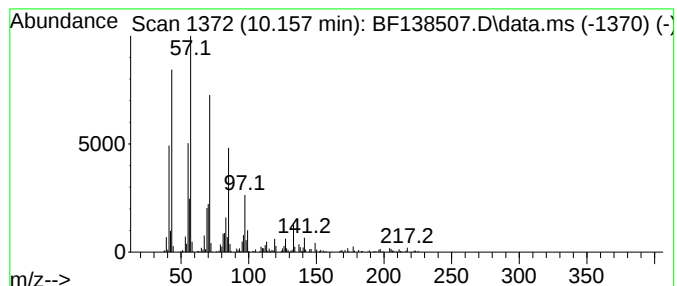
Instrument :  
BNA\_F  
ClientSampleId :  
SOIL-DRUMS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF070924.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Tetradecane, 1-iodo- Concentration Rank 7

| R.T.      | EstConc              | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|----------------------|--------|------------------|---------|-------------|------|
| 10.157    | 13.45 ng             | 543703 | Acenaphthene-d10 |         | 9.904       |      |
| Hit# of 5 | Tentative ID         |        | MW               | MolForm | CAS#        | Qual |
| 1         | Tetradecane, 1-iodo- |        | 324              | C14H29I | 019218-94-1 | 64   |
| 2         | Nonadecane           |        | 268              | C19H40  | 000629-92-5 | 58   |
| 3         | Octacosane           |        | 394              | C28H58  | 000630-02-4 | 58   |
| 4         | Tetracosane          |        | 338              | C24H50  | 000646-31-1 | 58   |
| 5         | Dodecane, 4-methyl-  |        | 184              | C13H28  | 006117-97-1 | 55   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF071024\  
Data File : BF138507.D  
Acq On : 10 Jul 2024 18:14  
Operator : CG\JU  
Sample : P3170-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOIL-DRUMS-1

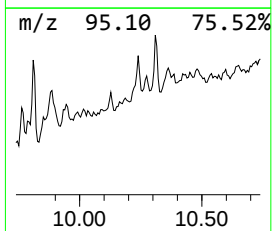
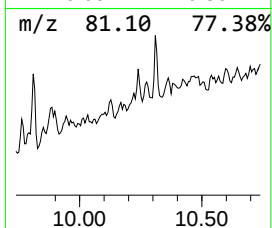
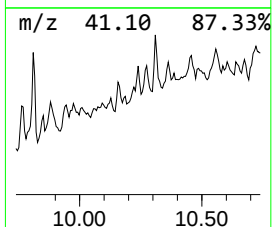
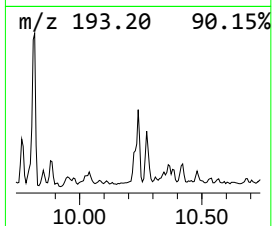
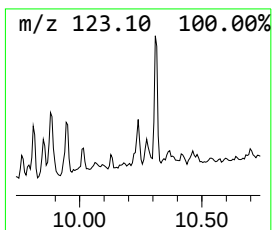
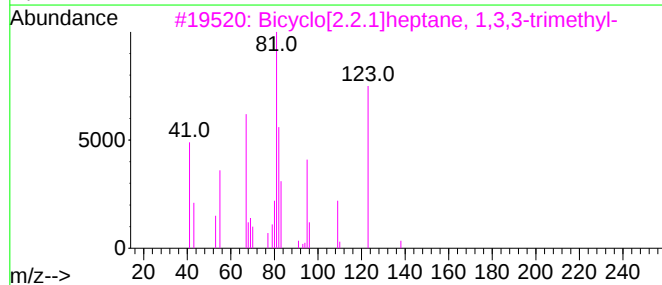
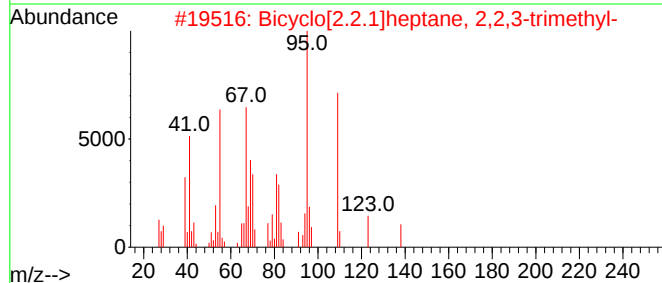
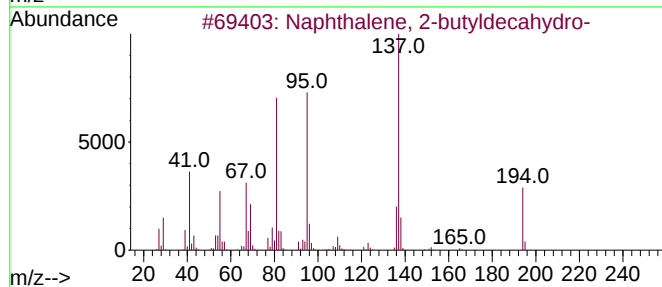
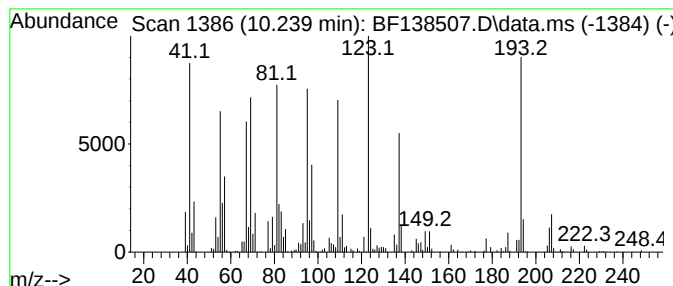
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 unknown10.239 Concentration Rank 10

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.  |
|--------|---------|--------|------------------|-------|
| 10.239 | 6.99 ng | 282697 | Acenaphthene-d10 | 9.904 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Naphthalene, 2-butyldecahydro-      | 194 | C14H26   | 006305-52-8 | 43   |
| 2    |    |   | Bicyclo[2.2.1]heptane, 2,2,3-tri... | 138 | C10H18   | 000473-19-8 | 38   |
| 3    |    |   | Bicyclo[2.2.1]heptane, 1,3,3-tri... | 138 | C10H18   | 006248-88-0 | 38   |
| 4    |    |   | Bicyclo[6.1.0]nonane, 9-bromo-9-... | 216 | C10H17Br | 059474-03-2 | 38   |
| 5    |    |   | 1H-Indene, octahydro-2,2,4,4,7,7... | 208 | C15H28   | 054832-83-6 | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF071024\  
Data File : BF138507.D  
Acq On : 10 Jul 2024 18:14  
Operator : CG\JU  
Sample : P3170-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOIL-DRUMS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF070924.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

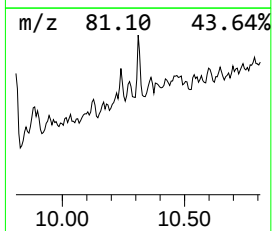
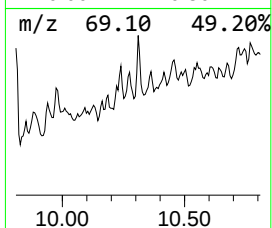
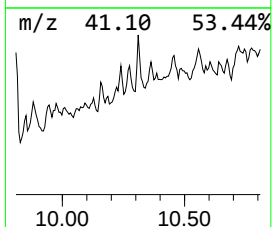
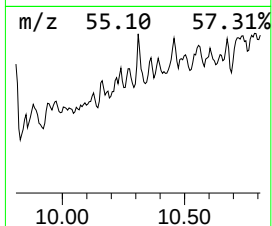
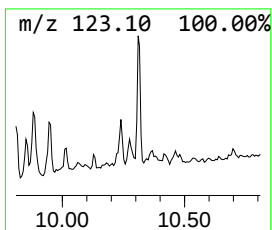
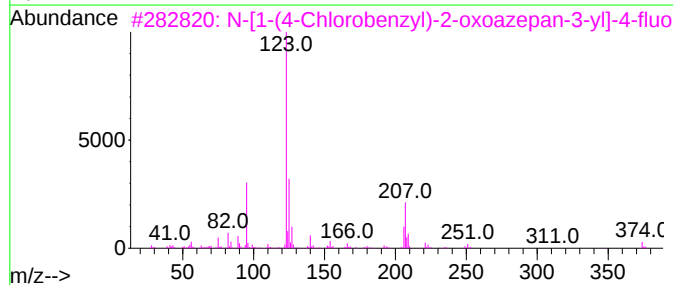
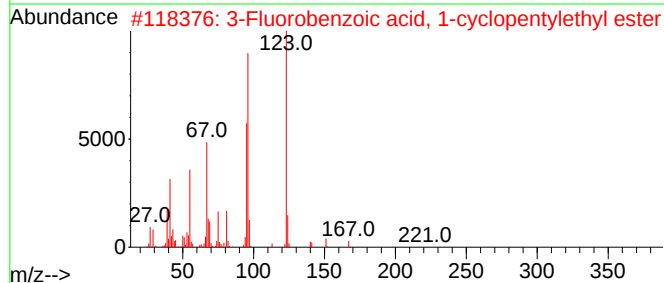
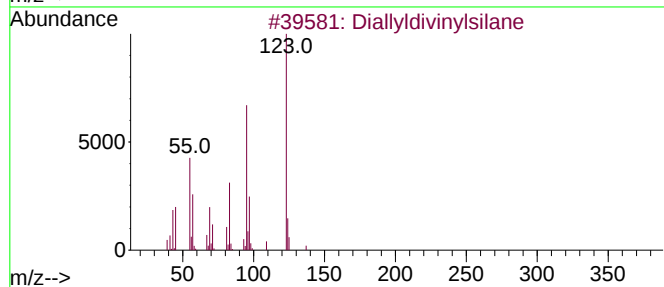
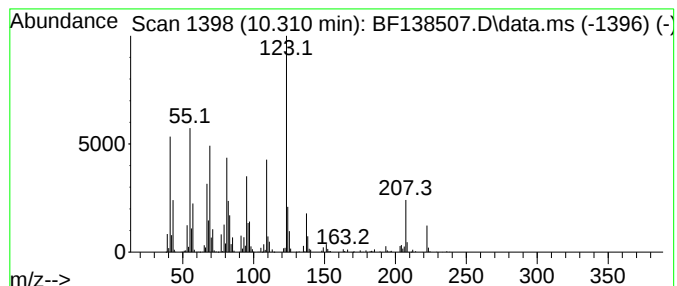
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 unknown10.310 Concentration Rank 5

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.  |
|--------|----------|--------|------------------|-------|
| 10.310 | 19.42 ng | 784943 | Acenaphthene-d10 | 9.904 |

| Hit# | of | 5 | Tentative ID                                             | MW  | MolForm       | CAS#         | Qual |
|------|----|---|----------------------------------------------------------|-----|---------------|--------------|------|
| 1    |    |   | Diallyldivinylsilane                                     | 164 | C10H16Si      | 119901-88-1  | 47   |
| 2    |    |   | 3-Fluorobenzoic acid, 1-cyclopentylethyl ester           | 236 | C14H17FO2     | 1000279-04-7 | 43   |
| 3    |    |   | N-[1-(4-Chlorobenzyl)-2-oxazepan-3-yl]-4-fluorobenzamide | 374 | C20H20ClFN2O2 | 1000481-70-9 | 43   |
| 4    |    |   | Pentanamide, N-(4-methoxyphenyl)-                        | 207 | C12H17NO2     | 1000306-89-3 | 38   |
| 5    |    |   | Benzamide, 3-fluoro-N-(2-phenylethyl)-                   | 313 | C20H24FNO     | 1000451-30-5 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF071024\  
Data File : BF138507.D  
Acq On : 10 Jul 2024 18:14  
Operator : CG\JU  
Sample : P3170-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOIL-DRUMS-1

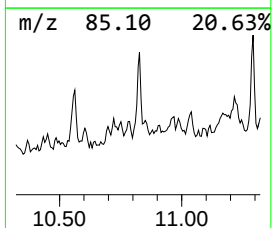
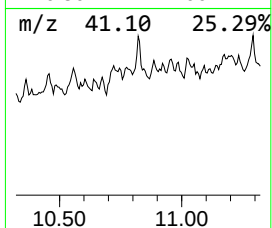
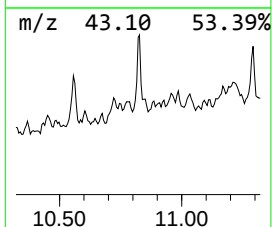
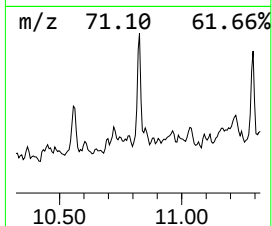
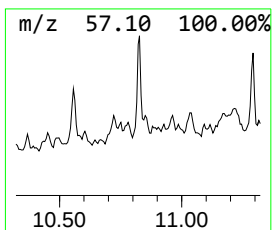
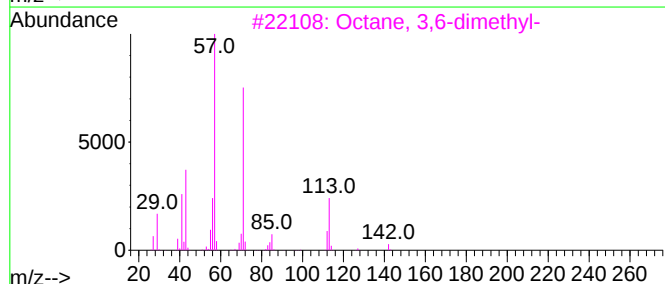
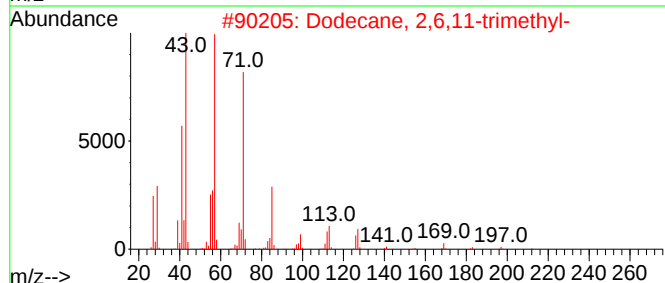
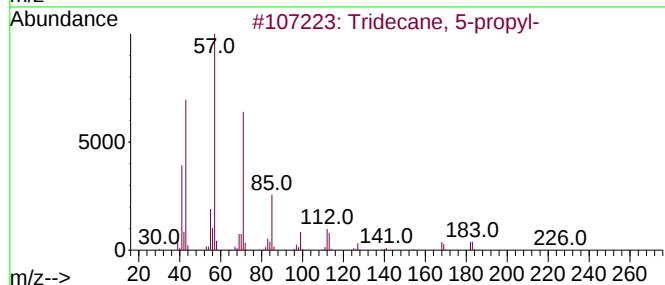
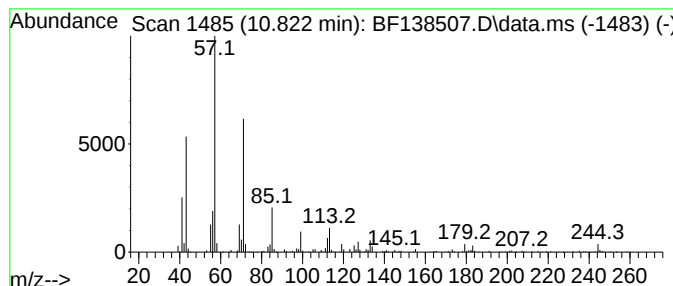
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Tridecane, 5-propyl- Concentration Rank 3

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 10.822 | 22.31 ng | 665566 | Phenanthrene-d10 | 11.398 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Tridecane, 5-propyl-        | 226 | C16H34  | 055045-11-9 | 89   |
| 2    |    |   | Dodecane, 2,6,11-trimethyl- | 212 | C15H32  | 031295-56-4 | 80   |
| 3    |    |   | Octane, 3,6-dimethyl-       | 142 | C10H22  | 015869-94-0 | 64   |
| 4    |    |   | Undecane, 3,6-dimethyl-     | 184 | C13H28  | 017301-28-9 | 58   |
| 5    |    |   | Tridecane, 3-ethyl-         | 212 | C15H32  | 013286-73-2 | 52   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF071024\  
Data File : BF138507.D  
Acq On : 10 Jul 2024 18:14  
Operator : CG\JU  
Sample : P3170-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOIL-DRUMS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF070924.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L  
TIC Integration Parameters: LSCINT.P

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |        |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|--------|------|
|                    |        |         |       |          | #                     | RT     | Resp   | Conc |
| Butane, 2-metho... | 2.175  | 83.4    | ng    | 2219510  | 1                     | 6.863  | 532119 | 20.0 |
| (S)-(+)-1,2-Pro... | 3.393  | 2.8     | ng    | 74040    | 1                     | 6.863  | 532119 | 20.0 |
| 2-Pentanone, 4-... | 5.087  | 6.4     | ng    | 169473   | 1                     | 6.863  | 532119 | 20.0 |
| Decahydro-1,1,4... | 9.298  | 6.6     | ng    | 268181   | 3                     | 9.904  | 808491 | 20.0 |
| unknown9.492       | 9.492  | 2.4     | ng    | 97795    | 3                     | 9.904  | 808491 | 20.0 |
| unknown9.604       | 9.604  | 19.4    | ng    | 783365   | 3                     | 9.904  | 808491 | 20.0 |
| 1H-Indene, octa... | 9.663  | 7.6     | ng    | 305453   | 3                     | 9.904  | 808491 | 20.0 |
| unknown9.763       | 9.763  | 19.9    | ng    | 804111   | 3                     | 9.904  | 808491 | 20.0 |
| unknown9.810       | 9.810  | 32.9    | ng    | 1329660  | 3                     | 9.904  | 808491 | 20.0 |
| unknown9.851       | 9.851  | 10.0    | ng    | 405118   | 3                     | 9.904  | 808491 | 20.0 |
| Tetradecane, 1-... | 10.157 | 13.4    | ng    | 543703   | 3                     | 9.904  | 808491 | 20.0 |
| unknown10.239      | 10.239 | 7.0     | ng    | 282697   | 3                     | 9.904  | 808491 | 20.0 |
| unknown10.310      | 10.310 | 19.4    | ng    | 784943   | 3                     | 9.904  | 808491 | 20.0 |
| Tridecane, 5-pr... | 10.822 | 22.3    | ng    | 665566   | 4                     | 11.398 | 596566 | 20.0 |