

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
 Data File : BF130866.D  
 Acq On : 29 Oct 2022 19:49  
 Operator : CG\JU  
 Sample : N5308-01  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PL-2-102622

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-BF102722.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF130866.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         | 2.246       | 19            | 27          | 33           | rVB      | 1061728        | 1362616       | 88.28%          | 7.738%        |
| 2         | 2.357       | 41            | 46          | 55           | rVB2     | 35527          | 52298         | 3.39%           | 0.297%        |
| 3         | 4.622       | 425           | 431         | 437          | rBV2     | 14724          | 20014         | 1.30%           | 0.114%        |
| 4         | 5.169       | 520           | 524         | 530          | rVB      | 226181         | 238573        | 15.46%          | 1.355%        |
| 5         | 5.557       | 585           | 590         | 593          | rBV      | 1619794        | 1543512       | 100.00%         | 8.765%        |
| 6         | 6.563       | 756           | 761         | 764          | rBV      | 1467628        | 1467925       | 95.10%          | 8.336%        |
| 7         | 6.945       | 822           | 826         | 830          | rVB      | 542934         | 419592        | 27.18%          | 2.383%        |
| 8         | 7.504       | 917           | 921         | 925          | rBV      | 941200         | 871878        | 56.49%          | 4.951%        |
| 9         | 8.128       | 1024          | 1027        | 1037         | rBV      | 36584          | 60958         | 3.95%           | 0.346%        |
| 10        | 8.228       | 1041          | 1044        | 1047         | rVV      | 514731         | 441932        | 28.63%          | 2.510%        |
| 11        | 8.816       | 1141          | 1144        | 1148         | rBV      | 30570          | 24950         | 1.62%           | 0.142%        |
| 12        | 9.304       | 1223          | 1227        | 1231         | rVB      | 1486288        | 1255569       | 81.34%          | 7.130%        |
| 13        | 9.381       | 1237          | 1240        | 1243         | rBV      | 52357          | 44022         | 2.85%           | 0.250%        |
| 14        | 9.569       | 1269          | 1272        | 1277         | rBV2     | 42443          | 42599         | 2.76%           | 0.242%        |
| 15        | 9.704       | 1291          | 1295        | 1297         | rBV3     | 31986          | 39019         | 2.53%           | 0.222%        |
| 16        | 9.851       | 1317          | 1320        | 1323         | rBV      | 30279          | 29931         | 1.94%           | 0.170%        |
| 17        | 9.910       | 1327          | 1330        | 1334         | rVB      | 50031          | 47136         | 3.05%           | 0.268%        |
| 18        | 9.992       | 1340          | 1344        | 1347         | rVB      | 550144         | 442032        | 28.64%          | 2.510%        |
| 19        | 10.233      | 1382          | 1385        | 1389         | rBV5     | 19640          | 25020         | 1.62%           | 0.142%        |
| 20        | 10.416      | 1411          | 1416        | 1418         | rVB      | 51050          | 50031         | 3.24%           | 0.284%        |
| 21        | 10.539      | 1434          | 1437        | 1441         | rBV2     | 23714          | 24810         | 1.61%           | 0.141%        |
| 22        | 10.633      | 1449          | 1453        | 1458         | rBV      | 30215          | 38147         | 2.47%           | 0.217%        |
| 23        | 10.780      | 1474          | 1478        | 1481         | rVB      | 963677         | 828113        | 53.65%          | 4.703%        |
| 24        | 10.886      | 1494          | 1496        | 1501         | rBV2     | 68001          | 105567        | 6.84%           | 0.599%        |
| 25        | 11.339      | 1570          | 1573        | 1575         | rBV      | 64522          | 51673         | 3.35%           | 0.293%        |
| 26        | 11.369      | 1575          | 1578        | 1583         | rVB2     | 51732          | 63745         | 4.13%           | 0.362%        |
| 27        | 11.416      | 1583          | 1586        | 1588         | rBV3     | 20192          | 24697         | 1.60%           | 0.140%        |
| 28        | 11.480      | 1594          | 1597        | 1600         | rBV      | 539234         | 493345        | 31.96%          | 2.802%        |
| 29        | 11.504      | 1600          | 1601        | 1604         | rVV      | 208955         | 145315        | 9.41%           | 0.825%        |
| 30        | 11.557      | 1607          | 1610        | 1613         | rVV      | 66494          | 54360         | 3.52%           | 0.309%        |
| 31        | 11.769      | 1643          | 1646        | 1649         | rBV      | 52188          | 51047         | 3.31%           | 0.290%        |
| 32        | 11.869      | 1660          | 1663        | 1666         | rVB      | 295914         | 217389        | 14.08%          | 1.235%        |
| 33        | 11.927      | 1671          | 1673        | 1677         | rBV5     | 23448          | 21718         | 1.41%           | 0.123%        |
| 34        | 11.963      | 1677          | 1679        | 1681         | rBV2     | 23116          | 24727         | 1.60%           | 0.140%        |
| 35        | 12.004      | 1681          | 1686        | 1693         | rVV2     | 466310         | 455194        | 29.49%          | 2.585%        |
| 36        | 12.098      | 1699          | 1702        | 1704         | rBV2     | 54543          | 51572         | 3.34%           | 0.293%        |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
 Data File : BF130866.D  
 Acq On : 29 Oct 2022 19:49  
 Operator : CG\JU  
 Sample : N5308-01  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PL-2-102622

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-BF102722.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 12.174 | 1712 | 1715 | 1720 | rBV  | 88789   | 95254   | 6.17%  | 0.541% |
| 38 | 12.457 | 1761 | 1763 | 1766 | rBV2 | 34541   | 41195   | 2.67%  | 0.234% |
| 39 | 12.533 | 1774 | 1776 | 1777 | rBV2 | 41931   | 36790   | 2.38%  | 0.209% |
| 40 | 12.557 | 1777 | 1780 | 1782 | rVV  | 1290619 | 1208955 | 78.32% | 6.865% |
| 41 | 12.580 | 1782 | 1784 | 1787 | rVB  | 1727354 | 1272911 | 82.47% | 7.229% |
| 42 | 12.663 | 1795 | 1798 | 1801 | rBV  | 338829  | 267536  | 17.33% | 1.519% |
| 43 | 12.698 | 1801 | 1804 | 1806 | rBV  | 642725  | 592504  | 38.39% | 3.365% |
| 44 | 12.716 | 1806 | 1807 | 1816 | rVV2 | 317123  | 399293  | 25.87% | 2.268% |
| 45 | 12.792 | 1818 | 1820 | 1824 | rVB4 | 66931   | 68769   | 4.46%  | 0.391% |
| 46 | 12.933 | 1838 | 1844 | 1850 | rBV  | 358588  | 428811  | 27.78% | 2.435% |
| 47 | 13.069 | 1864 | 1867 | 1876 | rVB  | 1183837 | 1118640 | 72.47% | 6.353% |
| 48 | 14.133 | 2043 | 2048 | 2050 | rBV2 | 428942  | 552689  | 35.81% | 3.139% |
| 49 | 14.227 | 2062 | 2064 | 2068 | rVB2 | 173231  | 159928  | 10.36% | 0.908% |
| 50 | 15.274 | 2240 | 2242 | 2247 | rVB  | 211850  | 234882  | 15.22% | 1.334% |

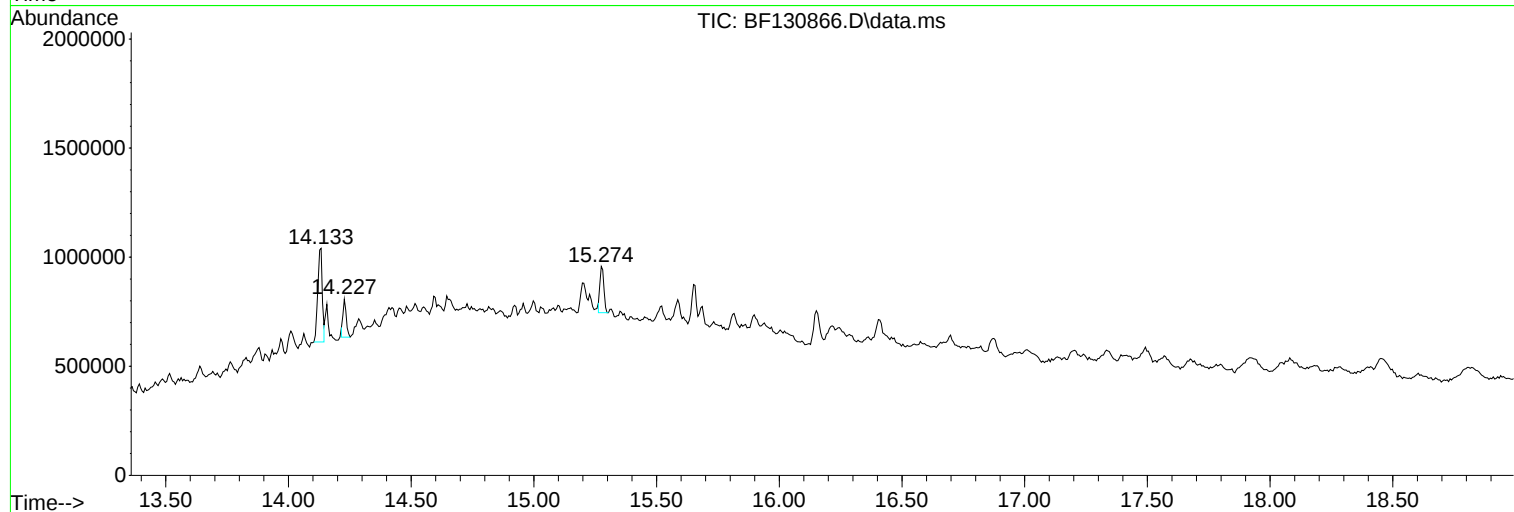
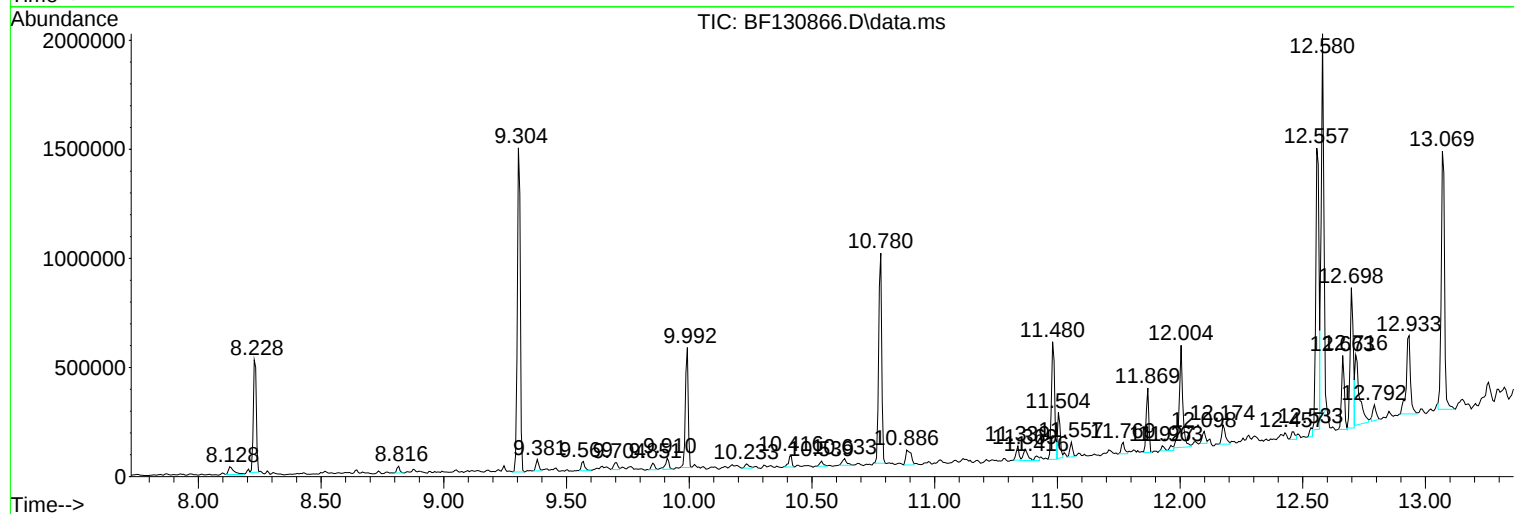
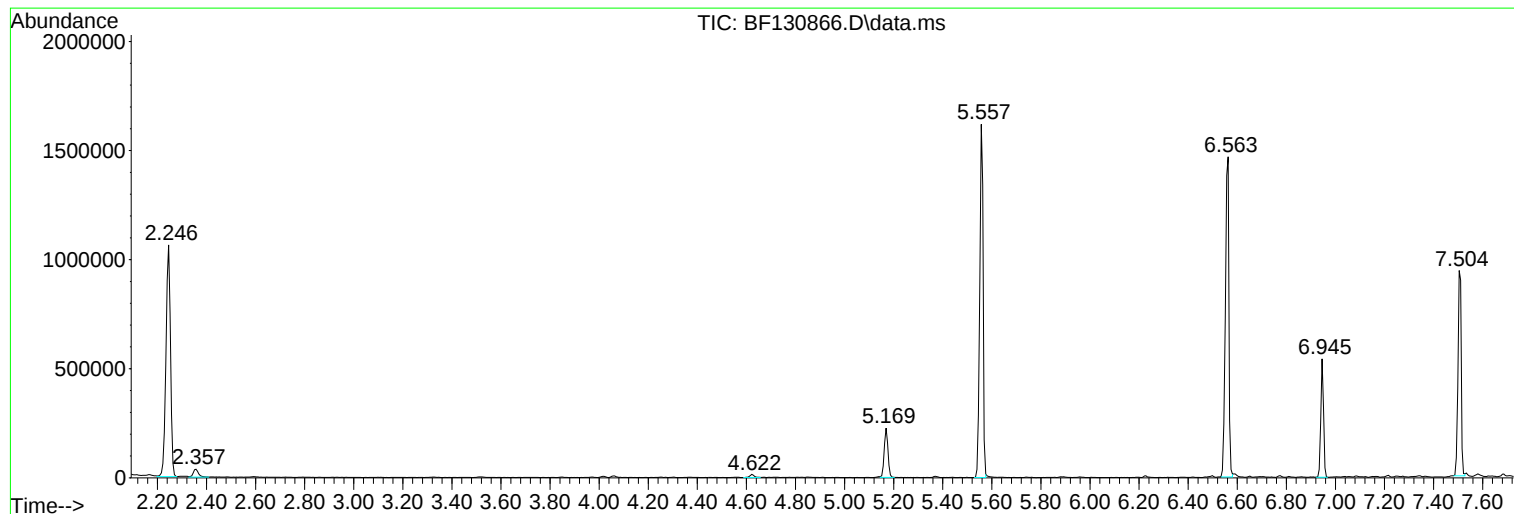
Sum of corrected areas: 17609183

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.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\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

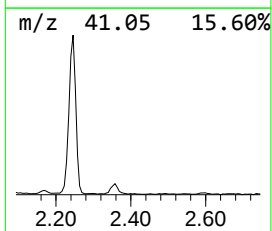
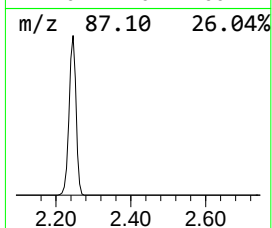
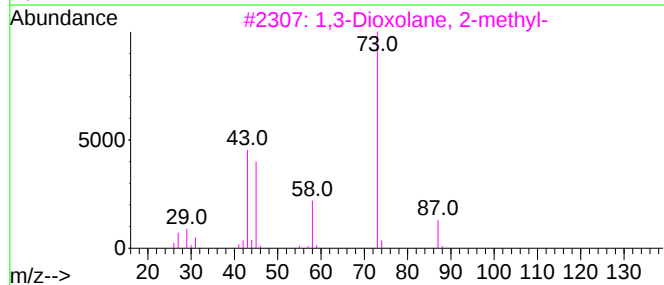
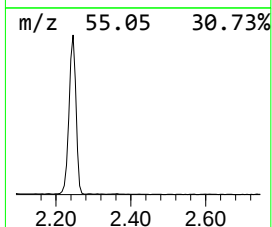
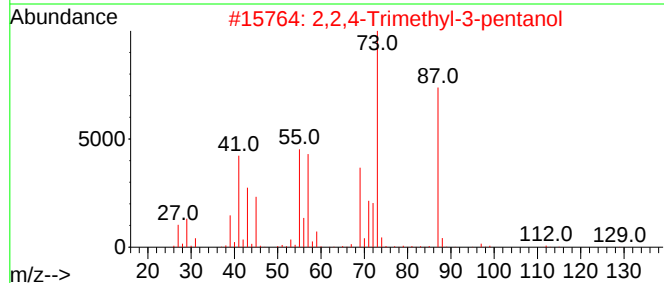
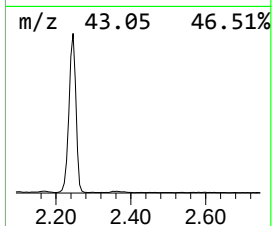
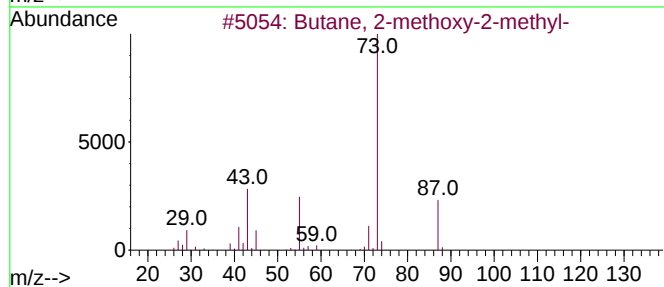
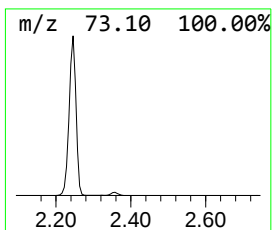
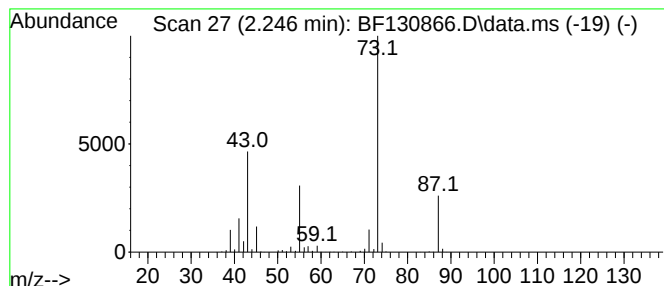
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

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

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 2.246 | 64.95 ng | 1362620 | 1,4-Dichlorobenzene-d4 | 6.945 |

| 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 | 40   |
| 3    |    |   | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 23   |
| 4    |    |   | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 17   |
| 5    |    |   | Borane, dimethoxy-          | 74  | C2H7BO2 | 004542-61-4 | 9    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

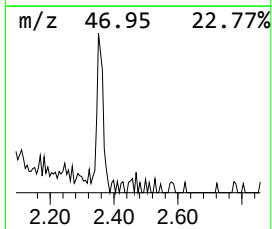
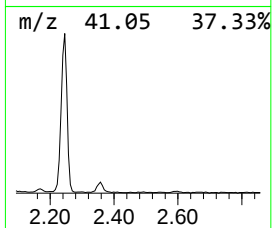
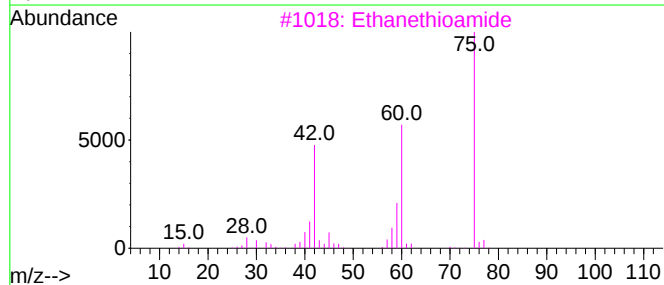
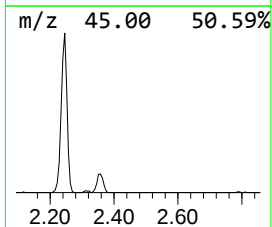
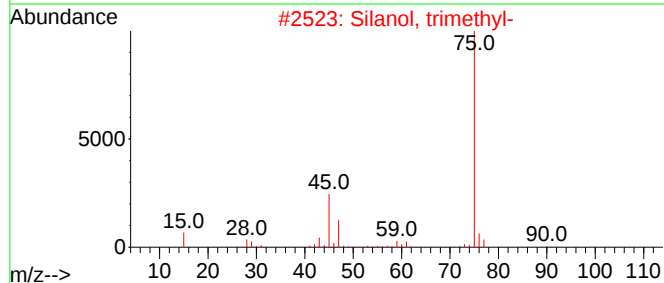
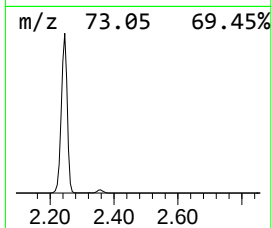
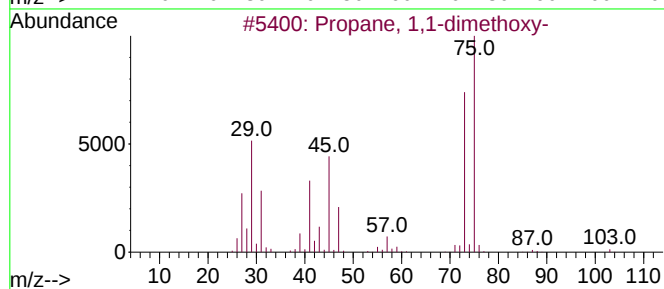
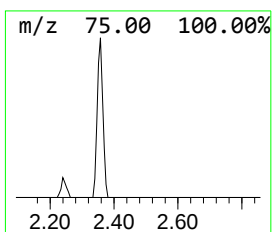
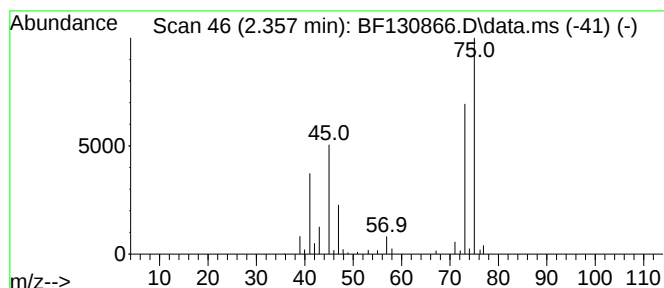
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 2 unknown2.357 Concentration Rank 13

| R.T.  | EstConc | Area  | Relative to ISTD       | R.T.  |
|-------|---------|-------|------------------------|-------|
| 2.357 | 2.49 ng | 52298 | 1,4-Dichlorobenzene-d4 | 6.945 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Propane, 1,1-dimethoxy-             | 104 | C5H12O2  | 004744-10-9 | 40   |
| 2    |    |   | Silanol, trimethyl-                 | 90  | C3H10OSi | 001066-40-6 | 40   |
| 3    |    |   | Ethanethioamide                     | 75  | C2H5NS   | 000062-55-5 | 9    |
| 4    |    |   | 2,3-Butanediol                      | 90  | C4H10O2  | 000513-85-9 | 9    |
| 5    |    |   | Trimethylene glycol monomethyl e... | 90  | C4H10O2  | 001589-49-7 | 9    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

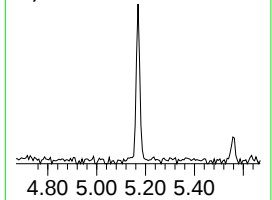
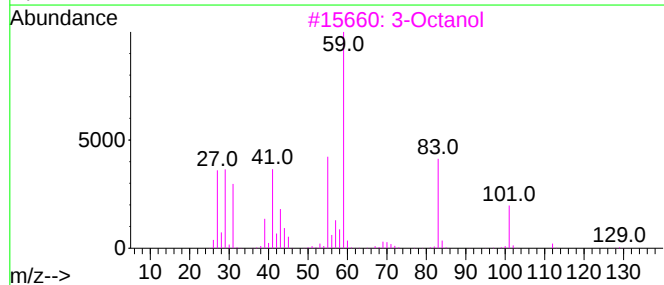
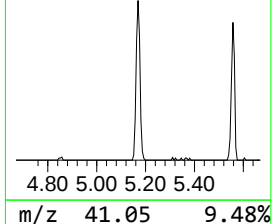
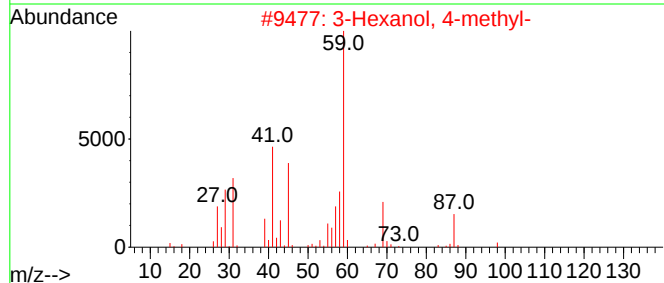
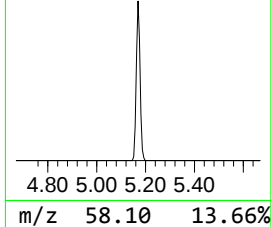
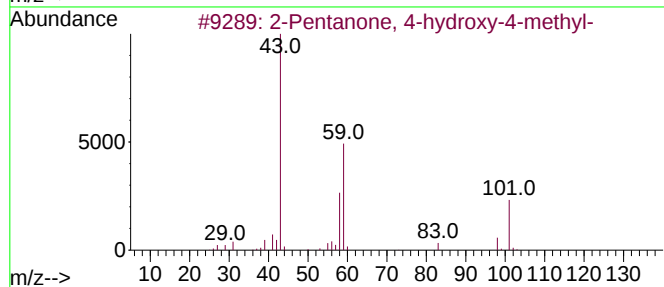
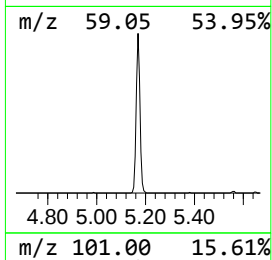
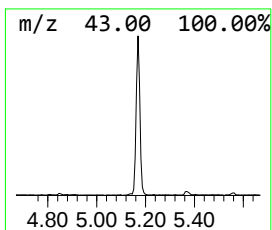
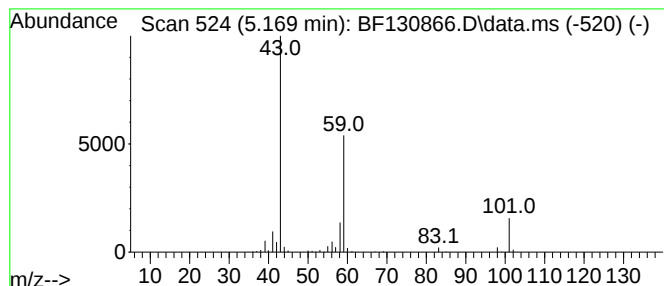
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 5.169 | 11.37 ng | 238573 | 1,4-Dichlorobenzene-d4 | 6.945 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|------|-------------------------------------|-----|----------|-------------|------|
| 1    |      | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 56   |
| 2    |      | 3-Hexanol, 4-methyl-                | 116 | C7H16O   | 000615-29-2 | 33   |
| 3    |      | 3-Octanol                           | 130 | C8H18O   | 000589-98-0 | 33   |
| 4    |      | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 32   |
| 5    |      | 1-Propen-2-ol, acetate              | 100 | C5H8O2   | 000108-22-5 | 12   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

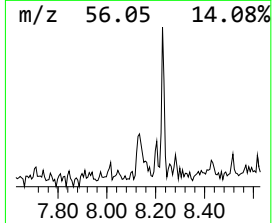
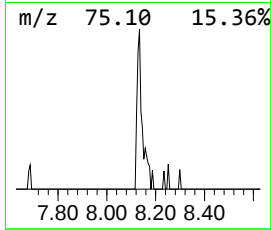
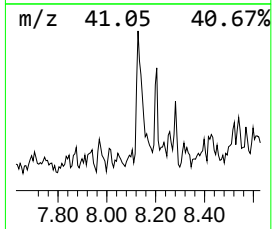
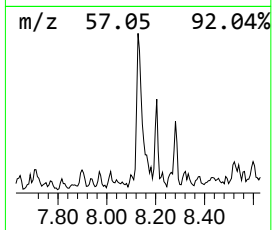
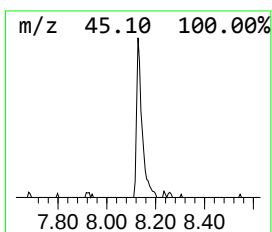
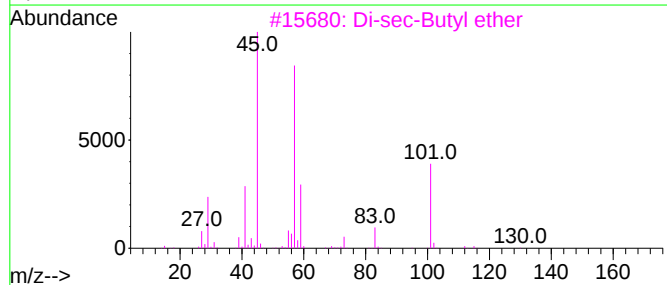
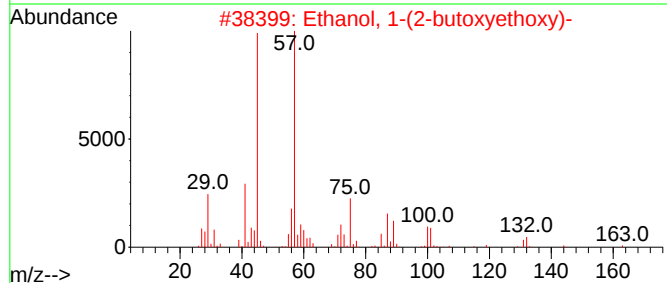
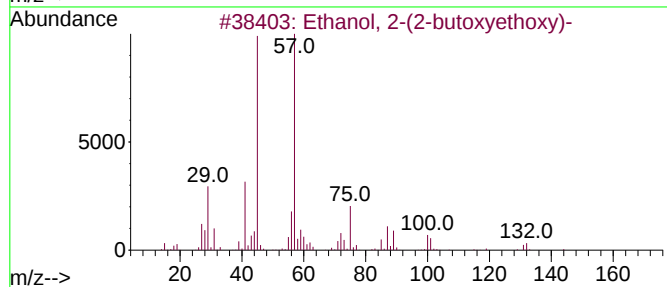
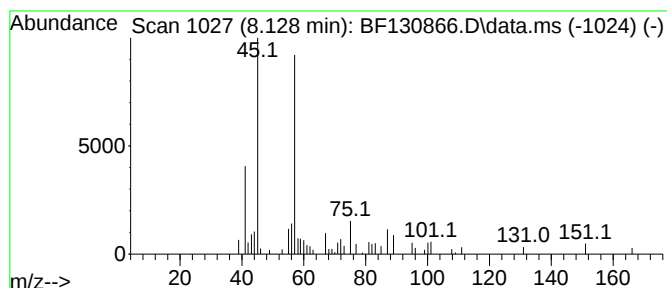
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 4 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 11

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 8.128 | 2.76 ng | 60958 | Naphthalene-d8   | 8.228 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Ethanol, 2-(2-butoxyethoxy)-         | 162 | C8H18O3  | 000112-34-5 | 86   |
| 2    |    |   | Ethanol, 1-(2-butoxyethoxy)-         | 162 | C8H18O3  | 054446-78-5 | 78   |
| 3    |    |   | Di-sec-Butyl ether                   | 130 | C8H18O   | 006863-58-7 | 53   |
| 4    |    |   | Butane, 1,1'-[ethylidenebis(oxy)...] | 174 | C10H22O2 | 000871-22-7 | 53   |
| 5    |    |   | Tetraethylene glycol                 | 194 | C8H18O5  | 000112-60-7 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

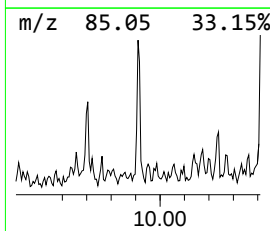
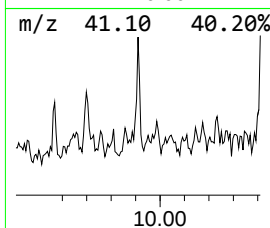
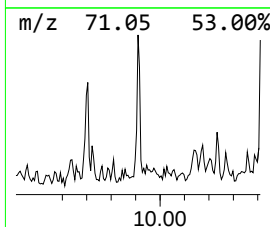
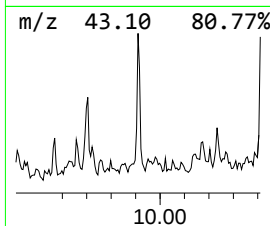
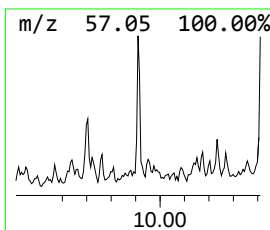
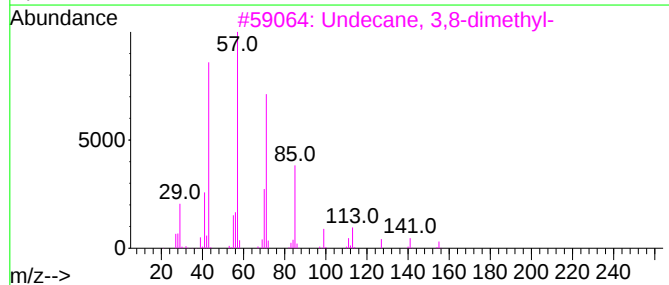
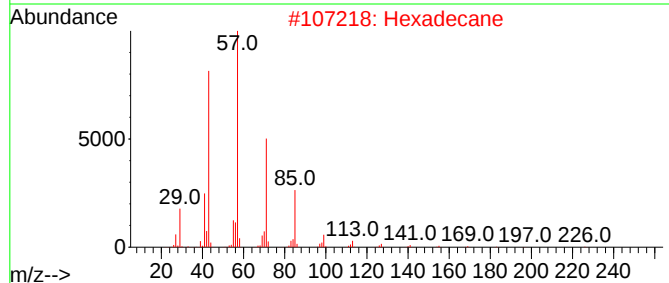
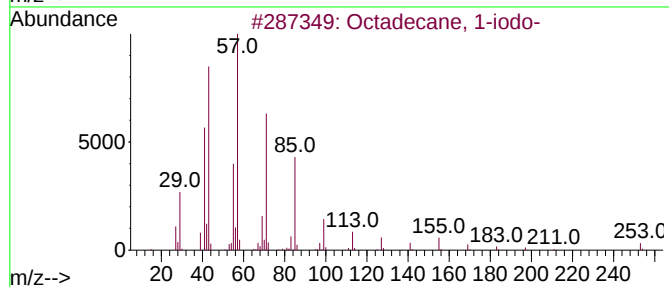
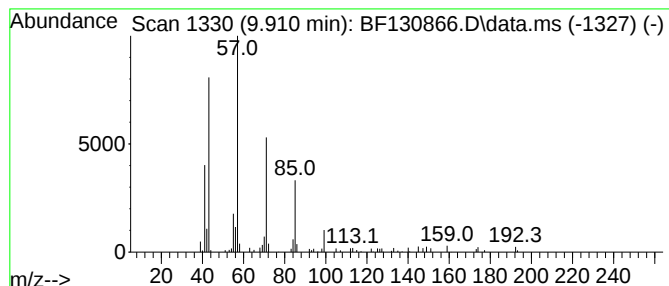
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 5 Octadecane, 1-iodo- Concentration Rank 15

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 9.910 | 2.13 ng | 47136 | Acenaphthene-d10 | 9.992 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm     | CAS#        | Qual |
|------|----|---|-----------------------------|-----|-------------|-------------|------|
| 1    |    |   | Octadecane, 1-iodo-         | 380 | C18H37I     | 000629-93-6 | 83   |
| 2    |    |   | Hexadecane                  | 226 | C16H34      | 000544-76-3 | 83   |
| 3    |    |   | Undecane, 3,8-dimethyl-     | 184 | C13H28      | 017301-30-3 | 83   |
| 4    |    |   | Silane, trichlorooctadecyl- | 386 | C18H37Cl3Si | 000112-04-9 | 78   |
| 5    |    |   | Dodecane, 1-iodo-           | 296 | C12H25I     | 004292-19-7 | 78   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

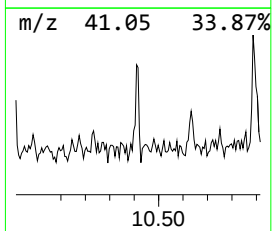
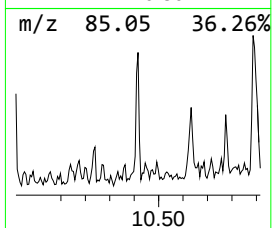
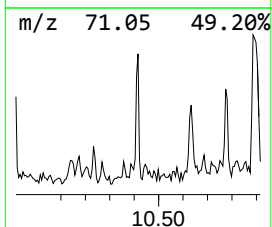
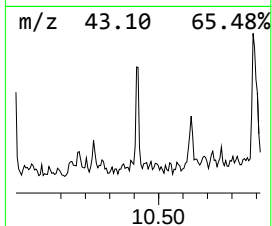
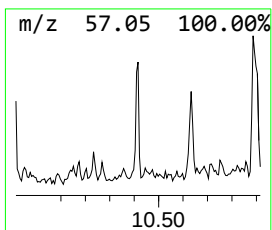
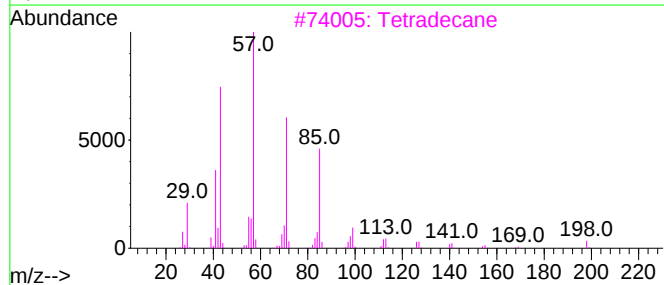
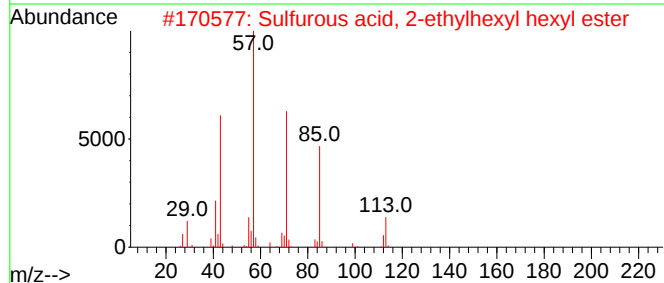
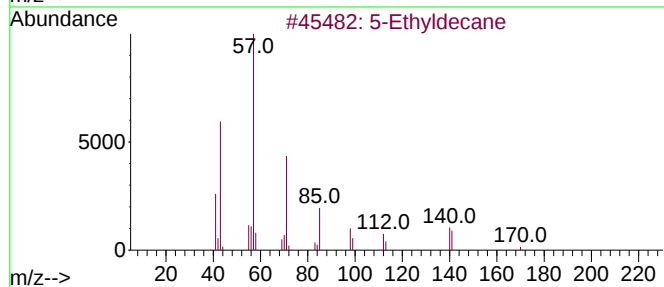
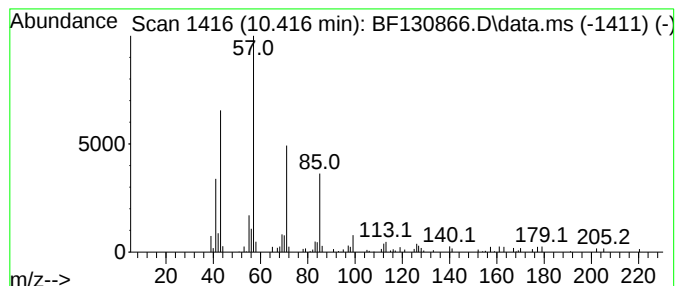
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 6 5-Ethyldecane Concentration Rank 14

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.  |
|--------|---------|-------|------------------|-------|
| 10.416 | 2.26 ng | 50031 | Acenaphthene-d10 | 9.992 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 5-Ethyldecane                       | 170 | C12H26    | 017302-36-2  | 83   |
| 2    |    |   | Sulfurous acid, 2-ethylhexyl hex... | 278 | C14H30O3S | 1000309-20-2 | 80   |
| 3    |    |   | Tetradecane                         | 198 | C14H30    | 000629-59-4  | 80   |
| 4    |    |   | Hexadecane                          | 226 | C16H34    | 000544-76-3  | 80   |
| 5    |    |   | Tetradecane, 1-iodo-                | 324 | C14H29I   | 019218-94-1  | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

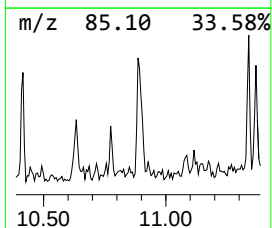
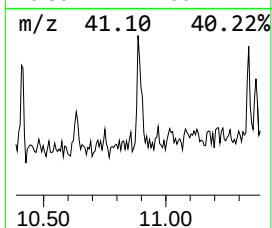
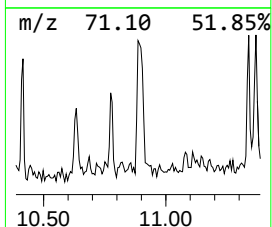
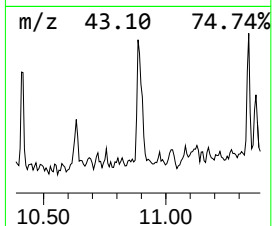
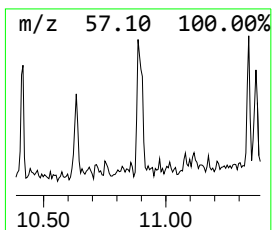
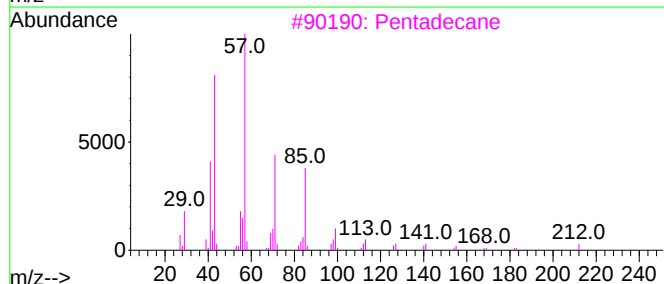
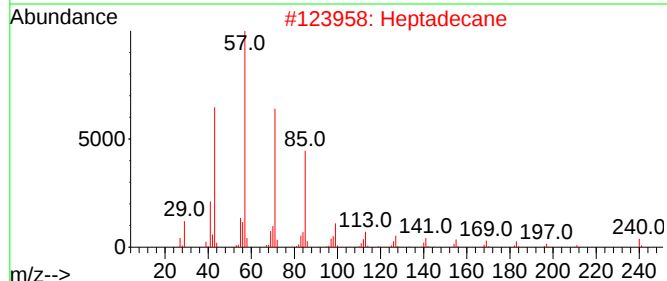
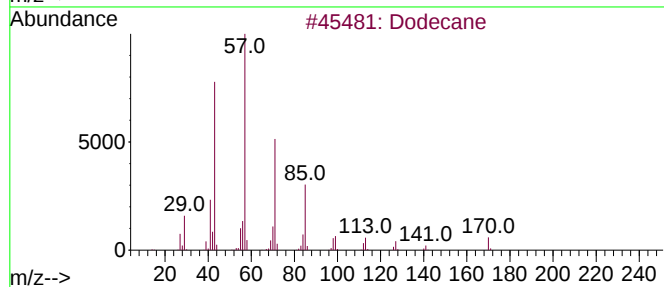
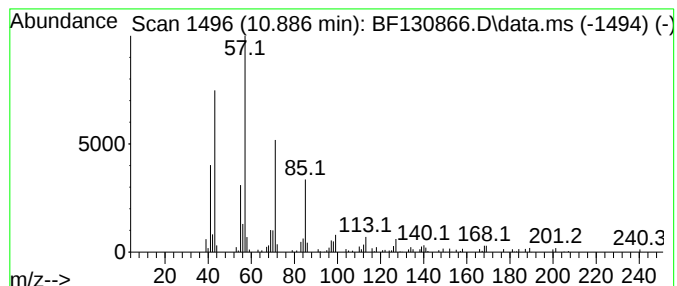
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 7 Dodecane Concentration Rank 8

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 10.886 | 4.28 ng | 105567 | Phenanthrene-d10 | 11.480 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Dodecane     | 170 | C12H26  | 000112-40-3 | 90   |
| 2    |    |   | Heptadecane  | 240 | C17H36  | 000629-78-7 | 90   |
| 3    |    |   | Pentadecane  | 212 | C15H32  | 000629-62-9 | 86   |
| 4    |    |   | Tetradecane  | 198 | C14H30  | 000629-59-4 | 86   |
| 5    |    |   | Tridecane    | 184 | C13H28  | 000629-50-5 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

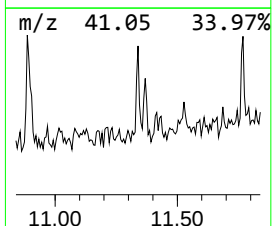
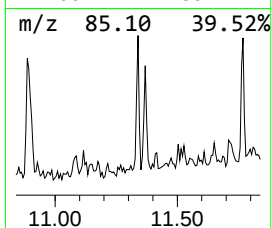
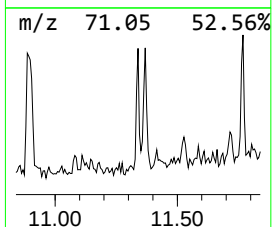
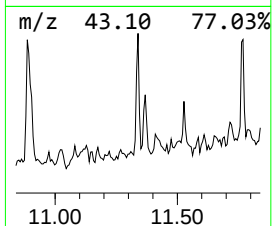
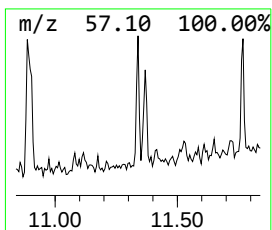
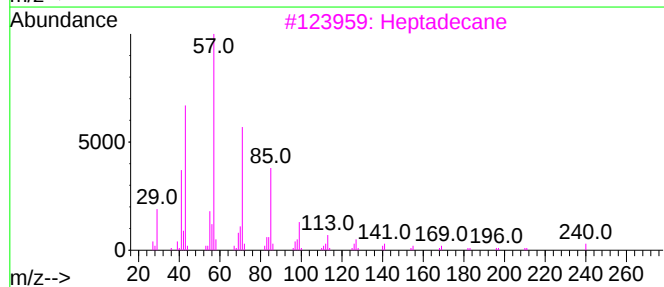
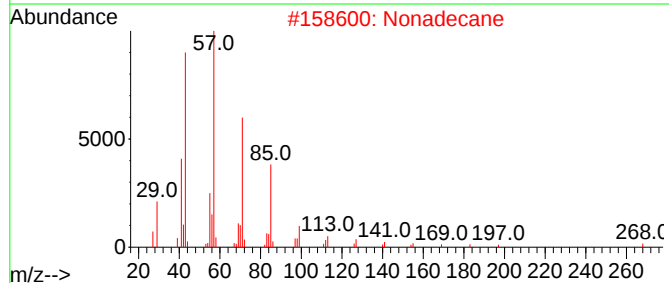
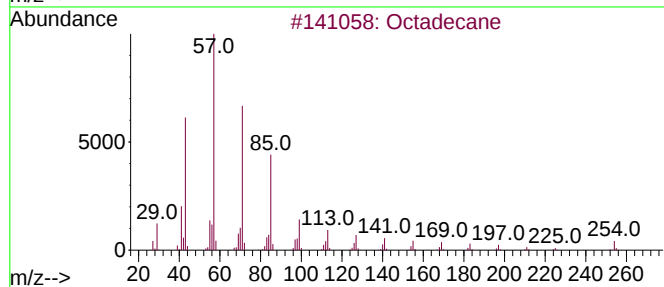
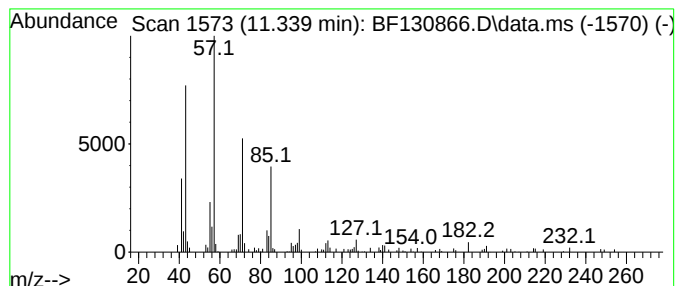
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 8 Octadecane Concentration Rank 16

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 11.339 | 2.09 ng | 51673 | Phenanthrene-d10 | 11.480 |

| Hit# | of | 5 | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------|-----|---------|-------------|------|
| 1    |    |   | Octadecane        | 254 | C18H38  | 000593-45-3 | 91   |
| 2    |    |   | Nonadecane        | 268 | C19H40  | 000629-92-5 | 83   |
| 3    |    |   | Heptadecane       | 240 | C17H36  | 000629-78-7 | 83   |
| 4    |    |   | Tetradecane       | 198 | C14H30  | 000629-59-4 | 80   |
| 5    |    |   | Dodecane, 1-iodo- | 296 | C12H25I | 004292-19-7 | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

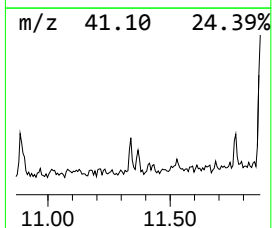
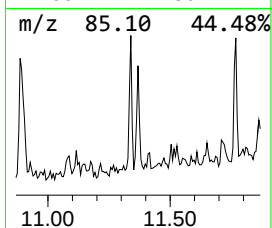
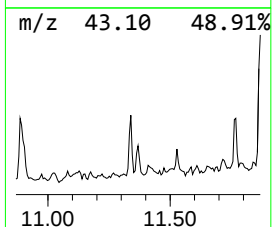
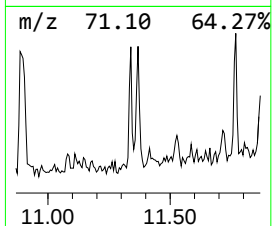
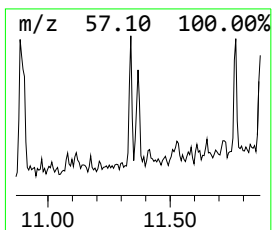
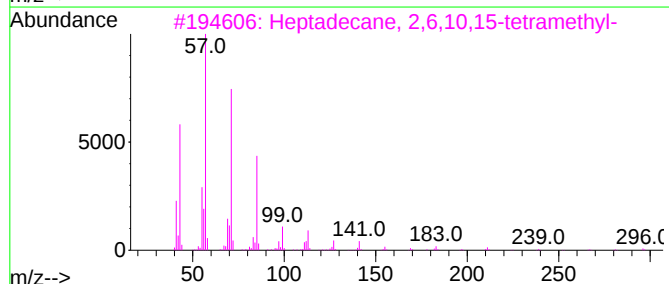
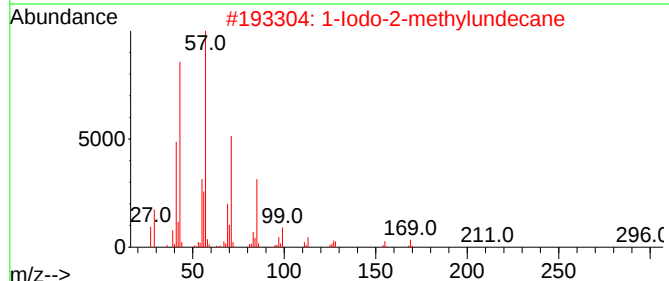
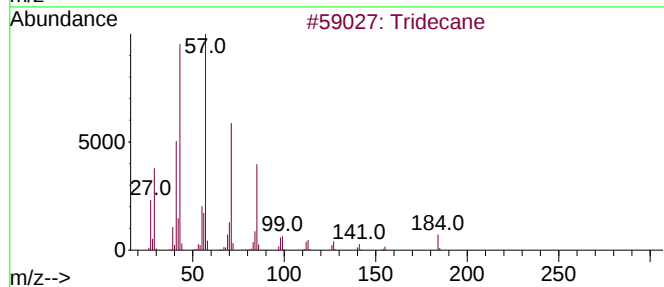
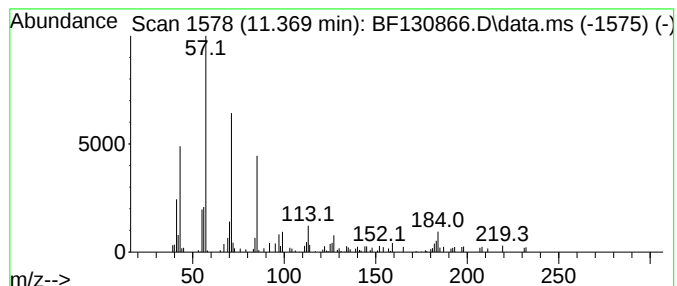
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*

Peak Number 9 Tridecane Concentration Rank 12

| R.T.      | EstConc                             | Area  | Relative to ISTD |         | R.T.        |      |
|-----------|-------------------------------------|-------|------------------|---------|-------------|------|
| 11.369    | 2.58 ng                             | 63745 | Phenanthrene-d10 |         | 11.480      |      |
| Hit# of 5 | Tentative ID                        |       | MW               | MolForm | CAS#        | Qual |
| 1         | Tridecane                           |       | 184              | C13H28  | 000629-50-5 | 90   |
| 2         | 1-Iodo-2-methylundecane             |       | 296              | C12H25I | 073105-67-6 | 72   |
| 3         | Heptadecane, 2,6,10,15-tetramethyl- |       | 296              | C21H44  | 054833-48-6 | 72   |
| 4         | Dodecane, 3-methyl-                 |       | 184              | C13H28  | 017312-57-1 | 64   |
| 5         | Dodecane, 2,7,10-trimethyl-         |       | 212              | C15H32  | 074645-98-0 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

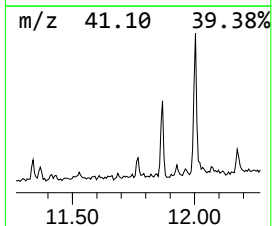
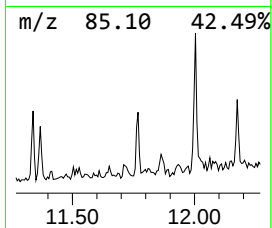
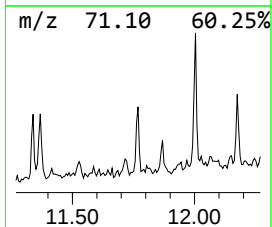
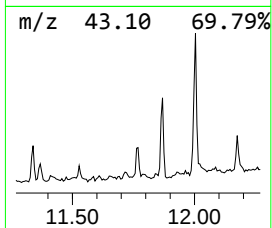
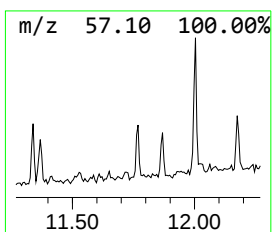
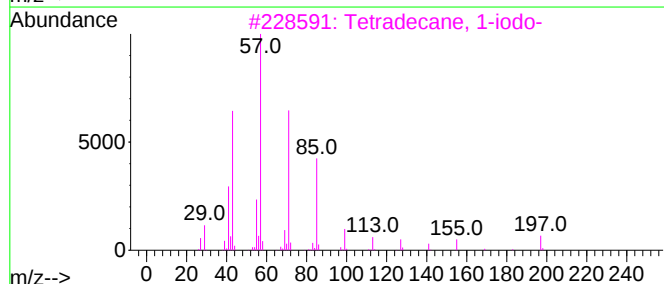
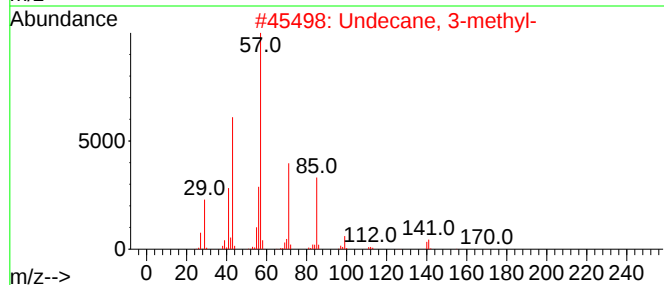
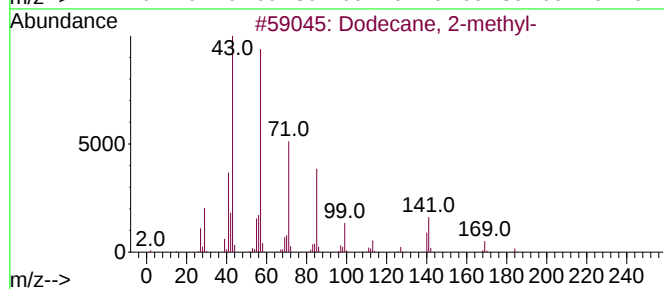
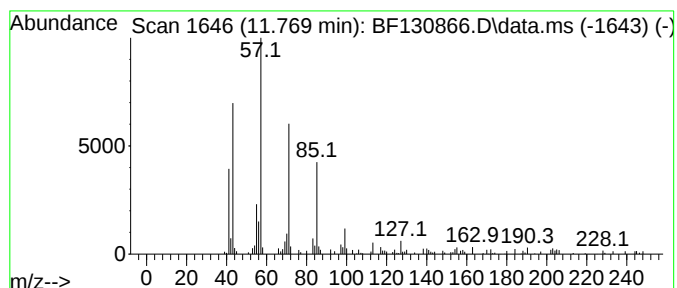
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*

Peak Number 10 Dodecane, 2-methyl- Concentration Rank 18

| R.T.      | EstConc               | Area  | Relative to ISTD | R.T.        |      |
|-----------|-----------------------|-------|------------------|-------------|------|
| 11.769    | 2.07 ng               | 51047 | Phenanthrene-d10 | 11.480      |      |
| Hit# of 5 | Tentative ID          | MW    | MolForm          | CAS#        | Qual |
| 1         | Dodecane, 2-methyl-   | 184   | C13H28           | 001560-97-0 | 80   |
| 2         | Undecane, 3-methyl-   | 170   | C12H26           | 001002-43-3 | 80   |
| 3         | Tetradecane, 1-iodo-  | 324   | C14H29I          | 019218-94-1 | 80   |
| 4         | Dodecane              | 170   | C12H26           | 000112-40-3 | 80   |
| 5         | Decane, 3,6-dimethyl- | 170   | C12H26           | 017312-53-7 | 74   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

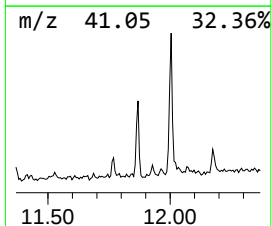
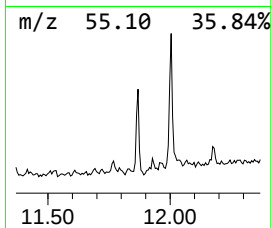
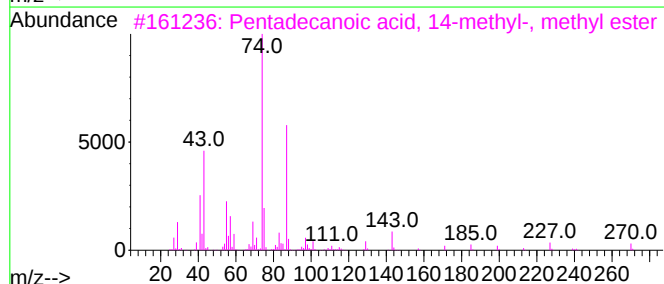
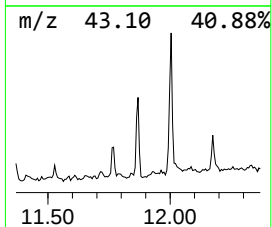
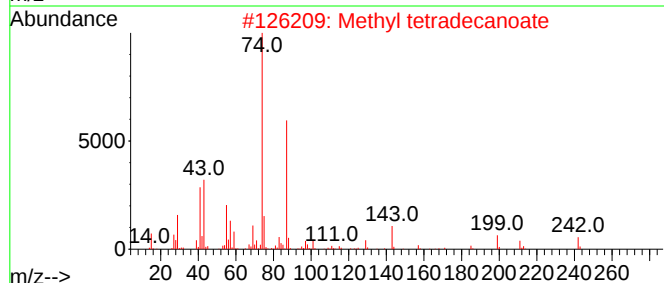
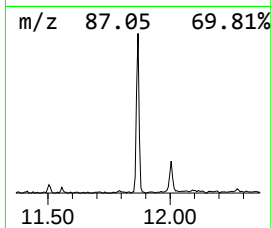
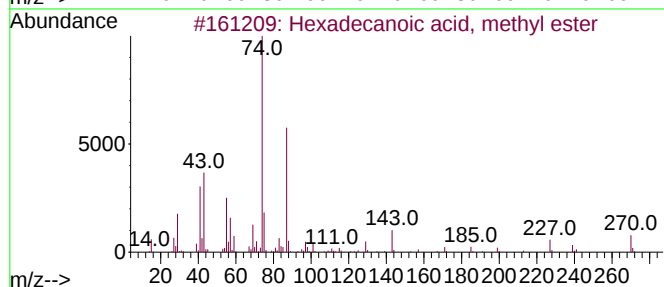
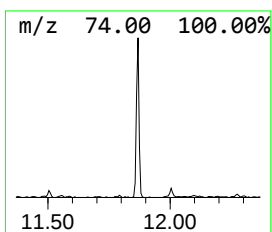
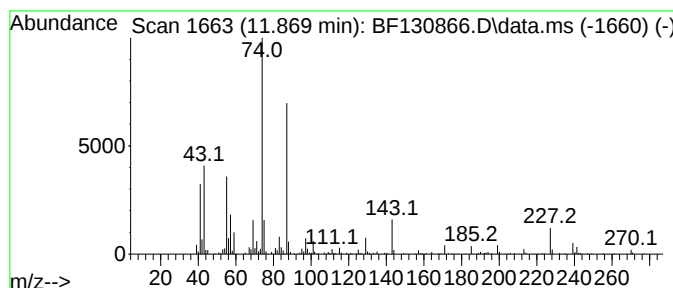
\*\*\*\*\*

Peak Number 11 Hexadecanoic acid, methyl e... Concentration Rank 6

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 11.869 | 8.81 ng | 217389 | Phenanthrene-d10 | 11.480 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|------|-------------------------------------|-----|----------|-------------|------|
| 1    |      | Hexadecanoic acid, methyl ester     | 270 | C17H34O2 | 000112-39-0 | 91   |
| 2    |      | Methyl tetradecanoate               | 242 | C15H30O2 | 000124-10-7 | 87   |
| 3    |      | Pentadecanoic acid, 14-methyl-, ... | 270 | C17H34O2 | 005129-60-2 | 86   |
| 4    |      | Pentadecanoic acid, methyl ester    | 256 | C16H32O2 | 007132-64-1 | 86   |
| 5    |      | Tridecanoic acid, methyl ester      | 228 | C14H28O2 | 001731-88-0 | 81   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

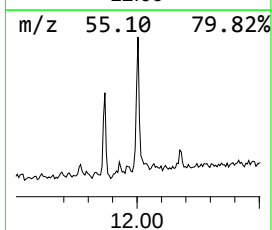
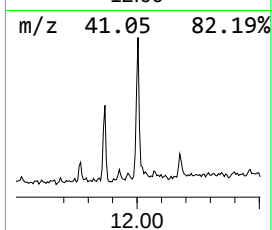
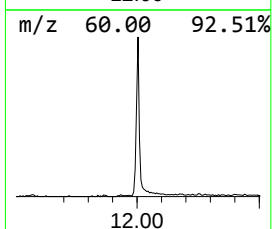
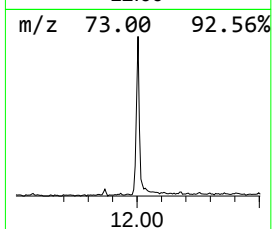
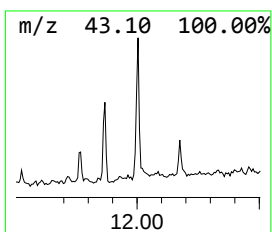
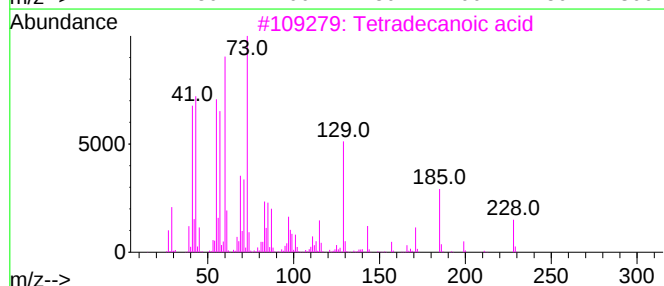
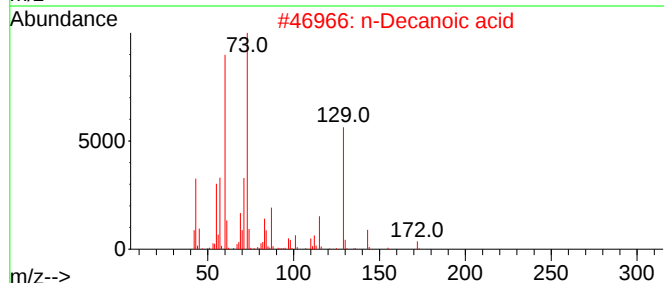
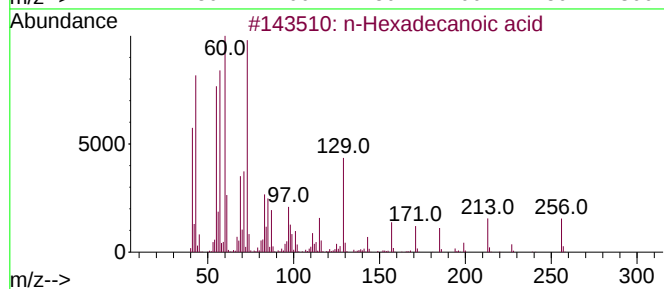
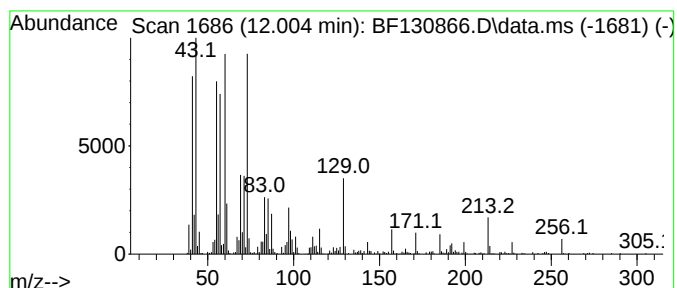
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*

Peak Number 12 n-Hexadecanoic acid Concentration Rank 4

| R.T.      | EstConc             | Area   | Relative to ISTD | R.T.        |      |
|-----------|---------------------|--------|------------------|-------------|------|
| 12.004    | 18.45 ng            | 455194 | Phenanthrene-d10 | 11.480      |      |
| Hit# of 5 | Tentative ID        | MW     | MolForm          | CAS#        | Qual |
| 1         | n-Hexadecanoic acid | 256    | C16H32O2         | 000057-10-3 | 98   |
| 2         | n-Decanoic acid     | 172    | C10H20O2         | 000334-48-5 | 91   |
| 3         | Tetradecanoic acid  | 228    | C14H28O2         | 000544-63-8 | 90   |
| 4         | Octadecanoic acid   | 284    | C18H36O2         | 000057-11-4 | 90   |
| 5         | Pentadecanoic acid  | 242    | C15H30O2         | 001002-84-2 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

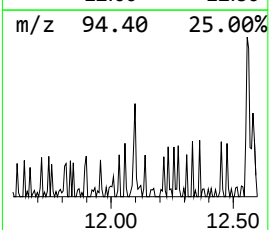
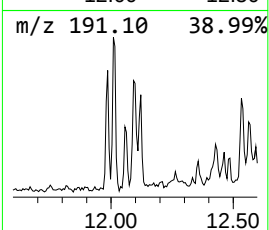
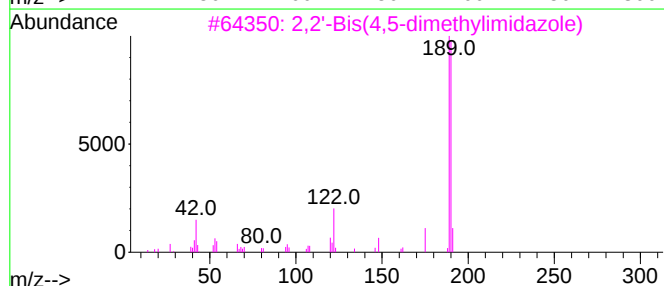
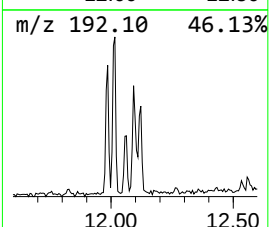
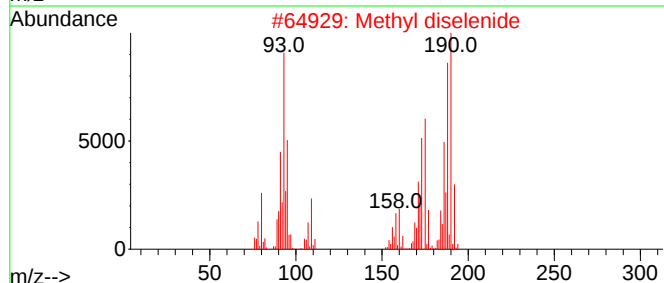
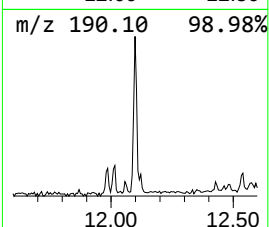
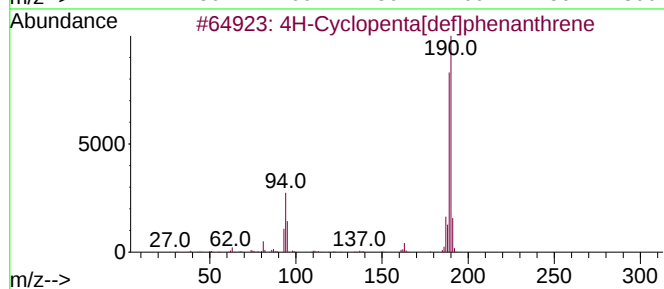
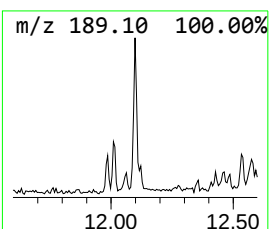
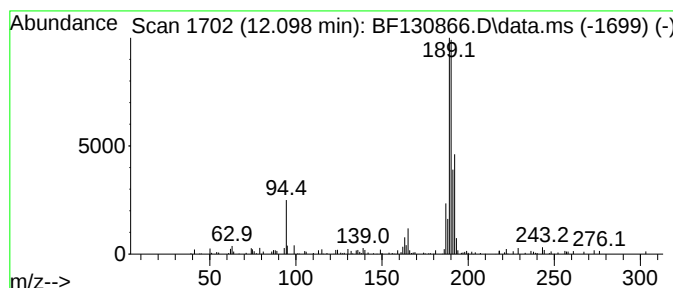
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*

Peak Number 13 4H-Cyclopenta[def]phenanthrene Concentration Rank 17

| R.T.      | EstConc                             | Area  | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|-------|------------------|--------------|------|
| 12.098    | 2.09 ng                             | 51572 | Phenanthrene-d10 | 11.480       |      |
| Hit# of 5 | Tentative ID                        | MW    | MolForm          | CAS#         | Qual |
| 1         | 4H-Cyclopenta[def]phenanthrene      | 190   | C15H10           | 000203-64-5  | 50   |
| 2         | Methyl diselenide                   | 190   | C2H6Se2          | 007101-31-7  | 45   |
| 3         | 2,2'-Bis(4,5-dimethylimidazole)     | 190   | C10H14N4         | 069286-06-2  | 43   |
| 4         | Carbonic acid, ethyl 2-formyl-4,... | 262   | C10H8Cl2O4       | 1000331-34-4 | 40   |
| 5         | 2-Amino-4,6-dichloropyrimidine-5... | 191   | C5H3Cl2N3O       | 005604-46-6  | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

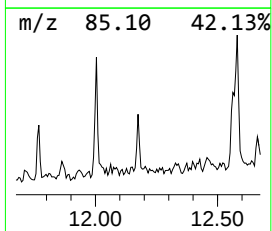
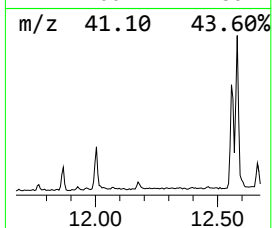
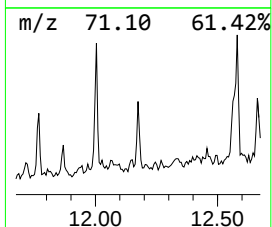
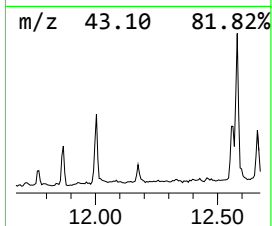
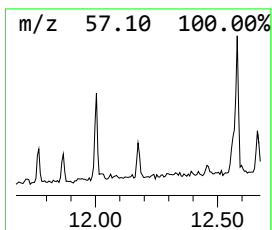
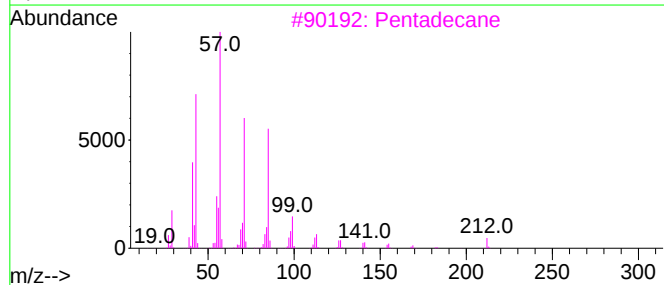
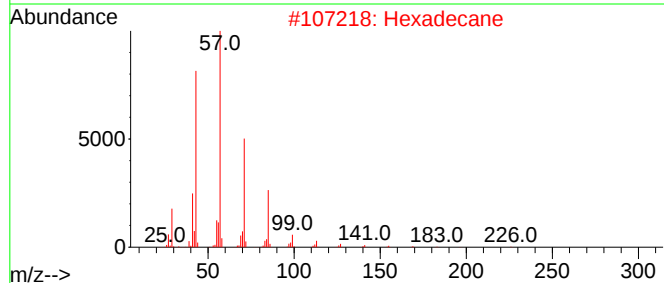
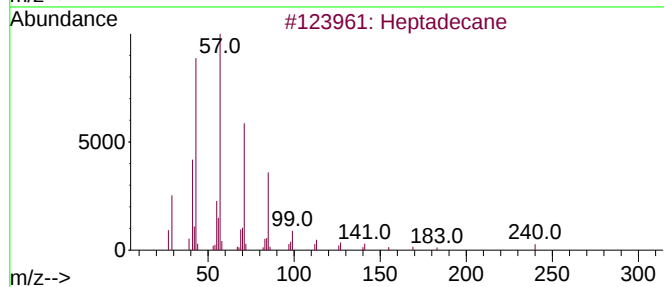
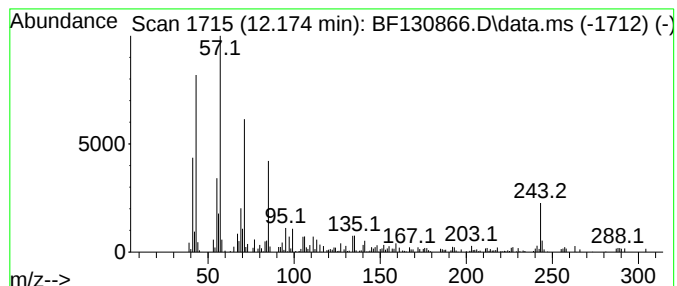
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 14 Heptadecane Concentration Rank 9

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 12.175 | 3.86 ng | 95254 | Phenanthrene-d10 | 11.480 |

| Hit# | of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------|--------------|-----|---------|-------------|------|
| 1    |      | Heptadecane  | 240 | C17H36  | 000629-78-7 | 86   |
| 2    |      | Hexadecane   | 226 | C16H34  | 000544-76-3 | 86   |
| 3    |      | Pentadecane  | 212 | C15H32  | 000629-62-9 | 62   |
| 4    |      | Nonadecane   | 268 | C19H40  | 000629-92-5 | 62   |
| 5    |      | Octacosane   | 394 | C28H58  | 000630-02-4 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

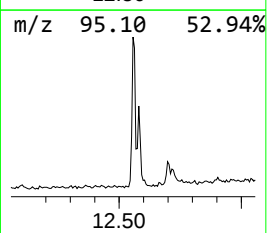
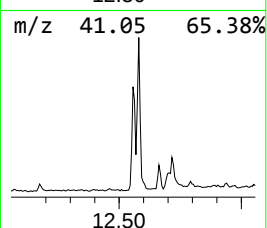
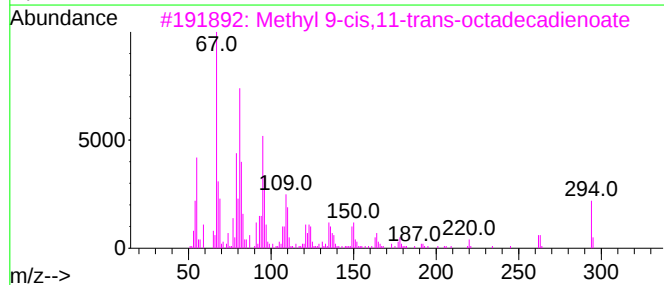
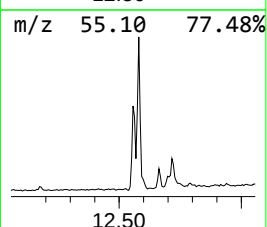
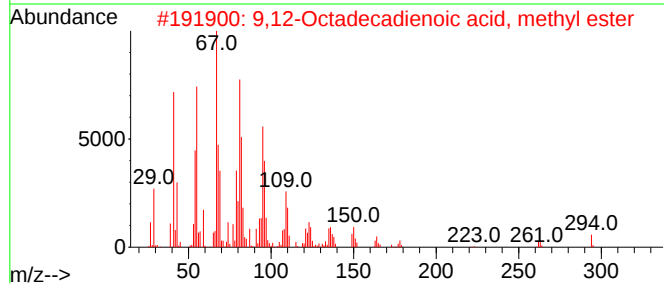
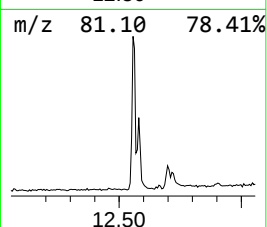
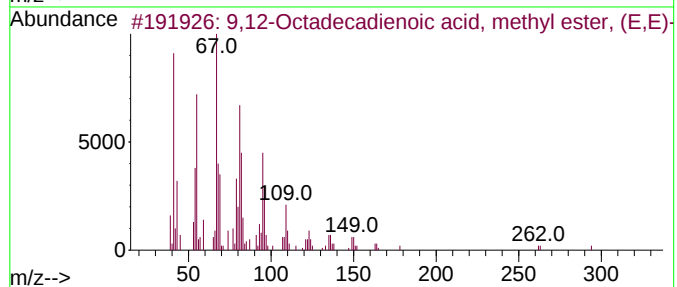
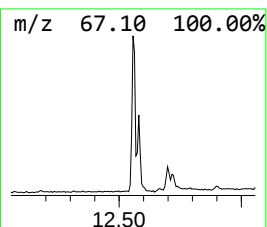
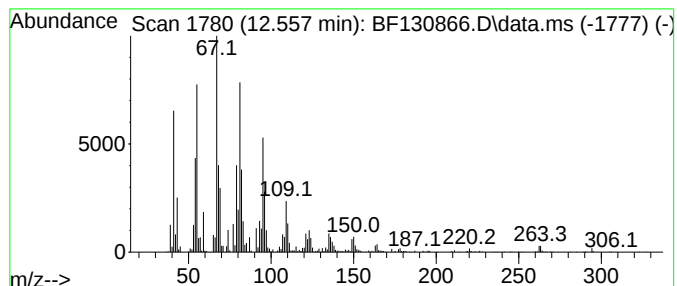
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*

Peak Number 15 9,12-Octadecadienoic acid, ... Concentration Rank 3

| R.T.   | EstConc  | Area                                | Relative to ISTD |          | R.T.        |      |
|--------|----------|-------------------------------------|------------------|----------|-------------|------|
| 12.557 | 49.01 ng | 1208960                             | Phenanthrene-d10 |          | 11.480      |      |
| Hit#   | of 5     | Tentative ID                        | MW               | MolForm  | CAS#        | Qual |
| 1      |          | 9,12-Octadecadienoic acid, methy... | 294              | C19H34O2 | 002566-97-4 | 99   |
| 2      |          | 9,12-Octadecadienoic acid, methy... | 294              | C19H34O2 | 002462-85-3 | 99   |
| 3      |          | Methyl 9-cis,11-trans-octadecadi... | 294              | C19H34O2 | 013058-52-1 | 99   |
| 4      |          | 9,12-Octadecadienoic acid (Z,Z)-... | 294              | C19H34O2 | 000112-63-0 | 99   |
| 5      |          | 10,13-Octadecadienoic acid, meth... | 294              | C19H34O2 | 056554-62-2 | 97   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

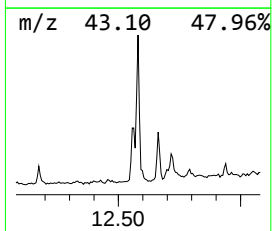
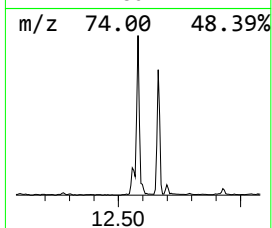
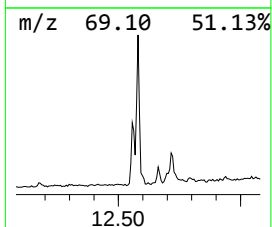
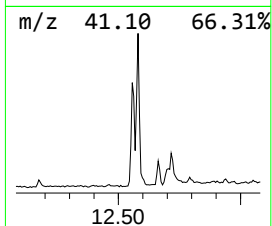
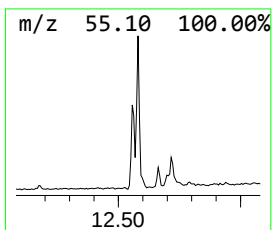
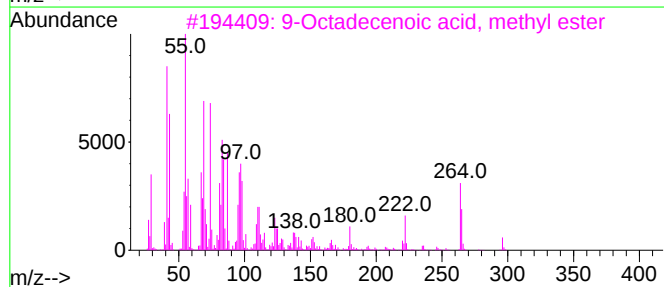
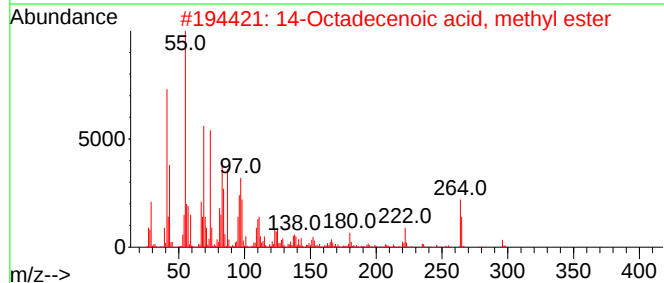
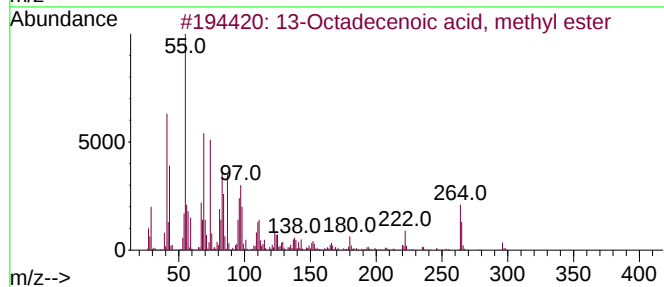
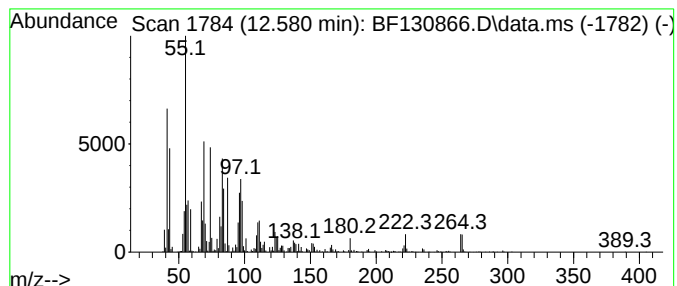
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*

Peak Number 16 13-Octadecenoic acid, methyl... Concentration Rank 2

| R.T.      | EstConc                            | Area    | Relative to ISTD |          | R.T.        |      |
|-----------|------------------------------------|---------|------------------|----------|-------------|------|
| 12.580    | 51.60 ng                           | 1272910 | Phenanthrene-d10 |          | 11.480      |      |
| Hit# of 5 | Tentative ID                       |         | MW               | MolForm  | CAS#        | Qual |
| 1         | 13-Octadecenoic acid, methyl ester |         | 296              | C19H36O2 | 056554-47-3 | 99   |
| 2         | 14-Octadecenoic acid, methyl ester |         | 296              | C19H36O2 | 056554-48-4 | 99   |
| 3         | 9-Octadecenoic acid, methyl ester  |         | 296              | C19H36O2 | 002462-84-2 | 98   |
| 4         | 16-Octadecenoic acid, methyl ester |         | 296              | C19H36O2 | 056554-49-5 | 96   |
| 5         | 8-Octadecenoic acid, methyl ester  |         | 296              | C19H36O2 | 002345-29-1 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

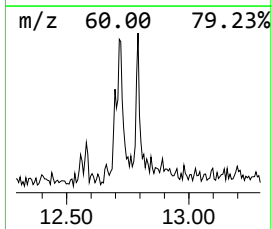
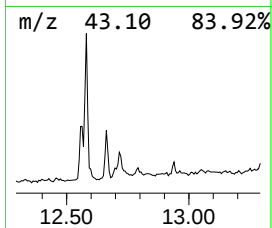
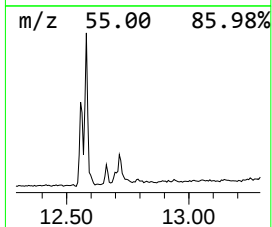
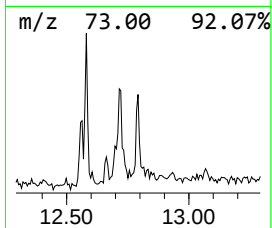
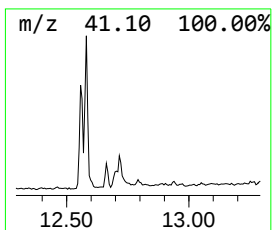
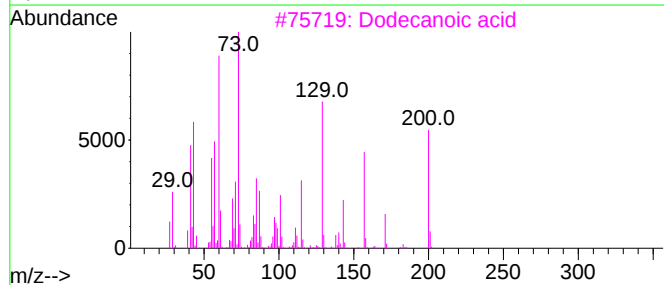
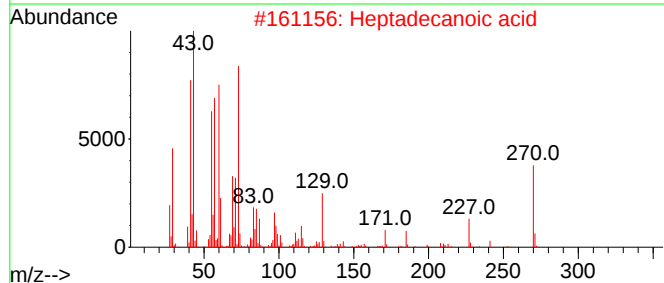
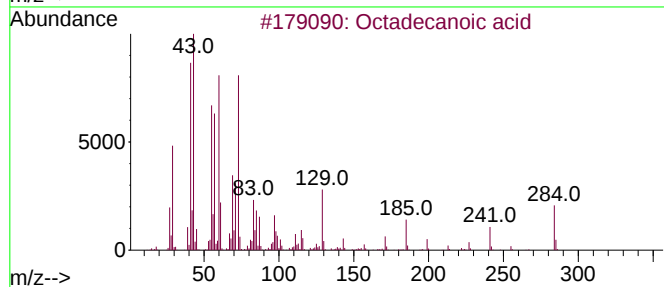
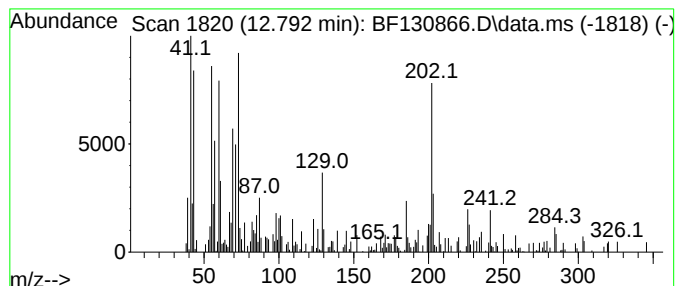
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 17 Octadecanoic acid Concentration Rank 10

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 12.792 | 2.79 ng | 68769 | Phenanthrene-d10 | 11.480 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Octadecanoic acid                   | 284 | C18H36O2 | 000057-11-4  | 50   |
| 2    |    |   | Heptadecanoic acid                  | 270 | C17H34O2 | 000506-12-7  | 30   |
| 3    |    |   | Dodecanoic acid                     | 200 | C12H24O2 | 000143-07-7  | 30   |
| 4    |    |   | Pyrene                              | 202 | C16H10   | 000129-00-0  | 25   |
| 5    |    |   | 4-Methoxy-naphthalene-1-carboxyl... | 202 | C12H10O3 | 1000317-85-4 | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
Data File : BF130866.D  
Acq On : 29 Oct 2022 19:49  
Operator : CG\JU  
Sample : N5308-01  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PL-2-102622

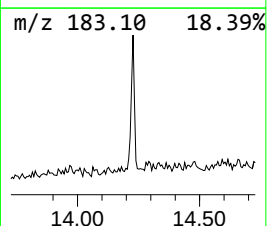
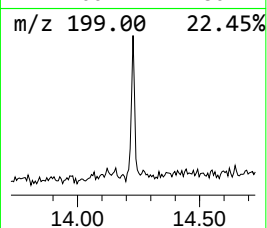
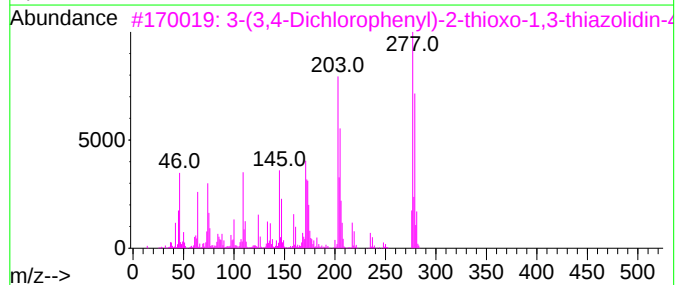
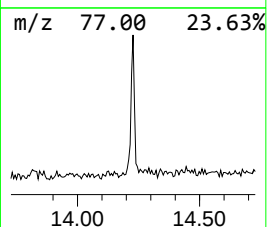
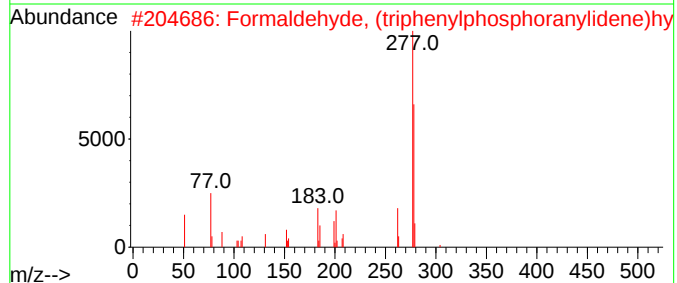
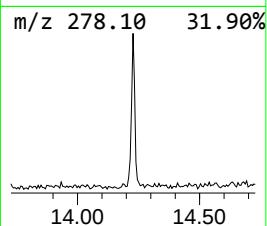
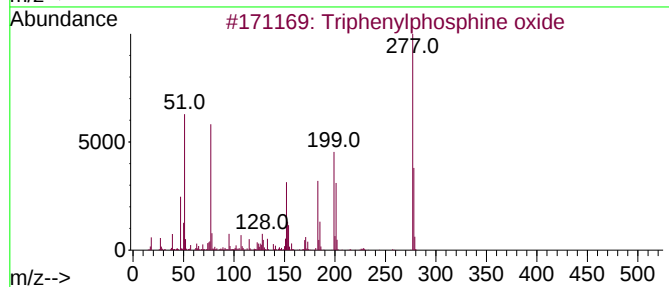
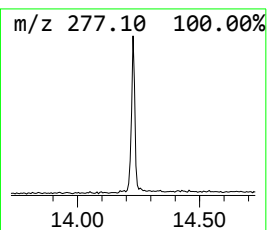
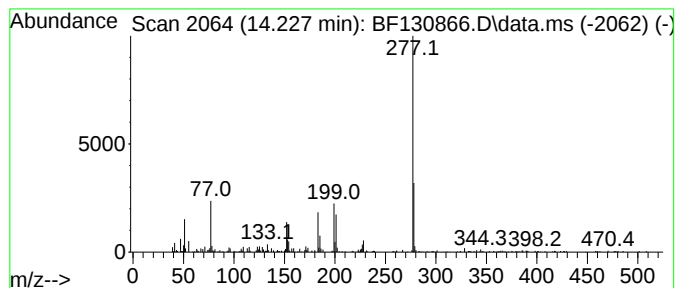
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 18 Triphenylphosphine oxide Concentration Rank 7

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 14.227 | 5.79 ng | 159928 | Chrysene-d12     | 14.133 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Triphenylphosphine oxide            | 278 | C18H15OP    | 000791-28-6  | 93   |
| 2    |    |   | Formaldehyde, (triphenylphosphor... | 304 | C19H17N2P   | 015990-54-2  | 50   |
| 3    |    |   | 3-(3,4-Dichlorophenyl)-2-thioxo-... | 277 | C9H5Cl2NOS2 | 017046-34-3  | 43   |
| 4    |    |   | Vanadium, cycloheptatrienyl-(pen... | 277 | C17H22V     | 1000155-20-5 | 38   |
| 5    |    |   | (E)-7-Benzylidene-1-azabicyclo[3... | 293 | C15H19N03S  | 1000483-83-4 | 37   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102922\  
 Data File : BF130866.D  
 Acq On : 29 Oct 2022 19:49  
 Operator : CG\JU  
 Sample : N5308-01  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PL-2-102622

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.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... | 2.246  | 65.0    | ng    | 1362620  | 1                     | 6.945  | 419592 | 20.0 |
| unknown2.357       | 2.357  | 2.5     | ng    | 52298    | 1                     | 6.945  | 419592 | 20.0 |
| 2-Pentanone, 4-... | 5.169  | 11.4    | ng    | 238573   | 1                     | 6.945  | 419592 | 20.0 |
| Ethanol, 2-(2-b... | 8.128  | 2.8     | ng    | 60958    | 2                     | 8.228  | 441932 | 20.0 |
| Octadecane, 1-i... | 9.910  | 2.1     | ng    | 47136    | 3                     | 9.992  | 442032 | 20.0 |
| 5-Ethyldecane      | 10.416 | 2.3     | ng    | 50031    | 3                     | 9.992  | 442032 | 20.0 |
| Dodecane           | 10.886 | 4.3     | ng    | 105567   | 4                     | 11.480 | 493345 | 20.0 |
| Octadecane         | 11.339 | 2.1     | ng    | 51673    | 4                     | 11.480 | 493345 | 20.0 |
| Tridecane          | 11.369 | 2.6     | ng    | 63745    | 4                     | 11.480 | 493345 | 20.0 |
| Dodecane, 2-met... | 11.769 | 2.1     | ng    | 51047    | 4                     | 11.480 | 493345 | 20.0 |
| Hexadecanoic ac... | 11.869 | 8.8     | ng    | 217389   | 4                     | 11.480 | 493345 | 20.0 |
| n-Hexadecanoic ... | 12.004 | 18.4    | ng    | 455194   | 4                     | 11.480 | 493345 | 20.0 |
| 4H-Cyclopenta[d... | 12.098 | 2.1     | ng    | 51572    | 4                     | 11.480 | 493345 | 20.0 |
| Heptadecane        | 12.175 | 3.9     | ng    | 95254    | 4                     | 11.480 | 493345 | 20.0 |
| 9,12-Octadecadi... | 12.557 | 49.0    | ng    | 1208960  | 4                     | 11.480 | 493345 | 20.0 |
| 13-Octadecenoic... | 12.580 | 51.6    | ng    | 1272910  | 4                     | 11.480 | 493345 | 20.0 |
| Octadecanoic acid  | 12.792 | 2.8     | ng    | 68769    | 4                     | 11.480 | 493345 | 20.0 |
| Triphenylphosph... | 14.227 | 5.8     | ng    | 159928   | 5                     | 14.133 | 552689 | 20.0 |