

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072722\  
 Data File : BF129383.D  
 Acq On : 27 Jul 2022 18:51  
 Operator : CG\JU  
 Sample : N3861-13  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 MW-21RR

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-BF072522.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF129383.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         | 3.263       | 161           | 166         | 176          | rVB      | 103044         | 180799        | 2.08%           | 0.349%        |
| 2         | 5.116       | 478           | 481         | 486          | rBV      | 164157         | 187616        | 2.16%           | 0.363%        |
| 3         | 5.504       | 543           | 547         | 551          | rVB      | 3936717        | 3593457       | 41.30%          | 6.944%        |
| 4         | 6.051       | 636           | 640         | 643          | rBV      | 301897         | 254577        | 2.93%           | 0.492%        |
| 5         | 6.340       | 686           | 689         | 693          | rVB      | 388662         | 294526        | 3.38%           | 0.569%        |
| 6         | 6.510       | 713           | 718         | 724          | rVB      | 2389209        | 2263609       | 26.01%          | 4.374%        |
| 7         | 6.881       | 777           | 781         | 784          | rBV      | 1951159        | 1597074       | 18.35%          | 3.086%        |
| 8         | 7.057       | 808           | 811         | 815          | rVB      | 952215         | 888425        | 10.21%          | 1.717%        |
| 9         | 7.128       | 819           | 823         | 826          | rVB      | 236227         | 213202        | 2.45%           | 0.412%        |
| 10        | 7.216       | 831           | 838         | 844          | rVB3     | 130783         | 191993        | 2.21%           | 0.371%        |
| 11        | 7.416       | 865           | 872         | 873          | rBV      | 259279         | 309079        | 3.55%           | 0.597%        |
| 12        | 7.451       | 873           | 878         | 881          | rVB      | 5129752        | 5602820       | 64.39%          | 10.827%       |
| 13        | 7.645       | 904           | 911         | 918          | rBV      | 493360         | 449091        | 5.16%           | 0.868%        |
| 14        | 7.834       | 940           | 943         | 948          | rVB      | 241150         | 244052        | 2.80%           | 0.472%        |
| 15        | 7.898       | 949           | 954         | 958          | rBV      | 351507         | 323597        | 3.72%           | 0.625%        |
| 16        | 8.163       | 991           | 999         | 1005         | rBV      | 2343291        | 2317841       | 26.64%          | 4.479%        |
| 17        | 8.975       | 1133          | 1137        | 1141         | rBV      | 379105         | 343574        | 3.95%           | 0.664%        |
| 18        | 9.245       | 1176          | 1183        | 1186         | rBV      | 8995019        | 8701208       | 100.00%         | 16.814%       |
| 19        | 9.504       | 1222          | 1227        | 1230         | rBV4     | 82208          | 91390         | 1.05%           | 0.177%        |
| 20        | 9.586       | 1236          | 1241        | 1245         | rBV5     | 88074          | 134483        | 1.55%           | 0.260%        |
| 21        | 9.781       | 1271          | 1274        | 1278         | rVB      | 106643         | 93846         | 1.08%           | 0.181%        |
| 22        | 9.922       | 1292          | 1298        | 1302         | rBV      | 2677621        | 2349329       | 27.00%          | 4.540%        |
| 23        | 10.539      | 1400          | 1403        | 1408         | rVB2     | 128059         | 138893        | 1.60%           | 0.268%        |
| 24        | 10.722      | 1427          | 1434        | 1437         | rBV      | 5215902        | 5505871       | 63.28%          | 10.639%       |
| 25        | 11.416      | 1548          | 1552        | 1555         | rVB      | 2273032        | 2054518       | 23.61%          | 3.970%        |
| 26        | 11.657      | 1589          | 1593        | 1598         | rBV6     | 65587          | 131334        | 1.51%           | 0.254%        |
| 27        | 11.792      | 1613          | 1616        | 1619         | rVB      | 315953         | 280938        | 3.23%           | 0.543%        |
| 28        | 11.933      | 1637          | 1640        | 1648         | rBV2     | 350136         | 346287        | 3.98%           | 0.669%        |
| 29        | 12.474      | 1728          | 1732        | 1734         | rBV2     | 268974         | 412592        | 4.74%           | 0.797%        |
| 30        | 12.504      | 1734          | 1737        | 1745         | rVB      | 766066         | 719141        | 8.26%           | 1.390%        |
| 31        | 12.592      | 1749          | 1752        | 1755         | rVB      | 263538         | 236341        | 2.72%           | 0.457%        |
| 32        | 12.651      | 1759          | 1762        | 1765         | rVB      | 447626         | 337184        | 3.88%           | 0.652%        |
| 33        | 12.804      | 1785          | 1788        | 1792         | rVB      | 110531         | 106002        | 1.22%           | 0.205%        |
| 34        | 13.004      | 1817          | 1822        | 1825         | rBV      | 7451851        | 7622596       | 87.60%          | 14.730%       |
| 35        | 13.439      | 1894          | 1896        | 1899         | rVB      | 124374         | 99093         | 1.14%           | 0.191%        |
| 36        | 13.916      | 1974          | 1977        | 1987         | rVB2     | 100181         | 201088        | 2.31%           | 0.389%        |

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Sample : N3861-13  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

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-BF072522.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |     |         |         |        |        |
|----|--------|------|------|------|-----|---------|---------|--------|--------|
| 37 | 14.057 | 1997 | 2001 | 2005 | rBV | 1571199 | 1396357 | 16.05% | 2.698% |
| 38 | 14.898 | 2141 | 2144 | 2150 | rVB | 274527  | 291639  | 3.35%  | 0.564% |
| 39 | 15.551 | 2250 | 2255 | 2264 | rVB | 1004475 | 1244581 | 14.30% | 2.405% |

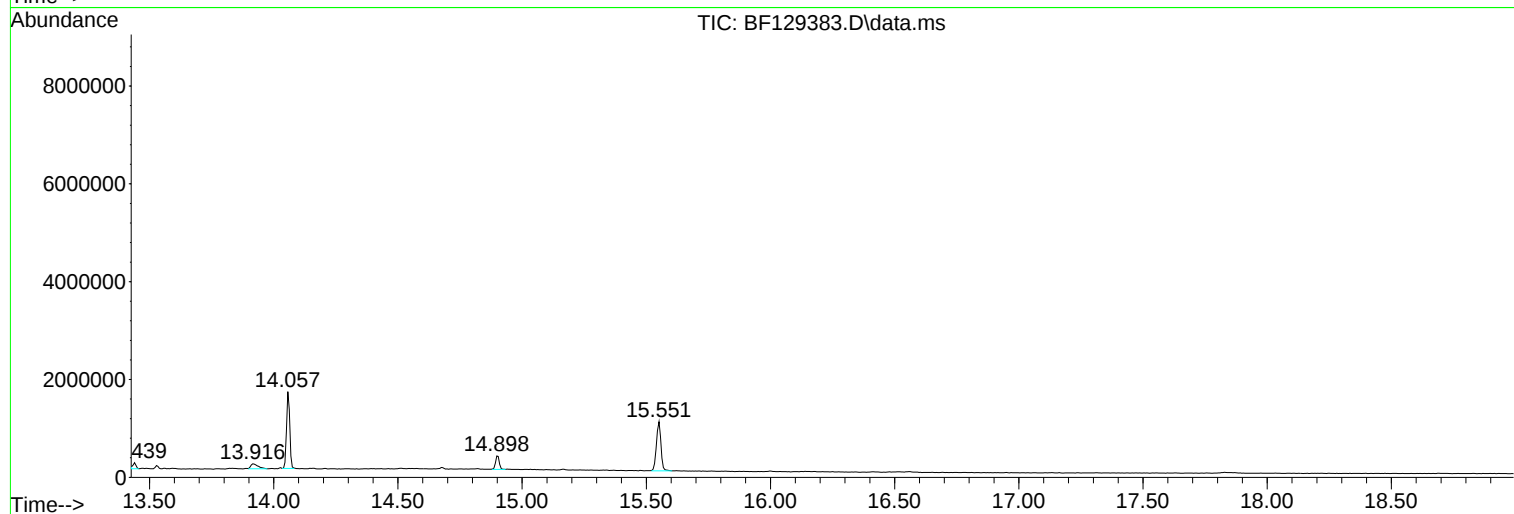
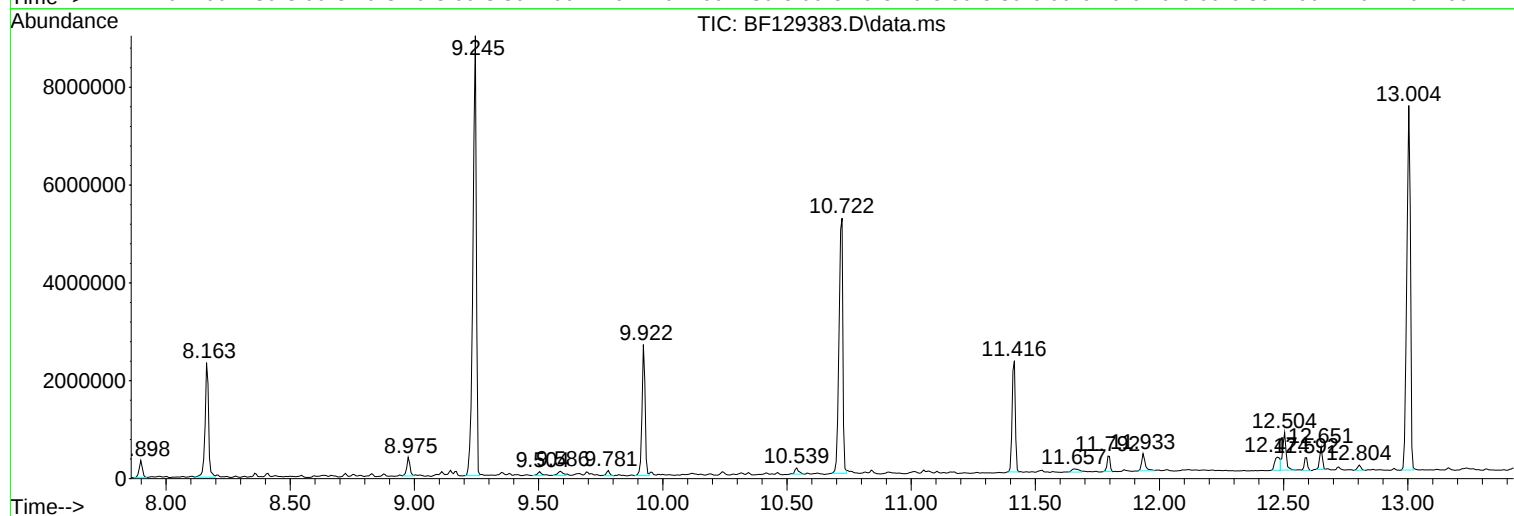
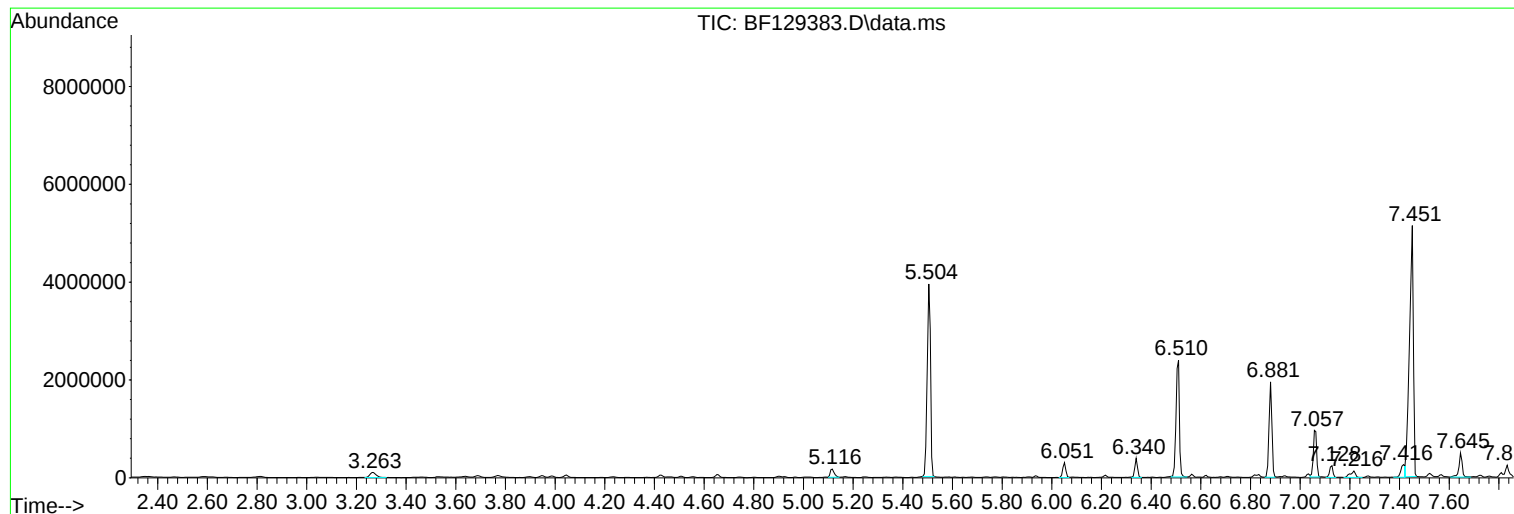
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Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072722\  
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Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF072522.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\BF072722\  
Data File : BF129383.D  
Acq On : 27 Jul 2022 18:51  
Operator : CG\JU  
Sample : N3861-13  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

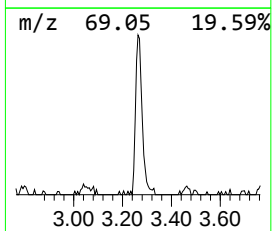
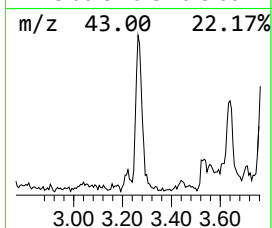
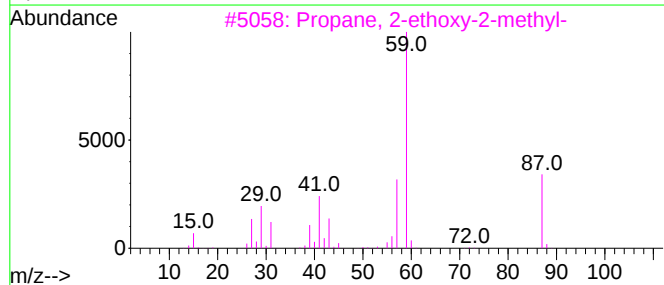
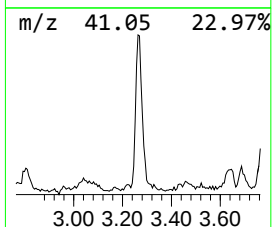
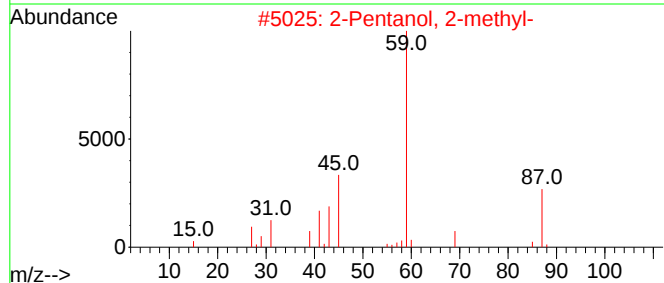
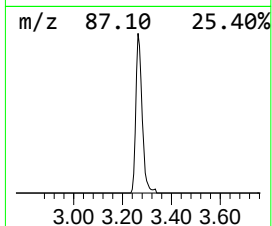
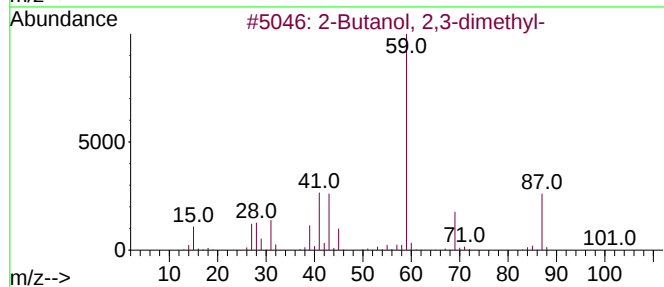
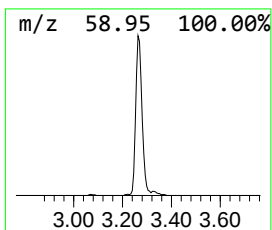
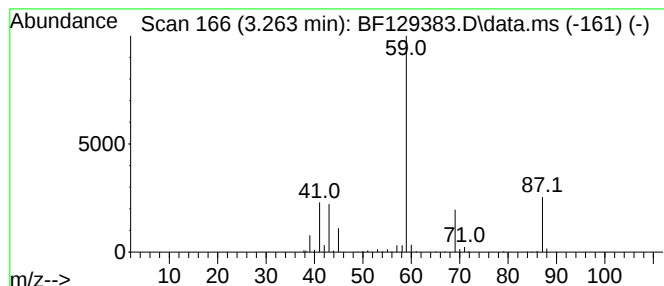
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 2-Butanol, 2,3-dimethyl- Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 3.263 | 2.26 ng | 180799 | 1,4-Dichlorobenzene-d4 | 6.881 |

| Hit# | of 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|------|-----------------------------|-----|---------|-------------|------|
| 1    |      | 2-Butanol, 2,3-dimethyl-    | 102 | C6H14O  | 000594-60-5 | 83   |
| 2    |      | 2-Pentanol, 2-methyl-       | 102 | C6H14O  | 000590-36-3 | 56   |
| 3    |      | Propane, 2-ethoxy-2-methyl- | 102 | C6H14O  | 000637-92-3 | 50   |
| 4    |      | 3-Heptanol                  | 116 | C7H16O  | 000589-82-2 | 40   |
| 5    |      | 3-Pentanol, 2,2-dimethyl-   | 116 | C7H16O  | 003970-62-5 | 40   |



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Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF072522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

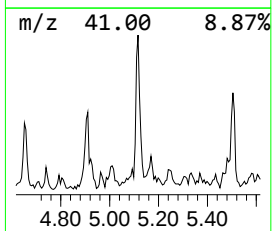
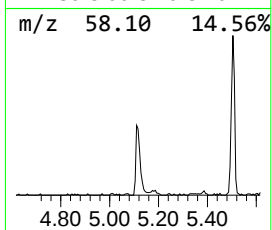
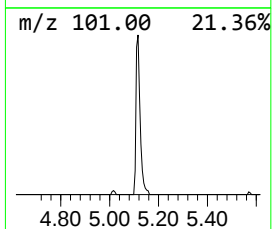
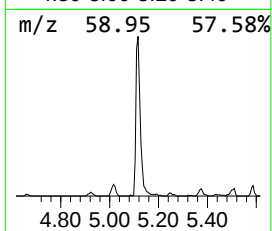
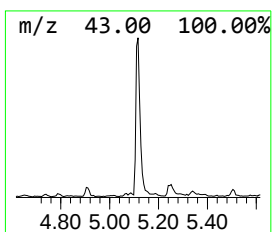
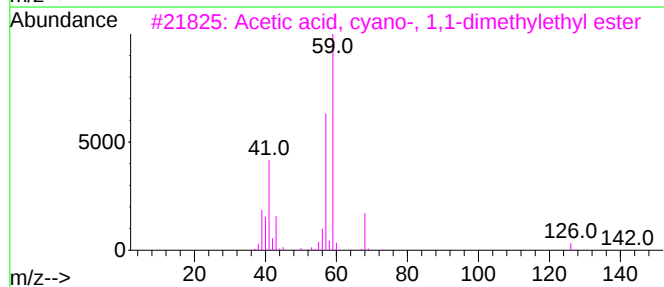
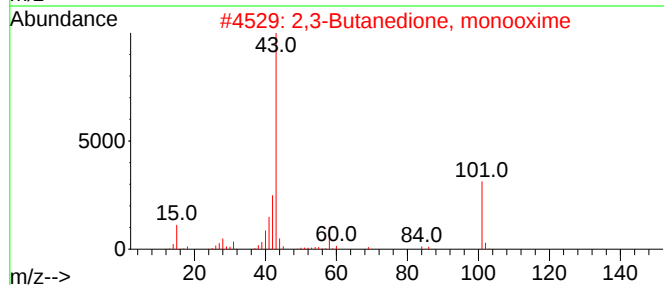
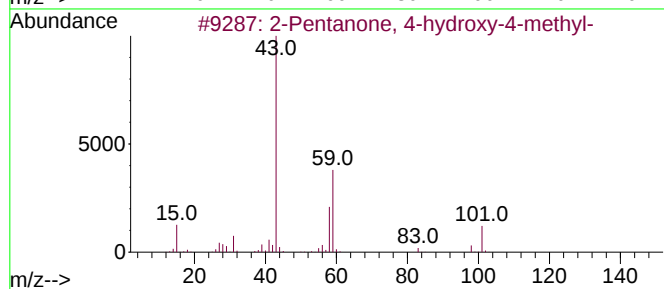
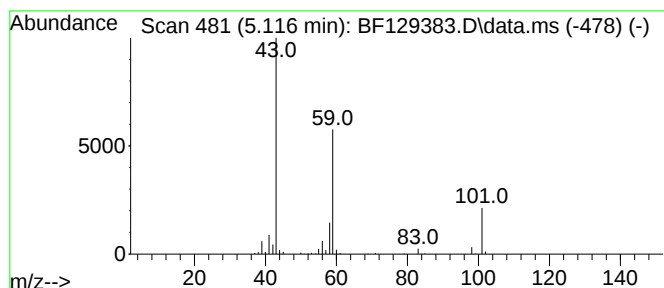
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 5.116 | 2.35 ng | 187616 | 1,4-Dichlorobenzene-d4 | 6.881 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 72   |
| 2    |    |   | 2,3-Butanedione, monooxime          | 101 | C4H7NO2  | 000057-71-6 | 17   |
| 3    |    |   | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 17   |
| 4    |    |   | 1-Propen-2-ol, acetate              | 100 | C5H8O2   | 000108-22-5 | 10   |
| 5    |    |   | Morpholine, 4-methyl-               | 101 | C5H11NO  | 000109-02-4 | 9    |



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Sample : N3861-13  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

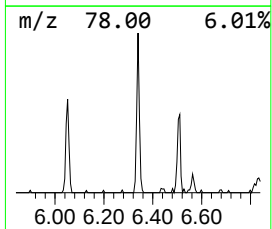
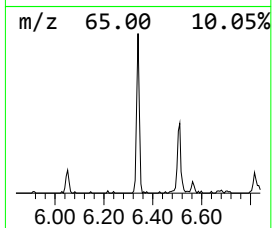
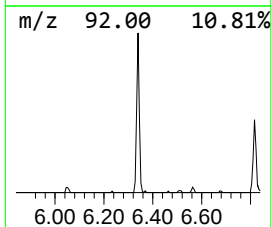
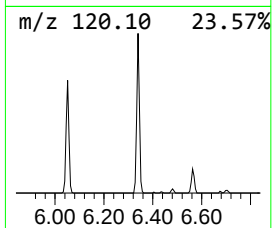
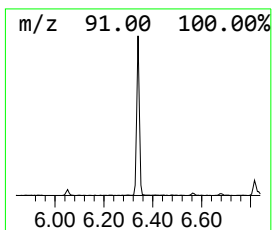
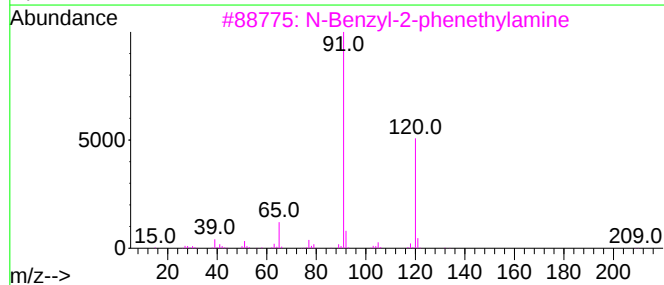
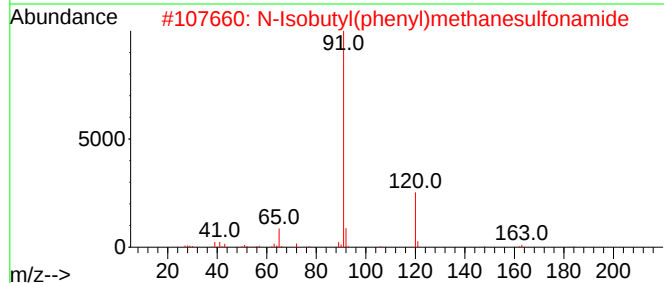
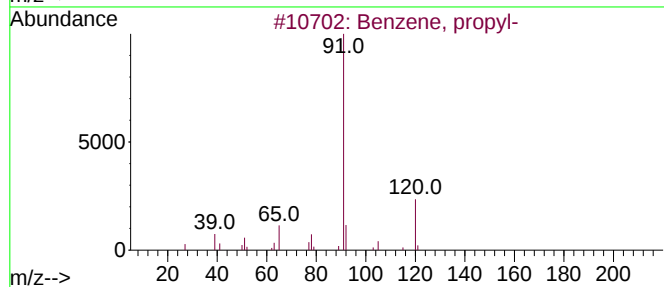
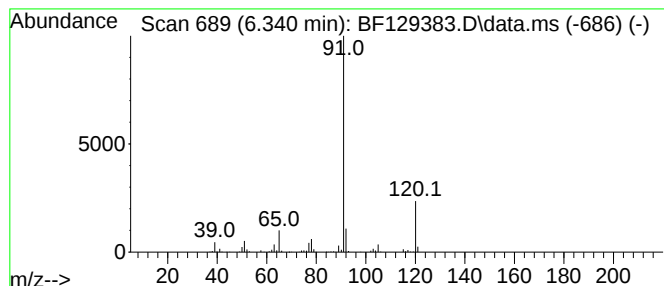
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF072522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 4 Benzene, propyl- Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 6.340 | 3.69 ng | 294526 | 1,4-Dichlorobenzene-d4 | 6.881 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|------------|-------------|------|
| 1    |    |   | Benzene, propyl-                    | 120 | C9H12      | 000103-65-1 | 94   |
| 2    |    |   | N-Isobutyl(phenyl)methanesulfona... | 227 | C11H17NO2S | 144615-94-1 | 72   |
| 3    |    |   | N-Benzyl-2-phenethylamine           | 211 | C15H17N    | 003647-71-0 | 72   |
| 4    |    |   | Phenethylamine, N-benzyl-p-chloro-  | 245 | C15H16ClN  | 013622-43-0 | 72   |
| 5    |    |   | 1,2-Ethanediamine, N,N'-bis(phen... | 240 | C16H20N2   | 000140-28-3 | 72   |



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Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

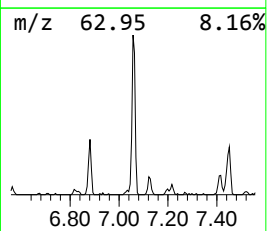
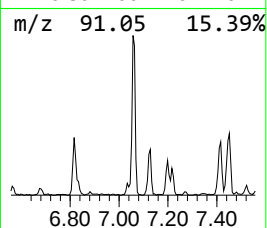
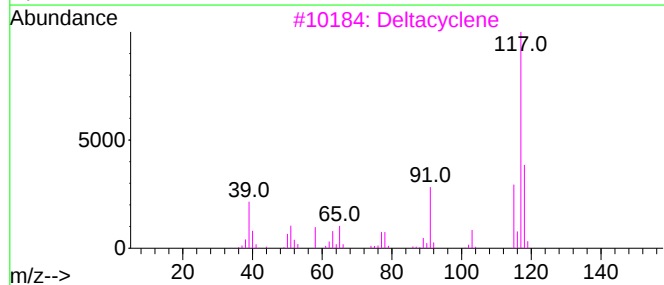
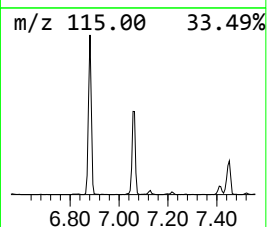
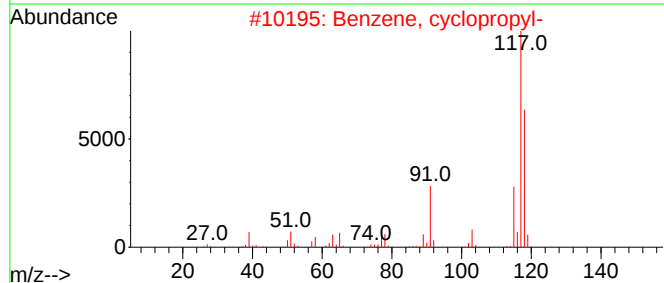
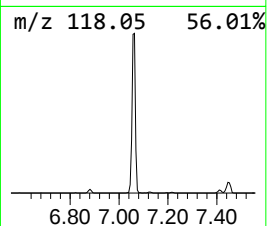
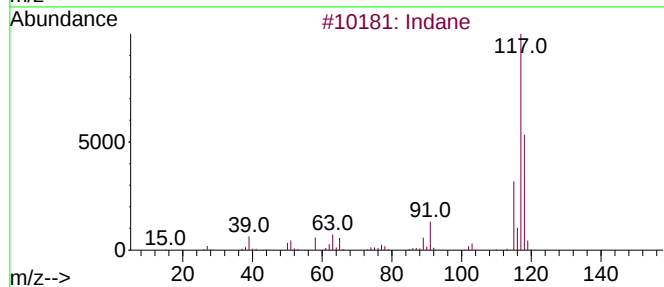
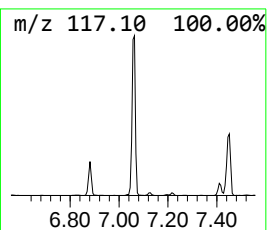
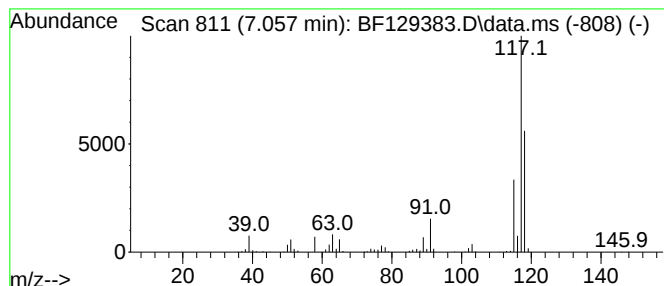
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF072522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 Indane Concentration Rank 1

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 7.057 | 11.13 ng | 888425 | 1,4-Dichlorobenzene-d4 | 6.881 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Indane                              | 118 | C9H10   | 000496-11-7  | 91   |
| 2    |    |   | Benzene, cyclopropyl-               | 118 | C9H10   | 000873-49-4  | 64   |
| 3    |    |   | Deltacyclene                        | 118 | C9H10   | 007785-10-6  | 59   |
| 4    |    |   | Benzene, 1-ethenyl-2-methyl-        | 118 | C9H10   | 000611-15-4  | 59   |
| 5    |    |   | Tetracyclo[3.3.1.0(2,8).0(4,6)]-... | 118 | C9H10   | 1000191-13-7 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072722\  
Data File : BF129383.D  
Acq On : 27 Jul 2022 18:51  
Operator : CG\JU  
Sample : N3861-13  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF072522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

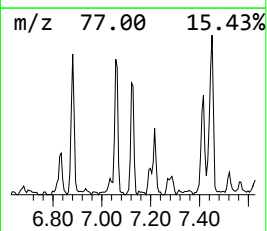
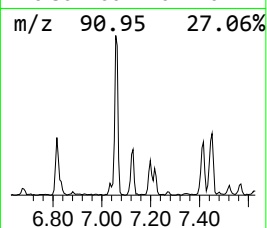
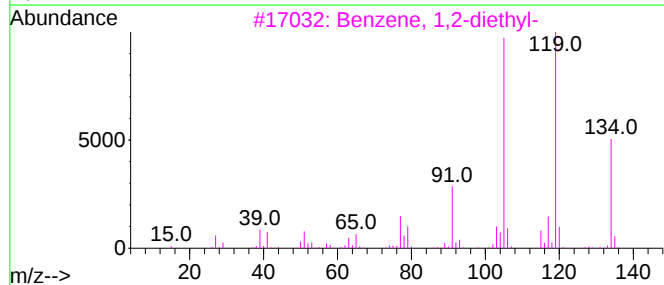
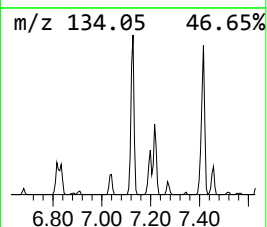
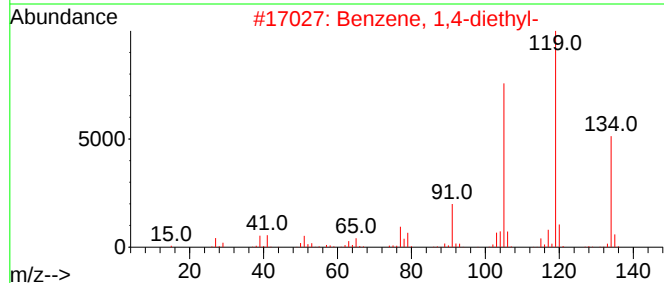
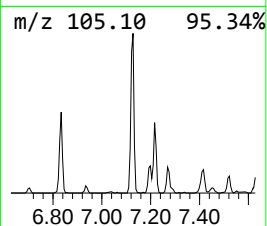
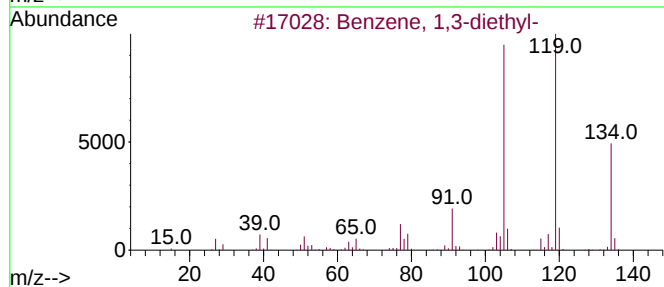
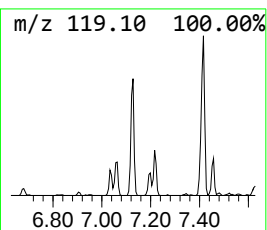
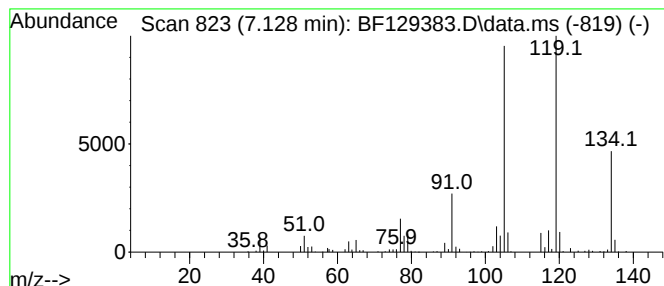
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 Benzene, 1,3-diethyl- Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 7.128 | 2.67 ng | 213202 | 1,4-Dichlorobenzene-d4 | 6.881 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,3-diethyl-          | 134 | C10H14  | 000141-93-5 | 95   |
| 2    |    |   | Benzene, 1,4-diethyl-          | 134 | C10H14  | 000105-05-5 | 94   |
| 3    |    |   | Benzene, 1,2-diethyl-          | 134 | C10H14  | 000135-01-3 | 93   |
| 4    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 91   |
| 5    |    |   | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072722\  
Data File : BF129383.D  
Acq On : 27 Jul 2022 18:51  
Operator : CG\JU  
Sample : N3861-13  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

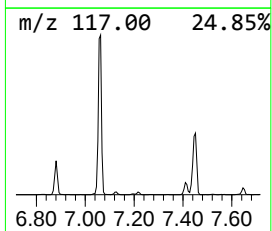
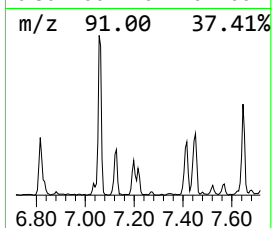
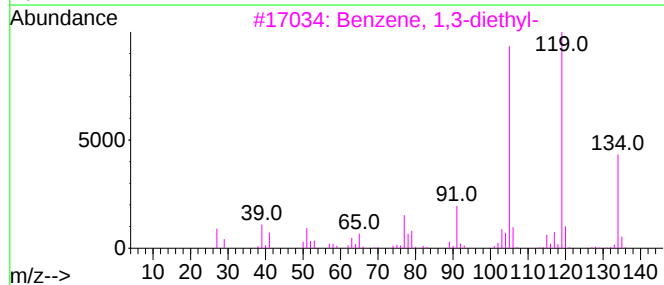
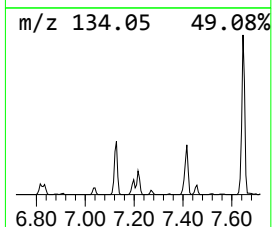
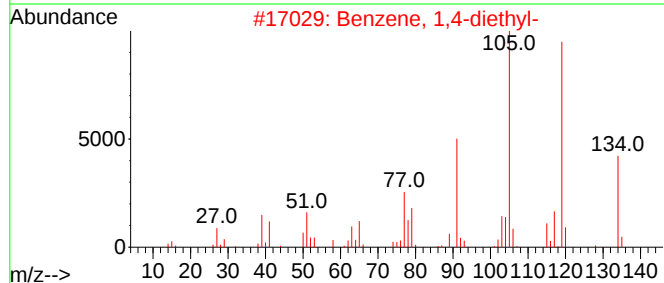
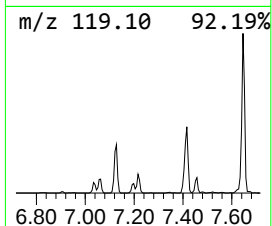
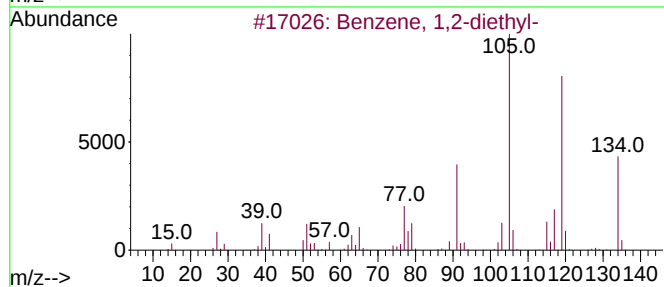
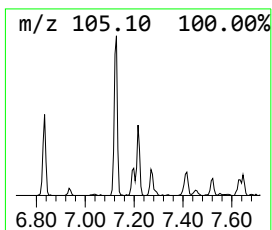
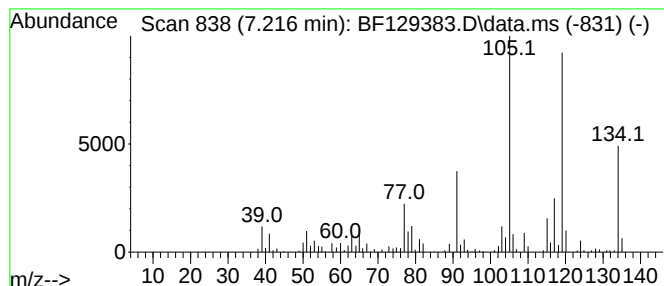
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 Benzene, 1,2-diethyl- Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 7.216 | 2.40 ng | 191993 | 1,4-Dichlorobenzene-d4 | 6.881 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2-diethyl-         | 134 | C10H14  | 000135-01-3 | 96   |
| 2    |    |   | Benzene, 1,4-diethyl-         | 134 | C10H14  | 000105-05-5 | 90   |
| 3    |    |   | Benzene, 1,3-diethyl-         | 134 | C10H14  | 000141-93-5 | 87   |
| 4    |    |   | 1,3,8-p-Menthatriene          | 134 | C10H14  | 018368-95-1 | 83   |
| 5    |    |   | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14  | 000527-53-7 | 81   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072722\  
Data File : BF129383.D  
Acq On : 27 Jul 2022 18:51  
Operator : CG\JU  
Sample : N3861-13  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

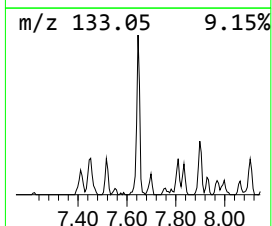
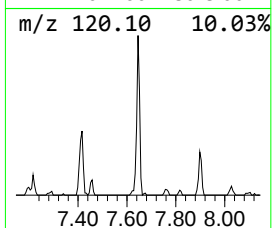
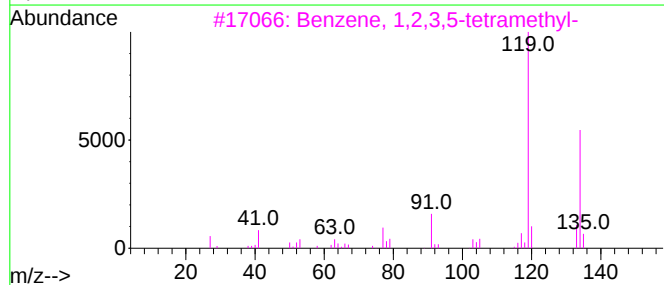
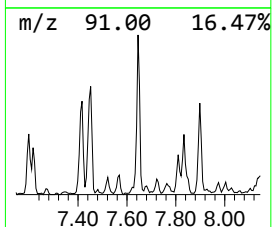
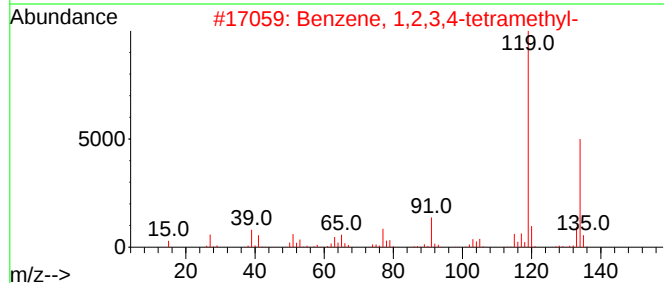
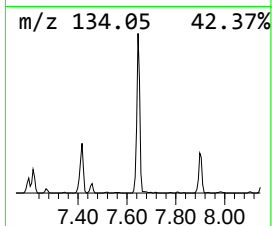
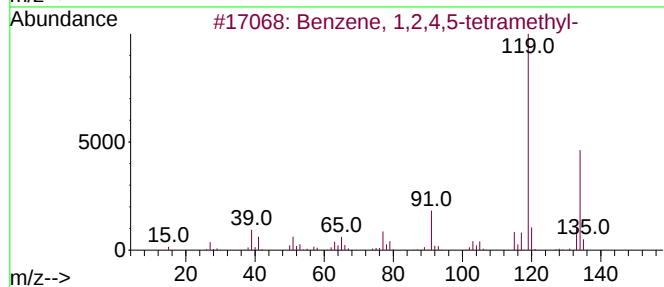
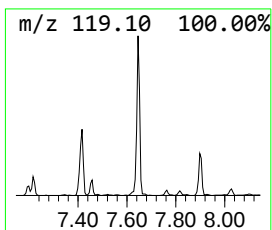
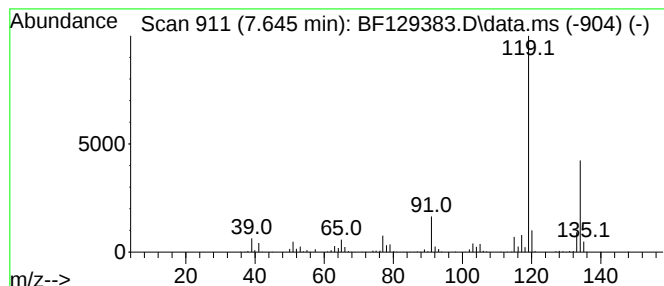
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 7.645 | 3.88 ng | 449091 | Naphthalene-d8   | 8.163 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,4,5-tetramethyl-       | 134 | C10H14  | 000095-93-2 | 97   |
| 2    |    |   | Benzene, 1,2,3,4-tetramethyl-       | 134 | C10H14  | 000488-23-3 | 97   |
| 3    |    |   | Benzene, 1,2,3,5-tetramethyl-       | 134 | C10H14  | 000527-53-7 | 96   |
| 4    |    |   | Benzene, 1-ethyl-2,4-dimethyl-      | 134 | C10H14  | 000874-41-9 | 94   |
| 5    |    |   | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072722\  
Data File : BF129383.D  
Acq On : 27 Jul 2022 18:51  
Operator : CG\JU  
Sample : N3861-13  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

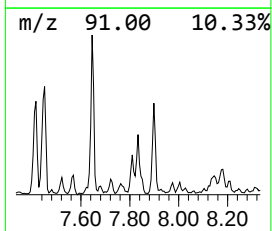
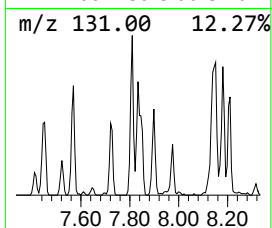
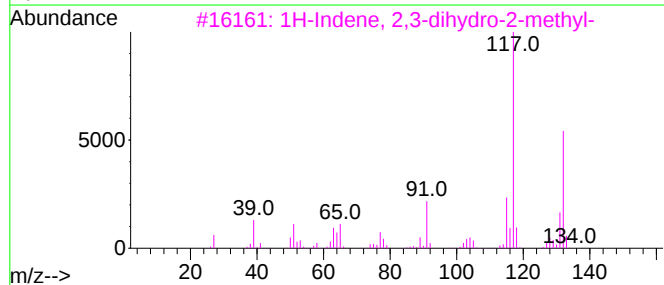
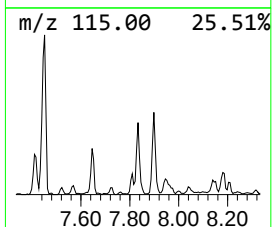
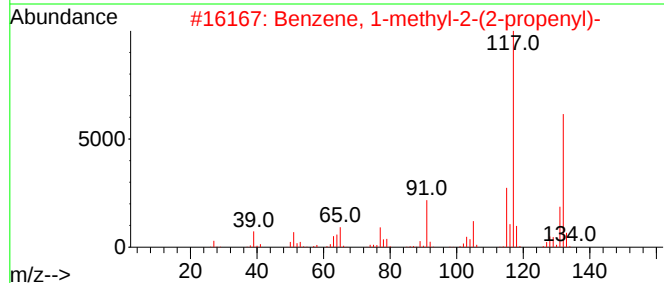
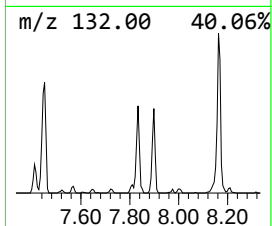
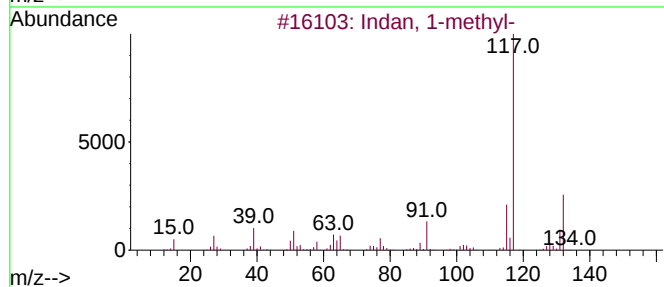
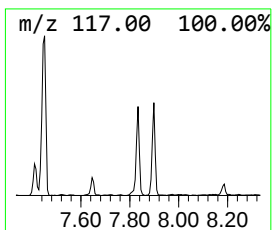
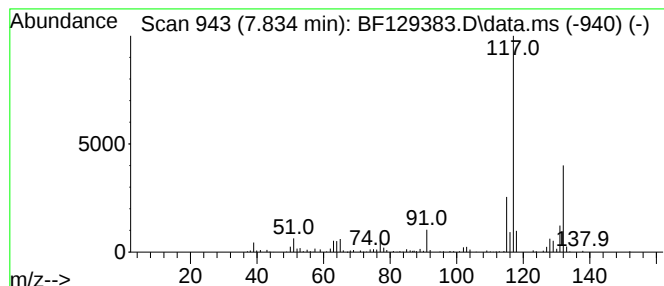
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 Indan, 1-methyl- Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 7.834 | 2.11 ng | 244052 | Naphthalene-d8   | 8.163 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Indan, 1-methyl-                  | 132 | C10H12  | 000767-58-8 | 90   |
| 2    |    |   | Benzene, 1-methyl-2-(2-propenyl)- | 132 | C10H12  | 001587-04-8 | 87   |
| 3    |    |   | 1H-Indene, 2,3-dihydro-2-methyl-  | 132 | C10H12  | 000824-63-5 | 87   |
| 4    |    |   | Benzene, 2-butenyl-               | 132 | C10H12  | 001560-06-1 | 87   |
| 5    |    |   | 1-Phenyl-1-butene                 | 132 | C10H12  | 000824-90-8 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072722\  
Data File : BF129383.D  
Acq On : 27 Jul 2022 18:51  
Operator : CG\JU  
Sample : N3861-13  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF072522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

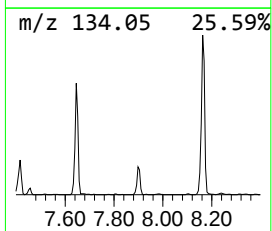
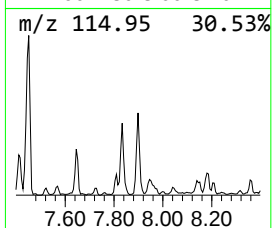
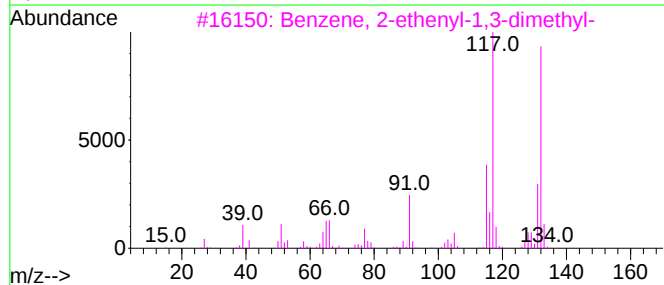
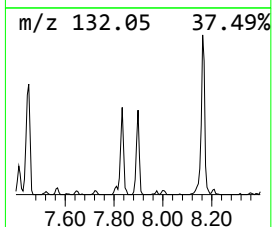
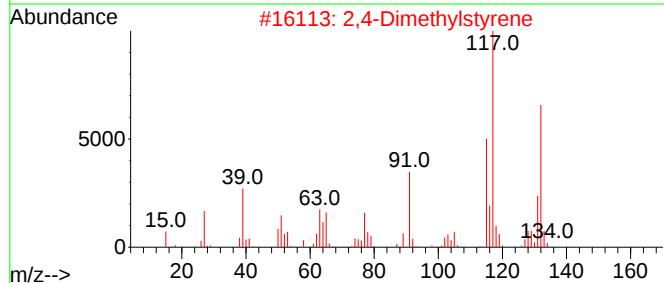
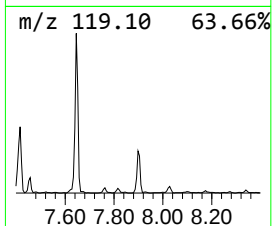
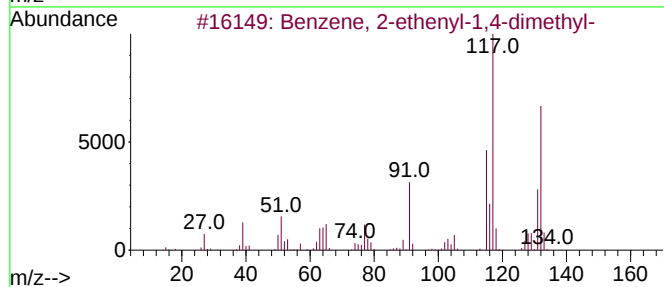
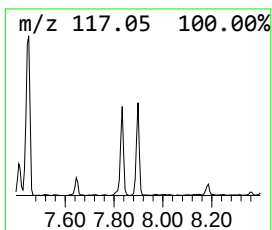
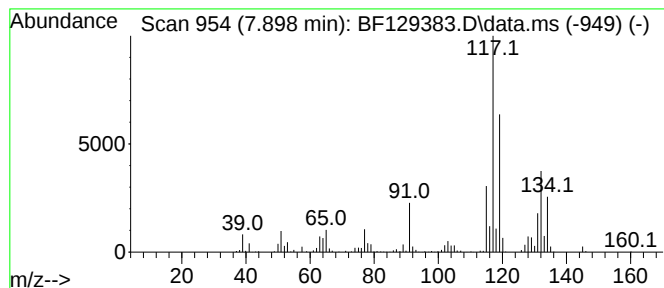
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Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 7.898 | 2.79 ng | 323597 | Naphthalene-d8   | 8.163 |

| Hit# | of 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|------|----------------------------------|-----|---------|-------------|------|
| 1    |      | Benzene, 2-ethenyl-1,4-dimethyl- | 132 | C10H12  | 002039-89-6 | 95   |
| 2    |      | 2,4-Dimethylstyrene              | 132 | C10H12  | 002234-20-0 | 95   |
| 3    |      | Benzene, 2-ethenyl-1,3-dimethyl- | 132 | C10H12  | 002039-90-9 | 86   |
| 4    |      | 1-Phenyl-1-butene                | 132 | C10H12  | 000824-90-8 | 70   |
| 5    |      | 1H-Indene, 2,3-dihydro-4-methyl- | 132 | C10H12  | 000824-22-6 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072722\  
Data File : BF129383.D  
Acq On : 27 Jul 2022 18:51  
Operator : CG\JU  
Sample : N3861-13  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

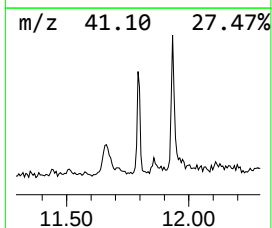
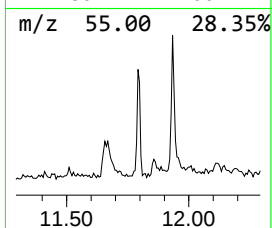
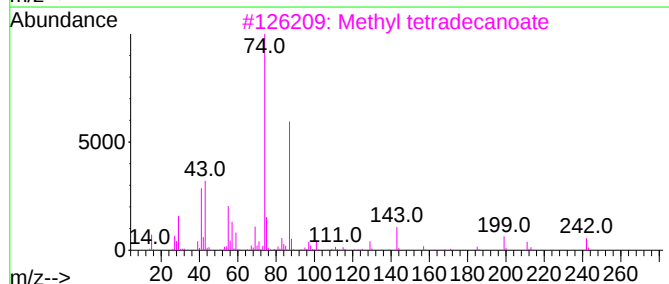
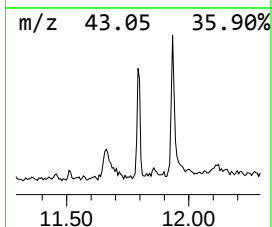
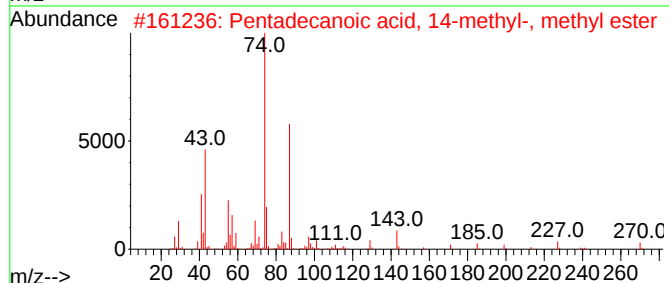
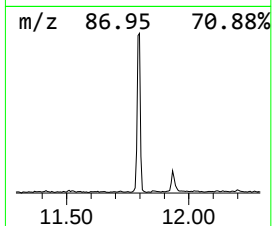
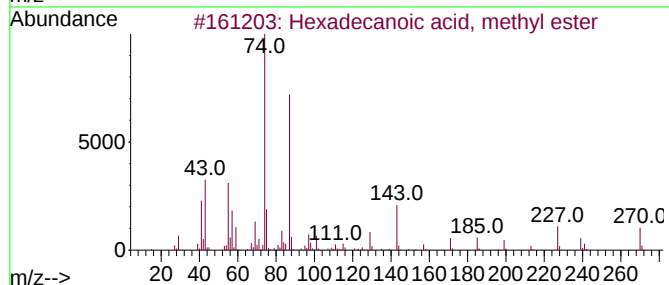
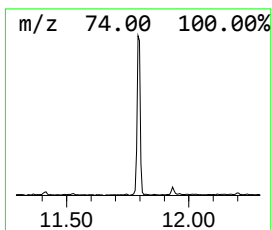
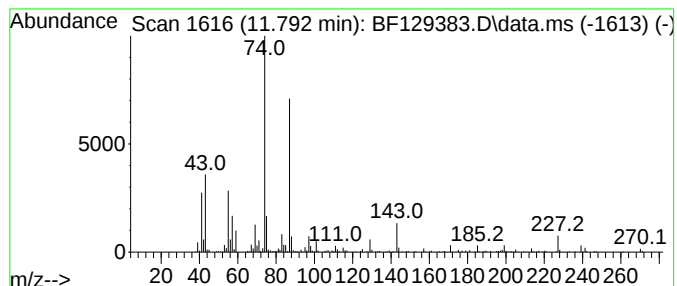
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 11 Hexadecanoic acid, methyl e... Concentration Rank 12

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 11.792    | 2.73 ng                             | 280938 | Phenanthrene-d10 | 11.416      |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Hexadecanoic acid, methyl ester     | 270    | C17H34O2         | 000112-39-0 | 94   |
| 2         | Pentadecanoic acid, 14-methyl-, ... | 270    | C17H34O2         | 005129-60-2 | 91   |
| 3         | Methyl tetradecanoate               | 242    | C15H30O2         | 000124-10-7 | 91   |
| 4         | Tridecanoic acid, 12-methyl-, me... | 242    | C15H30O2         | 005129-58-8 | 90   |
| 5         | Pentadecanoic acid, methyl ester    | 256    | C16H32O2         | 007132-64-1 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072722\  
Data File : BF129383.D  
Acq On : 27 Jul 2022 18:51  
Operator : CG\JU  
Sample : N3861-13  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

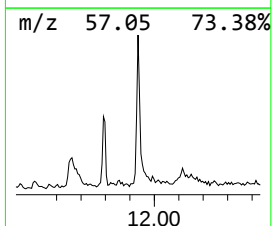
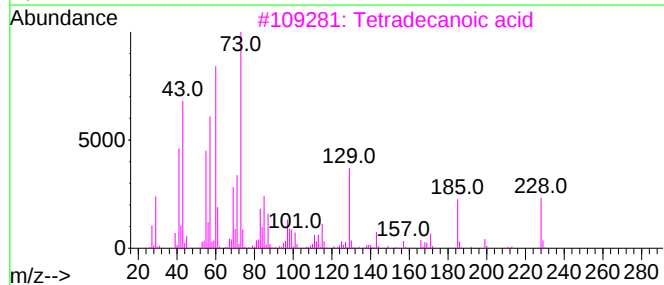
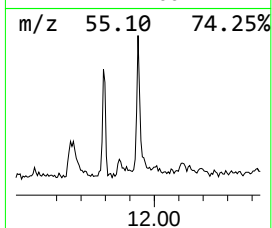
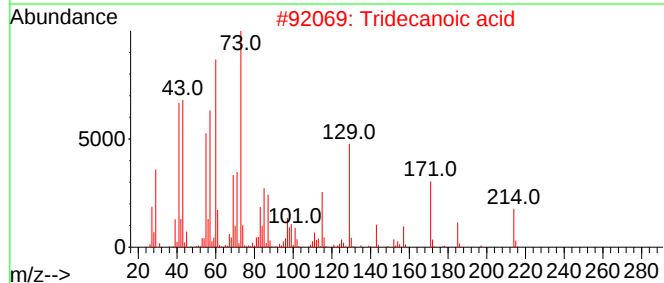
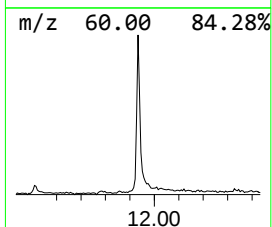
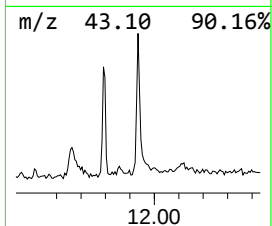
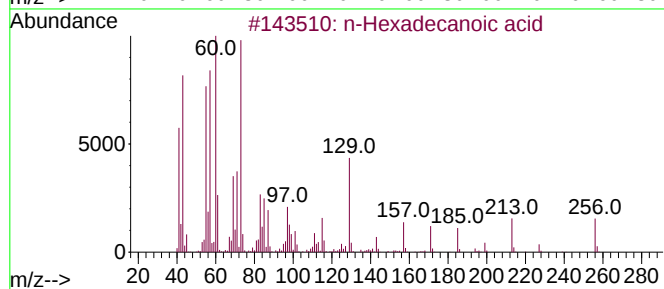
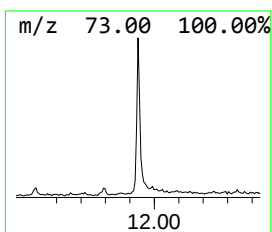
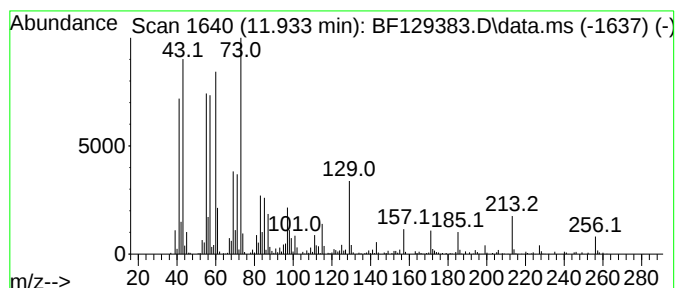
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 12 n-Hexadecanoic acid Concentration Rank 7

| R.T.      | EstConc             | Area   | Relative to ISTD | R.T.        |      |
|-----------|---------------------|--------|------------------|-------------|------|
| 11.933    | 3.37 ng             | 346287 | Phenanthrene-d10 | 11.416      |      |
| Hit# of 5 | Tentative ID        | MW     | MolForm          | CAS#        | Qual |
| 1         | n-Hexadecanoic acid | 256    | C16H32O2         | 000057-10-3 | 99   |
| 2         | Tridecanoic acid    | 214    | C13H26O2         | 000638-53-9 | 95   |
| 3         | Tetradecanoic acid  | 228    | C14H28O2         | 000544-63-8 | 93   |
| 4         | Pentadecanoic acid  | 242    | C15H30O2         | 001002-84-2 | 90   |
| 5         | Isopropyl palmitate | 298    | C19H38O2         | 000142-91-6 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072722\  
Data File : BF129383.D  
Acq On : 27 Jul 2022 18:51  
Operator : CG\JU  
Sample : N3861-13  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

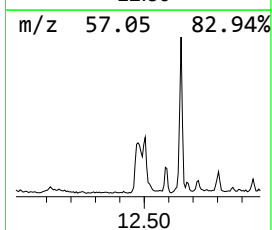
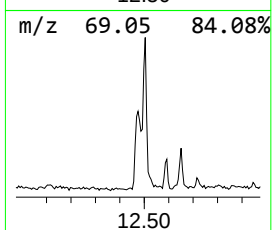
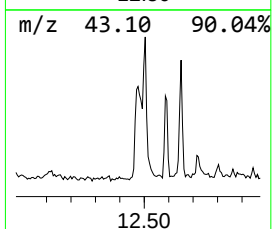
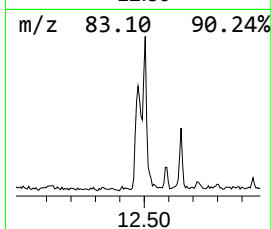
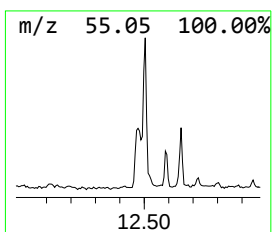
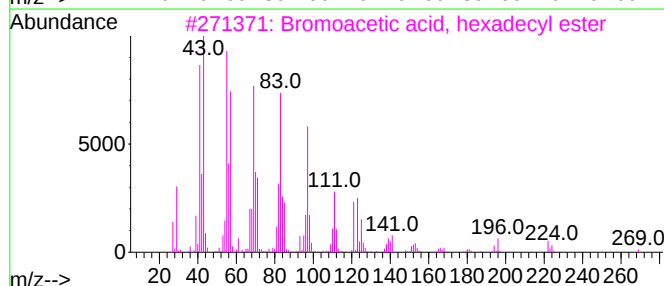
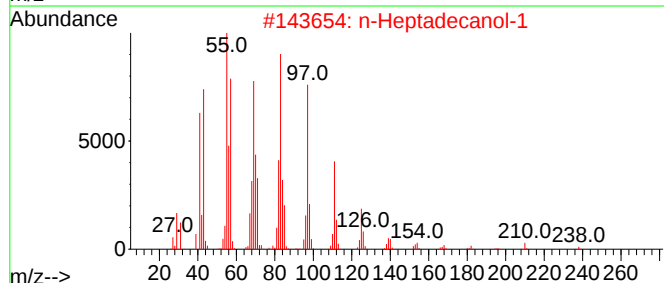
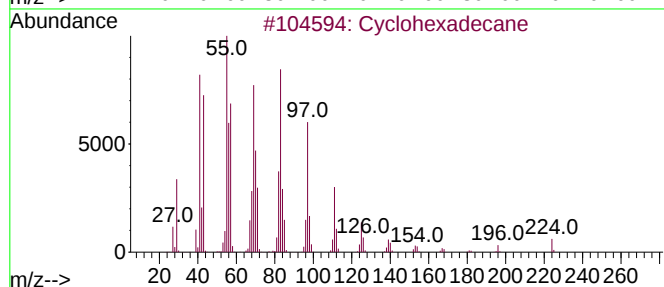
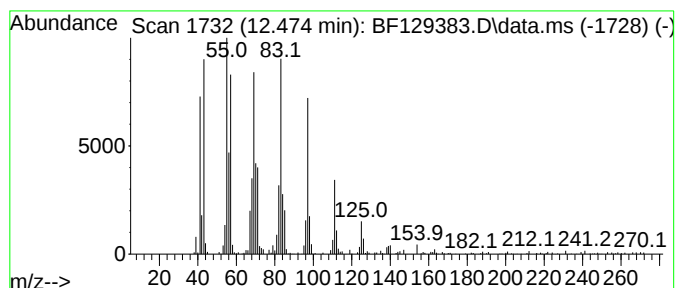
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 13 Cyclohexadecane Concentration Rank 4

| R.T.      | EstConc                           | Area   | Relative to ISTD | R.T.         |      |
|-----------|-----------------------------------|--------|------------------|--------------|------|
| 12.475    | 4.02 ng                           | 412592 | Phenanthrene-d10 | 11.416       |      |
| Hit# of 5 | Tentative ID                      | MW     | MolForm          | CAS#         | Qual |
| 1         | Cyclohexadecane                   | 224    | C16H32           | 000295-65-8  | 96   |
| 2         | n-Heptadecanol-1                  | 256    | C17H36O          | 001454-85-9  | 95   |
| 3         | Bromoacetic acid, hexadecyl ester | 362    | C18H35BrO2       | 005454-48-8  | 94   |
| 4         | 1-Octadecanol                     | 270    | C18H38O          | 000112-92-5  | 94   |
| 5         | Z-5-Nonadecene                    | 266    | C19H38           | 1000131-11-8 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072722\  
Data File : BF129383.D  
Acq On : 27 Jul 2022 18:51  
Operator : CG\JU  
Sample : N3861-13  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

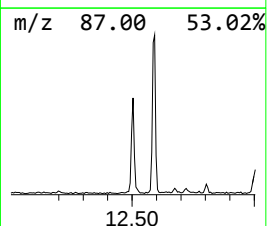
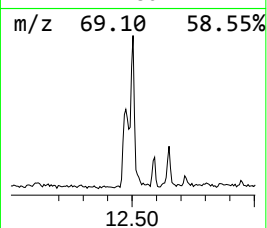
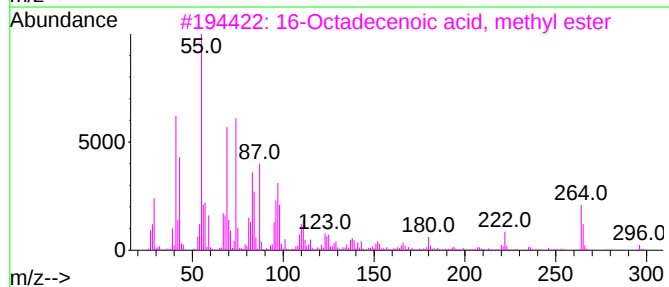
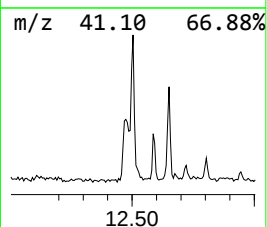
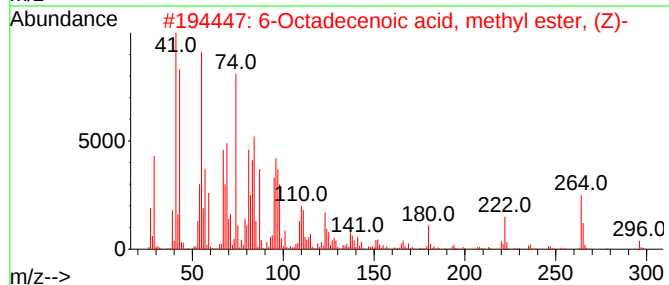
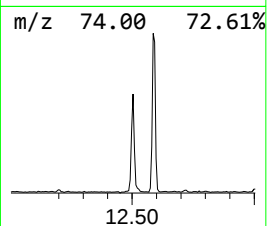
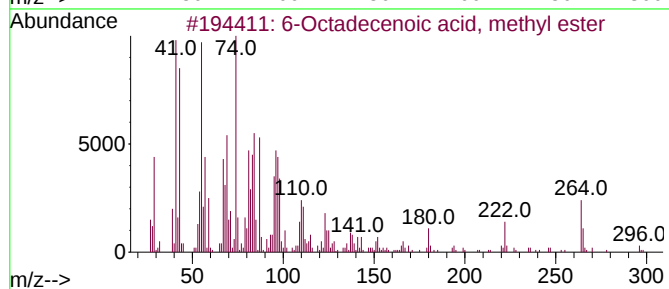
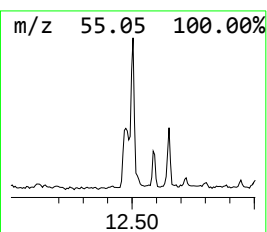
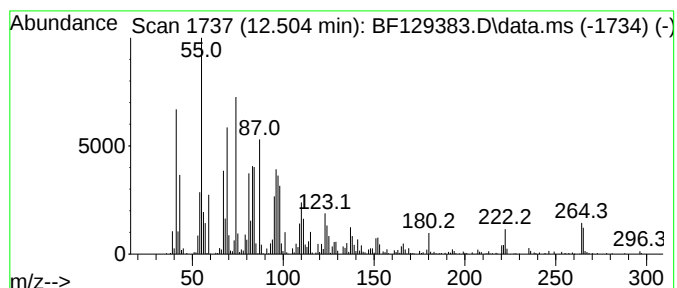
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 14 6-Octadecenoic acid, methyl... Concentration Rank 2

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 12.504    | 7.00 ng                             | 719141 | Phenanthrene-d10 | 11.416      |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | 6-Octadecenoic acid, methyl ester   | 296    | C19H36O2         | 052355-31-4 | 99   |
| 2         | 6-Octadecenoic acid, methyl este... | 296    | C19H36O2         | 002777-58-4 | 99   |
| 3         | 16-Octadecenoic acid, methyl ester  | 296    | C19H36O2         | 056554-49-5 | 98   |
| 4         | 15-Octadecenoic acid, methyl ester  | 296    | C19H36O2         | 004764-72-1 | 97   |
| 5         | 10-Octadecenoic acid, methyl ester  | 296    | C19H36O2         | 013481-95-3 | 97   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072722\  
Data File : BF129383.D  
Acq On : 27 Jul 2022 18:51  
Operator : CG\JU  
Sample : N3861-13  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

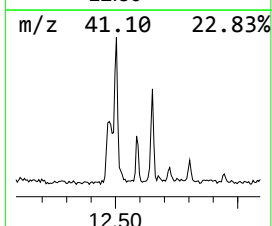
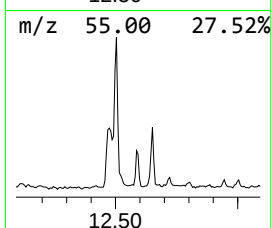
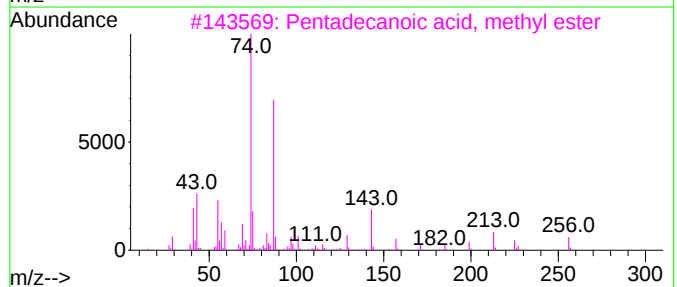
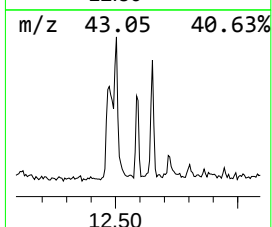
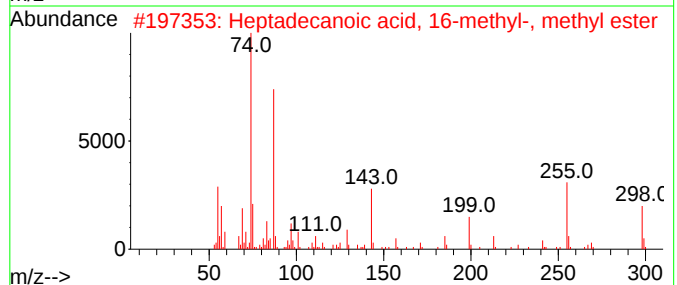
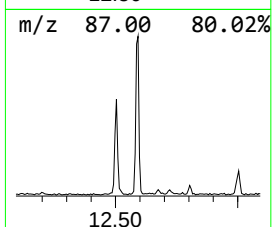
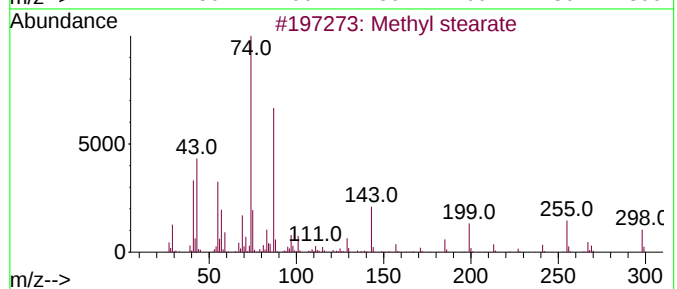
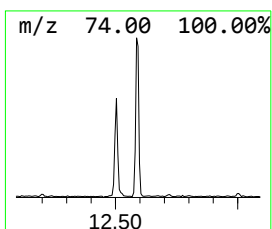
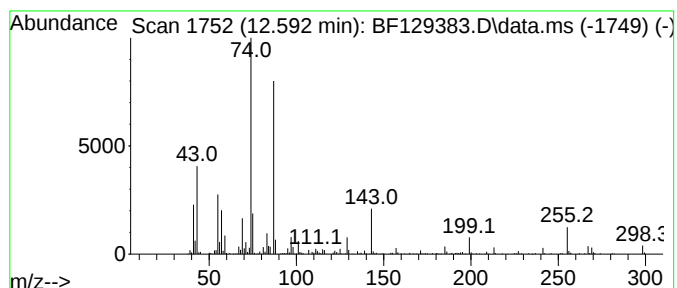
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 15 Methyl stearate Concentration Rank 16

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 12.592    | 2.30 ng                             | 236341 | Phenanthrene-d10 | 11.416      |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Methyl stearate                     | 298    | C19H38O2         | 000112-61-8 | 96   |
| 2         | Heptadecanoic acid, 16-methyl-, ... | 298    | C19H38O2         | 005129-61-3 | 93   |
| 3         | Pentadecanoic acid, methyl ester    | 256    | C16H32O2         | 007132-64-1 | 86   |
| 4         | Pentadecanoic acid, 14-methyl-, ... | 270    | C17H34O2         | 005129-60-2 | 83   |
| 5         | Hexadecanoic acid, methyl ester     | 270    | C17H34O2         | 000112-39-0 | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072722\  
Data File : BF129383.D  
Acq On : 27 Jul 2022 18:51  
Operator : CG\JU  
Sample : N3861-13  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

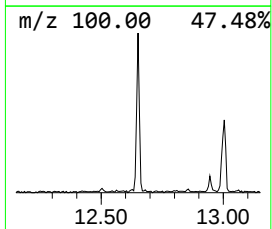
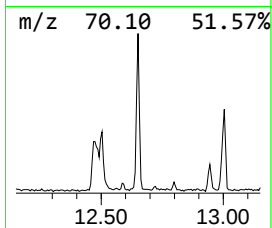
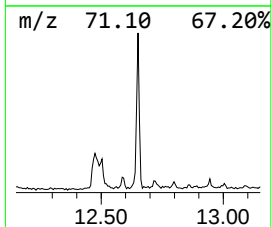
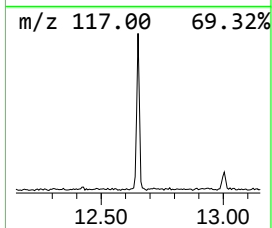
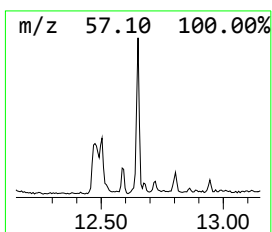
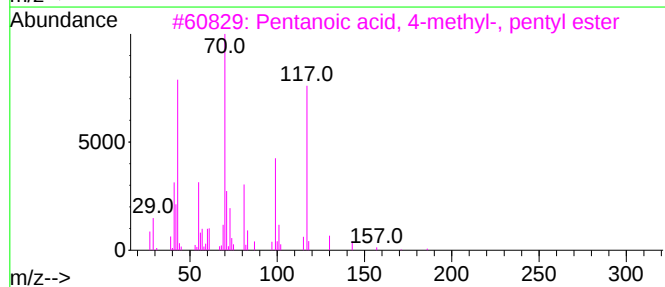
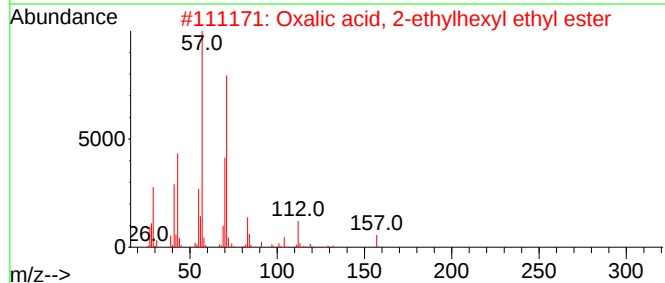
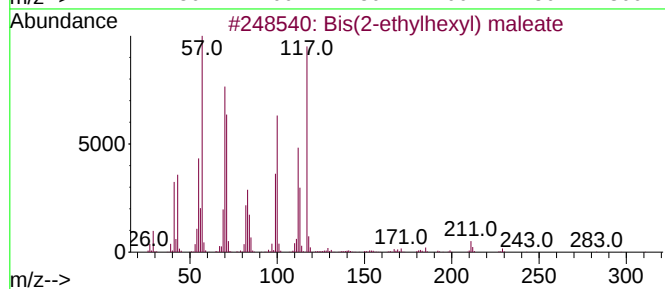
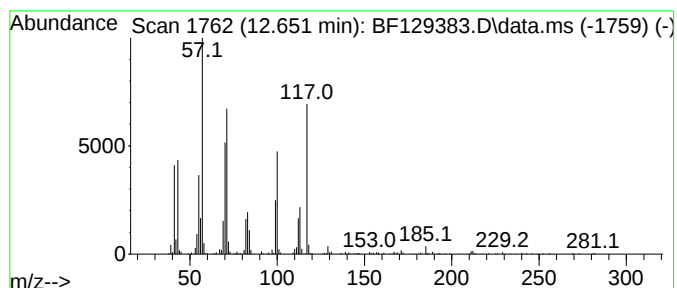
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 16 Bis(2-ethylhexyl) maleate Concentration Rank 8

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 12.651    | 3.28 ng                             | 337184 | Phenanthrene-d10 | 11.416       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | Bis(2-ethylhexyl) maleate           | 340    | C20H36O4         | 000142-16-5  | 80   |
| 2         | Oxalic acid, 2-ethylhexyl ethyl ... | 230    | C12H22O4         | 1000309-38-4 | 43   |
| 3         | Pentanoic acid, 4-methyl-, penty... | 186    | C11H22O2         | 025415-71-8  | 35   |
| 4         | 2-Ethyl-1-hexanol, pentafluoropr... | 276    | C11H17F5O2       | 959012-82-9  | 30   |
| 5         | Sulfurous acid, 2-ethylhexyl non... | 320    | C17H36O3S        | 1010309-19-2 | 27   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072722\  
Data File : BF129383.D  
Acq On : 27 Jul 2022 18:51  
Operator : CG\JU  
Sample : N3861-13  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

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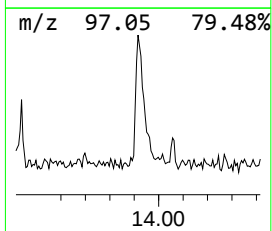
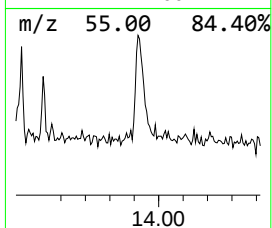
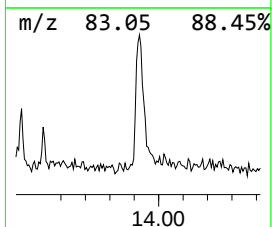
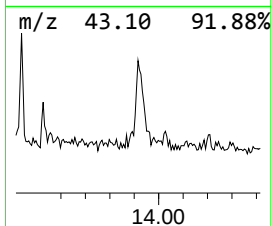
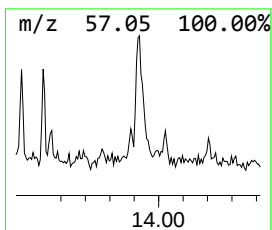
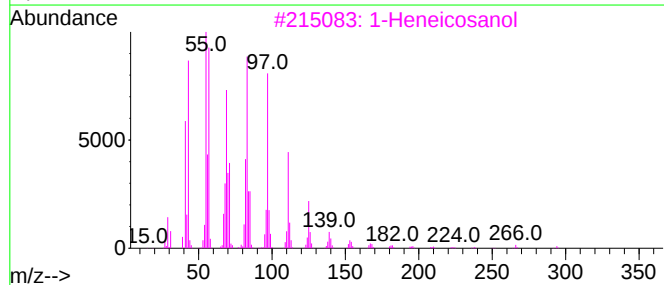
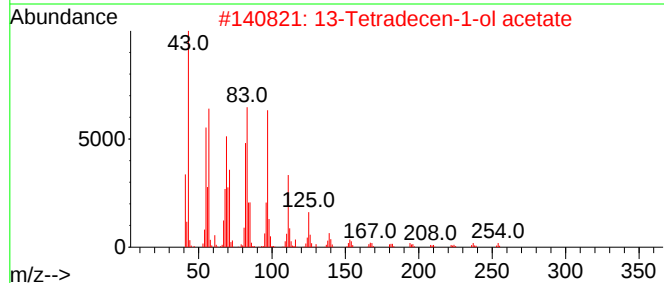
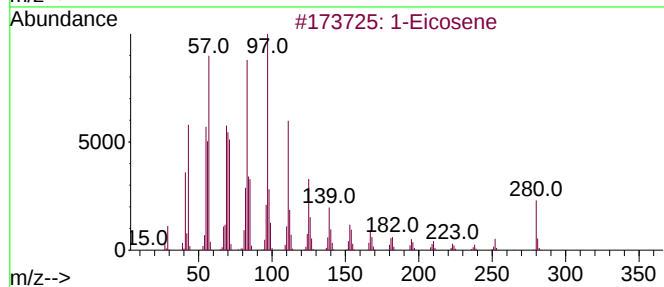
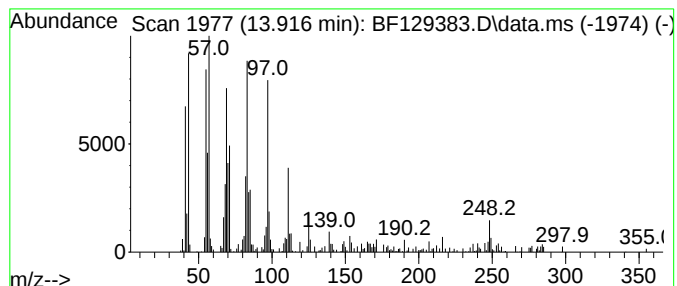
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TIC Integration Parameters: LSCINT.P

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Peak Number 17 1-Eicosene Concentration Rank 10

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 13.916 | 2.88 ng | 201088 | Chrysene-d12     | 14.057 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|------|-------------------------------------|-----|-------------|-------------|------|
| 1    |      | 1-Eicosene                          | 280 | C20H40      | 003452-07-1 | 97   |
| 2    |      | 13-Tetradecen-1-ol acetate          | 254 | C16H30O2    | 056221-91-1 | 94   |
| 3    |      | 1-Heneicosanol                      | 312 | C21H44O     | 015594-90-8 | 94   |
| 4    |      | Trichloroacetic acid, hexadecyl ... | 386 | C18H33Cl3O2 | 074339-54-1 | 94   |
| 5    |      | 1-Nonadecene                        | 266 | C19H38      | 018435-45-5 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072722\  
Data File : BF129383.D  
Acq On : 27 Jul 2022 18:51  
Operator : CG\JU  
Sample : N3861-13  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-21RR

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

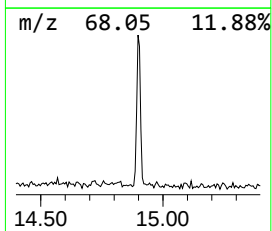
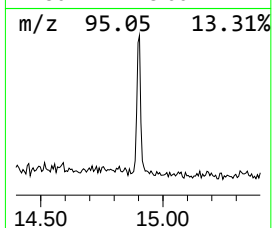
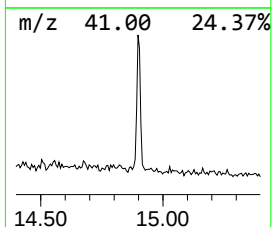
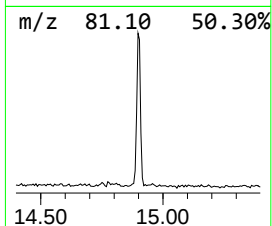
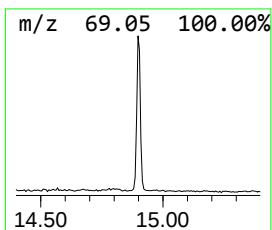
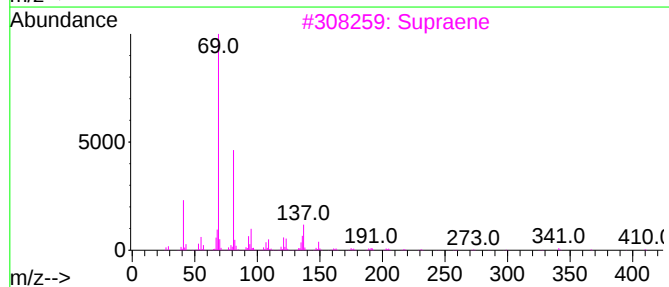
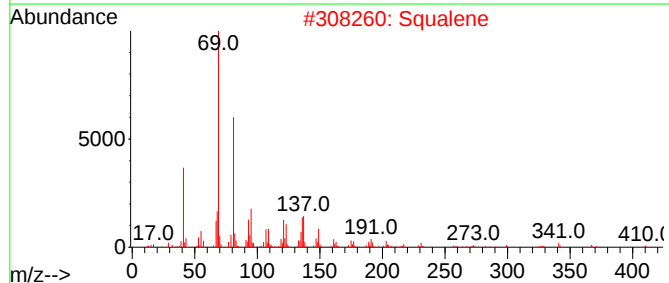
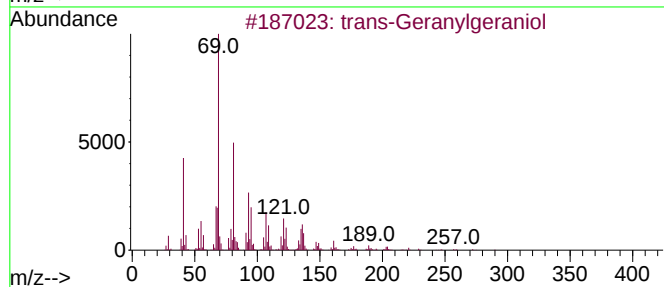
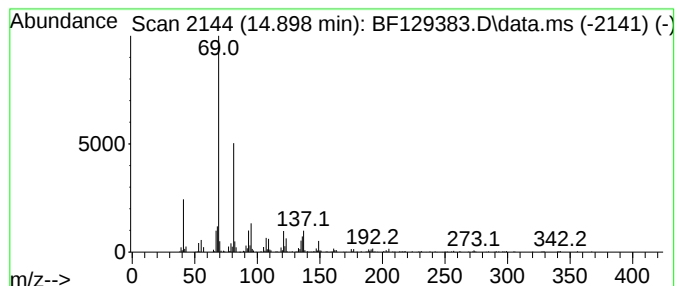
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 trans-Geranylgeraniol Concentration Rank 3

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 14.898 | 4.69 ng | 291639 | Perylene-d12     | 15.551 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|------|-------------------------------------|-----|----------|--------------|------|
| 1    |      | trans-Geranylgeraniol               | 290 | C20H34O  | 024034-73-9  | 92   |
| 2    |      | Squalene                            | 410 | C30H50   | 000111-02-4  | 90   |
| 3    |      | Supraene                            | 410 | C30H50   | 007683-64-9  | 87   |
| 4    |      | 5,9,13,17-Tetramethyl 4,8,12,16-... | 332 | C22H36O2 | 1000432-37-9 | 78   |
| 5    |      | 1,5,9-Undecatriene, 2,6,10-trime... | 192 | C14H24   | 062951-96-6  | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072722\  
 Data File : BF129383.D  
 Acq On : 27 Jul 2022 18:51  
 Operator : CG\JU  
 Sample : N3861-13  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 MW-21RR

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF072522.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 |
| 2-Butanol, 2,3-... | 3.263  | 2.3     | ng    | 180799   | 1                     | 6.881  | 1597070 | 20.0 |
| 2-Pentanone, 4-... | 5.116  | 2.4     | ng    | 187616   | 1                     | 6.881  | 1597070 | 20.0 |
| Benzene, propyl-   | 6.340  | 3.7     | ng    | 294526   | 1                     | 6.881  | 1597070 | 20.0 |
| Indane             | 7.057  | 11.1    | ng    | 888425   | 1                     | 6.881  | 1597070 | 20.0 |
| Benzene, 1,3-di... | 7.128  | 2.7     | ng    | 213202   | 1                     | 6.881  | 1597070 | 20.0 |
| Benzene, 1,2-di... | 7.216  | 2.4     | ng    | 191993   | 1                     | 6.881  | 1597070 | 20.0 |
| Benzene, 1,2,4,... | 7.645  | 3.9     | ng    | 449091   | 2                     | 8.163  | 2317840 | 20.0 |
| Indan, 1-methyl-   | 7.834  | 2.1     | ng    | 244052   | 2                     | 8.163  | 2317840 | 20.0 |
| Benzene, 2-ethe... | 7.898  | 2.8     | ng    | 323597   | 2                     | 8.163  | 2317840 | 20.0 |
| Hexadecanoic ac... | 11.792 | 2.7     | ng    | 280938   | 4                     | 11.416 | 2054520 | 20.0 |
| n-Hexadecanoic ... | 11.933 | 3.4     | ng    | 346287   | 4                     | 11.416 | 2054520 | 20.0 |
| Cyclohexadecane    | 12.475 | 4.0     | ng    | 412592   | 4                     | 11.416 | 2054520 | 20.0 |
| 6-Octadecenoic ... | 12.504 | 7.0     | ng    | 719141   | 4                     | 11.416 | 2054520 | 20.0 |
| Methyl stearate    | 12.592 | 2.3     | ng    | 236341   | 4                     | 11.416 | 2054520 | 20.0 |
| Bis(2-ethylhexy... | 12.651 | 3.3     | ng    | 337184   | 4                     | 11.416 | 2054520 | 20.0 |
| 1-Eicosene         | 13.916 | 2.9     | ng    | 201088   | 5                     | 14.057 | 1396360 | 20.0 |
| trans-Geranylge... | 14.898 | 4.7     | ng    | 291639   | 6                     | 15.551 | 1244580 | 20.0 |