

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF060821\  
 Data File : BF124178.D  
 Acq On : 8 Jun 2021 13:51  
 Operator : JU/CG  
 Sample : M2615-01  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 1245

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-BF060321.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF124178.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         | 2.116       | 3             | 5           | 11           | rVB      | 1563819        | 1469079       | 89.13%          | 14.924%       |
| 2         | 5.487       | 574           | 578         | 587          | rBB      | 535950         | 456450        | 27.69%          | 4.637%        |
| 3         | 6.498       | 746           | 750         | 757          | rBV      | 272911         | 265345        | 16.10%          | 2.696%        |
| 4         | 6.857       | 808           | 811         | 814          | rBV      | 263685         | 200336        | 12.15%          | 2.035%        |
| 5         | 7.422       | 902           | 907         | 910          | rBV      | 798896         | 799205        | 48.49%          | 8.119%        |
| 6         | 8.045       | 1009          | 1013        | 1026         | rBV      | 305987         | 290497        | 17.62%          | 2.951%        |
| 7         | 8.139       | 1026          | 1029        | 1036         | rVB      | 300463         | 308538        | 18.72%          | 3.134%        |
| 8         | 8.263       | 1045          | 1050        | 1057         | rVB      | 188282         | 157132        | 9.53%           | 1.596%        |
| 9         | 8.416       | 1074          | 1076        | 1080         | rVB      | 21984          | 20118         | 1.22%           | 0.204%        |
| 10        | 9.222       | 1208          | 1213        | 1216         | rBV      | 1720785        | 1518572       | 92.13%          | 15.427%       |
| 11        | 9.333       | 1228          | 1232        | 1236         | rVB      | 25367          | 22321         | 1.35%           | 0.227%        |
| 12        | 9.898       | 1324          | 1328        | 1331         | rBV      | 376016         | 321703        | 19.52%          | 3.268%        |
| 13        | 10.527      | 1430          | 1435        | 1438         | rBV      | 54451          | 53611         | 3.25%           | 0.545%        |
| 14        | 10.692      | 1458          | 1463        | 1466         | rBV      | 1163988        | 1085526       | 65.86%          | 11.028%       |
| 15        | 10.963      | 1506          | 1509        | 1512         | rBV      | 17509          | 17056         | 1.03%           | 0.173%        |
| 16        | 11.386      | 1577          | 1581        | 1585         | rVB      | 440211         | 353170        | 21.43%          | 3.588%        |
| 17        | 11.774      | 1644          | 1647        | 1650         | rVB2     | 53032          | 51363         | 3.12%           | 0.522%        |
| 18        | 11.910      | 1666          | 1670        | 1674         | rBV3     | 97042          | 102135        | 6.20%           | 1.038%        |
| 19        | 12.186      | 1714          | 1717        | 1719         | rBV      | 20023          | 17627         | 1.07%           | 0.179%        |
| 20        | 12.480      | 1765          | 1767        | 1770         | rVB4     | 25159          | 22264         | 1.35%           | 0.226%        |
| 21        | 12.698      | 1802          | 1804        | 1807         | rBV2     | 49380          | 43075         | 2.61%           | 0.438%        |
| 22        | 12.774      | 1813          | 1817        | 1822         | rVB7     | 15778          | 30038         | 1.82%           | 0.305%        |
| 23        | 12.827      | 1822          | 1826        | 1829         | rBV5     | 10584          | 17727         | 1.08%           | 0.180%        |
| 24        | 12.974      | 1846          | 1851        | 1854         | rBV      | 1742153        | 1648321       | 100.00%         | 16.745%       |
| 25        | 13.892      | 2003          | 2007        | 2014         | rBV6     | 29784          | 57040         | 3.46%           | 0.579%        |
| 26        | 14.027      | 2026          | 2030        | 2035         | rVB      | 291918         | 271743        | 16.49%          | 2.761%        |
| 27        | 15.509      | 2278          | 2282        | 2291         | rVB2     | 190864         | 243426        | 14.77%          | 2.473%        |

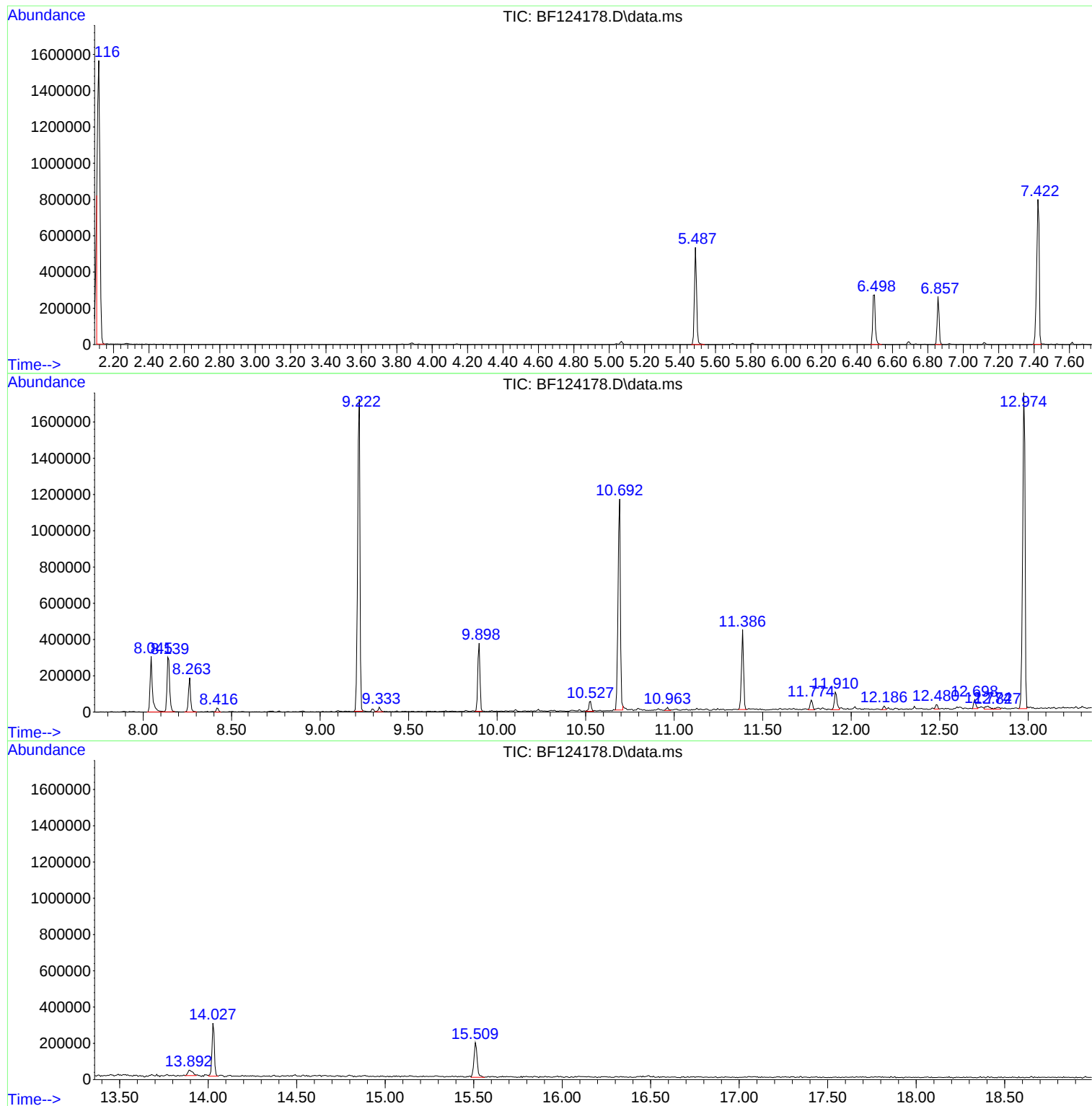
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF060321.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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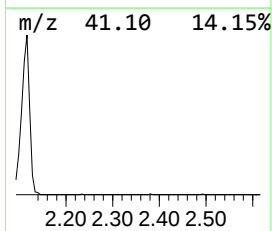
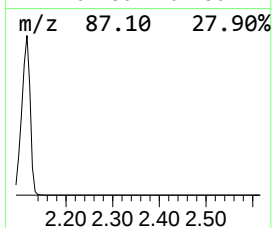
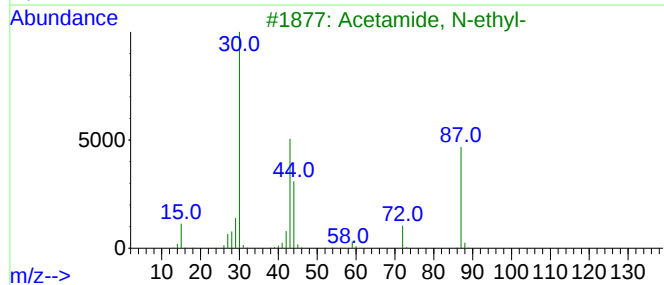
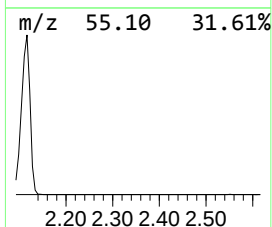
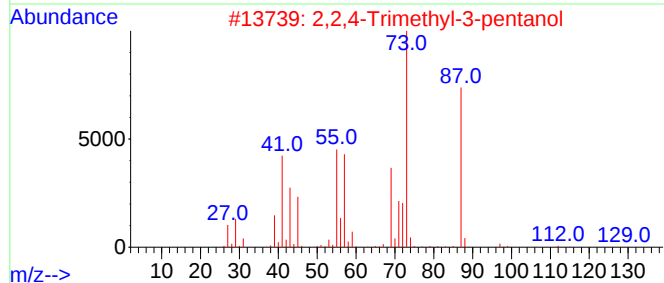
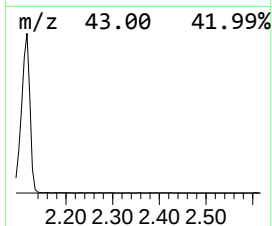
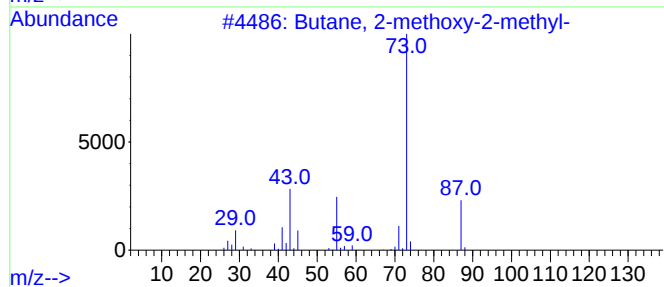
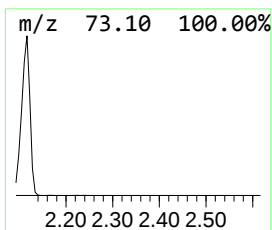
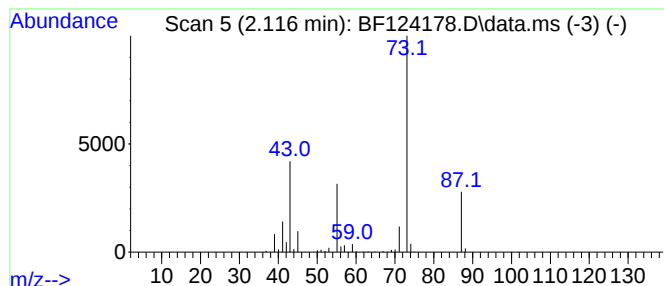
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF060321.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 2.116 | 146.66 ng | 1469080 | 1,4-Dichlorobenzene-d4 | 6.857 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 74   |
| 2    |    |   | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 28   |
| 3    |    |   | Acetamide, N-ethyl-         | 87  | C4H9NO  | 000625-50-3 | 25   |
| 4    |    |   | 2-Methyl-5-hexen-3-ol       | 114 | C7H14O  | 032815-70-6 | 23   |
| 5    |    |   | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 9    |



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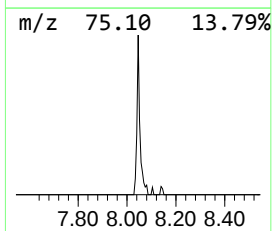
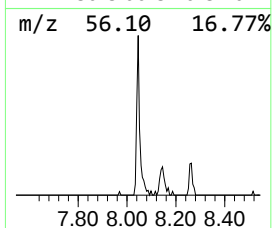
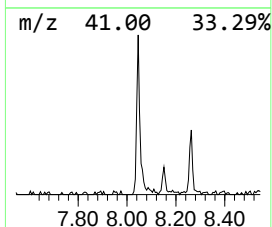
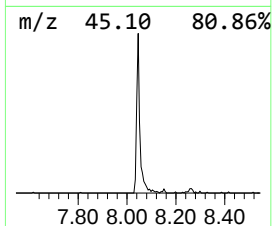
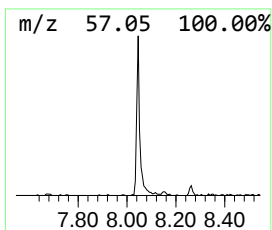
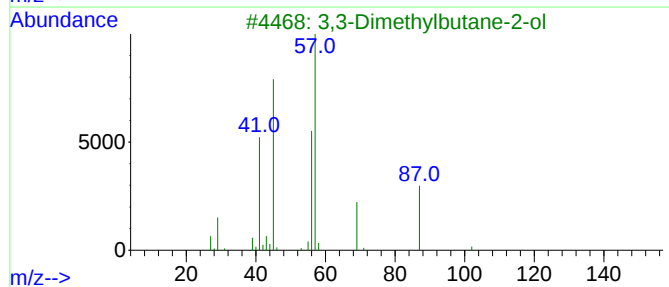
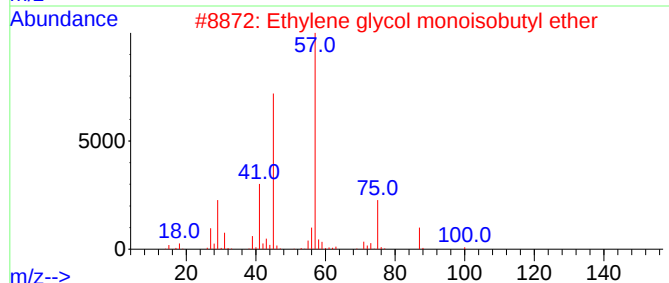
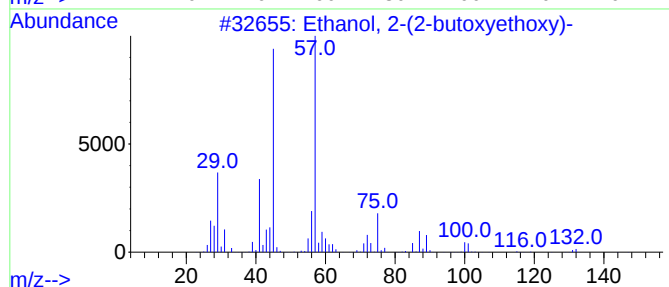
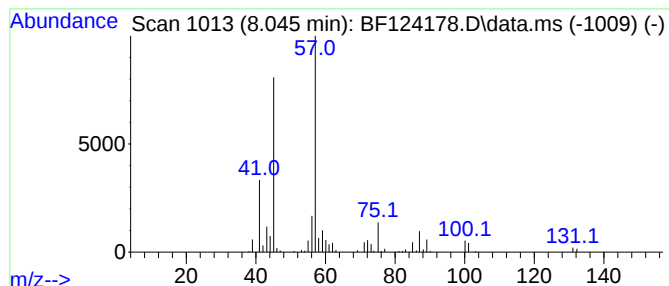
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 8.045 | 18.83 ng | 290497 | Naphthalene-d8   | 8.139 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Ethanol, 2-(2-butoxyethoxy)-       | 162 | C8H18O3 | 000112-34-5 | 90   |
| 2    |    |   | Ethylene glycol monoisobutyl ether | 118 | C6H14O2 | 004439-24-1 | 59   |
| 3    |    |   | 3,3-Dimethylbutane-2-ol            | 102 | C6H14O  | 000464-07-3 | 53   |
| 4    |    |   | Ethanol, 2-butoxy-                 | 118 | C6H14O2 | 000111-76-2 | 50   |
| 5    |    |   | 2-Propanol, 1-butoxy-              | 132 | C7H16O2 | 005131-66-8 | 45   |



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ALS Vial : 8 Sample Multiplier: 1

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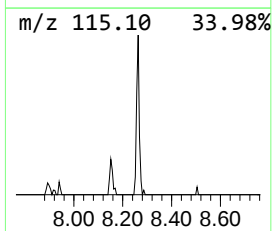
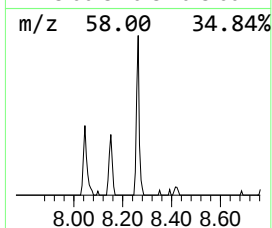
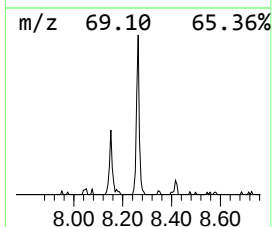
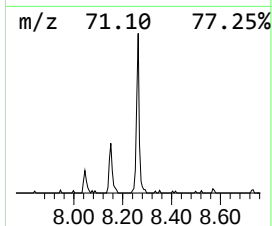
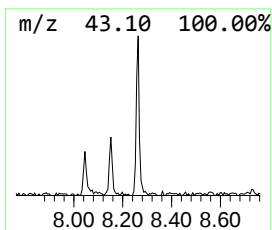
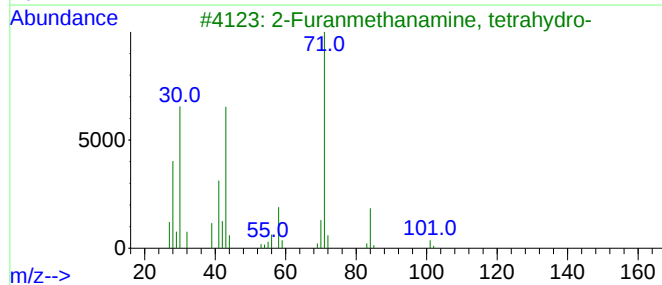
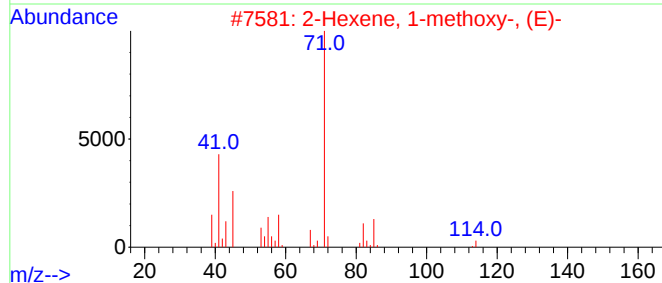
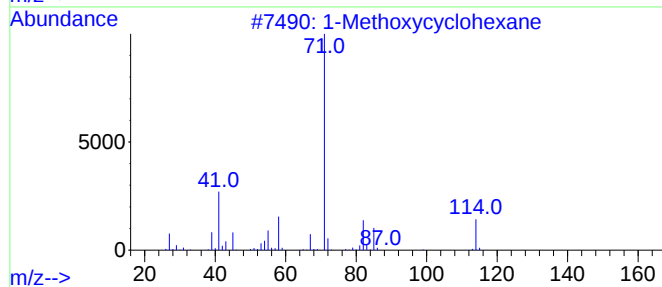
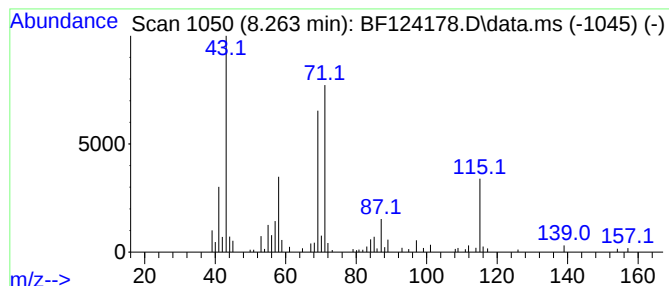
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TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 1-Methoxycyclohexane Concentration Rank 3

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 8.263 | 10.19 ng | 157132 | Naphthalene-d8   | 8.139 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1-Methoxycyclohexane            | 114 | C7H14O  | 000931-56-6 | 50   |
| 2    |    |   | 2-Hexene, 1-methoxy-, (E)-      | 114 | C7H14O  | 056052-83-6 | 47   |
| 3    |    |   | 2-Furanmethanamine, tetrahydro- | 101 | C5H11NO | 004795-29-3 | 47   |
| 4    |    |   | 3-Buten-2-ol, 2-methyl-         | 86  | C5H10O  | 000115-18-4 | 43   |
| 5    |    |   | 4-Heptanone                     | 114 | C7H14O  | 000123-19-3 | 43   |



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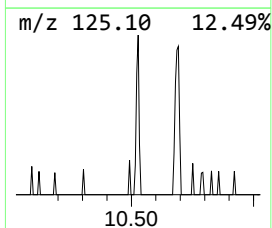
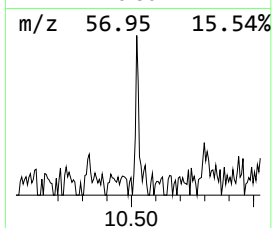
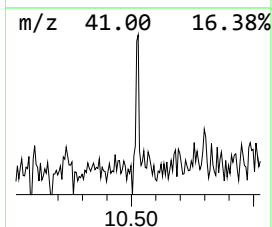
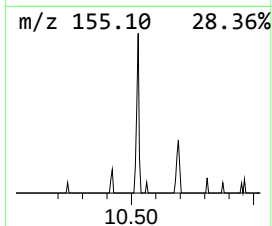
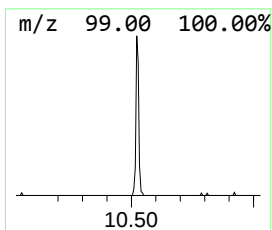
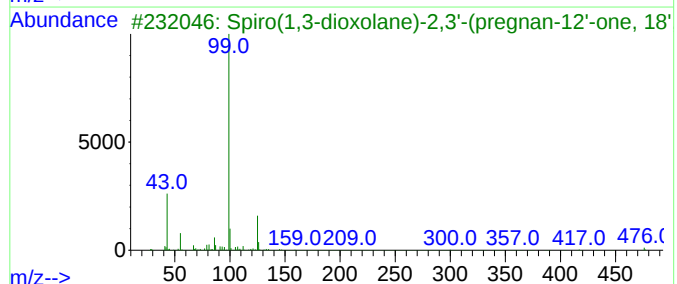
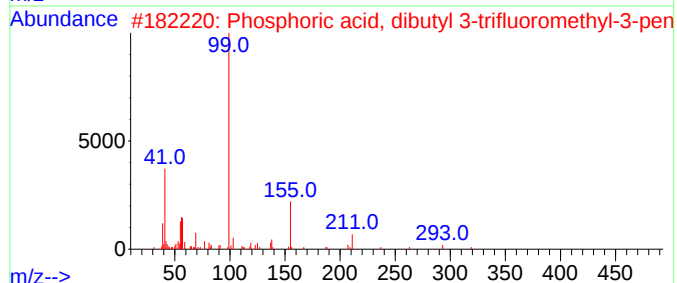
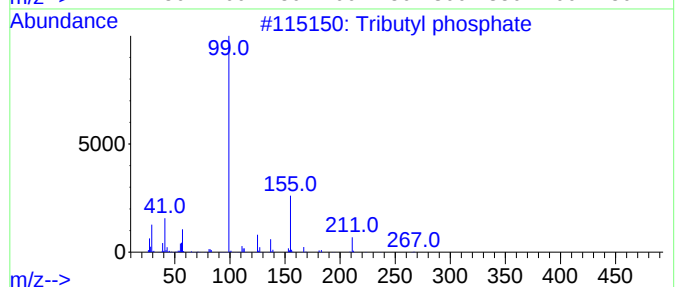
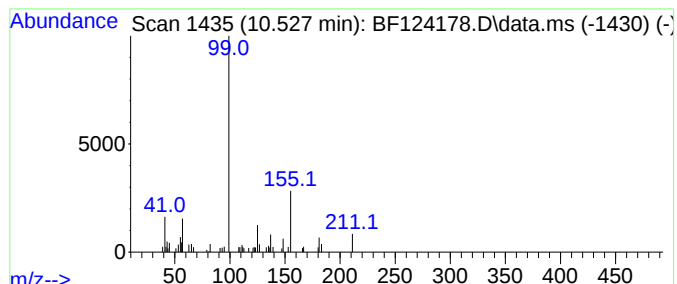
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Peak Number 4 Tributyl phosphate Concentration Rank 6

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.  |
|--------|---------|-------|------------------|-------|
| 10.528 | 3.33 ng | 53611 | Acenaphthene-d10 | 9.898 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Tributyl phosphate                  | 266 | C12H27O4P   | 000126-73-8  | 86   |
| 2    |    |   | Phosphoric acid, dibutyl 3-trifl... | 348 | C14H28F3O4P | 1000298-93-8 | 50   |
| 3    |    |   | Spiro(1,3-dioxolane)-2,3'-(pregn... | 476 | C27H40O7    | 1000194-70-1 | 47   |
| 4    |    |   | Phosphoric acid, dibutyl 1,1-dim... | 352 | C13H25F4O4P | 1000298-93-9 | 38   |
| 5    |    |   | Tricyclo[5.3.0.0(2,6)]decan-5-on... | 272 | C13H17ClO4  | 1000164-70-7 | 32   |



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

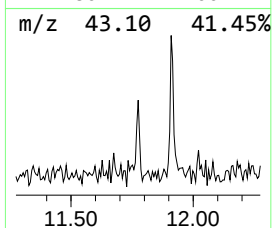
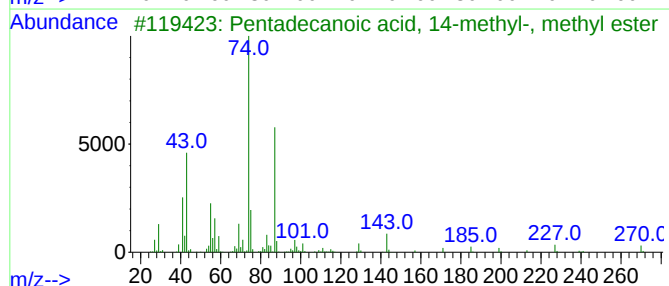
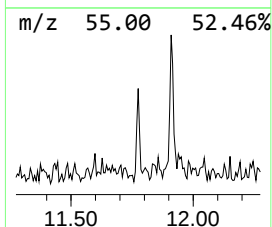
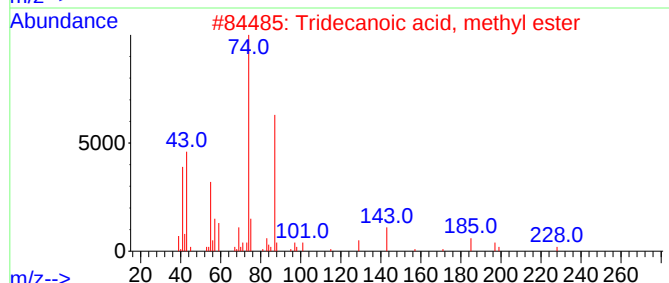
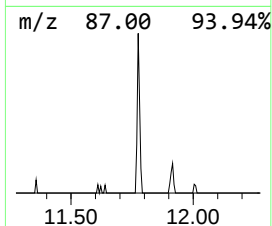
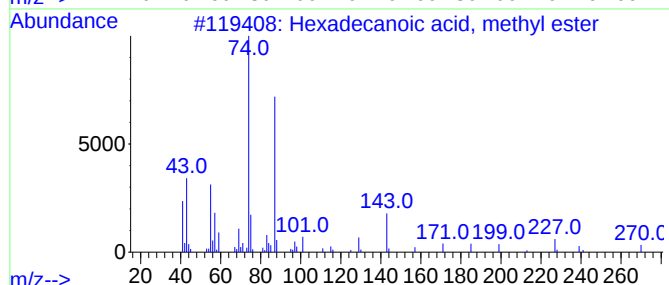
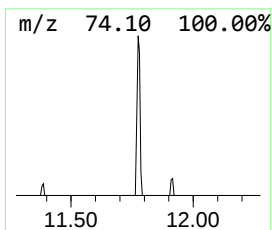
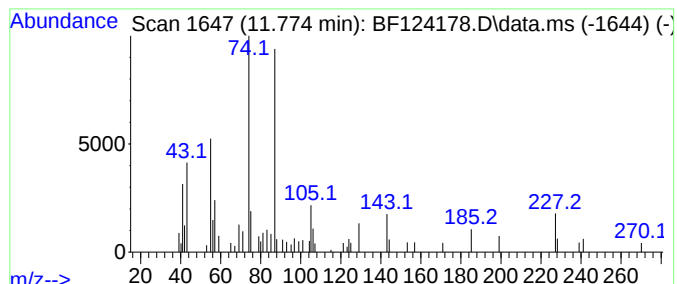
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Peak Number 5 Hexadecanoic acid, methyl e... Concentration Rank 7

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 11.774 | 2.91 ng | 51363 | Phenanthrene-d10 | 11.386 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Hexadecanoic acid, methyl ester     | 270 | C17H34O2 | 000112-39-0 | 90   |
| 2    |    |   | Tridecanoic acid, methyl ester      | 228 | C14H28O2 | 001731-88-0 | 64   |
| 3    |    |   | Pentadecanoic acid, 14-methyl-, ... | 270 | C17H34O2 | 005129-60-2 | 49   |
| 4    |    |   | Pentadecanoic acid, methyl ester    | 256 | C16H32O2 | 007132-64-1 | 43   |
| 5    |    |   | Methyl tetradecanoate               | 242 | C15H30O2 | 000124-10-7 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF060821\  
Data File : BF124178.D  
Acq On : 8 Jun 2021 13:51  
Operator : JU/CG  
Sample : M2615-01  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1245

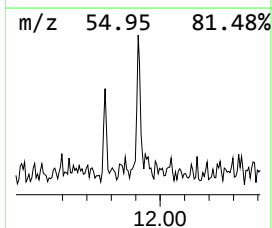
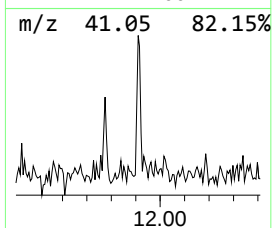
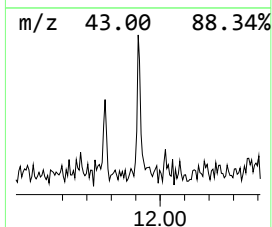
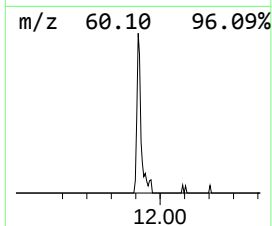
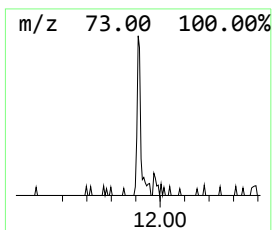
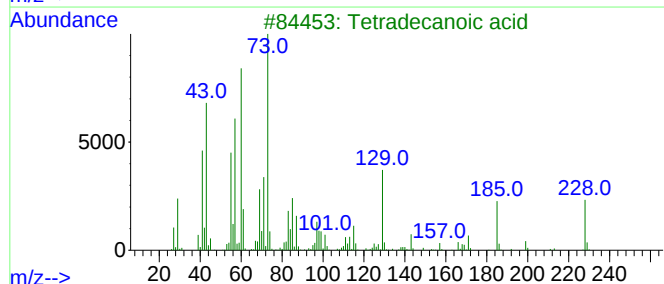
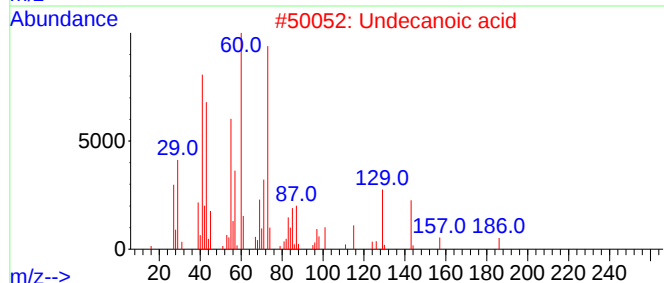
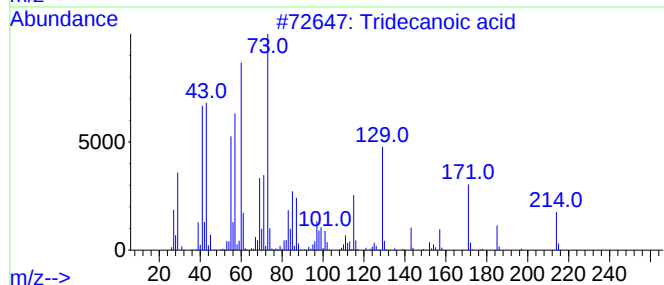
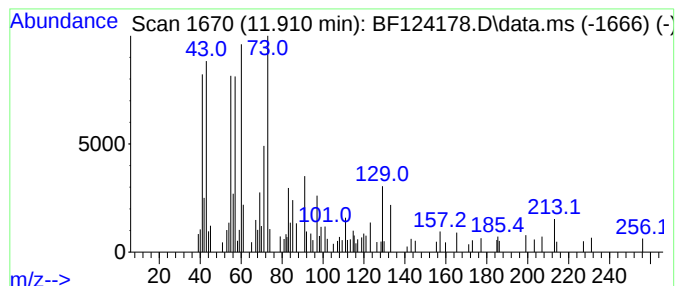
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 Tridecanoic acid Concentration Rank 4

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 11.910 | 5.78 ng | 102135 | Phenanthrene-d10 | 11.386 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm  | CAS#         | Qual |
|------|----|---|------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Tridecanoic acid                   | 214 | C13H26O2 | 000638-53-9  | 91   |
| 2    |    |   | Undecanoic acid                    | 186 | C11H22O2 | 000112-37-8  | 86   |
| 3    |    |   | Tetradecanoic acid                 | 228 | C14H28O2 | 000544-63-8  | 64   |
| 4    |    |   | n-PROPYL NONYL ETHER               | 186 | C12H26O  | 1000130-69-3 | 25   |
| 5    |    |   | Oxalic acid, allyl hexadecyl ester | 354 | C21H38O4 | 1000309-24-4 | 18   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF060821\  
Data File : BF124178.D  
Acq On : 8 Jun 2021 13:51  
Operator : JU/CG  
Sample : M2615-01  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1245

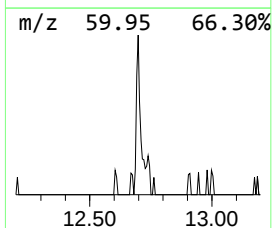
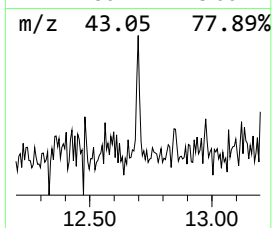
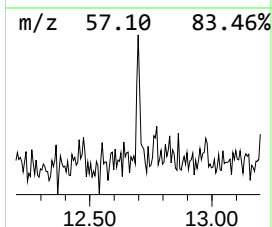
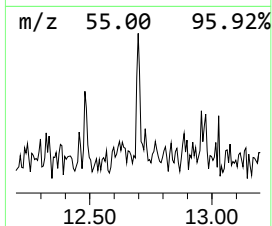
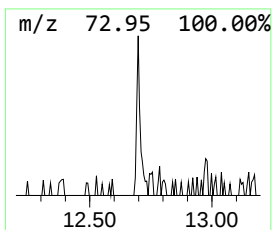
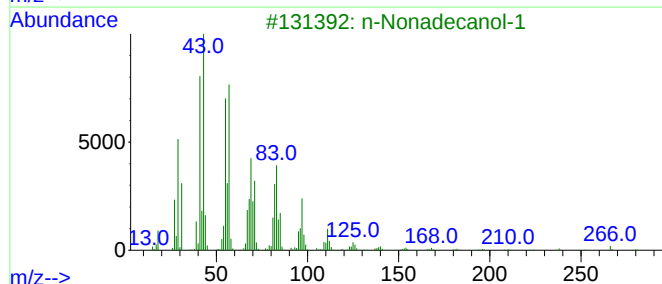
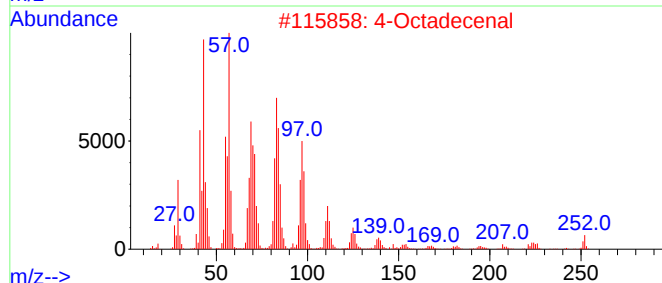
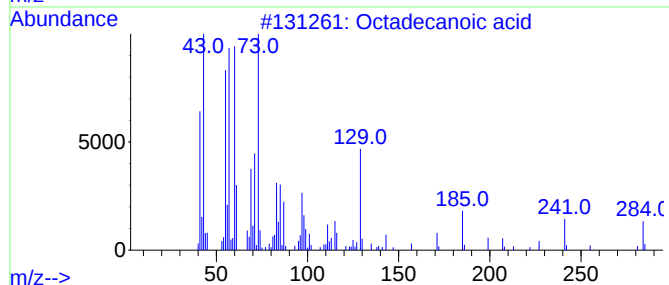
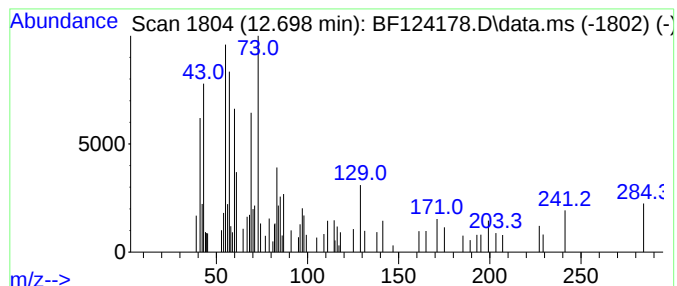
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 Octadecanoic acid Concentration Rank 8

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 12.698 | 2.44 ng | 43075 | Phenanthrene-d10 | 11.386 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm    | CAS#        | Qual |
|------|----|---|-----------------------------------|-----|------------|-------------|------|
| 1    |    |   | Octadecanoic acid                 | 284 | C18H36O2   | 000057-11-4 | 50   |
| 2    |    |   | 4-Octadecenal                     | 266 | C18H34O    | 056554-98-4 | 30   |
| 3    |    |   | n-Nonadecanol-1                   | 284 | C19H40O    | 001454-84-8 | 22   |
| 4    |    |   | 11-Tricosene                      | 322 | C23H46     | 052078-56-5 | 22   |
| 5    |    |   | Bromoacetic acid, hexadecyl ester | 362 | C18H35BrO2 | 005454-48-8 | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF060821\  
Data File : BF124178.D  
Acq On : 8 Jun 2021 13:51  
Operator : JU/CG  
Sample : M2615-01  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1245

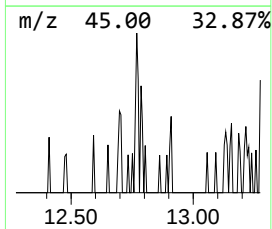
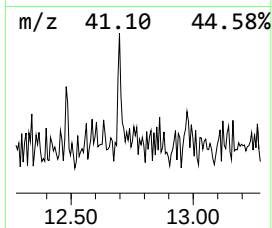
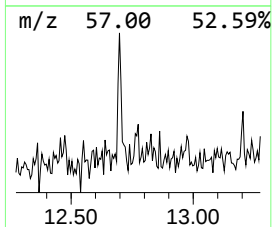
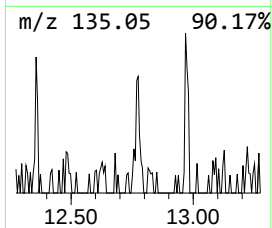
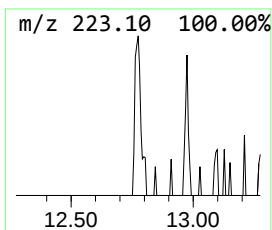
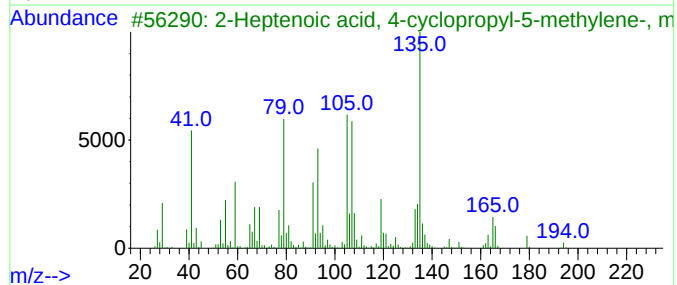
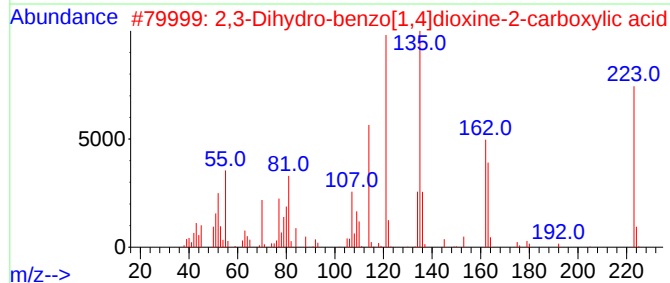
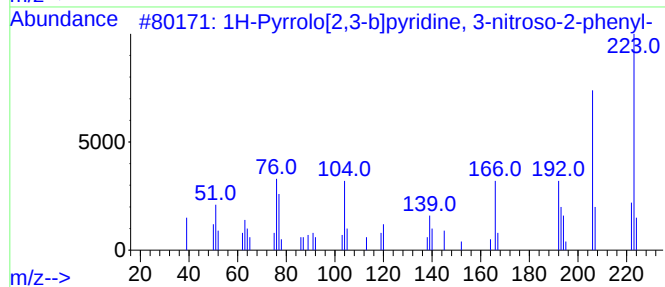
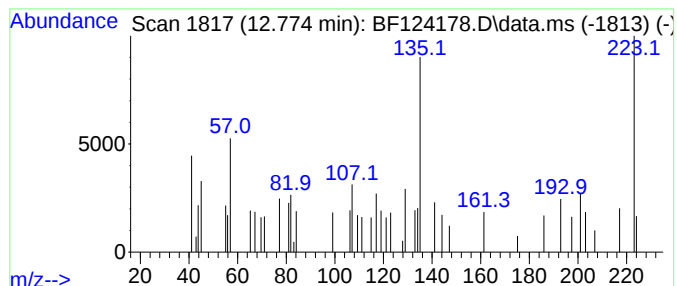
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 unknown12.774 Concentration Rank 9

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 12.774 | 2.21 ng | 30038 | Chrysene-d12     | 14.027 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 1H-Pyrrolo[2,3-b]pyridine, 3-nit... | 223 | C13H9N3O  | 023616-56-0  | 22   |
| 2    |    |   | 2,3-Dihydro-benzo[1,4]dioxine-2-... | 223 | C11H13NO4 | 1000297-29-5 | 17   |
| 3    |    |   | 2-Heptenoic acid, 4-cyclopropyl-... | 194 | C12H18O2  | 074793-23-0  | 16   |
| 4    |    |   | 2H-Indazol-3-amine, N-methyl-2-p... | 223 | C14H13N3  | 097990-17-5  | 14   |
| 5    |    |   | 6-Cyano-2,5-dimethoxy-4-nitroani... | 223 | C9H9N3O4  | 1000214-34-6 | 12   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF060821\  
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Acq On : 8 Jun 2021 13:51  
Operator : JU/CG  
Sample : M2615-01  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
1245

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

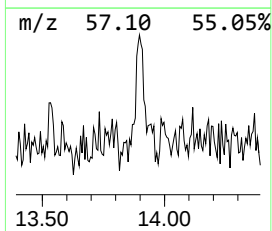
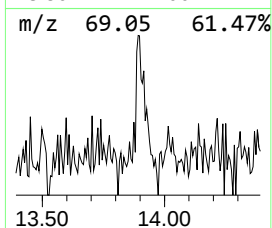
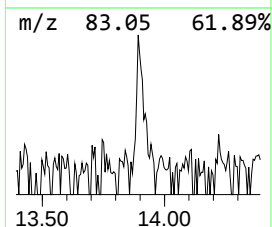
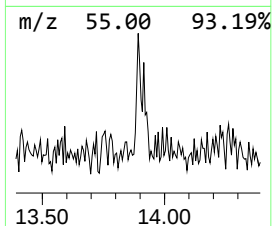
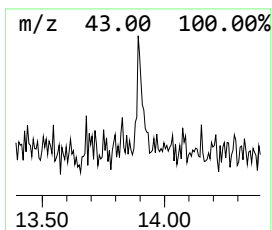
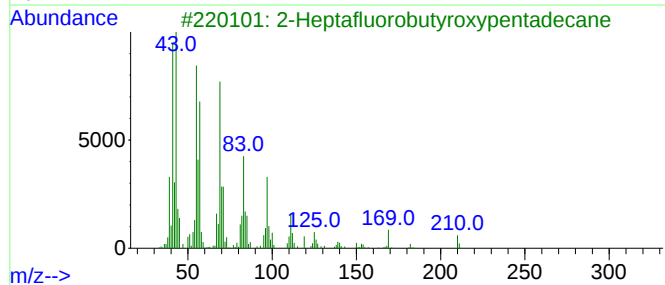
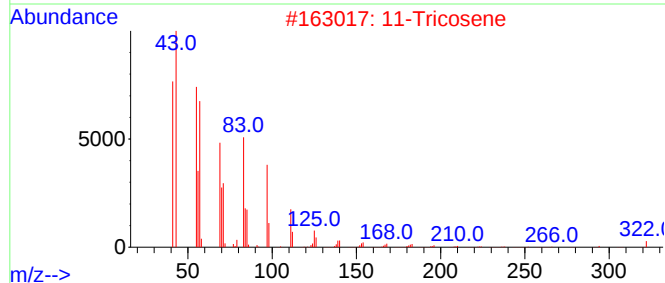
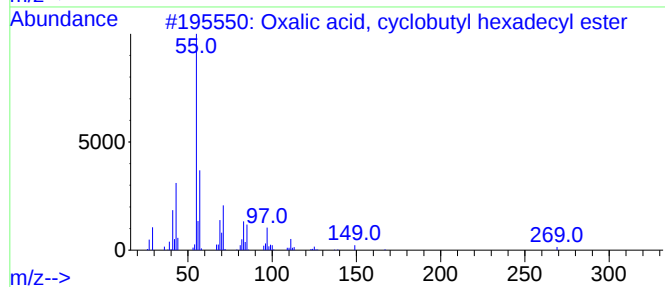
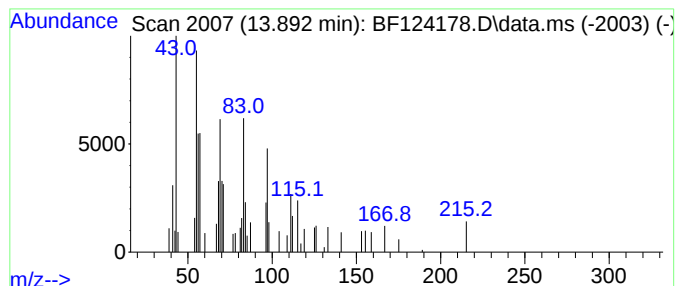
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Oxalic acid, cyclobutyl hex... Concentration Rank 5

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 13.892 | 4.20 ng | 57040 | Chrysene-d12     | 14.027 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Oxalic acid, cyclobutyl hexadecy... | 368 | C22H40O4   | 1000309-70-6 | 59   |
| 2    |    |   | 11-Tricosene                        | 322 | C23H46     | 052078-56-5  | 58   |
| 3    |    |   | 2-Heptafluorobutyroxy pentadecane   | 424 | C19H31F7O2 | 1000245-50-0 | 58   |
| 4    |    |   | 9-Eicosene, (E)-                    | 280 | C20H40     | 074685-29-3  | 53   |
| 5    |    |   | 1-Heptadecene                       | 238 | C17H34     | 006765-39-5  | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF060821\  
Data File : BF124178.D  
Acq On : 8 Jun 2021 13:51  
Operator : JU/CG  
Sample : M2615-01  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1245

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF060321.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |        |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|--------|------|
|                    |        |         |       |          | #                     | RT     | Resp   | Conc |
| Butane, 2-metho... | 2.116  | 146.7   | ng    | 1469080  | 1                     | 6.857  | 200336 | 20.0 |
| Ethanol, 2-(2-b... | 8.045  | 18.8    | ng    | 290497   | 2                     | 8.139  | 308538 | 20.0 |
| 1-Methoxycycloh... | 8.263  | 10.2    | ng    | 157132   | 2                     | 8.139  | 308538 | 20.0 |
| Tributyl phosphate | 10.528 | 3.3     | ng    | 53611    | 3                     | 9.898  | 321703 | 20.0 |
| Hexadecanoic ac... | 11.774 | 2.9     | ng    | 51363    | 4                     | 11.386 | 353170 | 20.0 |
| Tridecanoic acid   | 11.910 | 5.8     | ng    | 102135   | 4                     | 11.386 | 353170 | 20.0 |
| Octadecanoic acid  | 12.698 | 2.4     | ng    | 43075    | 4                     | 11.386 | 353170 | 20.0 |
| unknown12.774      | 12.774 | 2.2     | ng    | 30038    | 5                     | 14.027 | 271743 | 20.0 |
| Oxalic acid, cy... | 13.892 | 4.2     | ng    | 57040    | 5                     | 14.027 | 271743 | 20.0 |