

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082025\  
 Data File : BP025540.D  
 Acq On : 21 Aug 2025 14:56  
 Operator : CG/JU  
 Sample : Q2819-16 2X  
 Misc :  
 ALS Vial : 30 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 38M-S

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\_P\Methods\8270E-BP081425.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP025540.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         | 3.037       | 12            | 17          | 26           | rVB      | 502014         | 629629        | 15.96%          | 1.810%        |
| 2         | 4.943       | 336           | 341         | 349          | rVB      | 139823         | 188648        | 4.78%           | 0.542%        |
| 3         | 5.401       | 413           | 419         | 432          | rBV      | 1457751        | 2122903       | 53.80%          | 6.104%        |
| 4         | 6.966       | 676           | 685         | 695          | rBV      | 1439544        | 2245512       | 56.90%          | 6.456%        |
| 5         | 7.784       | 816           | 824         | 833          | rBV      | 833678         | 1332069       | 33.76%          | 3.830%        |
| 6         | 8.925       | 1011          | 1018        | 1030         | rBV      | 893673         | 1515648       | 38.41%          | 4.358%        |
| 7         | 10.554      | 1284          | 1295        | 1306         | rBV      | 1017726        | 1780748       | 45.12%          | 5.120%        |
| 8         | 13.019      | 1706          | 1714        | 1727         | rBV      | 2014663        | 3166961       | 80.25%          | 9.106%        |
| 9         | 14.248      | 1917          | 1923        | 1932         | rBV2     | 35979          | 71127         | 1.80%           | 0.205%        |
| 10        | 14.413      | 1942          | 1951        | 1964         | rBV      | 1441052        | 2121854       | 53.77%          | 6.101%        |
| 11        | 14.519      | 1964          | 1969        | 1980         | rVV      | 206835         | 307275        | 7.79%           | 0.883%        |
| 12        | 15.789      | 2179          | 2185        | 2191         | rBV      | 343469         | 478353        | 12.12%          | 1.375%        |
| 13        | 15.918      | 2201          | 2207        | 2216         | rBV      | 1576030        | 2245210       | 56.89%          | 6.456%        |
| 14        | 16.366      | 2279          | 2283        | 2287         | rBV2     | 33361          | 44716         | 1.13%           | 0.129%        |
| 15        | 16.771      | 2348          | 2352        | 2359         | rVB      | 60662          | 95616         | 2.42%           | 0.275%        |
| 16        | 16.948      | 2376          | 2382        | 2393         | rBV      | 578400         | 887922        | 22.50%          | 2.553%        |
| 17        | 17.054      | 2395          | 2400        | 2401         | rVV      | 60204          | 81794         | 2.07%           | 0.235%        |
| 18        | 17.077      | 2401          | 2404        | 2411         | rVB      | 95702          | 152192        | 3.86%           | 0.438%        |
| 19        | 17.207      | 2420          | 2426        | 2432         | rBV2     | 1587671        | 2276970       | 57.70%          | 6.547%        |
| 20        | 17.907      | 2542          | 2545        | 2549         | rBV5     | 21756          | 39519         | 1.00%           | 0.114%        |
| 21        | 18.148      | 2581          | 2586        | 2594         | rBV      | 432317         | 747628        | 18.95%          | 2.150%        |
| 22        | 18.801      | 2694          | 2697        | 2705         | rVB2     | 73741          | 135468        | 3.43%           | 0.390%        |
| 23        | 18.930      | 2715          | 2719        | 2726         | rBV      | 510090         | 703884        | 17.84%          | 2.024%        |
| 24        | 18.995      | 2727          | 2730        | 2732         | rBV3     | 104738         | 131742        | 3.34%           | 0.379%        |
| 25        | 19.330      | 2783          | 2787        | 2791         | rBV      | 171156         | 262256        | 6.65%           | 0.754%        |
| 26        | 19.471      | 2807          | 2811        | 2817         | rBV      | 241026         | 456637        | 11.57%          | 1.313%        |
| 27        | 19.624      | 2835          | 2837        | 2840         | rVB2     | 90035          | 87308         | 2.21%           | 0.251%        |
| 28        | 19.707      | 2848          | 2851        | 2855         | rVB      | 118426         | 148311        | 3.76%           | 0.426%        |
| 29        | 19.848      | 2872          | 2875        | 2878         | rBV      | 99877          | 137429        | 3.48%           | 0.395%        |
| 30        | 19.924      | 2883          | 2888        | 2893         | rVB      | 3025518        | 3946275       | 100.00%         | 11.347%       |
| 31        | 20.512      | 2985          | 2988        | 2991         | rVB      | 173733         | 195729        | 4.96%           | 0.563%        |
| 32        | 21.318      | 3121          | 3125        | 3130         | rBV2     | 282263         | 442607        | 11.22%          | 1.273%        |
| 33        | 21.571      | 3165          | 3168        | 3175         | rVB      | 214645         | 276494        | 7.01%           | 0.795%        |
| 34        | 21.653      | 3176          | 3182        | 3187         | rVB      | 1681660        | 2512315       | 63.66%          | 7.224%        |
| 35        | 25.030      | 3748          | 3756        | 3764         | rBV      | 995281         | 2810610       | 71.22%          | 8.081%        |

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Sample : Q2819-16 2X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
38M-S

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\_P\Methods\8270E-BP081425.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

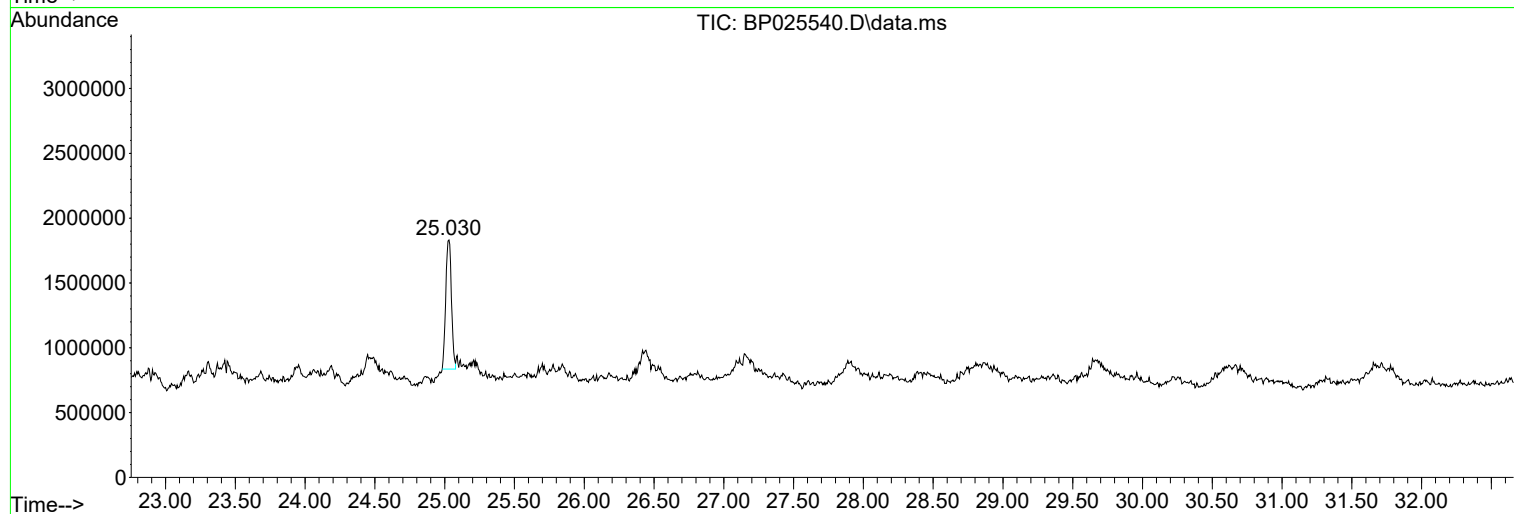
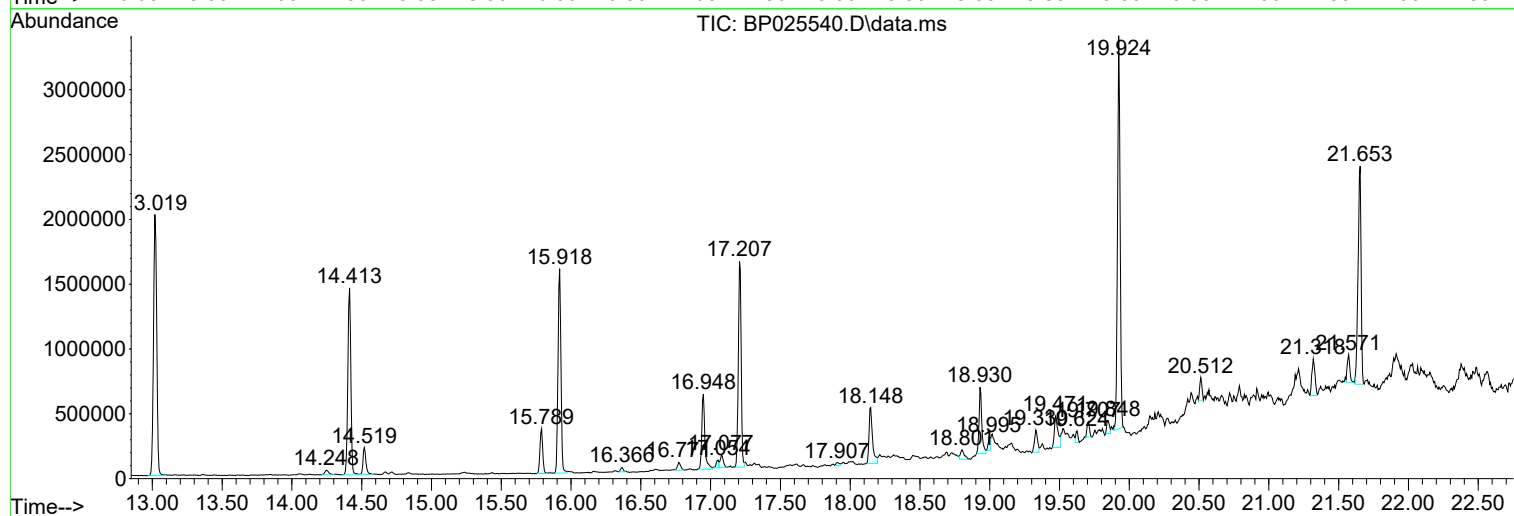
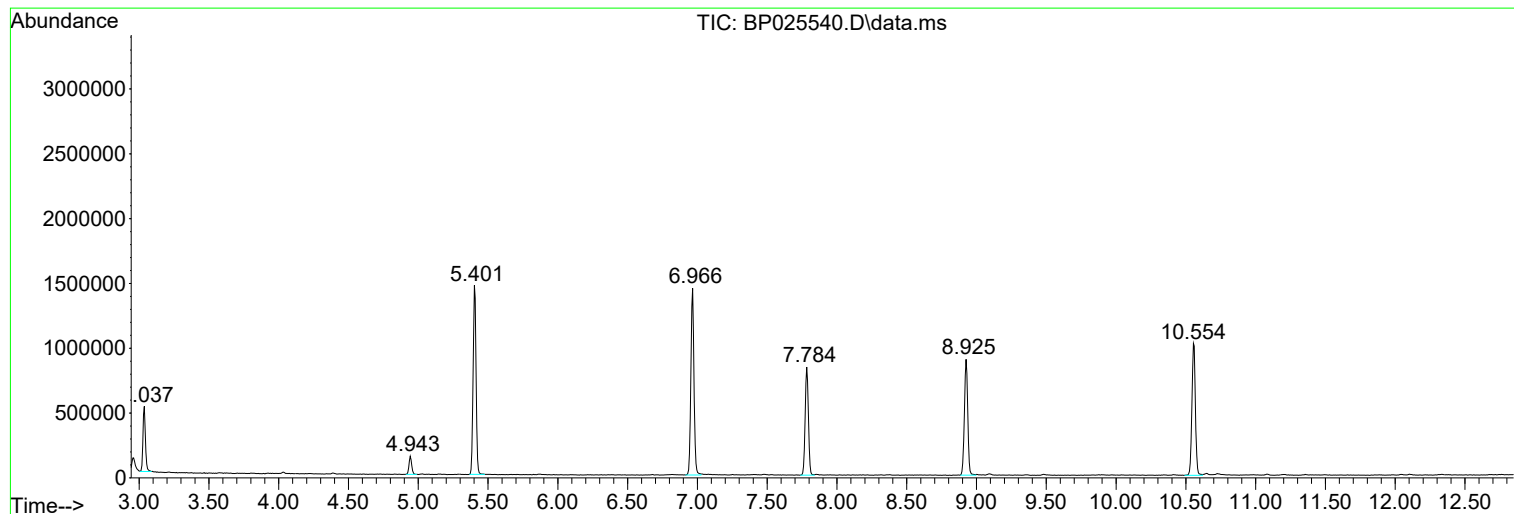
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP081425.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Sample : Q2819-16 2X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
38M-S

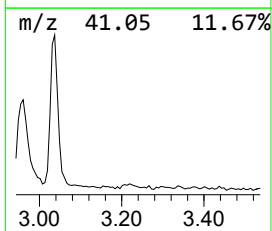
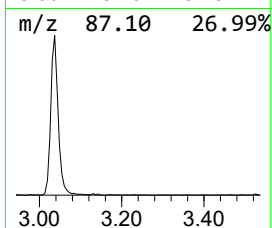
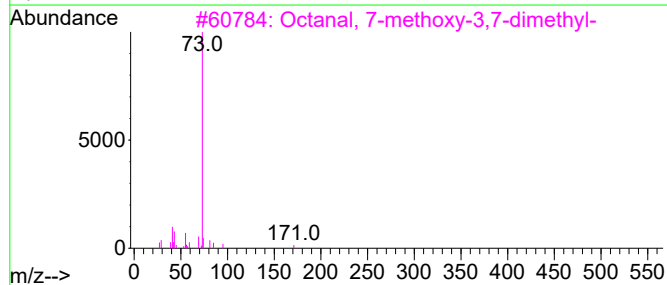
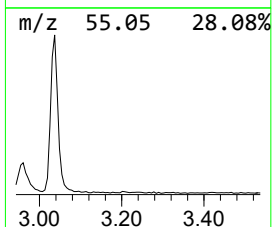
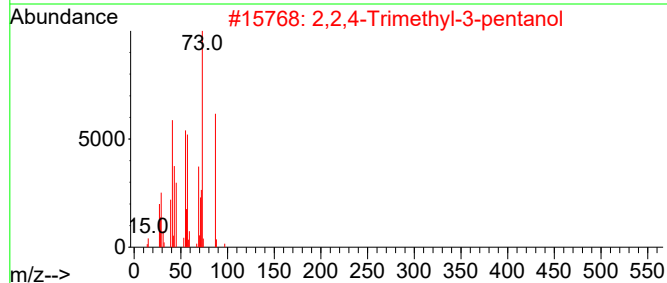
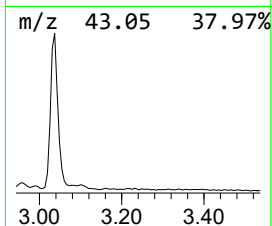
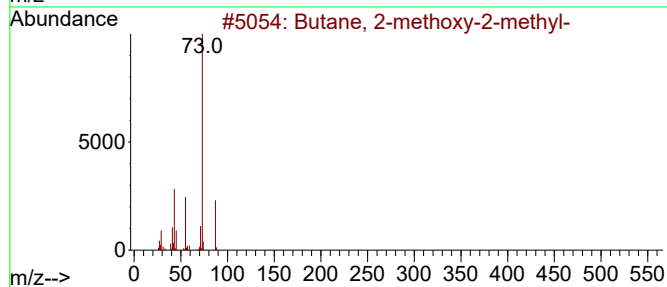
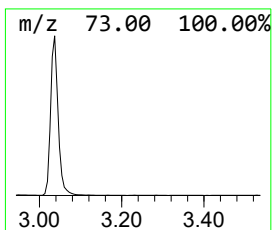
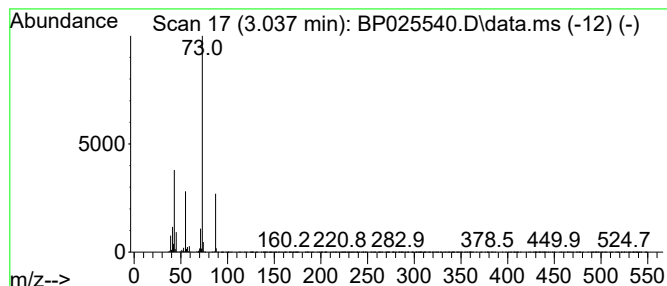
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP081425.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 3.037 | 9.45 ng | 629629 | 1,4-Dichlorobenzene-d4 | 7.784 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm  | CAS#        | Qual |
|------|----|---|----------------------------------|-----|----------|-------------|------|
| 1    |    |   | Butane, 2-methoxy-2-methyl-      | 102 | C6H14O   | 000994-05-8 | 83   |
| 2    |    |   | 2,2,4-Trimethyl-3-pentanol       | 130 | C8H18O   | 005162-48-1 | 40   |
| 3    |    |   | Octanal, 7-methoxy-3,7-dimethyl- | 186 | C11H22O2 | 003613-30-7 | 17   |
| 4    |    |   | Pentane, 3-methoxy-              | 102 | C6H14O   | 036839-67-5 | 12   |
| 5    |    |   | N-Ethylformamide                 | 73  | C3H7NO   | 000627-45-2 | 9    |



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Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
38M-S

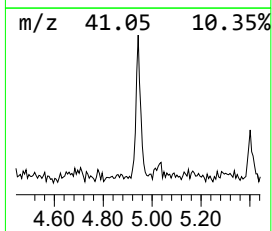
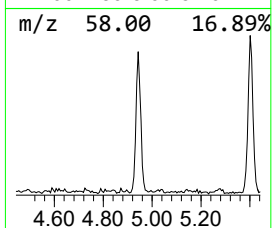
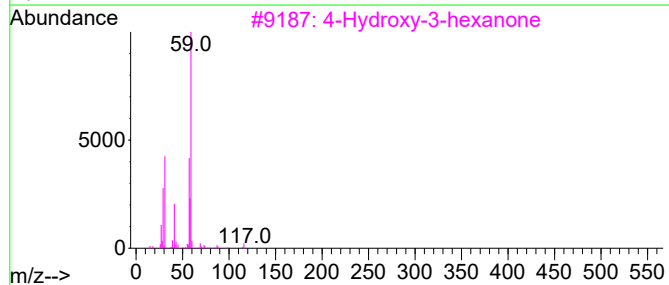
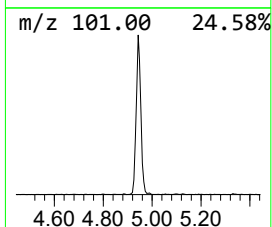
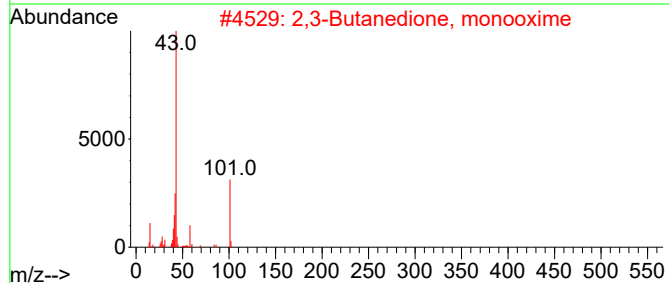
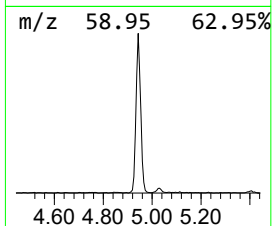
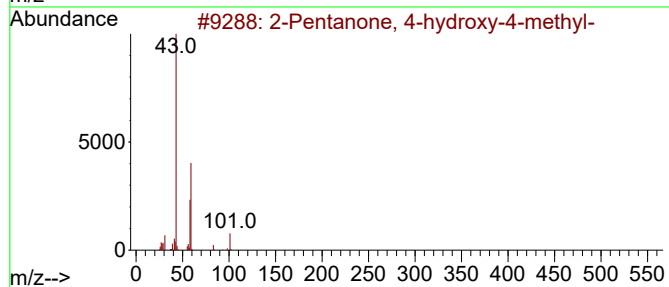
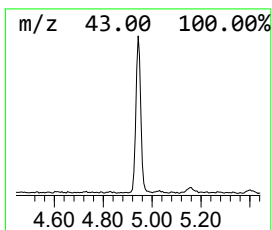
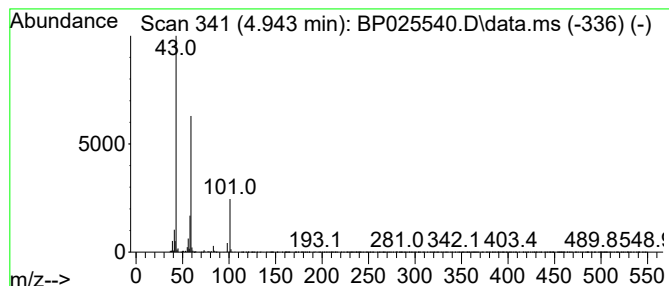
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP081425.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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 9

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 4.943 | 2.83 ng | 188648 | 1,4-Dichlorobenzene-d4 | 7.784 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-   | 116 | C6H12O2 | 000123-42-2 | 56   |
| 2    |    |   | 2,3-Butanedione, monooxime         | 101 | C4H7NO2 | 000057-71-6 | 25   |
| 3    |    |   | 4-Hydroxy-3-hexanone               | 116 | C6H12O2 | 004984-85-4 | 17   |
| 4    |    |   | 2,5,8,11-Tetraoxadodecane          | 178 | C8H18O4 | 000112-49-2 | 17   |
| 5    |    |   | (+)-4-Amino-4,5-dihydro-2(3H)-f... | 101 | C4H7NO2 | 016504-58-8 | 9    |



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ALS Vial : 30 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
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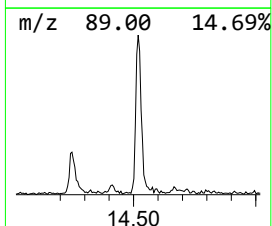
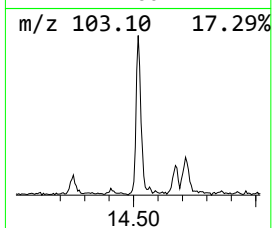
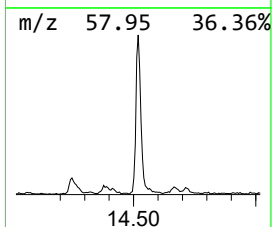
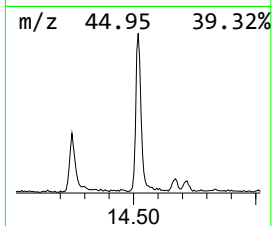
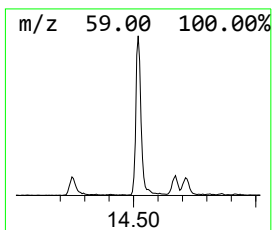
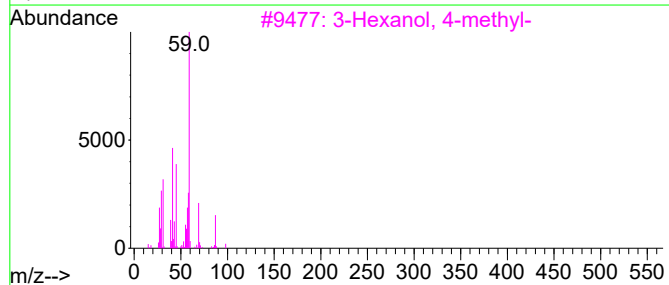
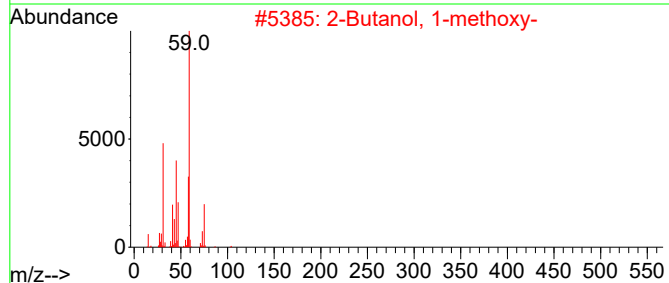
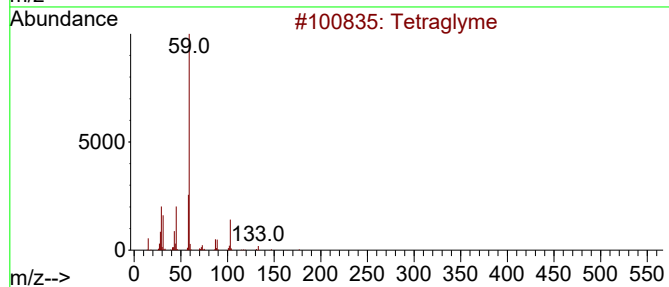
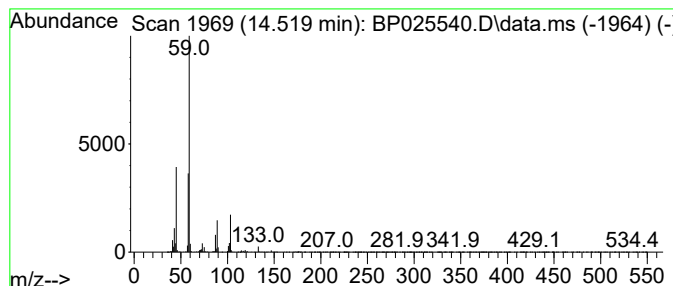
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP081425.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 3 Tetraglyme Concentration Rank 8

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 14.519 | 2.90 ng | 307275 | Acenaphthene-d10 | 14.413 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Tetraglyme                          | 222 | C10H22O5 | 000143-24-8  | 72   |
| 2    |    |   | 2-Butanol, 1-methoxy-               | 104 | C5H12O2  | 053778-73-7  | 64   |
| 3    |    |   | 3-Hexanol, 4-methyl-                | 116 | C7H16O   | 000615-29-2  | 64   |
| 4    |    |   | 2,5,8,11-Tetraoxadodecane           | 178 | C8H18O4  | 000112-49-2  | 59   |
| 5    |    |   | Methyl 2-(2-(2-methoxyethoxy)eth... | 192 | C8H16O5  | 1000366-88-2 | 56   |



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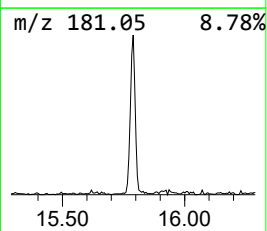
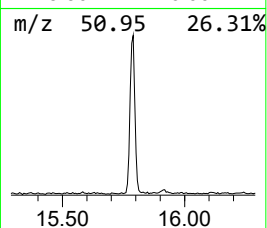
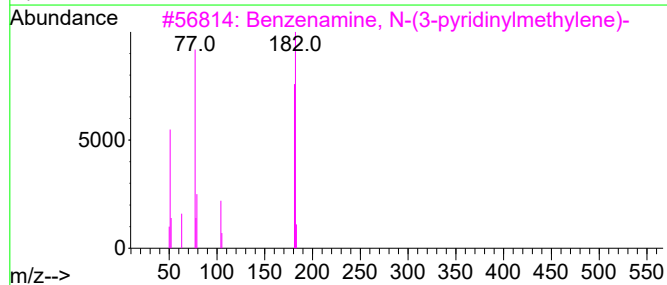
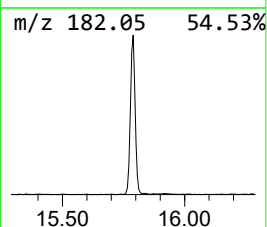
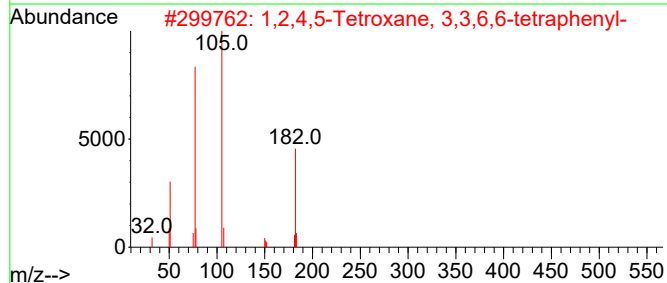
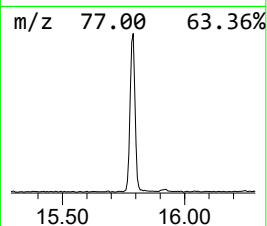
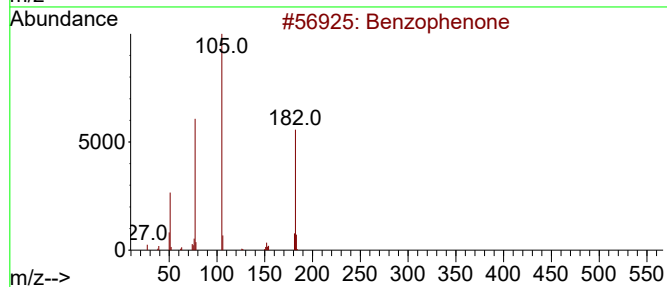
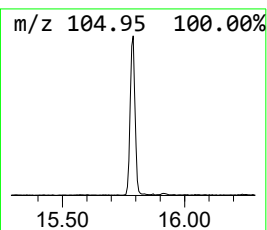
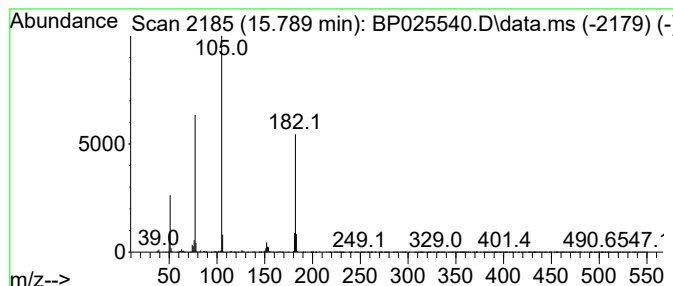
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP081425.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 4 Benzophenone Concentration Rank 5

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 15.789 | 4.51 ng | 478353 | Acenaphthene-d10 | 14.413 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | Benzophenone                        | 182 | C13H10O   | 000119-61-9 | 96   |
| 2    |    |   | 1,2,4,5-Tetroxane, 3,3,6,6-tetra... | 396 | C26H20O4  | 016204-36-7 | 64   |
| 3    |    |   | Benzenamine, N-(3-pyridinylmethy... | 182 | C12H10N2  | 029722-97-2 | 53   |
| 4    |    |   | 1,2,4-Trioxolane, 3,3,5-triphenyl-  | 304 | C20H16O3  | 023246-12-0 | 45   |
| 5    |    |   | 2,5-Dimethyl-7,7-diphenyl-3-aza-... | 279 | C18H17NO2 | 078950-91-1 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082025\  
Data File : BP025540.D  
Acq On : 21 Aug 2025 14:56  
Operator : CG/JU  
Sample : Q2819-16 2X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
38M-S

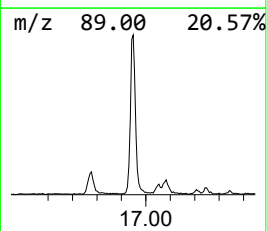
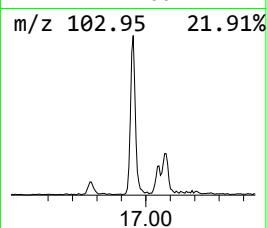
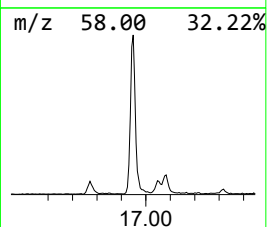
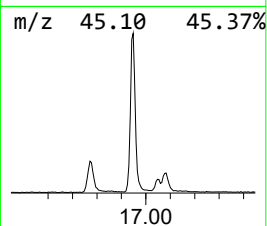
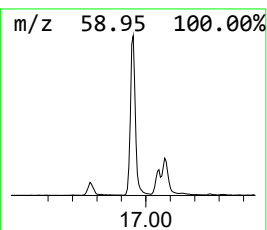
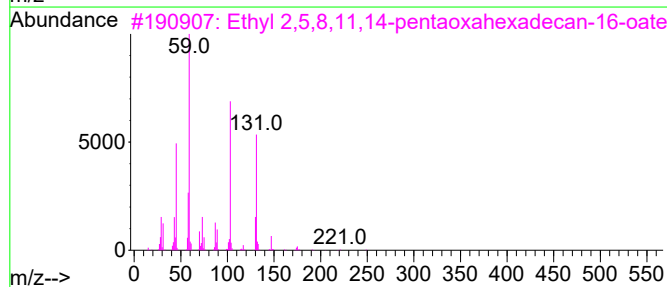
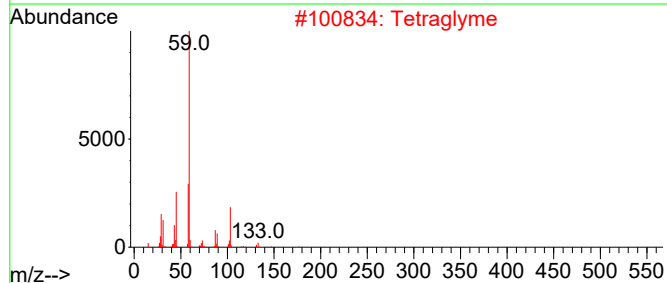
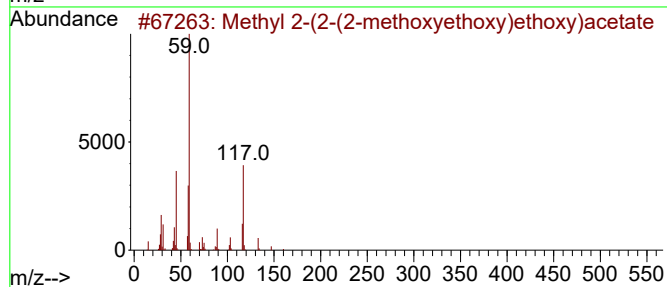
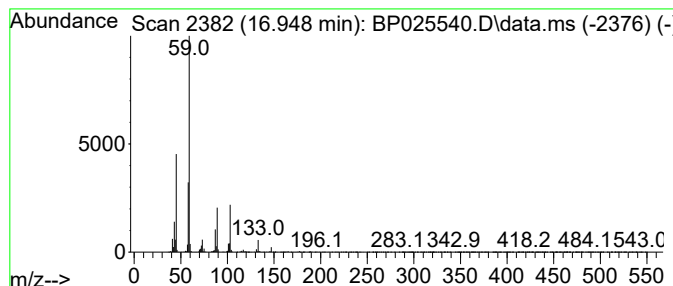
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 5 Methyl 2-(2-(2-methoxyethoxy)ethoxy)acetate Concentration Rank 2

| R.T.      | EstConc                                           | Area   | Relative to ISTD | R.T.         |      |
|-----------|---------------------------------------------------|--------|------------------|--------------|------|
| 16.948    | 7.80 ng                                           | 887922 | Phenanthrene-d10 | 17.207       |      |
| Hit# of 5 | Tentative ID                                      | MW     | MolForm          | CAS#         | Qual |
| 1         | Methyl 2-(2-(2-methoxyethoxy)ethoxy)ethoxyacetate | 192    | C8H16O5          | 1000366-88-2 | 64   |
| 2         | Tetraglyme                                        | 222    | C10H22O5         | 000143-24-8  | 64   |
| 3         | Ethyl 2,5,8,11,14-pentaoxahexadecane-16-oate      | 294    | C13H26O7         | 1000366-80-7 | 59   |
| 4         | 2,5,8,11-Tetraoxadodecane                         | 178    | C8H18O4          | 000112-49-2  | 59   |
| 5         | Hexaethylene glycol dimethyl ether                | 310    | C14H30O7         | 001072-40-8  | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082025\  
Data File : BP025540.D  
Acq On : 21 Aug 2025 14:56  
Operator : CG/JU  
Sample : Q2819-16 2X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
38M-S

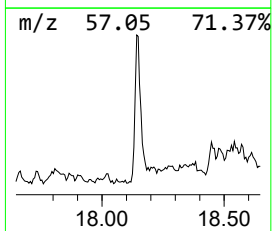
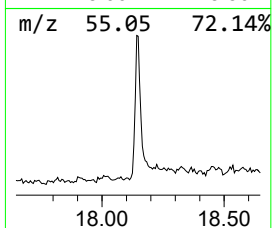
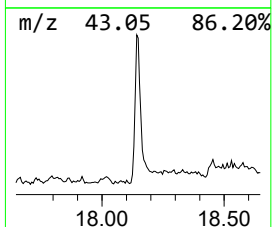
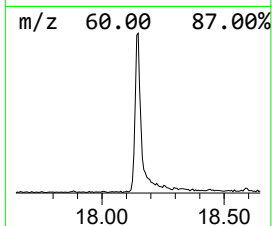
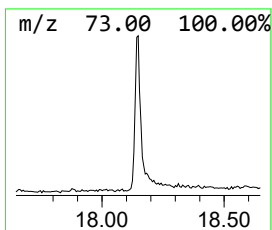
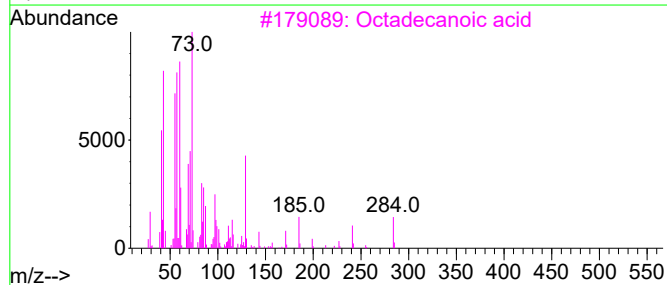
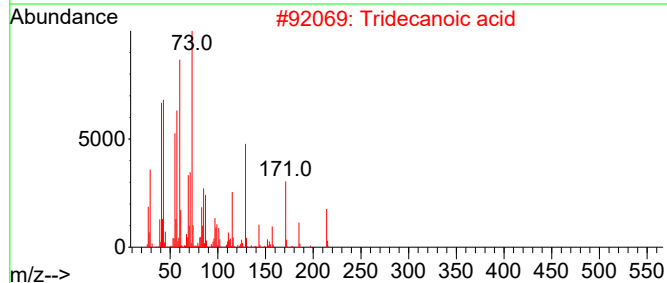
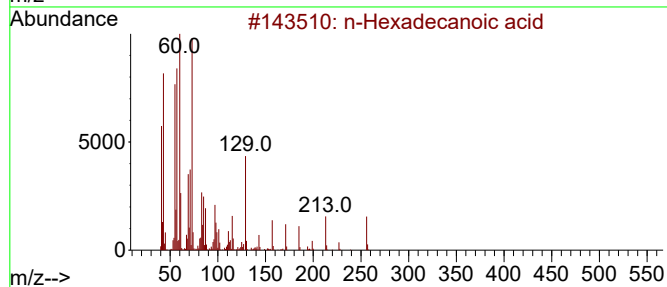
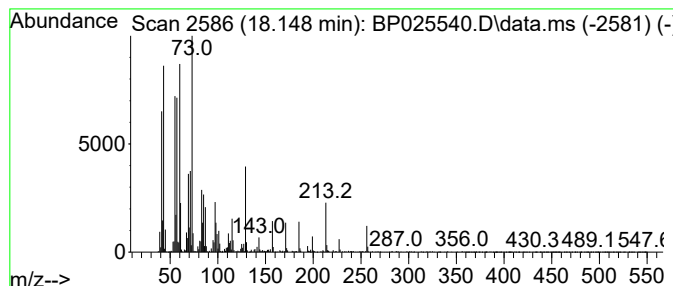
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 n-Hexadecanoic acid Concentration Rank 3

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 18.148 | 6.57 ng | 747628 | Phenanthrene-d10 | 17.207 |

| 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    |    |   | Octadecanoic acid   | 284 | C18H36O2 | 000057-11-4 | 93   |
| 4    |    |   | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 91   |
| 5    |    |   | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 91   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082025\  
Data File : BP025540.D  
Acq On : 21 Aug 2025 14:56  
Operator : CG/JU  
Sample : Q2819-16 2X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
38M-S

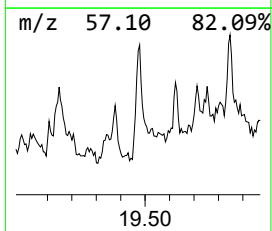
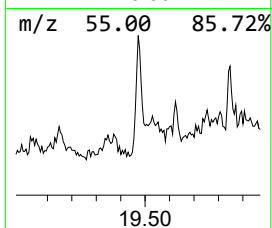
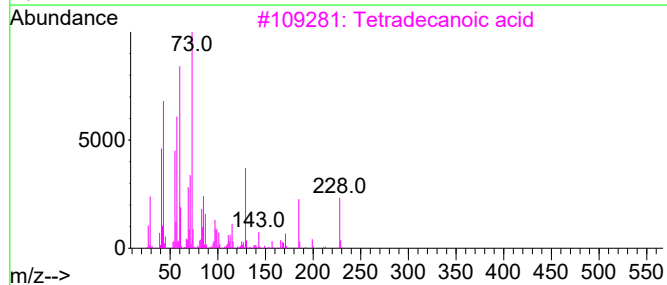
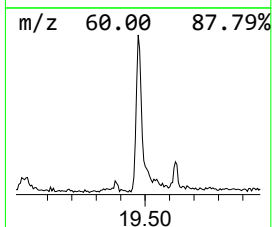
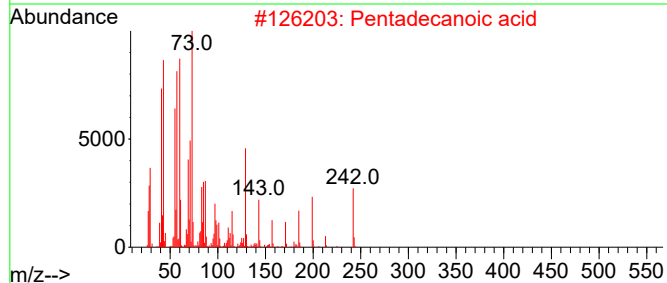
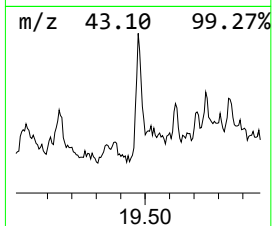
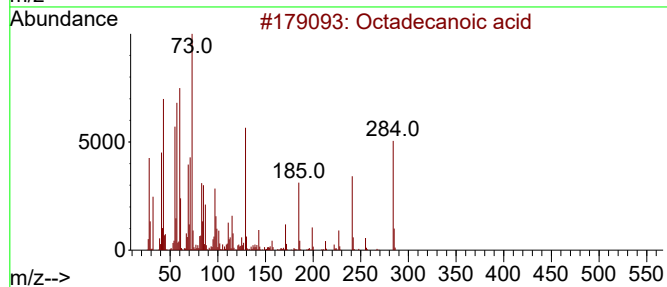
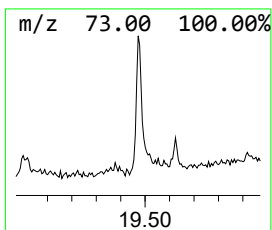
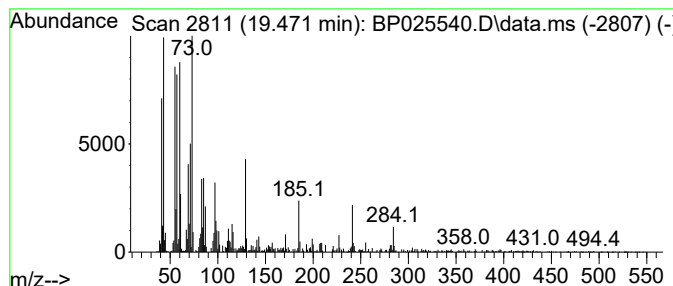
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 Octadecanoic acid Concentration Rank 6

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 19.471 | 3.64 ng | 456637 | Chrysene-d12     | 21.654 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|---------------------|-----|----------|-------------|------|
| 1    |    |   | Octadecanoic acid   | 284 | C18H36O2 | 000057-11-4 | 99   |
| 2    |    |   | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 93   |
| 3    |    |   | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 74   |
| 4    |    |   | n-Decanoic acid     | 172 | C10H20O2 | 000334-48-5 | 70   |
| 5    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082025\  
Data File : BP025540.D  
Acq On : 21 Aug 2025 14:56  
Operator : CG/JU  
Sample : Q2819-16 2X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
38M-S

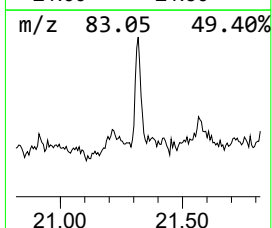
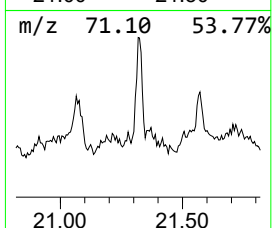
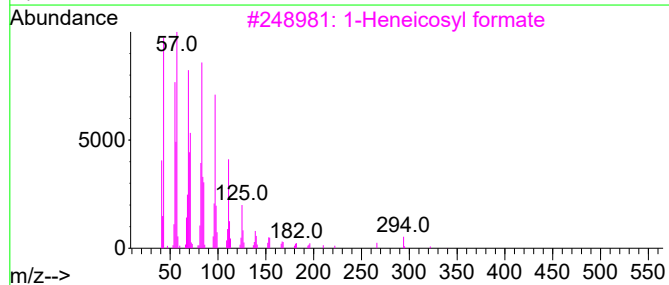
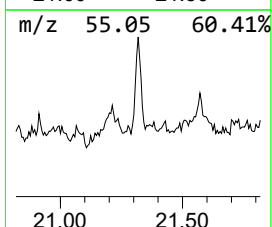
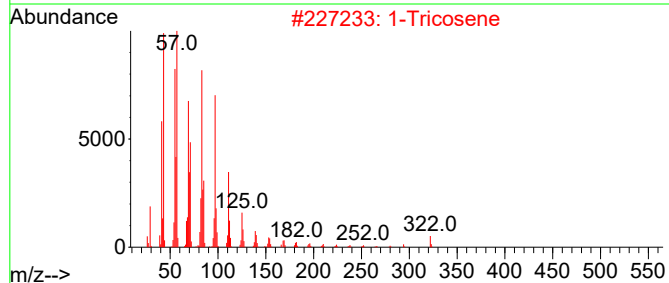
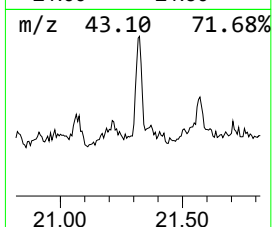
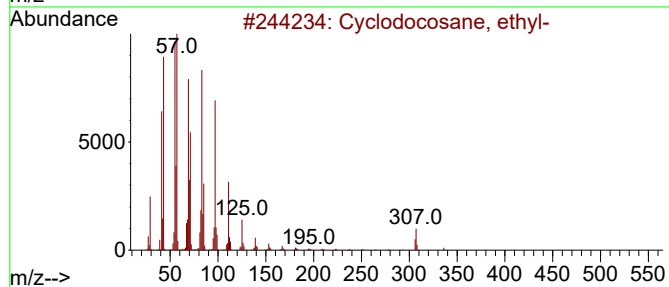
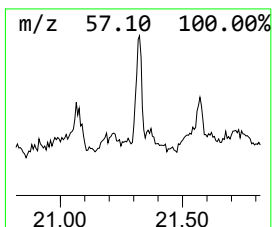
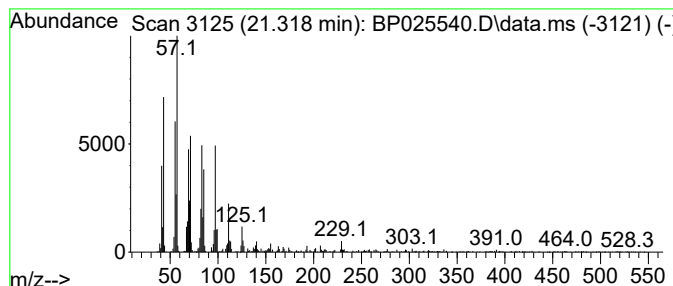
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 Cyclodocosane, ethyl- Concentration Rank 7

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 21.318 | 3.52 ng | 442607 | Chrysene-d12     | 21.654 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Cyclodocosane, ethyl-               | 336 | C24H48     | 1000151-22-6 | 97   |
| 2    |    |   | 1-Tricosene                         | 322 | C23H46     | 018835-32-0  | 92   |
| 3    |    |   | 1-Heneicosyl formate                | 340 | C22H44O2   | 077899-03-7  | 90   |
| 4    |    |   | Oxalic acid, isobutyl hexadecyl ... | 370 | C22H42O4   | 1000309-38-1 | 87   |
| 5    |    |   | Hexacosyl heptafluorobutyrate       | 578 | C30H53F7O2 | 1000351-83-3 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082025\  
Data File : BP025540.D  
Acq On : 21 Aug 2025 14:56  
Operator : CG/JU  
Sample : Q2819-16 2X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
38M-S

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP081425.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

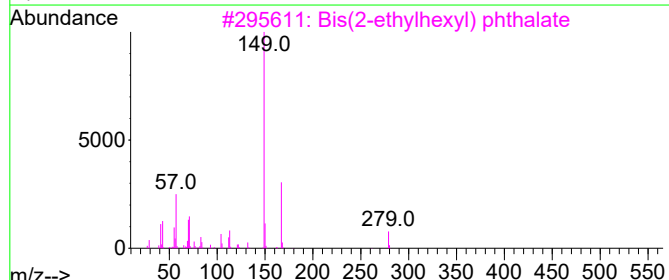
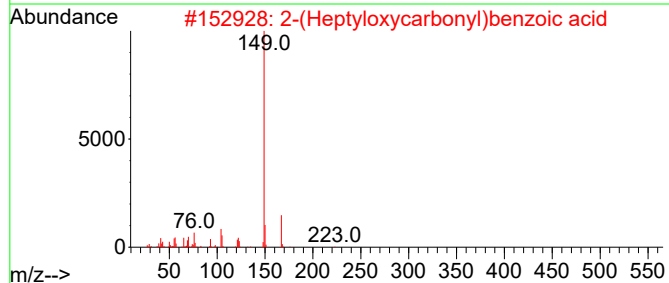
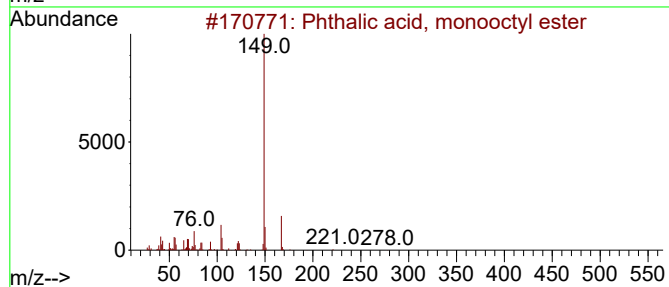
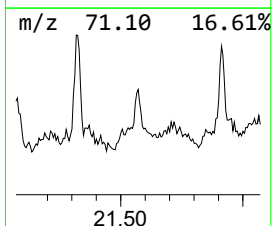
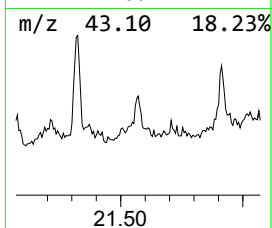
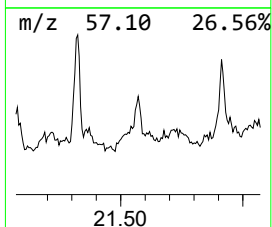
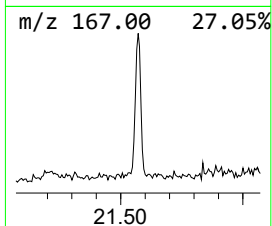
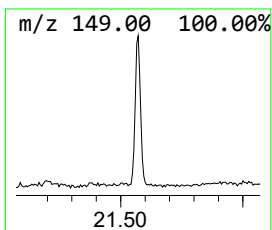
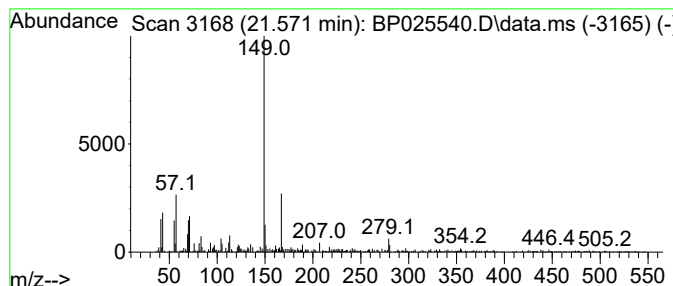
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Phthalic acid, monoethyl ester Concentration Rank 11

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 21.571 | 2.20 ng | 276494 | Chrysene-d12     | 21.654 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Phthalic acid, monoethyl ester      | 278 | C16H22O4 | 005393-19-1  | 81   |
| 2    |    |   | 2-(Heptyloxycarbonyl)benzoic acid   | 264 | C15H20O4 | 024539-58-0  | 74   |
| 3    |    |   | Bis(2-ethylhexyl) phthalate         | 390 | C24H38O4 | 000117-81-7  | 74   |
| 4    |    |   | Phthalic acid, isohexyl 4-methyl... | 334 | C20H30O4 | 1010356-87-5 | 72   |
| 5    |    |   | Phthalic acid, 2,4,4-trimethylpe... | 432 | C27H44O4 | 1000377-77-7 | 68   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082025\  
Data File : BP025540.D  
Acq On : 21 Aug 2025 14:56  
Operator : CG/JU  
Sample : Q2819-16 2X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
38M-S

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP081425.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
|                    |        |         |       |          | #                     | RT     | Resp    | Conc |
| Butane, 2-metho... | 3.037  | 9.4     | ng    | 629629   | 1                     | 7.784  | 1332070 | 20.0 |
| 2-Pentanone, 4-... | 4.943  | 2.8     | ng    | 188648   | 1                     | 7.784  | 1332070 | 20.0 |
| Tetraglyme         | 14.519 | 2.9     | ng    | 307275   | 3                     | 14.413 | 2121850 | 20.0 |
| Benzophenone       | 15.789 | 4.5     | ng    | 478353   | 3                     | 14.413 | 2121850 | 20.0 |
| Methyl 2-(2-(2-... | 16.948 | 7.8     | ng    | 887922   | 4                     | 17.207 | 2276970 | 20.0 |
| n-Hexadecanoic ... | 18.148 | 6.6     | ng    | 747628   | 4                     | 17.207 | 2276970 | 20.0 |
| 2-(2-(2-(2-(2-...  | 18.930 | 6.2     | ng    | 703884   | 4                     | 17.207 | 2276970 | 20.0 |
| Octadecanoic acid  | 19.471 | 3.6     | ng    | 456637   | 5                     | 21.654 | 2512320 | 20.0 |
| Cyclodocosane, ... | 21.318 | 3.5     | ng    | 442607   | 5                     | 21.654 | 2512320 | 20.0 |
| Phthalic acid, ... | 21.571 | 2.2     | ng    | 276494   | 5                     | 21.654 | 2512320 | 20.0 |