

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080521\  
 Data File : BP006523.D  
 Acq On : 06 Aug 2021 00:54  
 Operator : CG/JU  
 Sample : M3281-05  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 TWP02

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-BP080321.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP006523.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         | 3.546       | 74            | 78          | 85           | rBV2     | 183773         | 250048        | 1.61%           | 0.288%        |
| 2         | 4.287       | 199           | 204         | 214          | rBV      | 638226         | 913869        | 5.89%           | 1.054%        |
| 3         | 5.357       | 380           | 386         | 393          | rBV      | 1834359        | 2591776       | 16.70%          | 2.989%        |
| 4         | 6.875       | 640           | 644         | 648          | rBV      | 125323         | 199225        | 1.28%           | 0.230%        |
| 5         | 6.934       | 648           | 654         | 661          | rVB      | 1188756        | 1820006       | 11.73%          | 2.099%        |
| 6         | 7.757       | 787           | 794         | 800          | rVB      | 960002         | 1510987       | 9.73%           | 1.742%        |
| 7         | 8.075       | 837           | 848         | 853          | rBV      | 4274654        | 6948854       | 44.77%          | 8.013%        |
| 8         | 8.122       | 853           | 856         | 865          | rVB      | 1129720        | 1664862       | 10.73%          | 1.920%        |
| 9         | 8.916       | 983           | 991         | 1001         | rVB      | 3558275        | 5880407       | 37.89%          | 6.781%        |
| 10        | 9.722       | 1123          | 1128        | 1134         | rBV      | 168553         | 275714        | 1.78%           | 0.318%        |
| 11        | 10.328      | 1220          | 1231        | 1246         | rBV      | 835044         | 1310221       | 8.44%           | 1.511%        |
| 12        | 10.540      | 1259          | 1267        | 1270         | rBV      | 1271436        | 2125663       | 13.70%          | 2.451%        |
| 13        | 10.592      | 1270          | 1276        | 1283         | rVB      | 4944073        | 8457973       | 54.49%          | 9.753%        |
| 14        | 10.704      | 1288          | 1295        | 1304         | rVB2     | 272440         | 532253        | 3.43%           | 0.614%        |
| 15        | 12.422      | 1578          | 1587        | 1598         | rVB      | 239290         | 425224        | 2.74%           | 0.490%        |
| 16        | 13.010      | 1679          | 1687        | 1694         | rBV      | 7927744        | 11678950      | 75.24%          | 13.467%       |
| 17        | 13.222      | 1718          | 1723        | 1730         | rVB3     | 121691         | 190571        | 1.23%           | 0.220%        |
| 18        | 14.381      | 1914          | 1920        | 1926         | rBV      | 1819532        | 2619323       | 16.88%          | 3.020%        |
| 19        | 14.445      | 1926          | 1931        | 1938         | rVV      | 1052452        | 1557347       | 10.03%          | 1.796%        |
| 20        | 14.510      | 1938          | 1942        | 1947         | rVV3     | 92764          | 158673        | 1.02%           | 0.183%        |
| 21        | 14.575      | 1947          | 1953        | 1963         | rVB      | 169206         | 286309        | 1.84%           | 0.330%        |
| 22        | 14.781      | 1983          | 1988        | 1998         | rVB      | 370277         | 554807        | 3.57%           | 0.640%        |
| 23        | 14.875      | 1998          | 2004        | 2010         | rBV      | 153183         | 230158        | 1.48%           | 0.265%        |
| 24        | 15.433      | 2091          | 2099        | 2104         | rVB      | 445073         | 678536        | 4.37%           | 0.782%        |
| 25        | 15.575      | 2117          | 2123        | 2131         | rBV5     | 137520         | 326865        | 2.11%           | 0.377%        |
| 26        | 15.733      | 2146          | 2150        | 2159         | rVB4     | 81795          | 180030        | 1.16%           | 0.208%        |
| 27        | 15.875      | 2168          | 2174        | 2181         | rBV      | 3971324        | 5458111       | 35.17%          | 6.294%        |
| 28        | 16.098      | 2209          | 2212        | 2219         | rVB4     | 102400         | 201592        | 1.30%           | 0.232%        |
| 29        | 16.680      | 2308          | 2311        | 2319         | rVB4     | 113877         | 192752        | 1.24%           | 0.222%        |
| 30        | 17.133      | 2382          | 2388        | 2393         | rBV      | 2023271        | 2882531       | 18.57%          | 3.324%        |
| 31        | 17.174      | 2393          | 2395        | 2400         | rVB      | 199623         | 235983        | 1.52%           | 0.272%        |
| 32        | 17.257      | 2405          | 2409        | 2417         | rVB2     | 112067         | 241882        | 1.56%           | 0.279%        |
| 33        | 17.545      | 2453          | 2458        | 2463         | rBV      | 205393         | 296978        | 1.91%           | 0.342%        |
| 34        | 17.610      | 2465          | 2469        | 2480         | rVB2     | 124926         | 226552        | 1.46%           | 0.261%        |
| 35        | 17.822      | 2502          | 2505        | 2514         | rVB      | 165072         | 233404        | 1.50%           | 0.269%        |
| 36        | 18.045      | 2538          | 2543        | 2552         | rBV      | 571605         | 836825        | 5.39%           | 0.965%        |

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Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TWP02

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-BP080321.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |          |          |         |         |
|----|--------|------|------|------|------|----------|----------|---------|---------|
| 37 | 19.151 | 2727 | 2731 | 2737 | rVB  | 141546   | 195158   | 1.26%   | 0.225%  |
| 38 | 19.280 | 2749 | 2753 | 2761 | rBV  | 428239   | 573312   | 3.69%   | 0.661%  |
| 39 | 19.710 | 2820 | 2826 | 2832 | rBV  | 13109923 | 15521286 | 100.00% | 17.897% |
| 40 | 20.957 | 3035 | 3038 | 3044 | rBV2 | 207129   | 244382   | 1.57%   | 0.282%  |
| 41 | 21.233 | 3080 | 3085 | 3091 | rBV  | 2361065  | 2954324  | 19.03%  | 3.407%  |
| 42 | 23.557 | 3474 | 3480 | 3489 | rVB  | 1583996  | 3060583  | 19.72%  | 3.529%  |

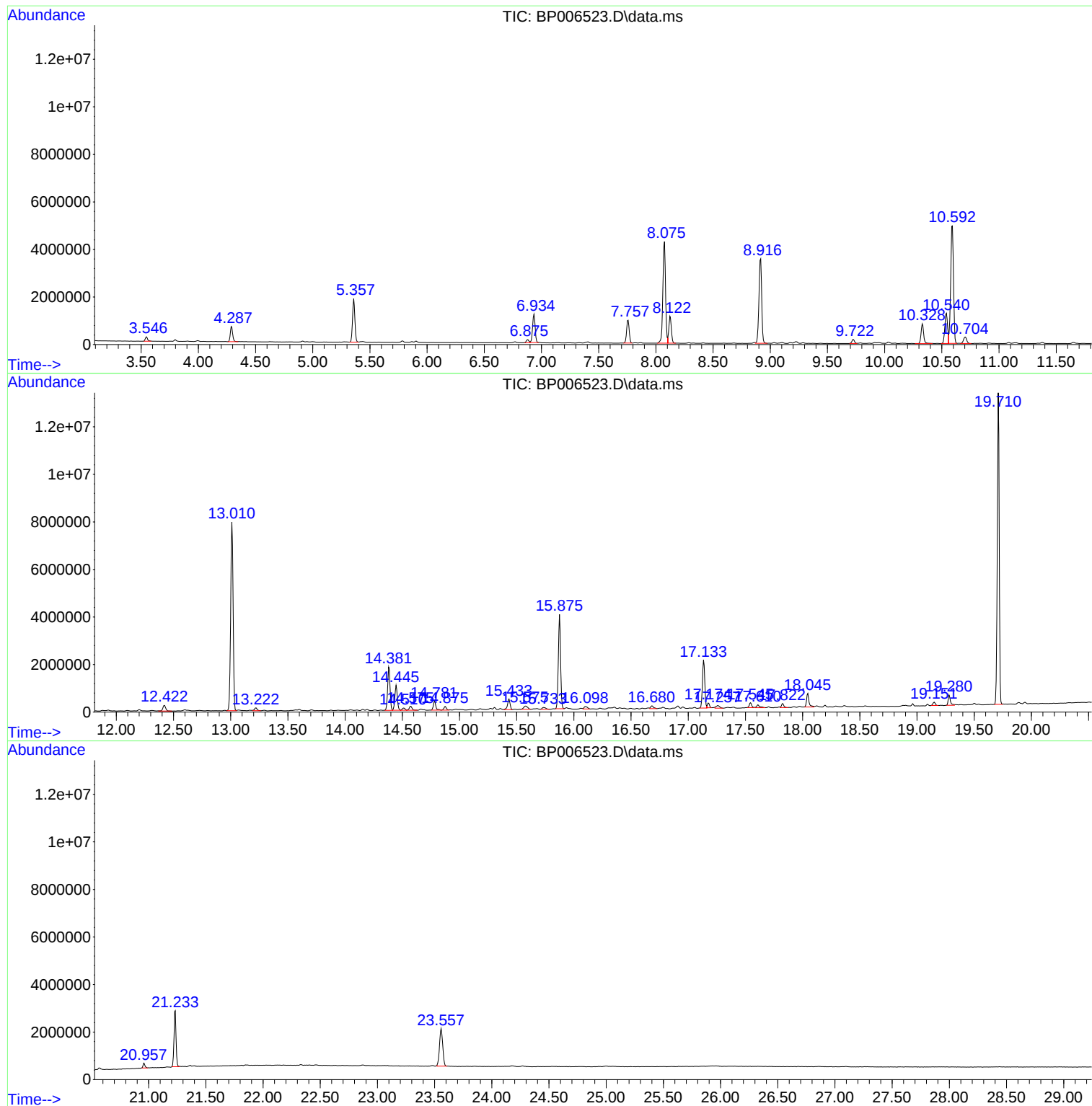
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Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080521\  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP080321.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080521\  
Data File : BP006523.D  
Acq On : 06 Aug 2021 00:54  
Operator : CG/JU  
Sample : M3281-05  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TWP02

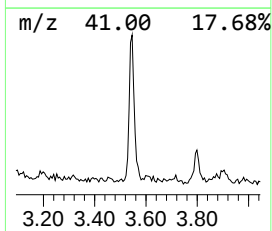
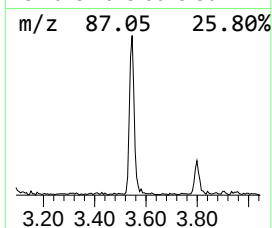
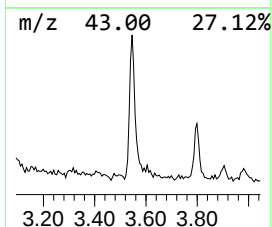
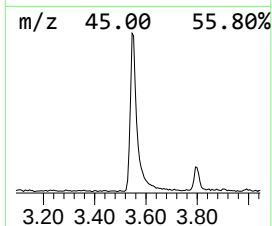
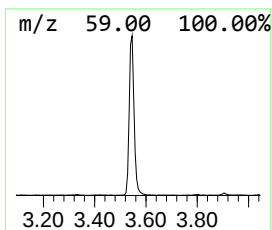
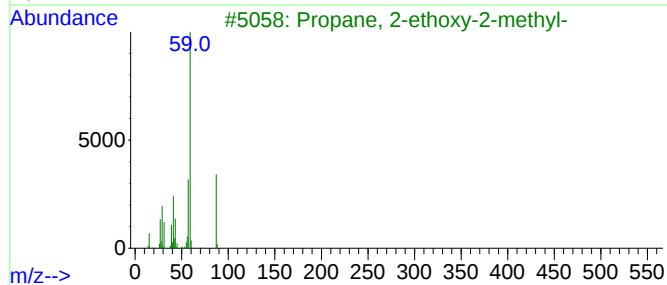
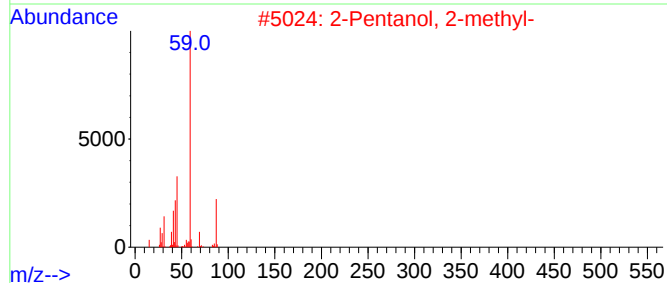
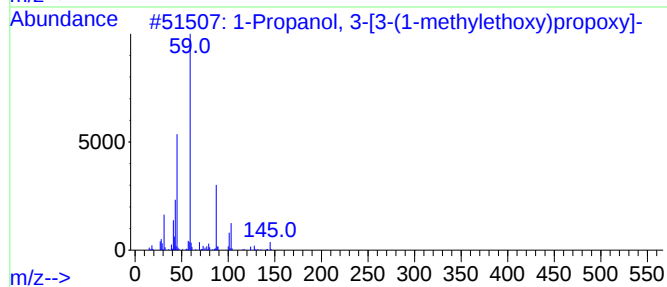
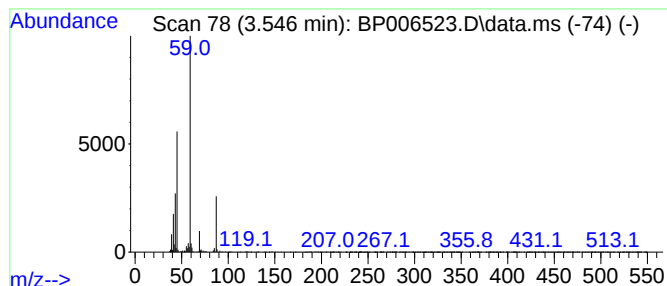
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 1-Propanol, 3-[3-(1-methyle... Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 3.546 | 3.31 ng | 250048 | 1,4-Dichlorobenzene-d4 | 7.757 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1-Propanol, 3-[3-(1-methylethoxy... | 176 | C9H20O3      | 054518-03-5 | 78      |      |      |
| 2    | 2-Pentanol, 2-methyl-               | 102 | C6H14O       | 000590-36-3 | 72      |      |      |
| 3    | Propane, 2-ethoxy-2-methyl-         | 102 | C6H14O       | 000637-92-3 | 50      |      |      |
| 4    | 2-Butanol, 2,3-dimethyl-            | 102 | C6H14O       | 000594-60-5 | 50      |      |      |
| 5    | 2-Butanone, 3-ethoxy-3-methyl-      | 130 | C7H14O2      | 036687-99-7 | 47      |      |      |



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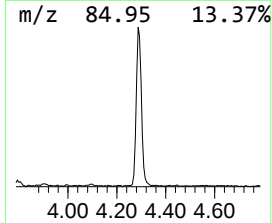
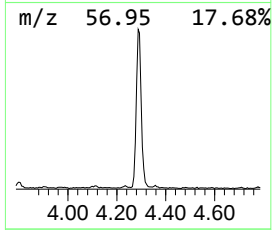
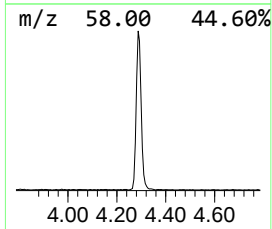
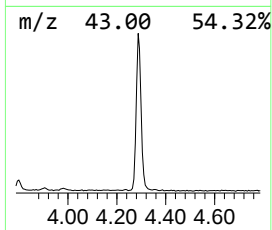
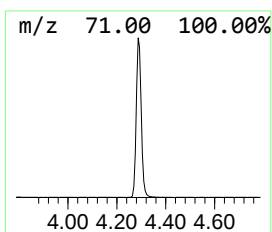
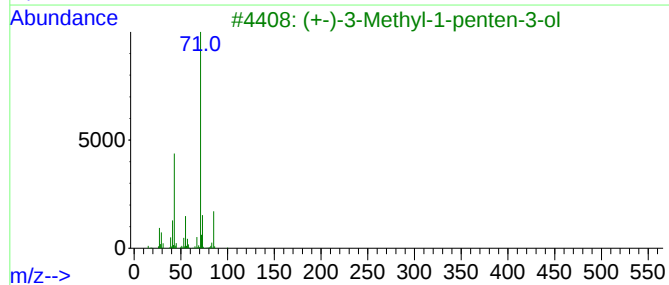
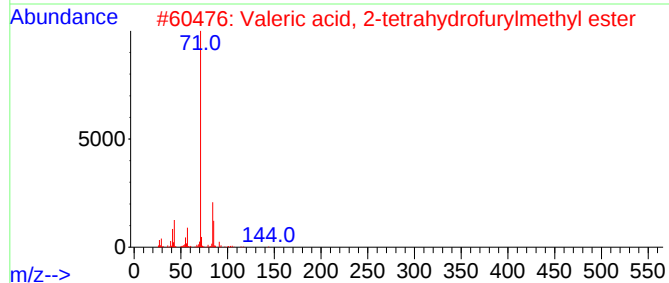
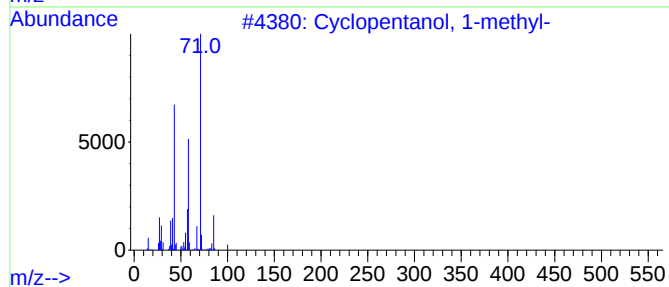
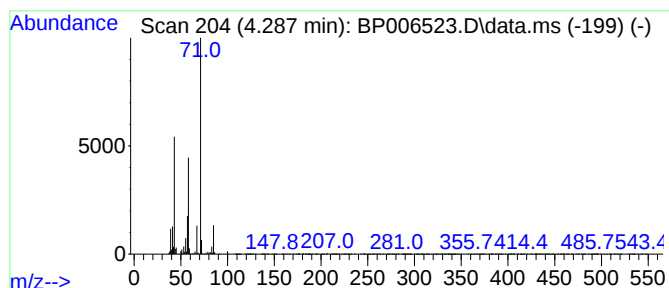
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP080321.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 2 Cyclopentanol, 1-methyl- Concentration Rank 4

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 4.287 | 12.10 ng | 913869 | 1,4-Dichlorobenzene-d4 | 7.757 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Cyclopentanol, 1-methyl-            | 100 | C6H12O     | 001462-03-9  | 91   |
| 2    |    |   | Valeric acid, 2-tetrahydrofurylm... | 186 | C10H18O3   | 005451-86-5  | 50   |
| 3    |    |   | (+)-3-Methyl-1-penten-3-ol          | 100 | C6H12O     | 1010238-75-4 | 40   |
| 4    |    |   | 1-Penten-3-ol, 2-methyl-            | 100 | C6H12O     | 002088-07-5  | 36   |
| 5    |    |   | Diethylmalonic acid, monochlorid... | 262 | C12H19ClO4 | 1000370-65-0 | 28   |



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Instrument :  
BNA\_P  
ClientSampleId :  
TWP02

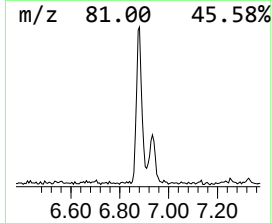
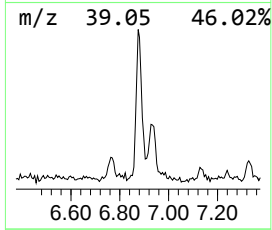
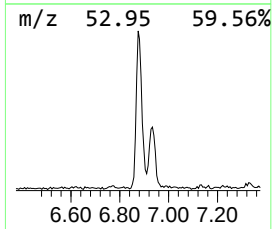
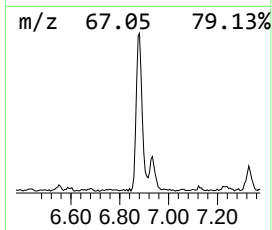
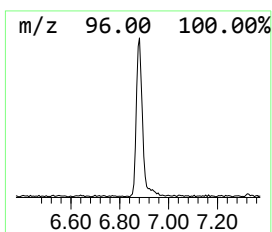
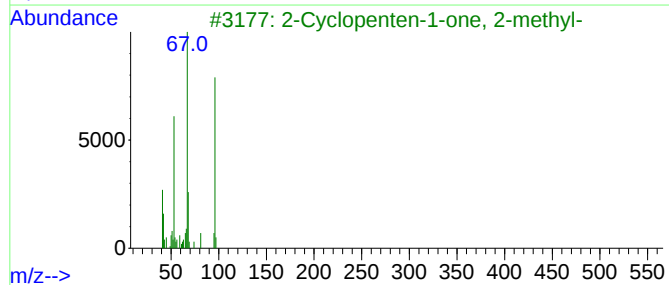
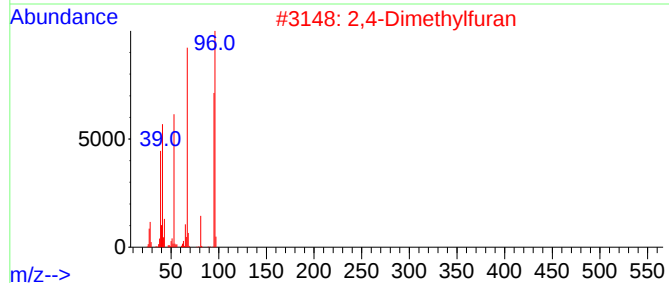
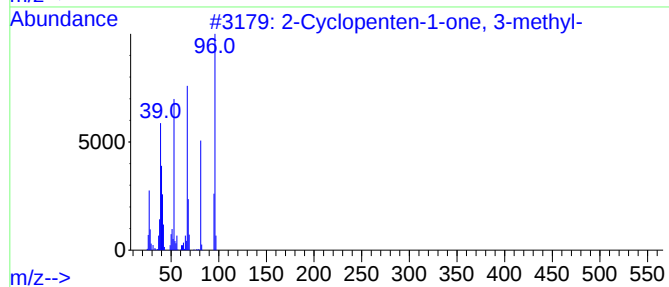
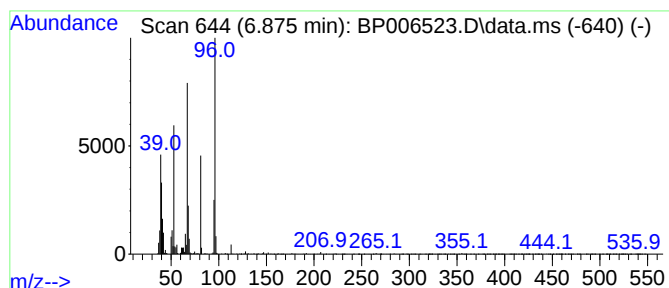
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP080321.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 3 2-Cyclopenten-1-one, 3-methyl- Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 6.875 | 2.64 ng | 199225 | 1,4-Dichlorobenzene-d4 | 7.757 |

| Hit# | of | 5 | Tentative ID                    | MW | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------|----|---------|-------------|------|
| 1    |    |   | 2-Cyclopenten-1-one, 3-methyl-  | 96 | C6H8O   | 002758-18-1 | 96   |
| 2    |    |   | 2,4-Dimethylfuran               | 96 | C6H8O   | 003710-43-8 | 87   |
| 3    |    |   | 2-Cyclopenten-1-one, 2-methyl-  | 96 | C6H8O   | 001120-73-6 | 87   |
| 4    |    |   | Cyclopentene                    | 68 | C5H8    | 000142-29-0 | 60   |
| 5    |    |   | 4-Vinyl-4,5-dihydro-3H-pyrazole | 96 | C5H8N2  | 063438-33-5 | 53   |



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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP080321.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

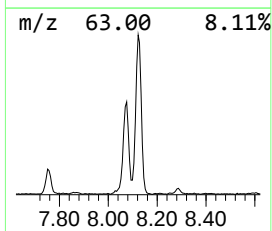
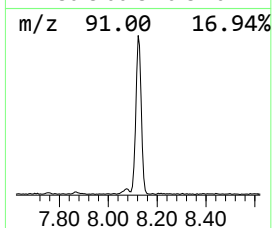
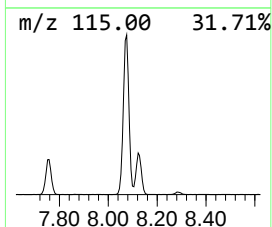
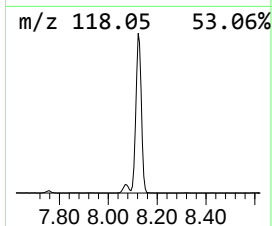
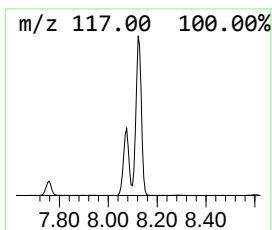
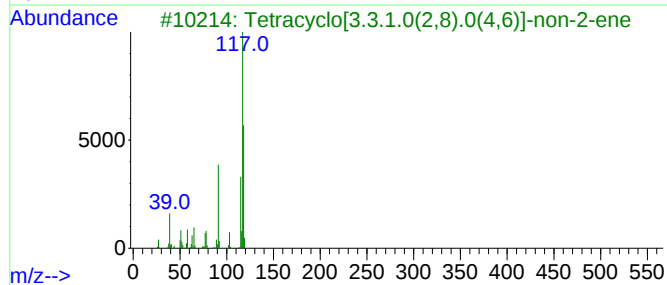
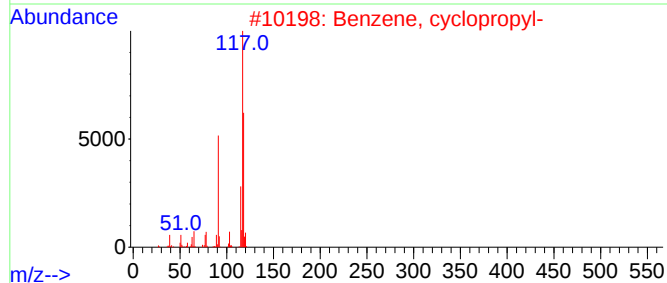
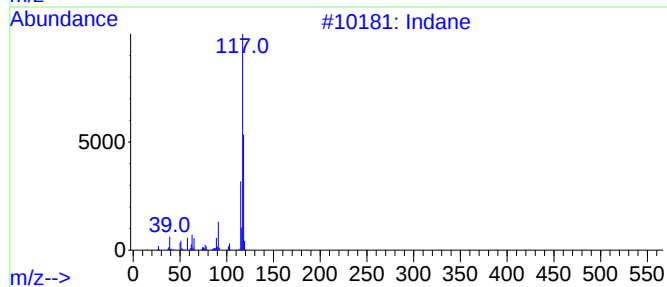
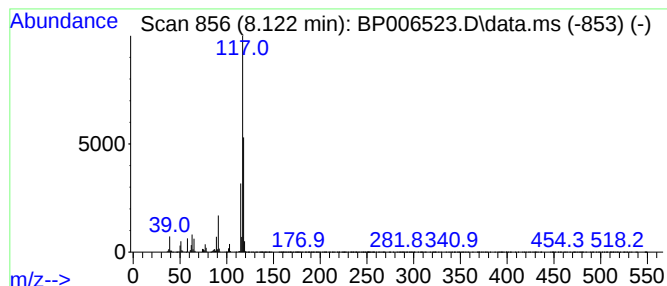
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Indane Concentration Rank 2

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 8.122 | 22.04 ng | 1664860 | 1,4-Dichlorobenzene-d4 | 7.757 |

| Hit# of 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|-----------|-------------------------------------|-----|---------|--------------|------|
| 1         | Indane                              | 118 | C9H10   | 000496-11-7  | 94   |
| 2         | Benzene, cyclopropyl-               | 118 | C9H10   | 000873-49-4  | 83   |
| 3         | Tetracyclo[3.3.1.0(2,8).0(4,6)]-... | 118 | C9H10   | 1000191-13-7 | 80   |
| 4         | Deltacyclene                        | 118 | C9H10   | 007785-10-6  | 74   |
| 5         | Benzene, 1-propenyl-                | 118 | C9H10   | 000637-50-3  | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080521\  
Data File : BP006523.D  
Acq On : 06 Aug 2021 00:54  
Operator : CG/JU  
Sample : M3281-05  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TWP02

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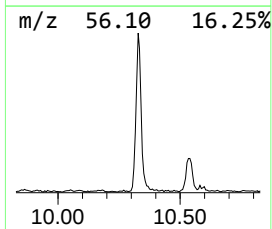
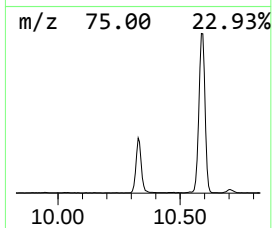
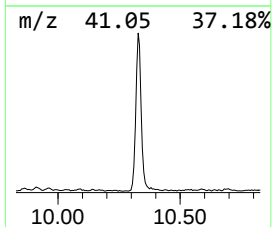
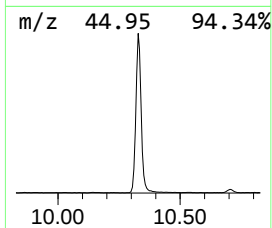
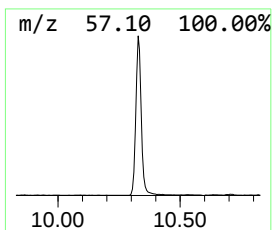
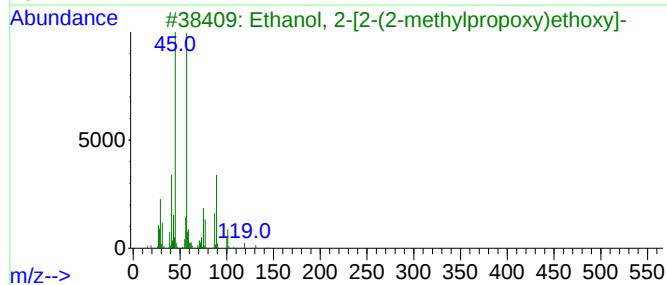
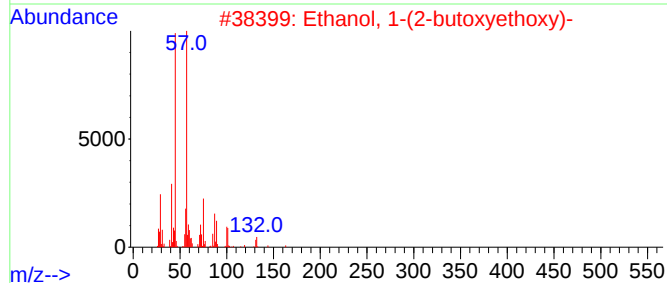
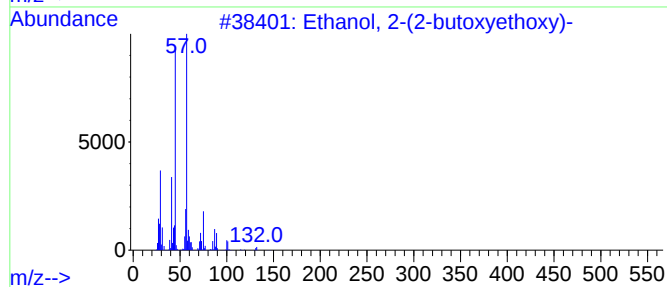
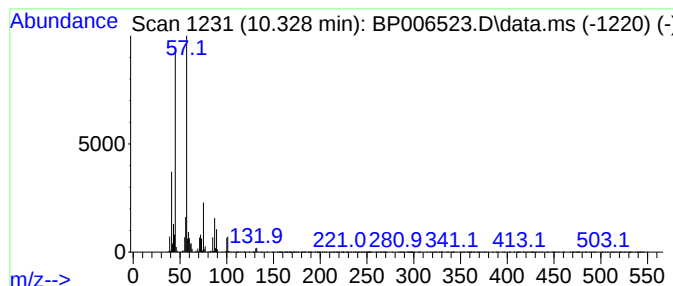
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 3

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 10.328 | 12.33 ng | 1310220 | Naphthalene-d8   | 10.540 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Ethanol, 2-(2-butoxyethoxy)-        | 162 | C8H18O3 | 000112-34-5 | 90   |
| 2    |    |   | Ethanol, 1-(2-butoxyethoxy)-        | 162 | C8H18O3 | 054446-78-5 | 90   |
| 3    |    |   | Ethanol, 2-[2-(2-methylpropoxy)e... | 162 | C8H18O3 | 018912-80-6 | 72   |
| 4    |    |   | Ethylene glycol monoisobutyl ether  | 118 | C6H14O2 | 004439-24-1 | 64   |
| 5    |    |   | Di-sec-Butyl ether                  | 130 | C8H18O  | 006863-58-7 | 50   |



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Data File : BP006523.D  
Acq On : 06 Aug 2021 00:54  
Operator : CG/JU  
Sample : M3281-05  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TWP02

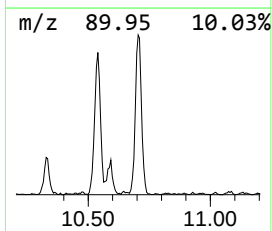
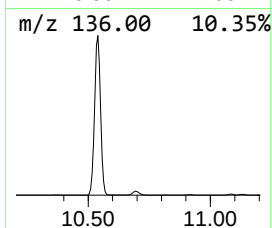
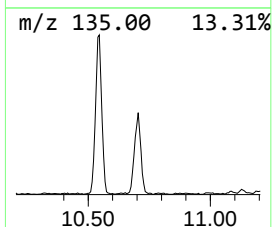
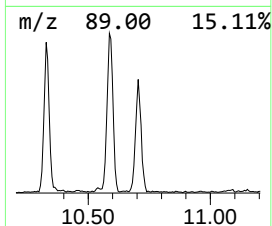
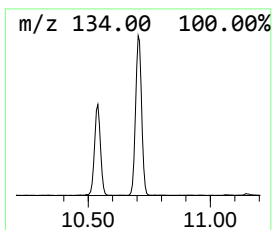
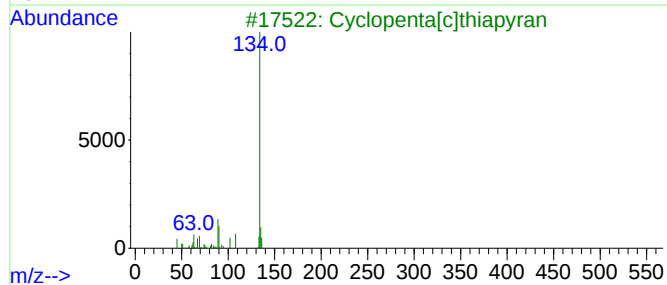
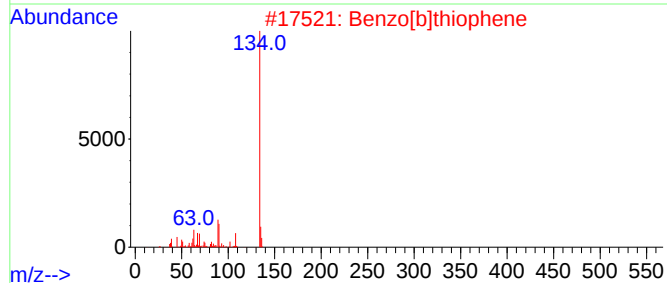
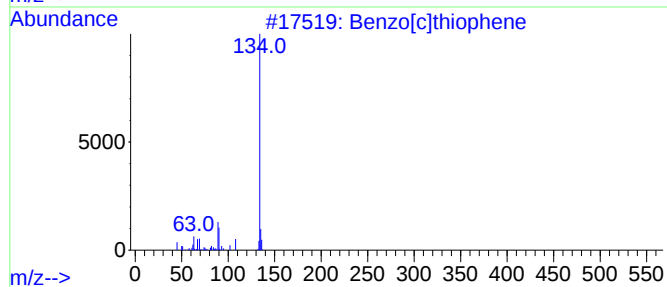
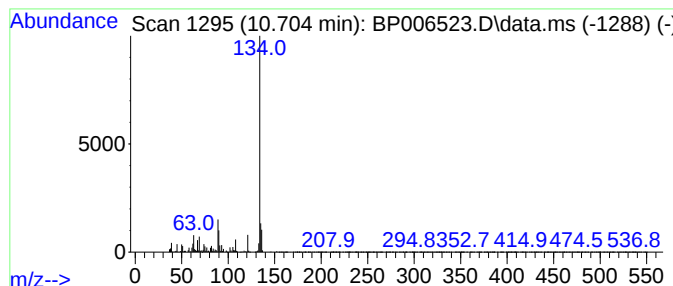
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 Benzo[c]thiophene Concentration Rank 6

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 10.704 | 5.01 ng | 532253 | Naphthalene-d8   | 10.540 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|------------|-------------|------|
| 1    |    |   | Benzo[c]thiophene                   | 134 | C8H6S      | 000270-82-6 | 94   |
| 2    |    |   | Benzo[b]thiophene                   | 134 | C8H6S      | 000095-15-8 | 93   |
| 3    |    |   | Cyclopenta[c]thiapyran              | 134 | C8H6S      | 000270-63-3 | 90   |
| 4    |    |   | Thiazolidin-4-one, 5-benzylideno... | 287 | C13H9N3OS2 | 351062-07-2 | 72   |
| 5    |    |   | [1]Benzothieno[2',3':3,4]cyclobu... | 246 | C13H10O3S  | 034002-19-2 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080521\  
Data File : BP006523.D  
Acq On : 06 Aug 2021 00:54  
Operator : CG/JU  
Sample : M3281-05  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TWP02

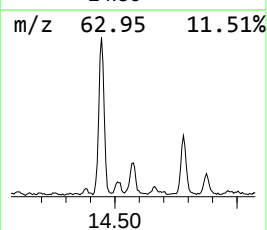
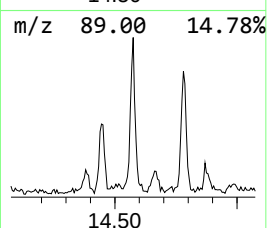
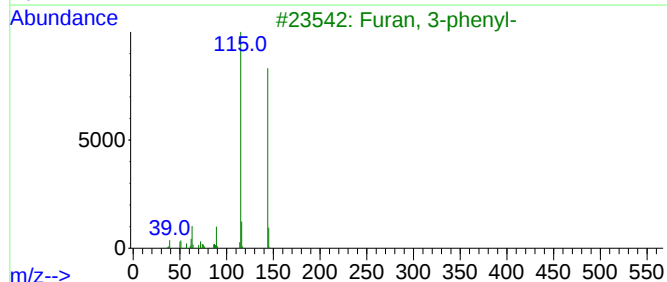
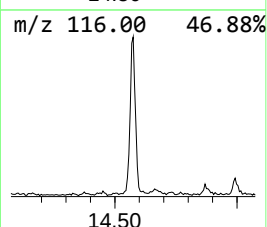
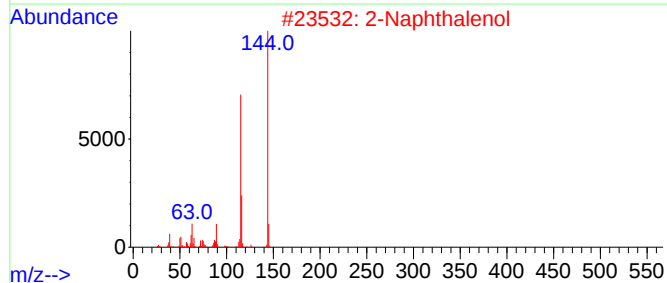
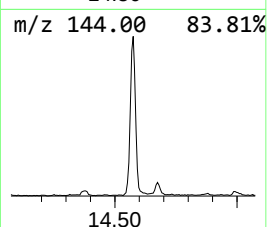
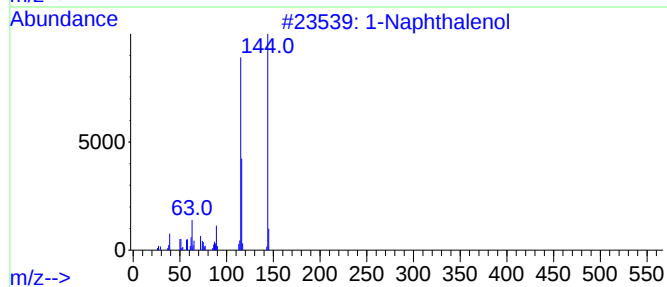
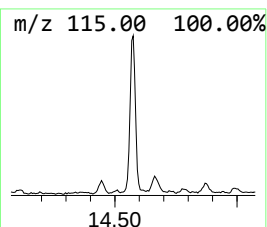
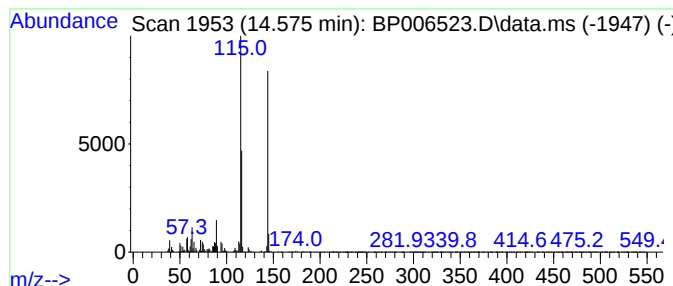
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 1-Naphthalenol Concentration Rank 11

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 14.575 | 2.19 ng | 286309 | Acenaphthene-d10 | 14.381 |

| Hit# | of 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|------|------------------------|-----|---------|-------------|------|
| 1    |      | 1-Naphthalenol         | 144 | C10H8O  | 000090-15-3 | 96   |
| 2    |      | 2-Naphthalenol         | 144 | C10H8O  | 000135-19-3 | 90   |
| 3    |      | Furan, 3-phenyl-       | 144 | C10H8O  | 013679-41-9 | 81   |
| 4    |      | Benzofuran, 2-ethenyl- | 144 | C10H8O  | 007522-79-4 | 81   |
| 5    |      | Cinnoline, 3-methyl-   | 144 | C9H8N2  | 017372-78-0 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080521\  
Data File : BP006523.D  
Acq On : 06 Aug 2021 00:54  
Operator : CG/JU  
Sample : M3281-05  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TWP02

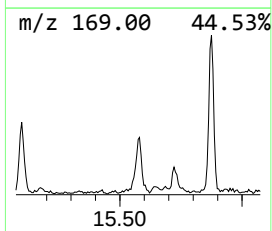
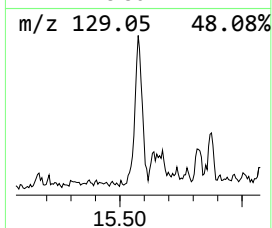
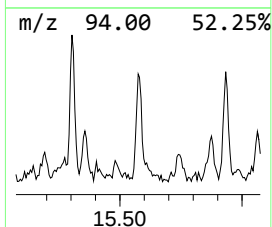
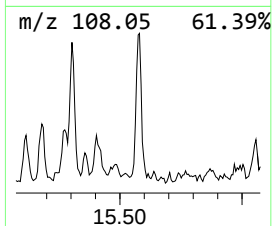
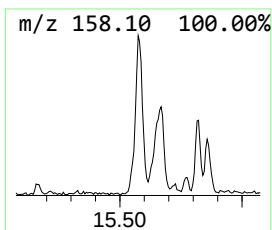
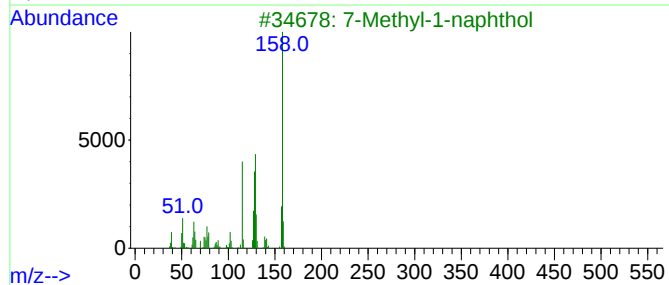
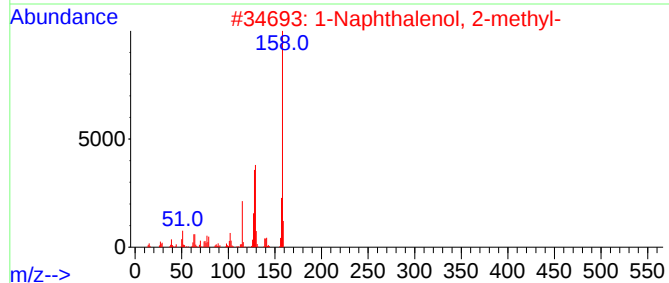
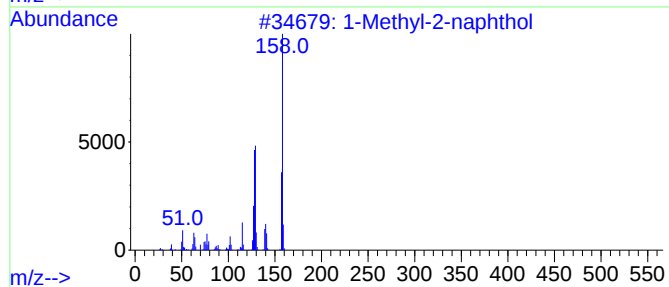
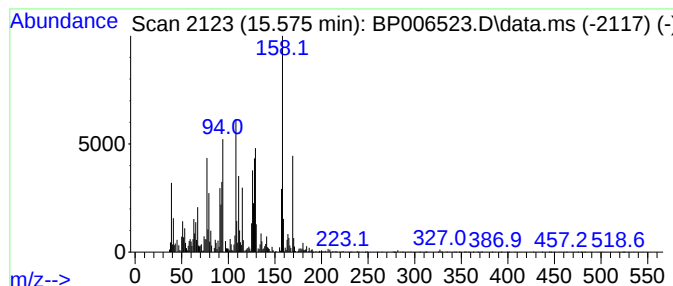
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 1-Methyl-2-naphthol Concentration Rank 10

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 15.575 | 2.50 ng | 326865 | Acenaphthene-d10 | 14.381 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm  | CAS#        | Qual |
|------|----|---|------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 1-Methyl-2-naphthol                | 158 | C11H10O  | 001076-26-2 | 55   |
| 2    |    |   | 1-Naphthalenol, 2-methyl-          | 158 | C11H10O  | 007469-77-4 | 50   |
| 3    |    |   | 7-Methyl-1-naphthol                | 158 | C11H10O  | 006939-33-9 | 49   |
| 4    |    |   | Cinnoline, 3,4-dimethyl-           | 158 | C10H10N2 | 003929-83-7 | 30   |
| 5    |    |   | Benzene, 1,3-bis(1-methylethenyl)- | 158 | C12H14   | 003748-13-8 | 25   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080521\  
Data File : BP006523.D  
Acq On : 06 Aug 2021 00:54  
Operator : CG/JU  
Sample : M3281-05  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TWP02

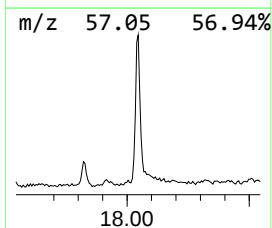
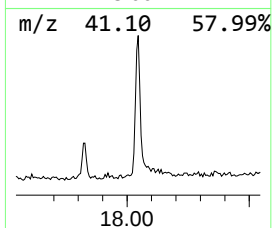
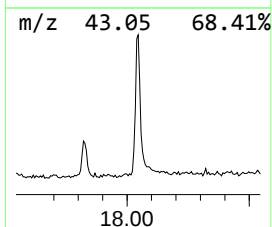
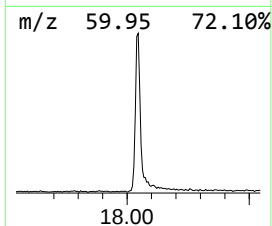
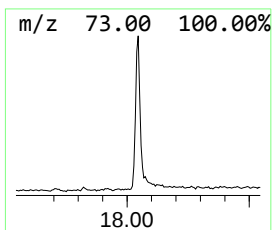
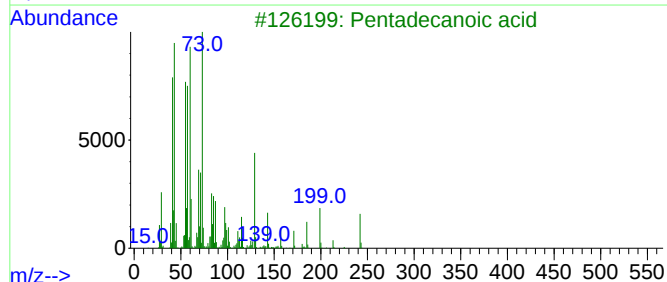
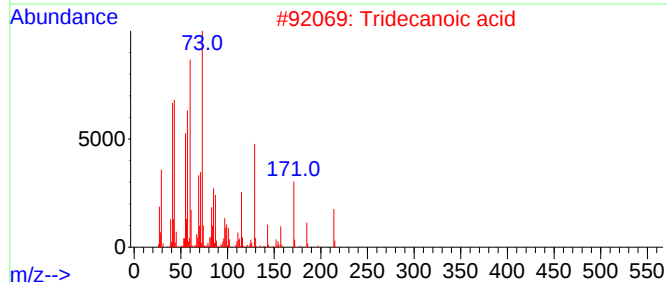
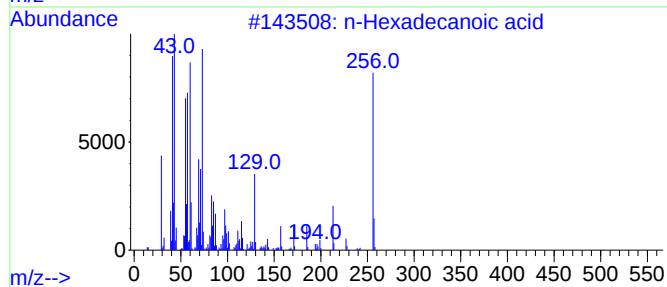
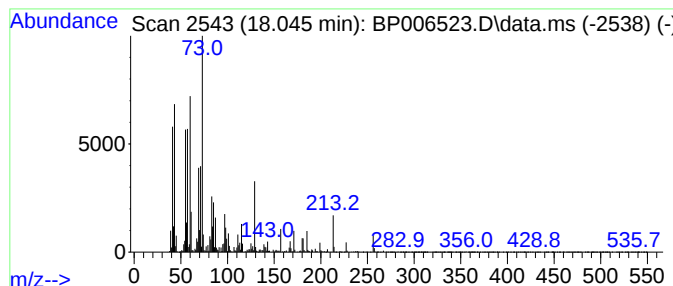
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 n-Hexadecanoic acid Concentration Rank 5

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 18.045 | 5.81 ng | 836825 | Phenanthrene-d10 | 17.133 |

| 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 | 92   |
| 3    |    |   | Pentadecanoic acid       | 242 | C15H30O2 | 001002-84-2 | 58   |
| 4    |    |   | 8-Methylnonanoic acid    | 172 | C10H20O2 | 005963-14-4 | 38   |
| 5    |    |   | Ethanol, 2-(dodecyloxy)- | 230 | C14H30O2 | 004536-30-5 | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080521\  
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Acq On : 06 Aug 2021 00:54  
Operator : CG/JU  
Sample : M3281-05  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TWP02

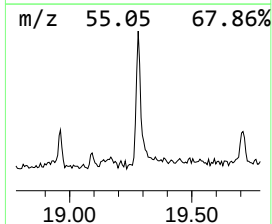
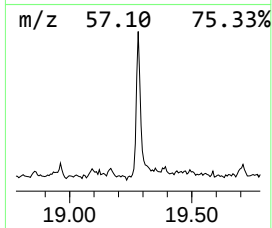
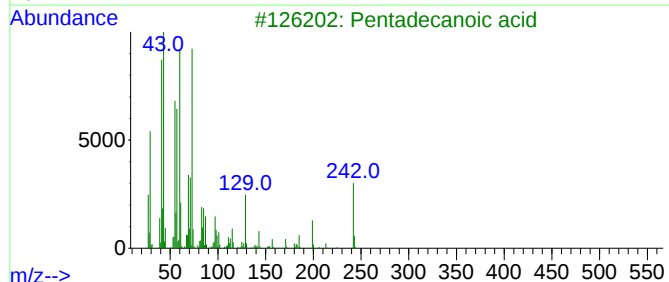
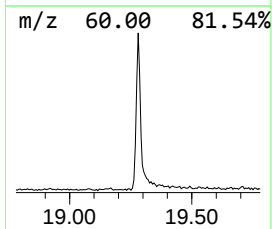
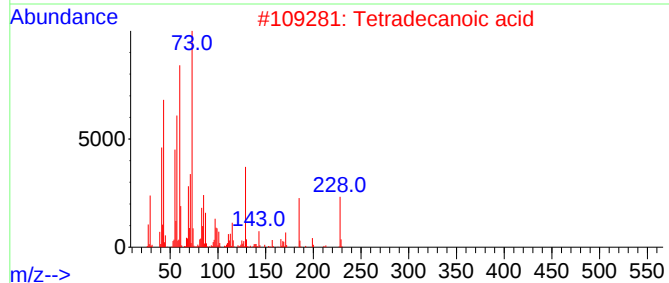
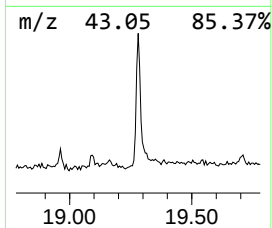
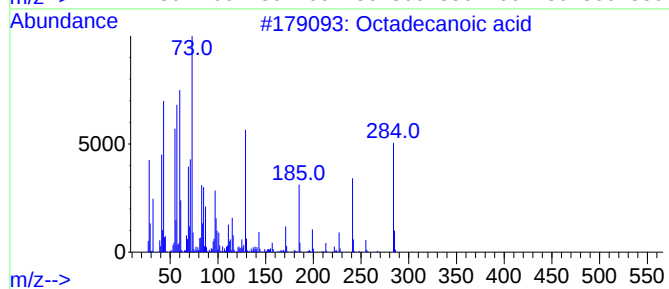
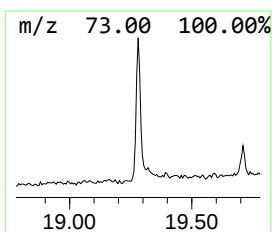
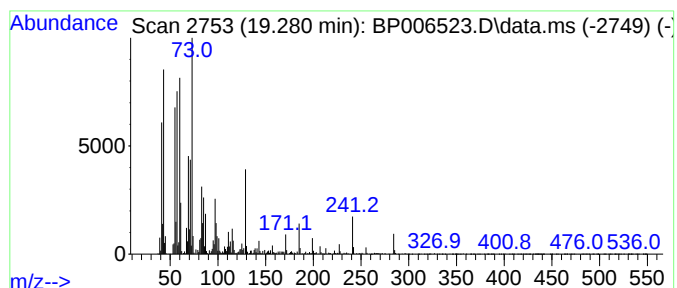
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 11 Octadecanoic acid Concentration Rank 7

| R.T.      | EstConc             | Area   | Relative to ISTD | R.T.        |      |
|-----------|---------------------|--------|------------------|-------------|------|
| 19.280    | 3.88 ng             | 573312 | Chrysene-d12     | 21.233      |      |
| Hit# of 5 | Tentative ID        | MW     | MolForm          | CAS#        | Qual |
| 1         | Octadecanoic acid   | 284    | C18H36O2         | 000057-11-4 | 99   |
| 2         | Tetradecanoic acid  | 228    | C14H28O2         | 000544-63-8 | 93   |
| 3         | Pentadecanoic acid  | 242    | C15H30O2         | 001002-84-2 | 90   |
| 4         | n-Hexadecanoic acid | 256    | C16H32O2         | 000057-10-3 | 74   |
| 5         | Tridecanoic acid    | 214    | C13H26O2         | 000638-53-9 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080521\  
Data File : BP006523.D  
Acq On : 06 Aug 2021 00:54  
Operator : CG/JU  
Sample : M3281-05  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TWP02

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP080321.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 |
| 1-Propanol, 3-[...] | 3.546  | 3.3     | ng    | 250048   | 1                     | 7.757  | 1510990 | 20.0 |
| Cyclopentanol, ...  | 4.287  | 12.1    | ng    | 913869   | 1                     | 7.757  | 1510990 | 20.0 |
| 2-Cyclopenten-1...  | 6.875  | 2.6     | ng    | 199225   | 1                     | 7.757  | 1510990 | 20.0 |
| Indane              | 8.122  | 22.0    | ng    | 1664860  | 1                     | 7.757  | 1510990 | 20.0 |
| Ethanol, 2-(2-b...  | 10.328 | 12.3    | ng    | 1310220  | 2                     | 10.540 | 2125660 | 20.0 |
| Benzo[c]thiophene   | 10.704 | 5.0     | ng    | 532253   | 2                     | 10.540 | 2125660 | 20.0 |
| 1-Naphthalenol      | 14.575 | 2.2     | ng    | 286309   | 3                     | 14.381 | 2619320 | 20.0 |
| 1-Methyl-2-naph...  | 15.575 | 2.5     | ng    | 326865   | 3                     | 14.381 | 2619320 | 20.0 |
| n-Hexadecanoic ...  | 18.045 | 5.8     | ng    | 836825   | 4                     | 17.133 | 2882530 | 20.0 |
| Octadecanoic acid   | 19.280 | 3.9     | ng    | 573312   | 5                     | 21.233 | 2954320 | 20.0 |