

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
 Data File : BP009541.D  
 Acq On : 24 Mar 2022 21:49  
 Operator : CG/JU  
 Sample : N2090-03  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 MW-48

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-BP031722.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP009541.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         | 3.022       | 3             | 6           | 26           | rVB      | 1486453        | 2096920       | 10.12%          | 1.792%        |
| 2         | 4.246       | 209           | 214         | 219          | rBV      | 206373         | 315863        | 1.52%           | 0.270%        |
| 3         | 4.305       | 219           | 224         | 233          | rVB      | 232261         | 379964        | 1.83%           | 0.325%        |
| 4         | 5.375       | 400           | 406         | 423          | rBV      | 2286726        | 4171921       | 20.14%          | 3.565%        |
| 5         | 6.269       | 552           | 558         | 570          | rVB      | 384522         | 680884        | 3.29%           | 0.582%        |
| 6         | 6.758       | 633           | 641         | 654          | rBV      | 500379         | 886708        | 4.28%           | 0.758%        |
| 7         | 6.963       | 666           | 676         | 686          | rBV      | 1335906        | 2680064       | 12.94%          | 2.290%        |
| 8         | 7.769       | 806           | 813         | 823          | rBV      | 884227         | 1586570       | 7.66%           | 1.356%        |
| 9         | 8.146       | 870           | 877         | 889          | rVB      | 1652601        | 2897407       | 13.99%          | 2.476%        |
| 10        | 8.269       | 889           | 898         | 904          | rBV      | 223383         | 388640        | 1.88%           | 0.332%        |
| 11        | 8.881       | 996           | 1002        | 1005         | rBV      | 550083         | 937054        | 4.52%           | 0.801%        |
| 12        | 8.928       | 1005          | 1010        | 1027         | rVB2     | 3080032        | 6559992       | 31.67%          | 5.606%        |
| 13        | 9.399       | 1077          | 1090        | 1098         | rBV      | 529512         | 984851        | 4.75%           | 0.842%        |
| 14        | 9.816       | 1157          | 1161        | 1174         | rVB      | 154105         | 332360        | 1.60%           | 0.284%        |
| 15        | 9.969       | 1176          | 1187        | 1197         | rBV2     | 958199         | 1953594       | 9.43%           | 1.670%        |
| 16        | 10.563      | 1273          | 1288        | 1294         | rBV      | 1284013        | 2583002       | 12.47%          | 2.207%        |
| 17        | 10.616      | 1294          | 1297        | 1304         | rVV3     | 165937         | 327965        | 1.58%           | 0.280%        |
| 18        | 10.681      | 1304          | 1308        | 1315         | rVB      | 124414         | 221315        | 1.07%           | 0.189%        |
| 19        | 11.034      | 1356          | 1368        | 1376         | rVB3     | 130396         | 298129        | 1.44%           | 0.255%        |
| 20        | 11.146      | 1381          | 1387        | 1393         | rBV2     | 176411         | 366107        | 1.77%           | 0.313%        |
| 21        | 11.481      | 1439          | 1444        | 1454         | rVB2     | 133170         | 307793        | 1.49%           | 0.263%        |
| 22        | 11.681      | 1467          | 1478        | 1488         | rVB4     | 161542         | 448754        | 2.17%           | 0.384%        |
| 23        | 12.122      | 1547          | 1553        | 1559         | rBV2     | 205343         | 383707        | 1.85%           | 0.328%        |
| 24        | 12.251      | 1569          | 1575        | 1587         | rVB9     | 101148         | 379053        | 1.83%           | 0.324%        |
| 25        | 12.451      | 1599          | 1609        | 1615         | rBV      | 1362246        | 2430546       | 11.73%          | 2.077%        |
| 26        | 13.040      | 1701          | 1709        | 1721         | rBV      | 8714745        | 14045097      | 67.80%          | 12.003%       |
| 27        | 13.469      | 1778          | 1782        | 1789         | rVB5     | 142611         | 286823        | 1.38%           | 0.245%        |
| 28        | 13.740      | 1822          | 1828        | 1833         | rBV2     | 378459         | 768004        | 3.71%           | 0.656%        |
| 29        | 13.975      | 1864          | 1868        | 1871         | rBV      | 147438         | 228786        | 1.10%           | 0.196%        |
| 30        | 14.139      | 1891          | 1896        | 1903         | rVB      | 313361         | 556654        | 2.69%           | 0.476%        |
| 31        | 14.257      | 1912          | 1916        | 1926         | rVB2     | 185637         | 352165        | 1.70%           | 0.301%        |
| 32        | 14.416      | 1937          | 1943        | 1949         | rVB      | 1929755        | 3071383       | 14.83%          | 2.625%        |
| 33        | 14.481      | 1950          | 1954        | 1960         | rVB3     | 156859         | 246619        | 1.19%           | 0.211%        |
| 34        | 15.916      | 2192          | 2198        | 2211         | rBV      | 6856959        | 10936525      | 52.80%          | 9.346%        |
| 35        | 17.181      | 2408          | 2413        | 2422         | rVB      | 2249190        | 3426745       | 16.54%          | 2.929%        |
| 36        | 18.098      | 2565          | 2569        | 2574         | rBV2     | 442790         | 628357        | 3.03%           | 0.537%        |

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Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

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-BP031722.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |          |          |         |         |
|----|--------|------|------|------|------|----------|----------|---------|---------|
| 37 | 19.763 | 2846 | 2852 | 2859 | rBV  | 13255461 | 18278856 | 88.24%  | 15.621% |
| 38 | 21.033 | 3065 | 3068 | 3075 | rVB  | 301041   | 442229   | 2.13%   | 0.378%  |
| 39 | 21.292 | 3108 | 3112 | 3119 | rBV  | 2937365  | 3874491  | 18.70%  | 3.311%  |
| 40 | 21.427 | 3129 | 3135 | 3155 | rVB  | 13410079 | 20714461 | 100.00% | 17.703% |
| 41 | 21.710 | 3180 | 3183 | 3186 | rBV  | 272897   | 298001   | 1.44%   | 0.255%  |
| 42 | 23.686 | 3513 | 3519 | 3530 | rVB2 | 1994843  | 4258001  | 20.56%  | 3.639%  |

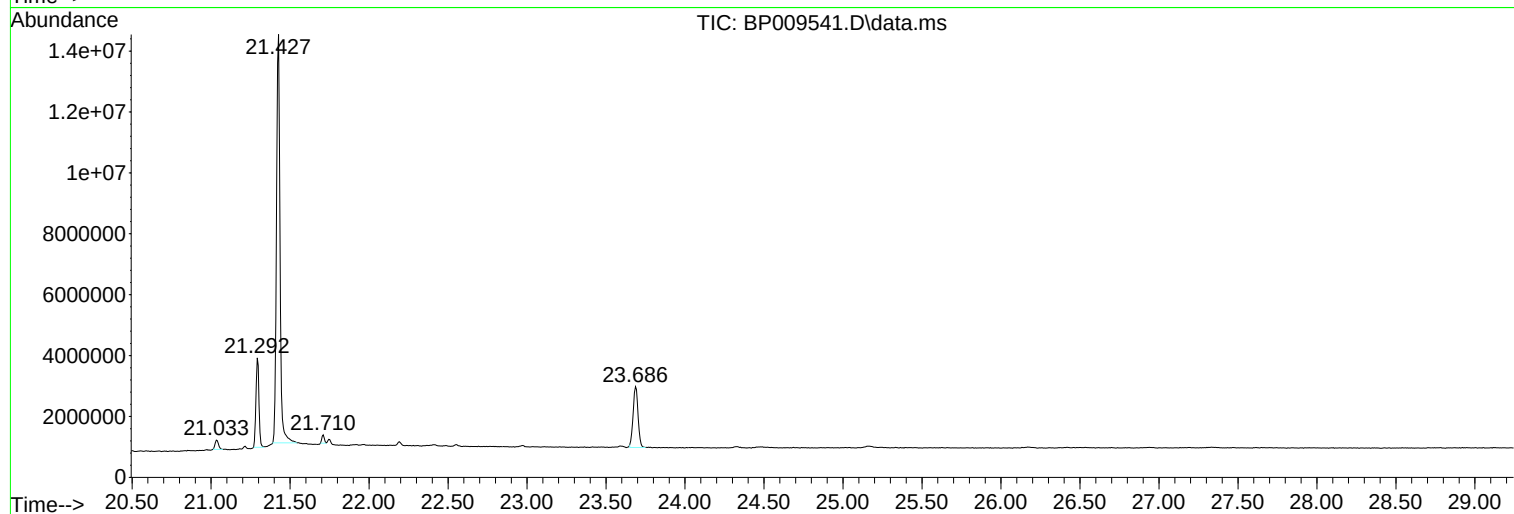
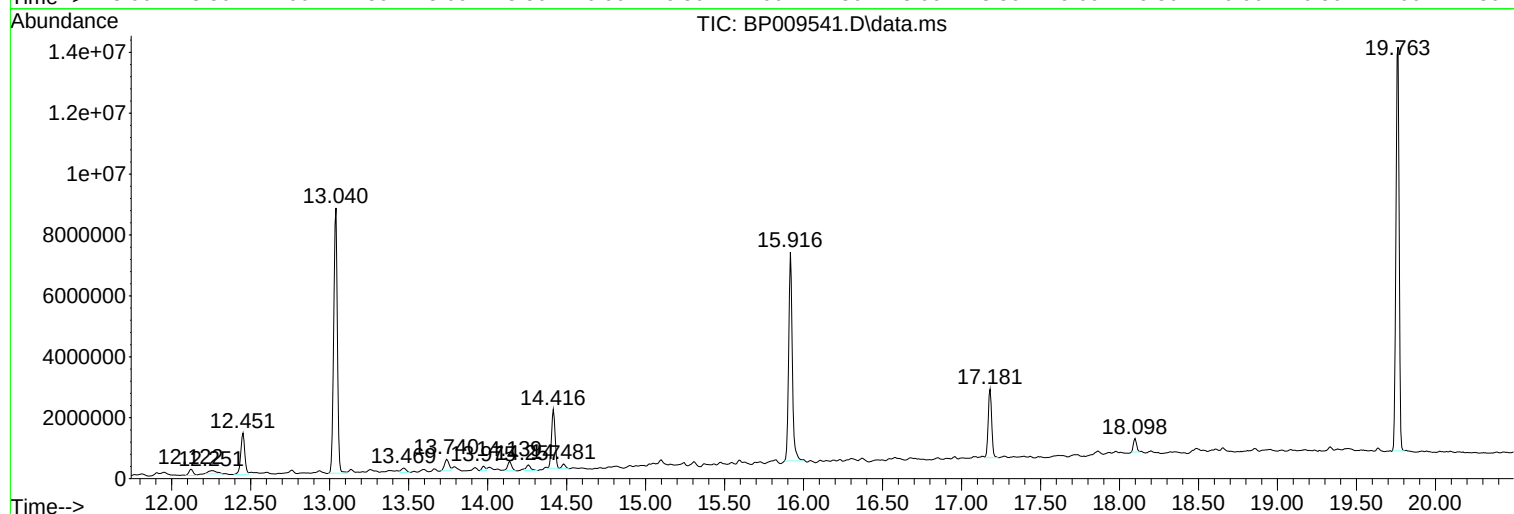
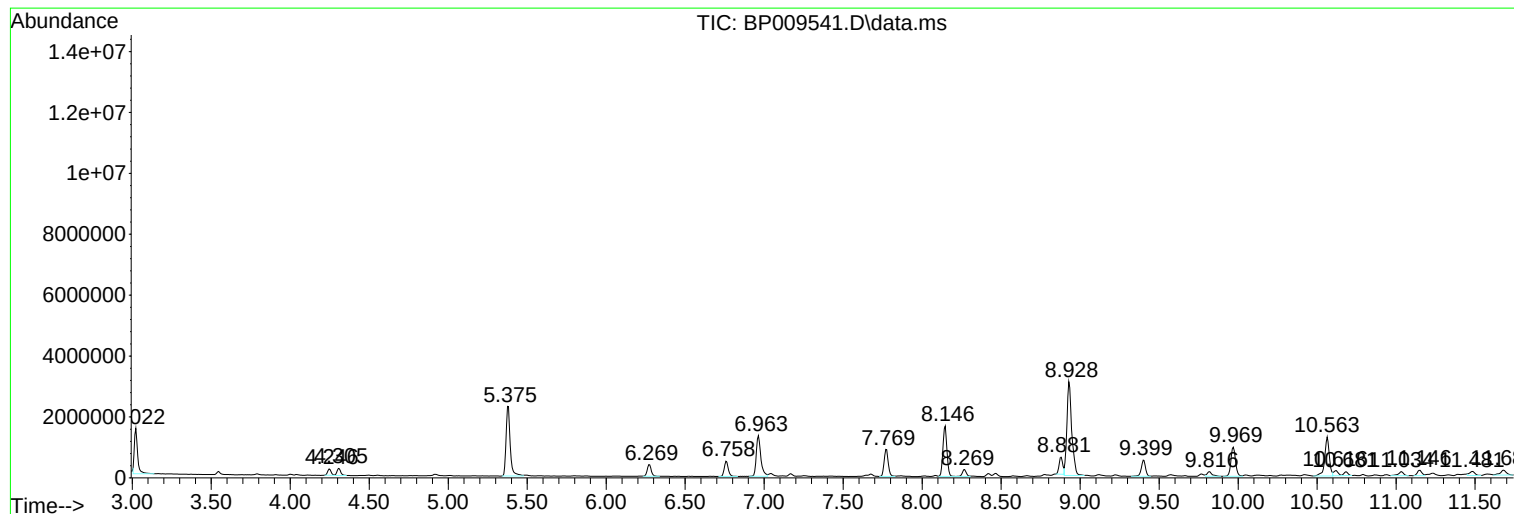
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Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
Data File : BP009541.D  
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Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.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\BP032422\  
Data File : BP009541.D  
Acq On : 24 Mar 2022 21:49  
Operator : CG/JU  
Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

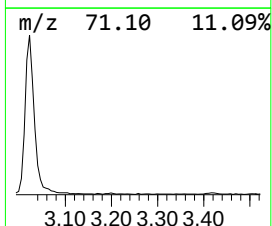
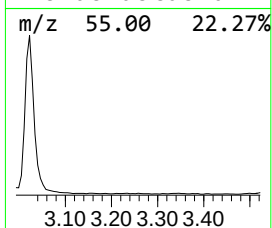
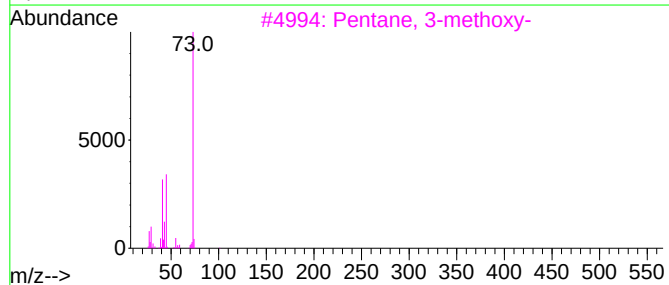
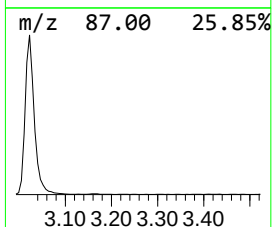
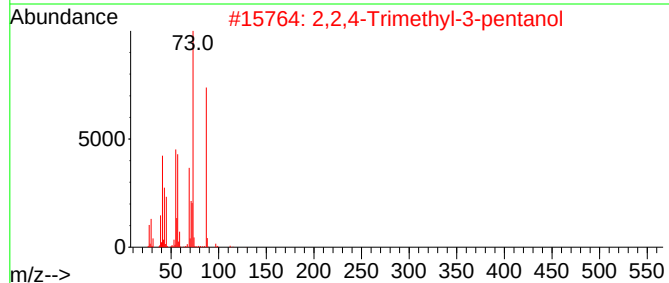
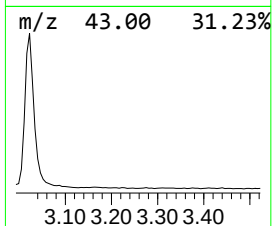
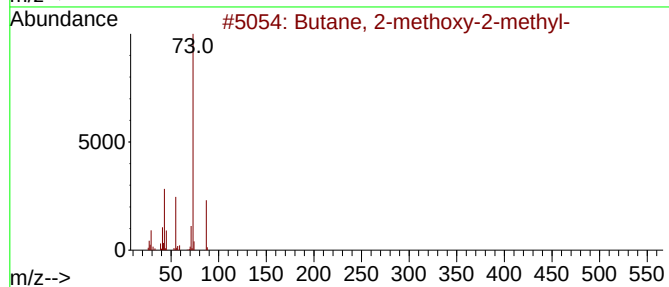
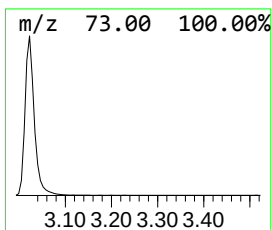
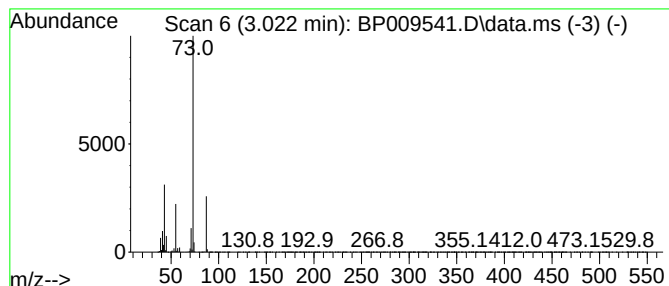
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 3.022 | 26.43 ng | 2096920 | 1,4-Dichlorobenzene-d4 | 7.769 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 90   |
| 2    |    |   | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 43   |
| 3    |    |   | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 23   |
| 4    |    |   | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 9    |
| 5    |    |   | N-Ethylformamide            | 73  | C3H7NO  | 000627-45-2 | 9    |



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Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

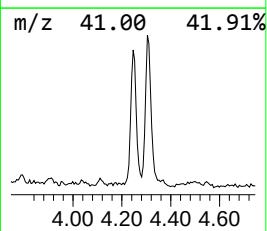
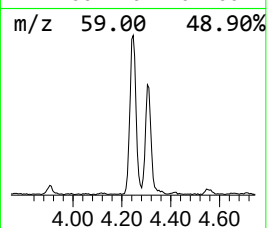
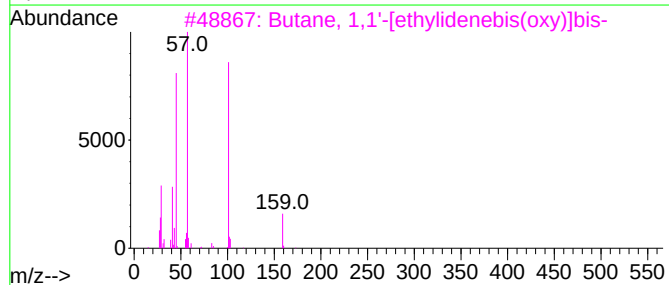
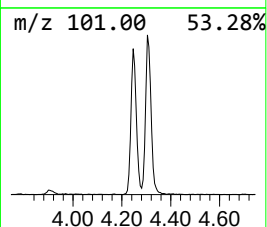
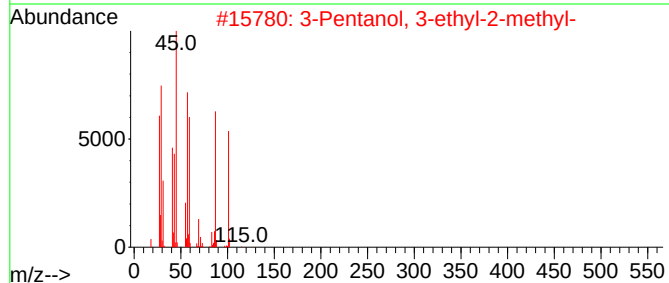
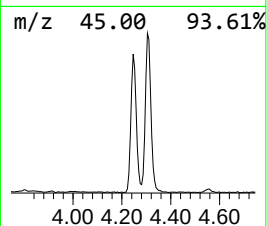
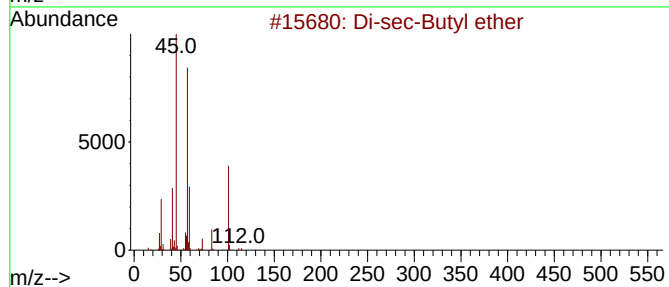
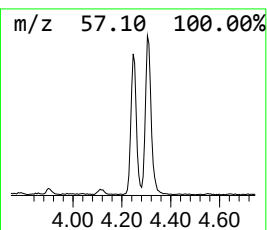
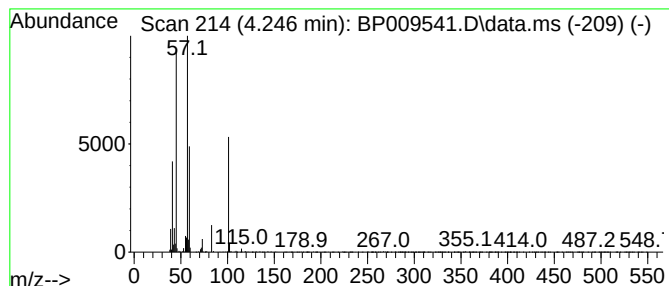
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 2 Di-sec-Butyl ether Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 4.246 | 3.98 ng | 315863 | 1,4-Dichlorobenzene-d4 | 7.769 |

| Hit# | of 5 | Tentative ID                         | MW  | MolForm  | CAS#         | Qual |
|------|------|--------------------------------------|-----|----------|--------------|------|
| 1    |      | Di-sec-Butyl ether                   | 130 | C8H18O   | 006863-58-7  | 83   |
| 2    |      | 3-Pentanol, 3-ethyl-2-methyl-        | 130 | C8H18O   | 000597-05-7  | 56   |
| 3    |      | Butane, 1,1'-[ethylidenebis(oxy)]... | 174 | C10H22O2 | 000871-22-7  | 53   |
| 4    |      | Butane, 1-(1-methylpropoxy)-         | 130 | C8H18O   | 000999-65-5  | 47   |
| 5    |      | DL-Lactamide, butyl ether            | 145 | C7H15NO2 | 1000452-56-5 | 42   |



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Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

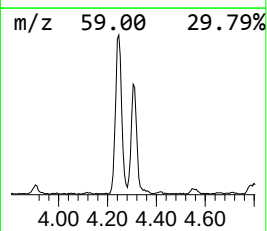
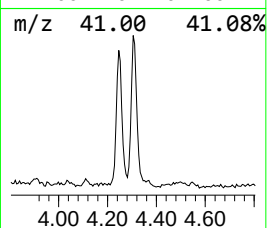
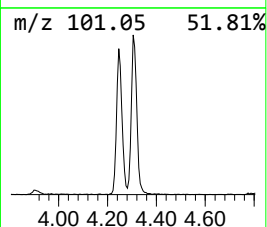
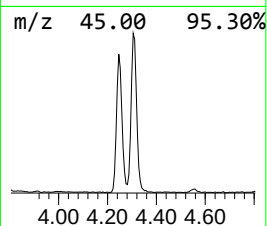
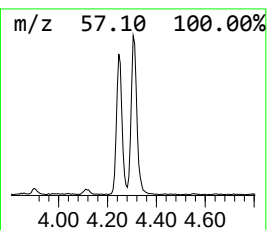
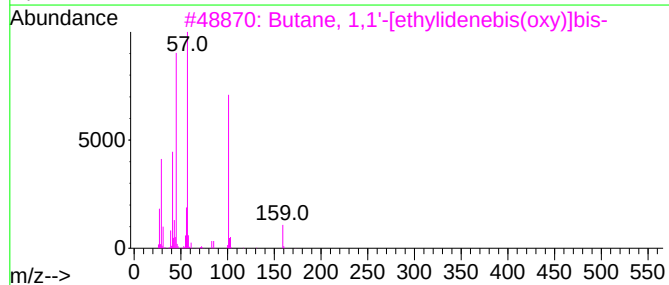
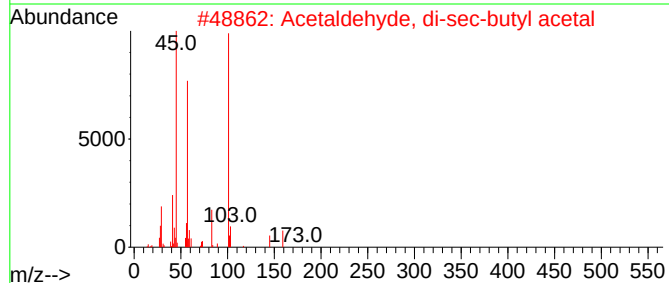
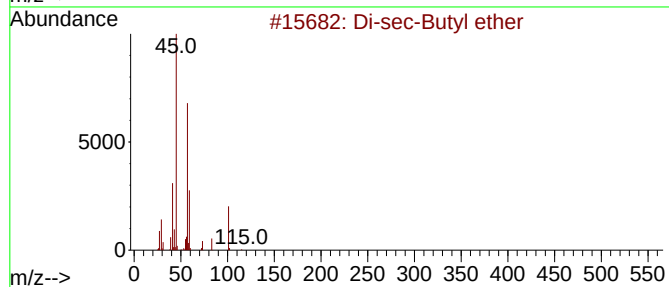
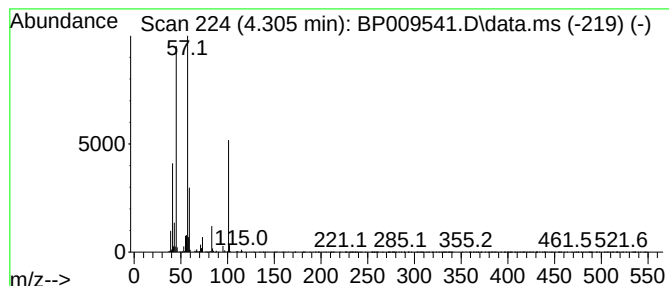
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 Acetaldehyde, di-sec-butyl ... Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 4.305 | 4.79 ng | 379964 | 1,4-Dichlorobenzene-d4 | 7.769 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm  | CAS#         | Qual |
|------|----|---|--------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Di-sec-Butyl ether                   | 130 | C8H18O   | 006863-58-7  | 83   |
| 2    |    |   | Acetaldehyde, di-sec-butyl acetal    | 174 | C10H22O2 | 005314-41-0  | 64   |
| 3    |    |   | Butane, 1,1'-[ethylidenebis(oxy)]... | 174 | C10H22O2 | 000871-22-7  | 59   |
| 4    |    |   | Butane, 1-(1-methylpropoxy)-         | 130 | C8H18O   | 000999-65-5  | 53   |
| 5    |    |   | DL-Lactamide, butyl ether            | 145 | C7H15NO2 | 1000452-56-5 | 45   |



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Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

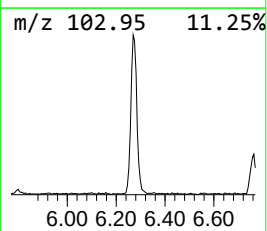
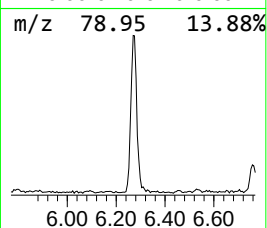
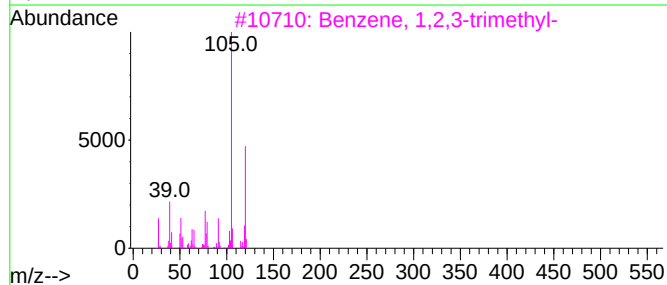
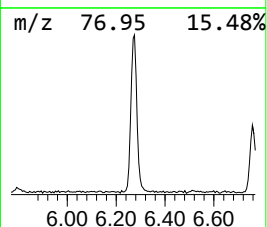
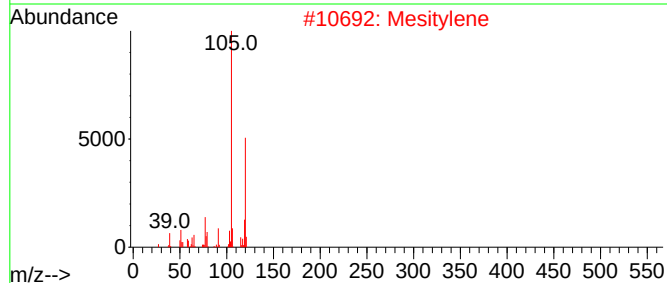
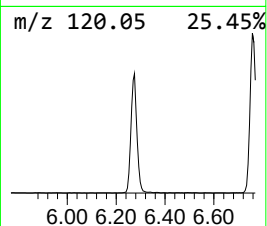
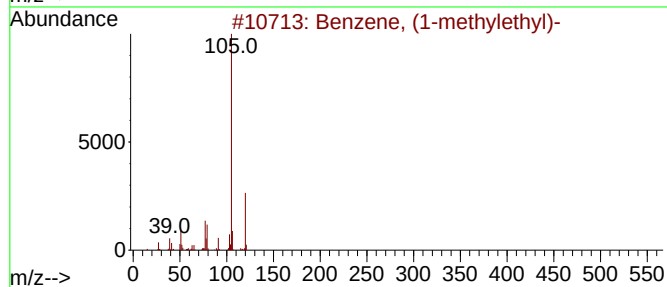
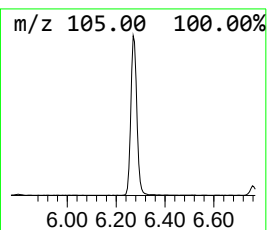
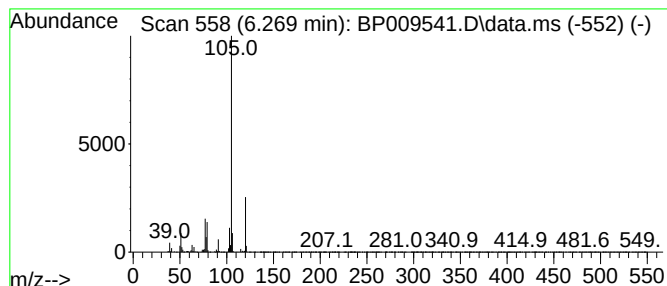
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 4 Benzene, (1-methylethyl)- Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 6.269 | 8.58 ng | 680884 | 1,4-Dichlorobenzene-d4 | 7.769 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, (1-methylethyl)-  | 120 | C9H12   | 000098-82-8 | 94   |
| 2    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 91   |
| 3    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |
| 4    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 87   |
| 5    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
Data File : BP009541.D  
Acq On : 24 Mar 2022 21:49  
Operator : CG/JU  
Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

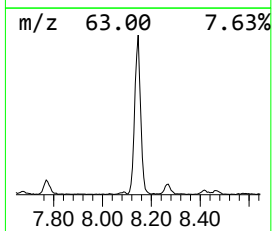
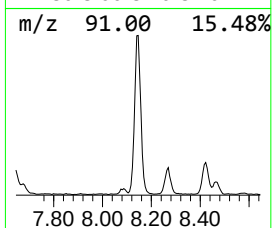
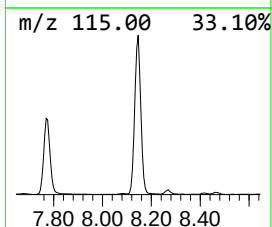
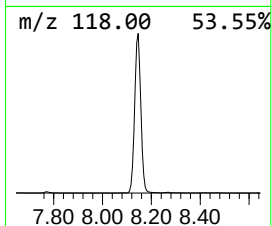
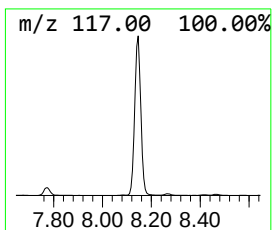
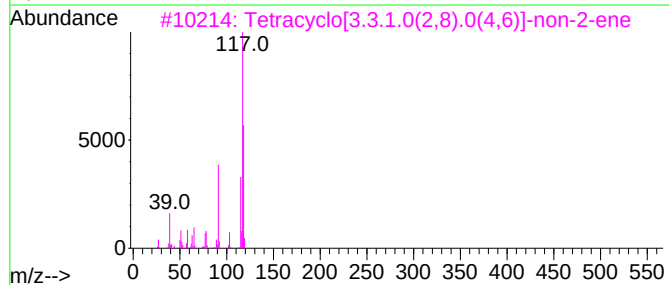
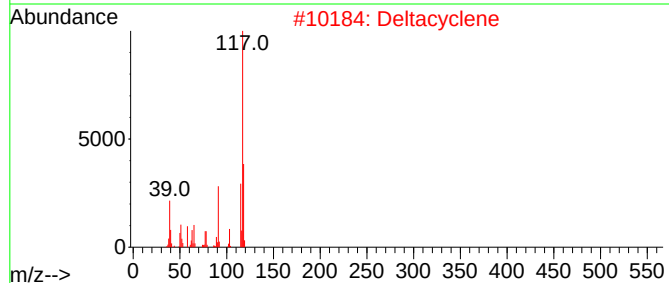
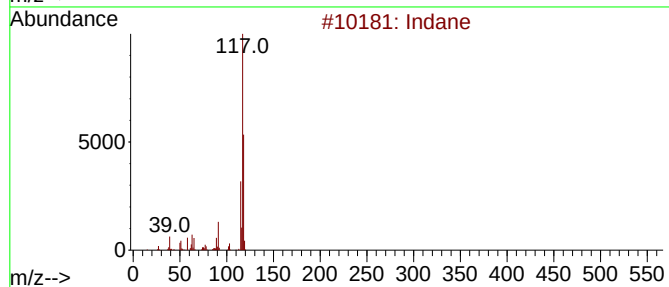
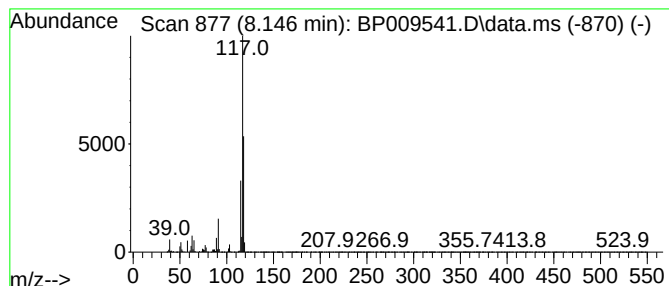
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 6 Indane Concentration Rank 2

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 8.146 | 36.52 ng | 2897410 | 1,4-Dichlorobenzene-d4 | 7.769 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Indane                              | 118 | C9H10   | 000496-11-7  | 95   |
| 2    |    |   | Deltacyclene                        | 118 | C9H10   | 007785-10-6  | 72   |
| 3    |    |   | Tetracyclo[3.3.1.0(2,8).0(4,6)]-... | 118 | C9H10   | 1000191-13-7 | 64   |
| 4    |    |   | Benzene, cyclopropyl-               | 118 | C9H10   | 000873-49-4  | 64   |
| 5    |    |   | Benzene, 1,1'-(1,5-hexadiene-1,6... | 234 | C18H18  | 004439-45-6  | 59   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
Data File : BP009541.D  
Acq On : 24 Mar 2022 21:49  
Operator : CG/JU  
Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

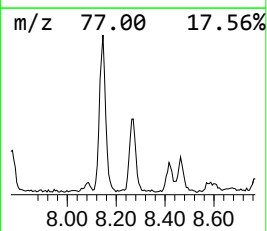
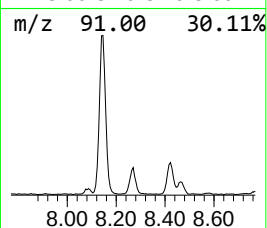
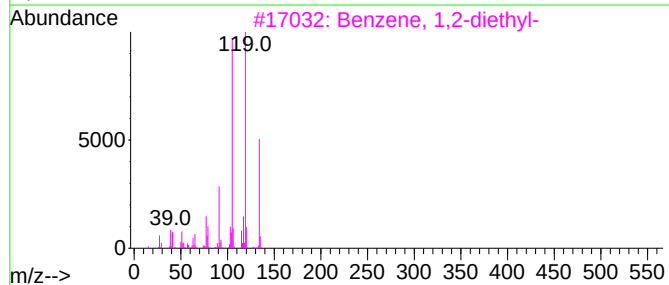
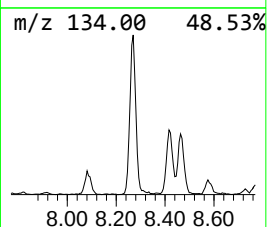
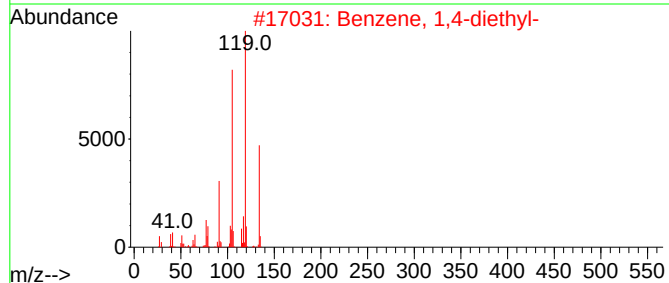
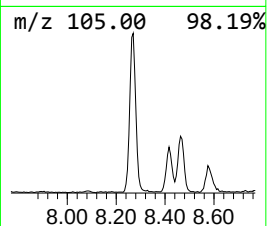
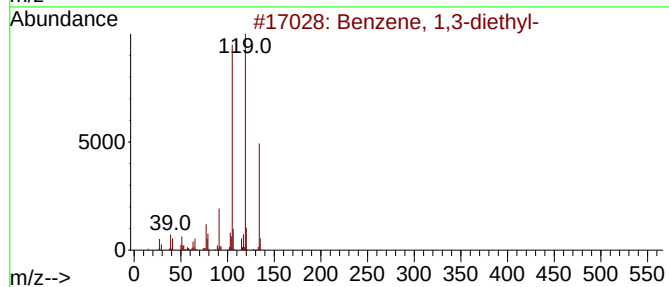
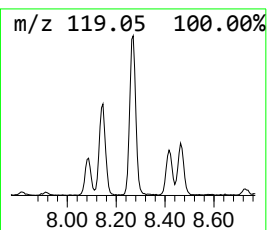
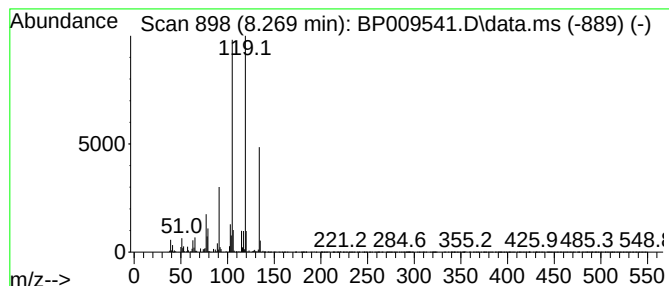
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 Benzene, 1,3-diethyl- Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 8.269 | 4.90 ng | 388640 | 1,4-Dichlorobenzene-d4 | 7.769 |

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



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
Data File : BP009541.D  
Acq On : 24 Mar 2022 21:49  
Operator : CG/JU  
Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.M  
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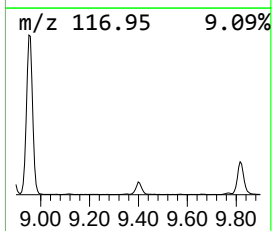
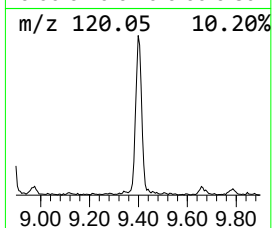
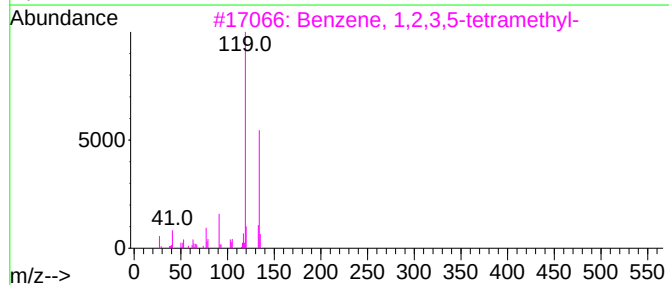
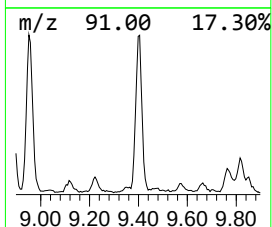
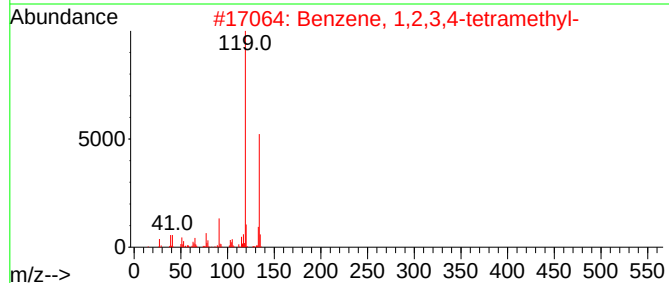
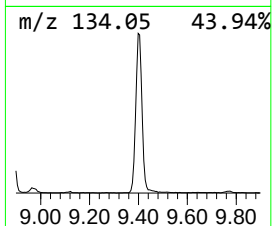
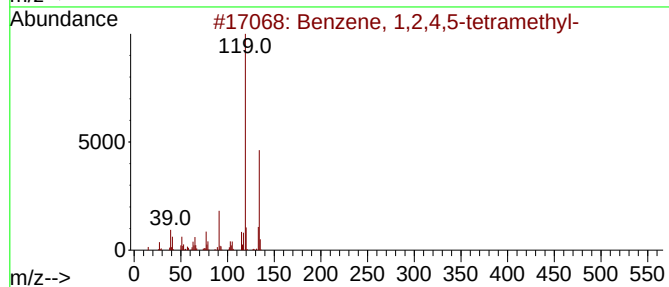
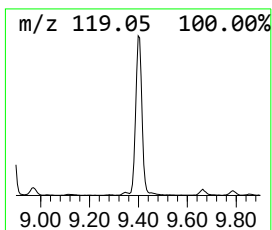
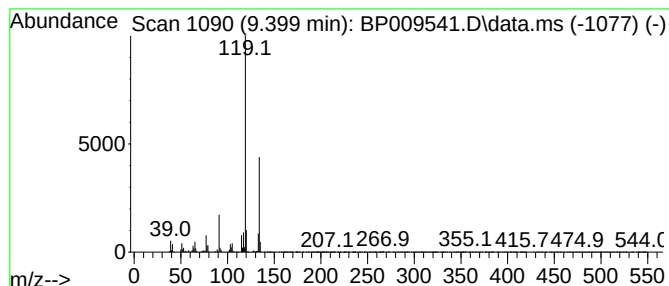
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.   |
|-------|---------|--------|------------------|--------|
| 9.399 | 7.63 ng | 984851 | Naphthalene-d8   | 10.563 |

| Hit# | of 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|------|--------------------------------|-----|---------|-------------|------|
| 1    |      | Benzene, 1,2,4,5-tetramethyl-  | 134 | C10H14  | 000095-93-2 | 97   |
| 2    |      | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 95   |
| 3    |      | Benzene, 1,2,3,5-tetramethyl-  | 134 | C10H14  | 000527-53-7 | 95   |
| 4    |      | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 95   |
| 5    |      | Benzene, 1-ethyl-2,4-dimethyl- | 134 | C10H14  | 000874-41-9 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
Data File : BP009541.D  
Acq On : 24 Mar 2022 21:49  
Operator : CG/JU  
Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
MW-48

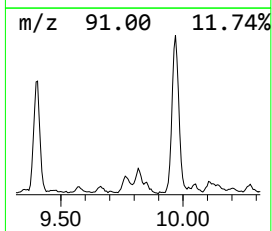
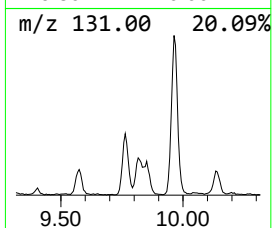
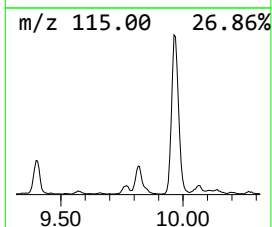
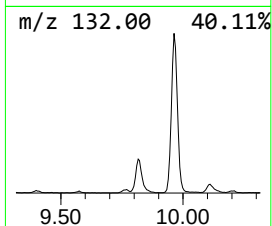
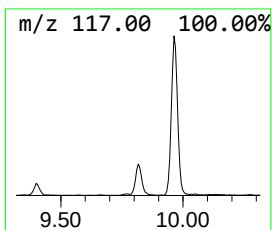
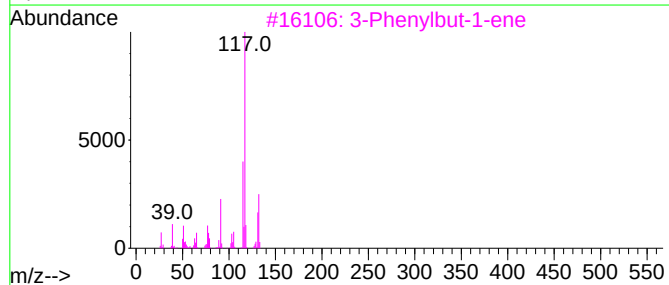
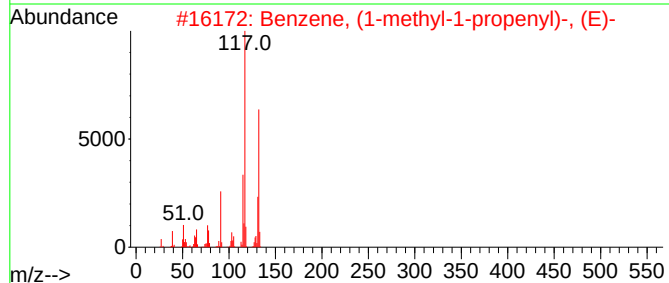
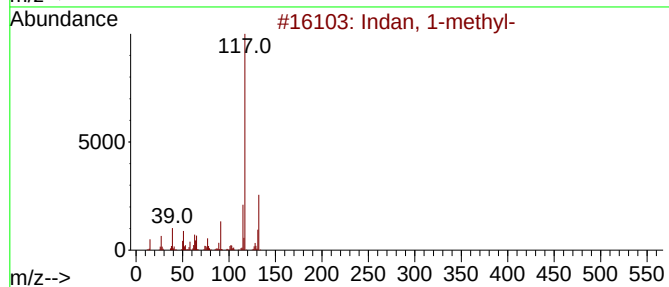
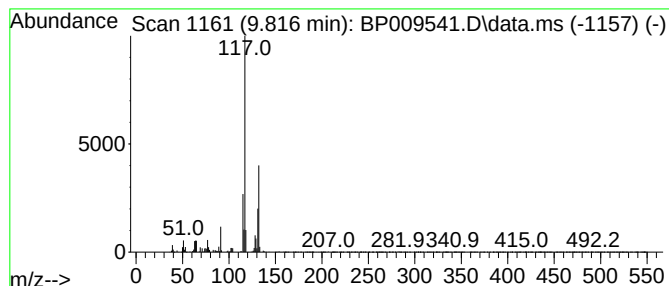
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.   |
|-------|---------|--------|------------------|--------|
| 9.816 | 2.57 ng | 332360 | Naphthalene-d8   | 10.563 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Indan, 1-methyl-                    | 132 | C10H12  | 000767-58-8 | 90   |
| 2    |    |   | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000768-00-3 | 87   |
| 3    |    |   | 3-Phenylbut-1-ene                   | 132 | C10H12  | 000934-10-1 | 87   |
| 4    |    |   | 1H-Indene, 2,3-dihydro-5-methyl-    | 132 | C10H12  | 000874-35-1 | 86   |
| 5    |    |   | Benzene, 1-ethenyl-4-ethyl-         | 132 | C10H12  | 003454-07-7 | 81   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
Data File : BP009541.D  
Acq On : 24 Mar 2022 21:49  
Operator : CG/JU  
Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
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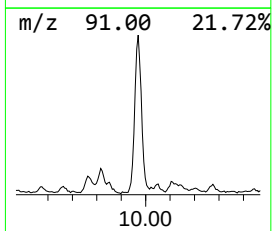
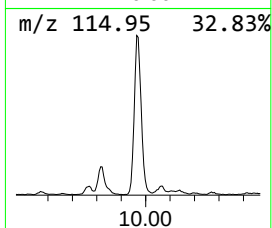
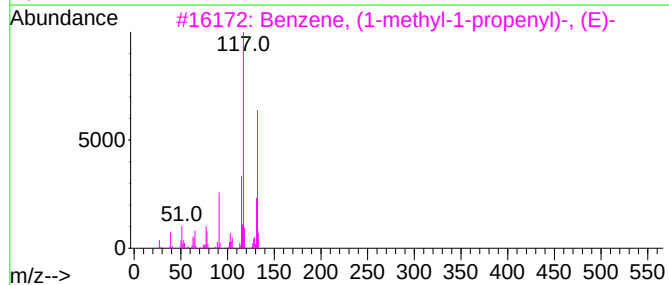
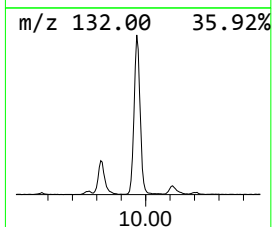
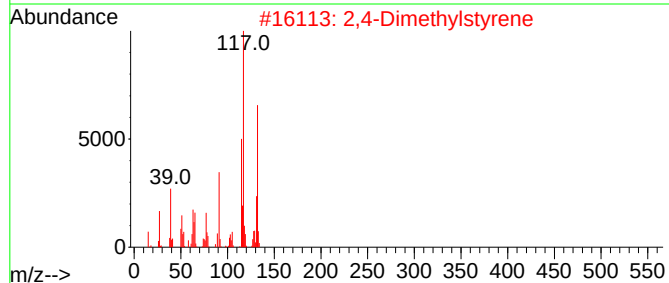
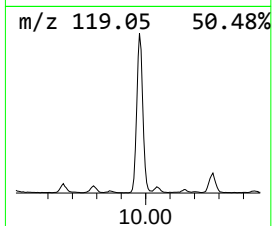
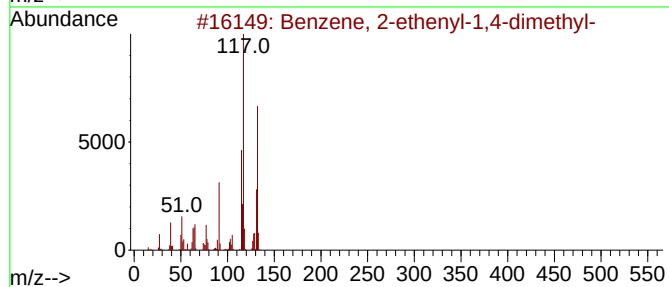
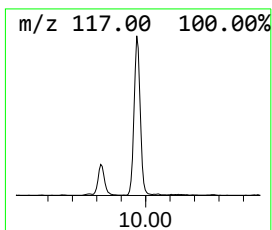
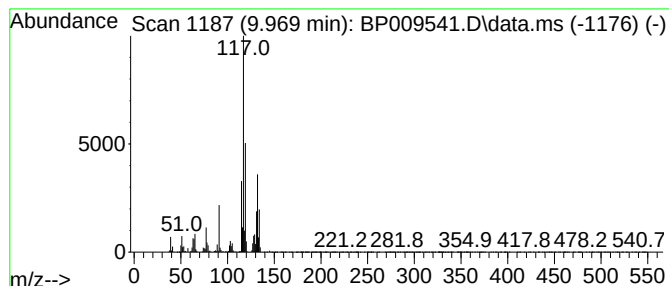
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.   |
|-------|----------|---------|------------------|--------|
| 9.969 | 15.13 ng | 1953590 | Naphthalene-d8   | 10.563 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Benzene, 2-ethenyl-1,4-dimethyl-    | 132 | C10H12  | 002039-89-6 | 95   |
| 2    |      | 2,4-Dimethylstyrene                 | 132 | C10H12  | 002234-20-0 | 89   |
| 3    |      | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000768-00-3 | 70   |
| 4    |      | Benzene, 2-ethenyl-1,3-dimethyl-    | 132 | C10H12  | 002039-90-9 | 70   |
| 5    |      | 3-Phenylbut-1-ene                   | 132 | C10H12  | 000934-10-1 | 70   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
Data File : BP009541.D  
Acq On : 24 Mar 2022 21:49  
Operator : CG/JU  
Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

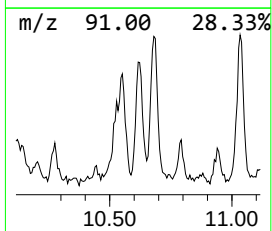
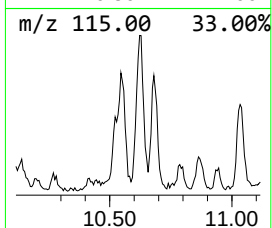
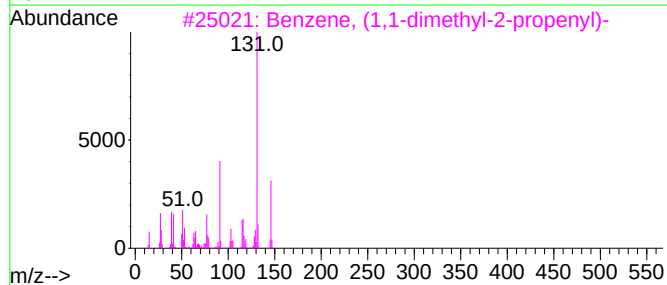
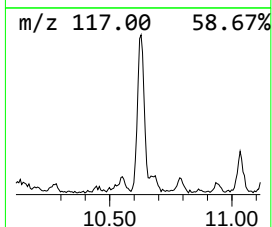
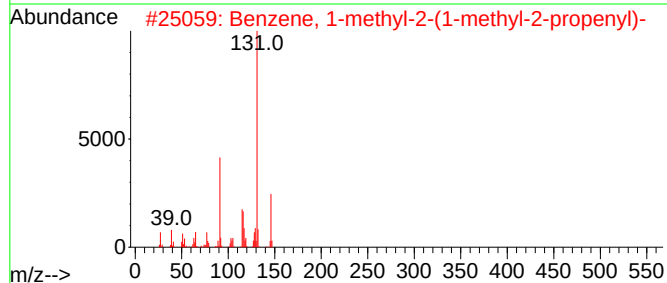
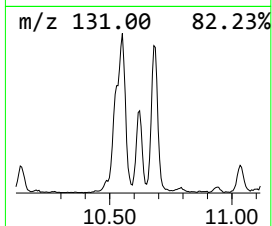
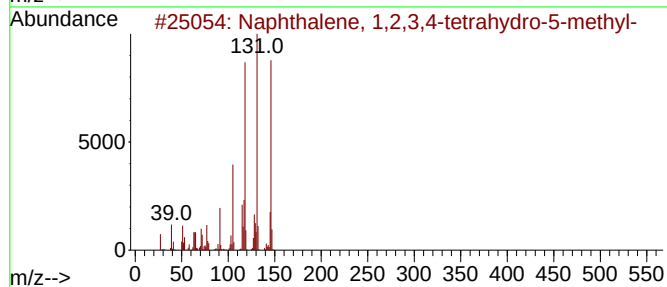
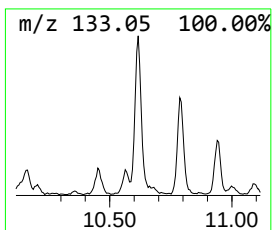
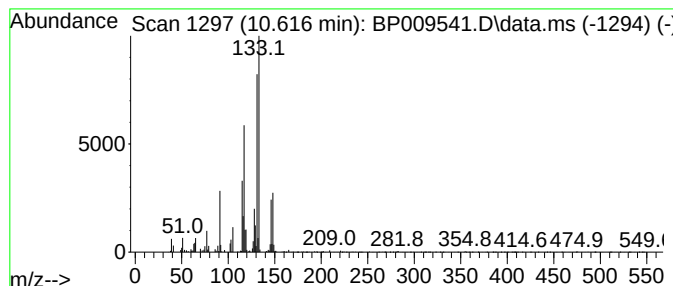
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 11 unknown10.616 Concentration Rank 19

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 10.616 | 2.54 ng | 327965 | Naphthalene-d8   | 10.563 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 002809-64-5 | 49   |
| 2    |    |   | Benzene, 1-methyl-2-(1-methyl-2-... | 146 | C11H14  | 097664-19-2 | 46   |
| 3    |    |   | Benzene, (1,1-dimethyl-2-propenyl)- | 146 | C11H14  | 018321-36-3 | 42   |
| 4    |    |   | Benzene, (1-methyl-1-butenyl)-      | 146 | C11H14  | 053172-84-2 | 38   |
| 5    |    |   | Benzene, 1-methyl-4-(1-methyl-2-... | 146 | C11H14  | 097664-18-1 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
Data File : BP009541.D  
Acq On : 24 Mar 2022 21:49  
Operator : CG/JU  
Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

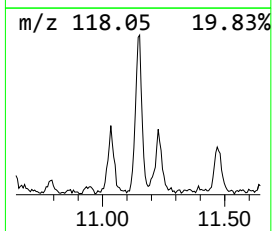
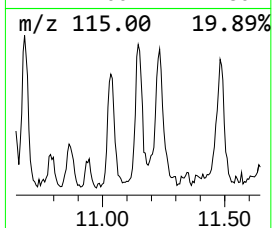
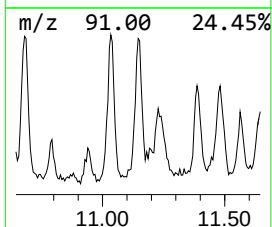
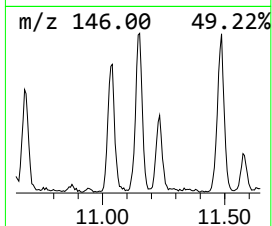
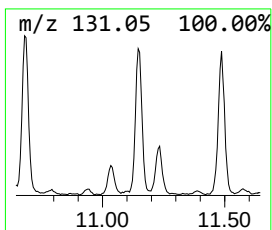
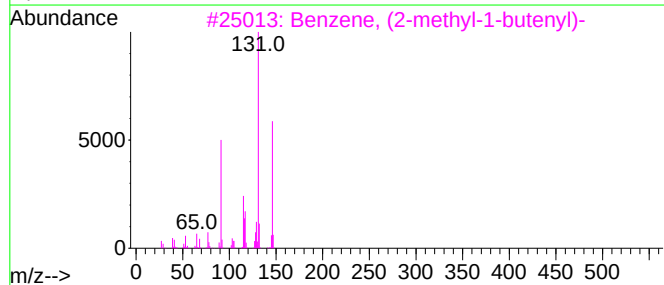
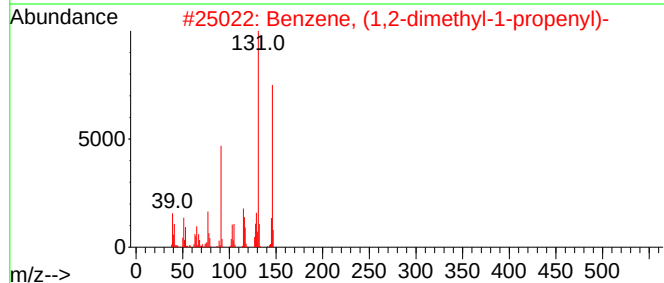
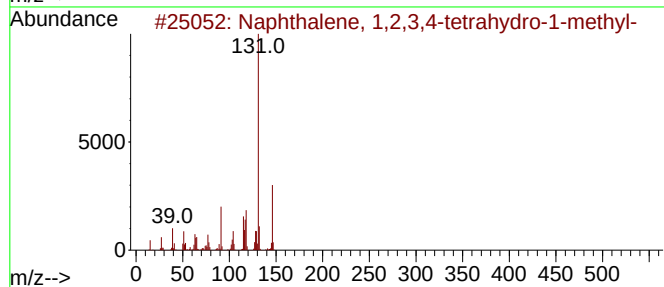
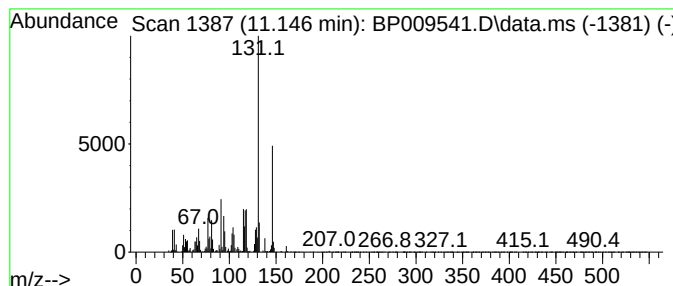
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 12 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 17

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 11.146    | 2.83 ng                             | 366107 | Naphthalene-d8   | 10.563      |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Naphthalene, 1,2,3,4-tetrahydro-... | 146    | C11H14           | 001559-81-5 | 90   |
| 2         | Benzene, (1,2-dimethyl-1-propenyl)- | 146    | C11H14           | 000769-57-3 | 90   |
| 3         | Benzene, (2-methyl-1-butenyl)-      | 146    | C11H14           | 056253-64-6 | 83   |
| 4         | 1H-Indene, 2,3-dihydro-1,3-dimet... | 146    | C11H14           | 004175-53-5 | 76   |
| 5         | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146    | C11H14           | 006682-71-9 | 76   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
Data File : BP009541.D  
Acq On : 24 Mar 2022 21:49  
Operator : CG/JU  
Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

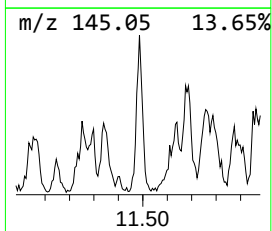
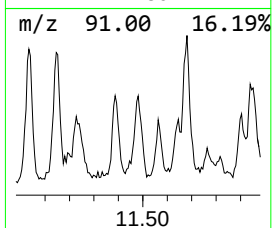
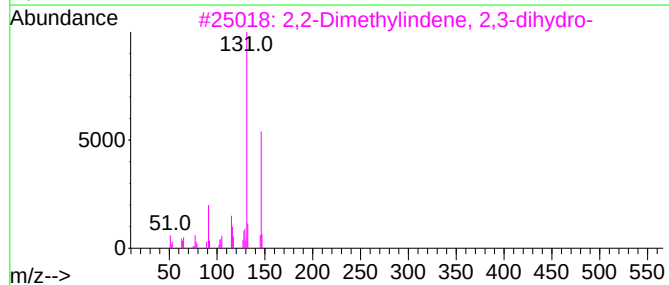
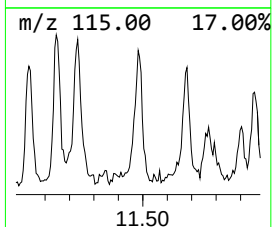
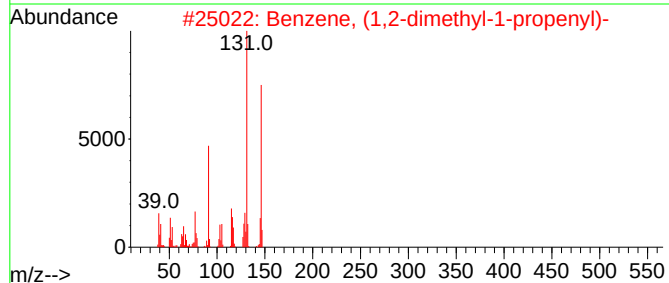
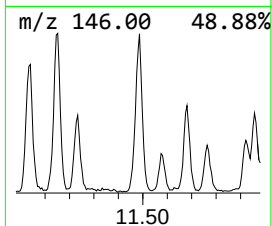
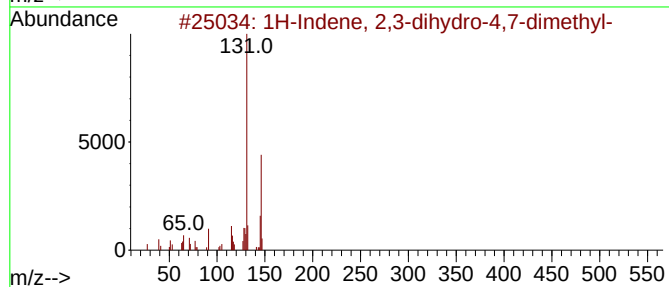
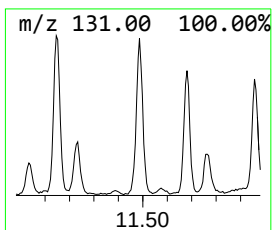
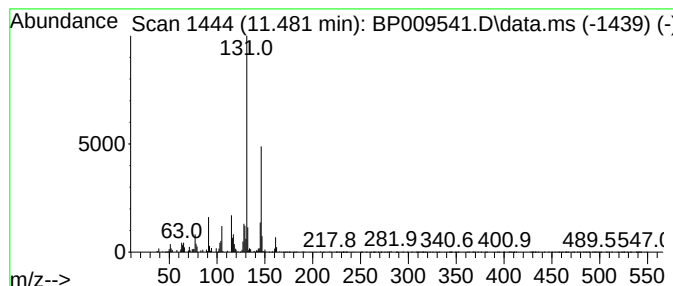
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 13 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 20

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 11.481 | 2.38 ng | 307793 | Naphthalene-d8   | 10.563 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 97   |
| 2    |    |   | Benzene, (1,2-dimethyl-1-propenyl)- | 146 | C11H14  | 000769-57-3 | 94   |
| 3    |    |   | 2,2-Dimethylindene, 2,3-dihydro-    | 146 | C11H14  | 020836-11-7 | 94   |
| 4    |    |   | Benzene, 1-methyl-3-(1-methyl-2-... | 146 | C11H14  | 052161-57-6 | 91   |
| 5    |    |   | Benzene, 1-methyl-4-(1-methyl-2-... | 146 | C11H14  | 097664-18-1 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
Data File : BP009541.D  
Acq On : 24 Mar 2022 21:49  
Operator : CG/JU  
Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

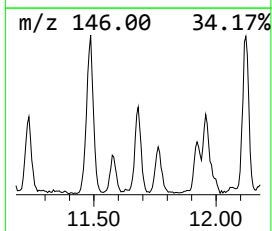
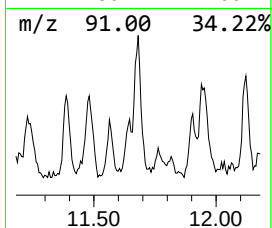
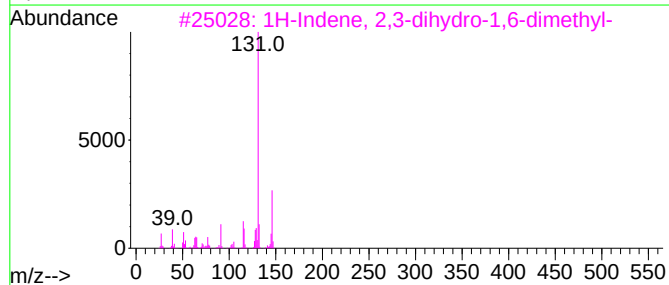
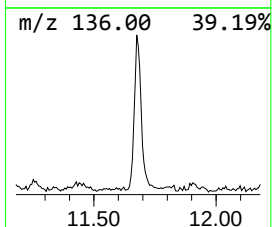
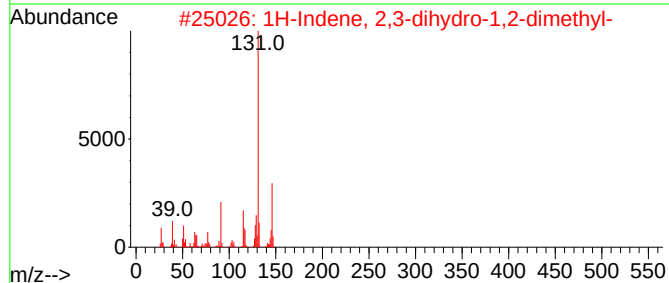
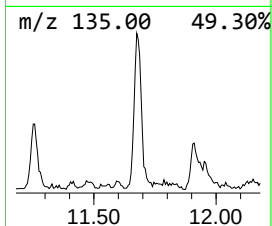
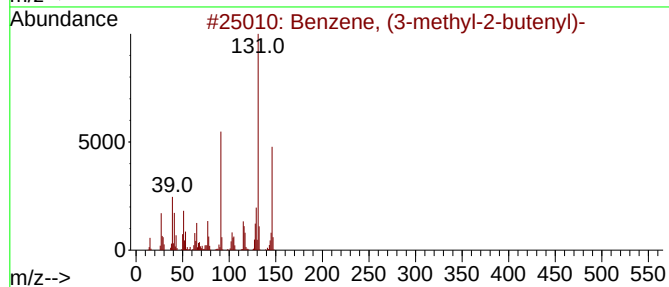
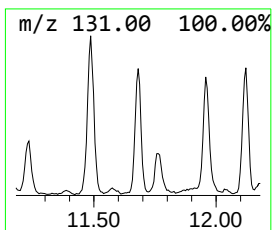
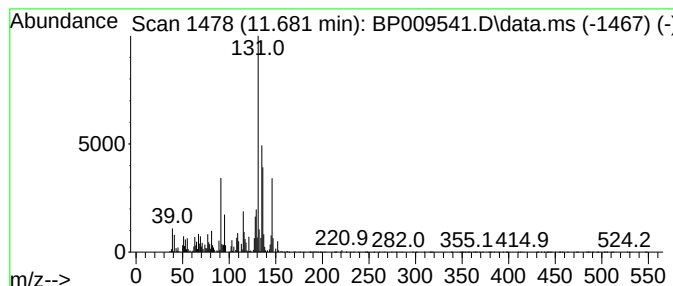
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Benzene, (3-methyl-2-butenyl)- Concentration Rank 14

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 11.681 | 3.47 ng | 448754 | Naphthalene-d8   | 10.563 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, (3-methyl-2-butenyl)-      | 146 | C11H14  | 004489-84-3 | 78   |
| 2    |    |   | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 | C11H14  | 017057-82-8 | 70   |
| 3    |    |   | 1H-Indene, 2,3-dihydro-1,6-dimet... | 146 | C11H14  | 017059-48-2 | 55   |
| 4    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 55   |
| 5    |    |   | Benzene, 2-ethenyl-1,3,5-trimethyl- | 146 | C11H14  | 000769-25-5 | 55   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
Data File : BP009541.D  
Acq On : 24 Mar 2022 21:49  
Operator : CG/JU  
Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

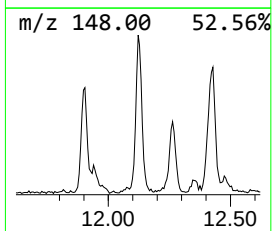
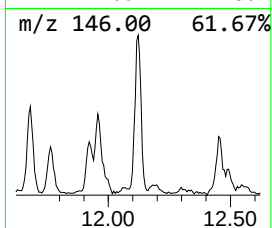
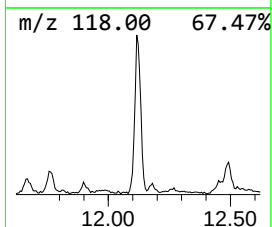
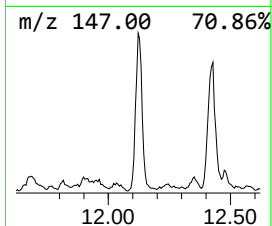
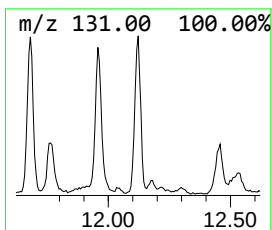
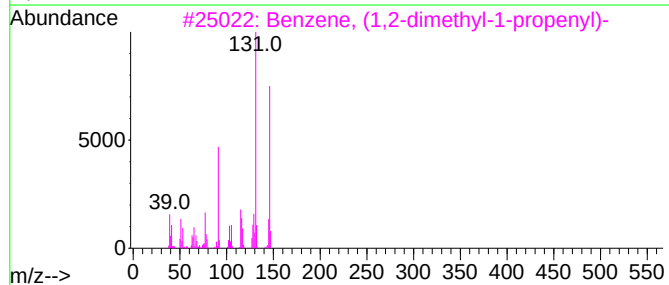
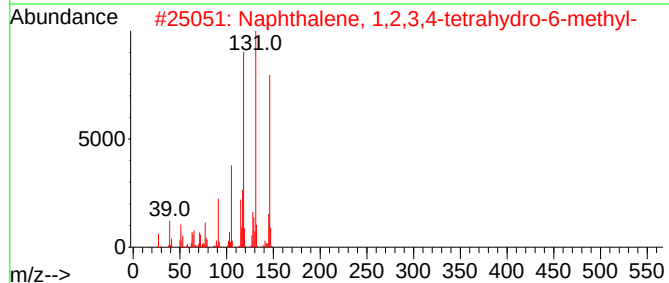
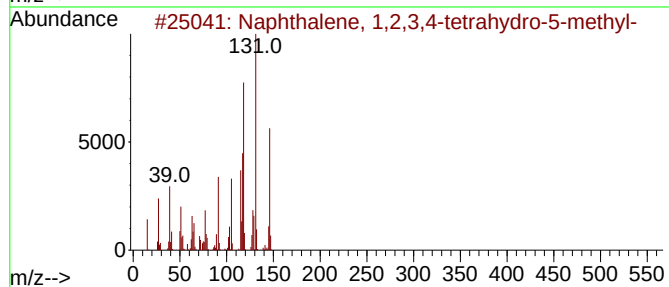
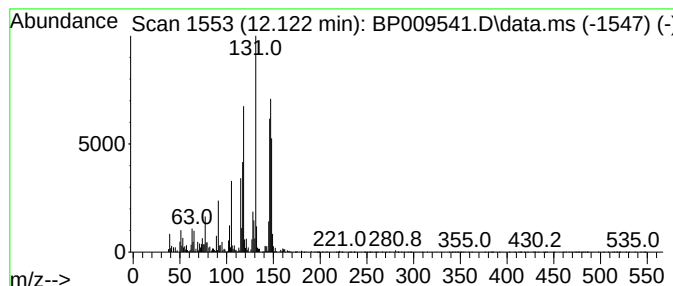
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 15

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 12.122 | 2.97 ng | 383707 | Naphthalene-d8   | 10.563 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 002809-64-5  | 94   |
| 2    |    |   | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 001680-51-9  | 91   |
| 3    |    |   | Benzene, (1,2-dimethyl-1-propenyl)- | 146 | C11H14  | 000769-57-3  | 86   |
| 4    |    |   | 2,3,4,5,6,7-Hexahydro-1H-cyclope... | 146 | C11H14  | 1010189-31-0 | 58   |
| 5    |    |   | Benzene, (1-methyl-1-butenyl)-      | 146 | C11H14  | 053172-84-2  | 55   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
Data File : BP009541.D  
Acq On : 24 Mar 2022 21:49  
Operator : CG/JU  
Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

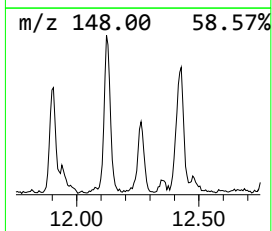
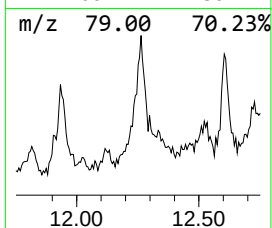
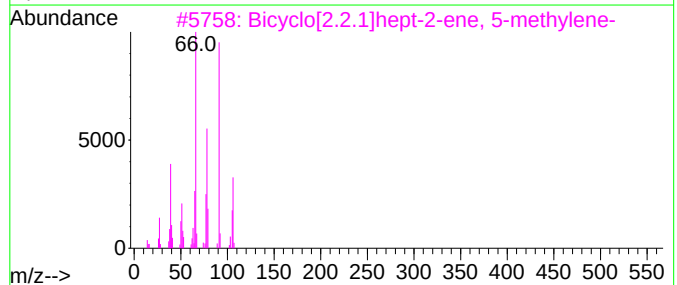
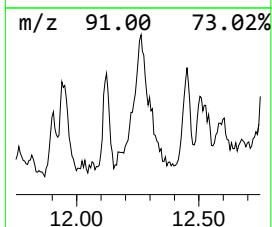
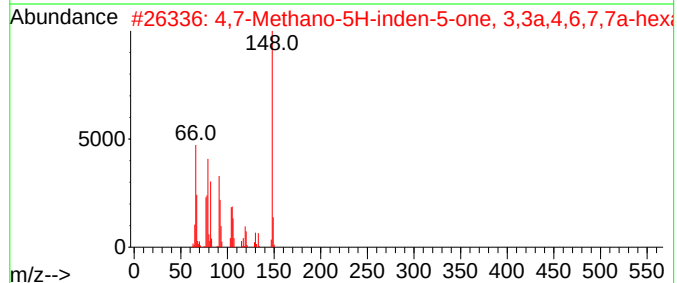
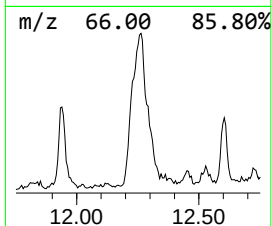
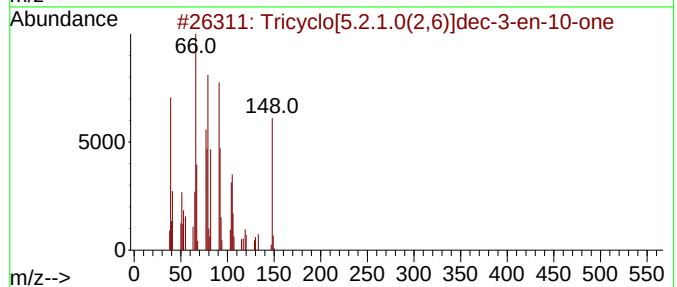
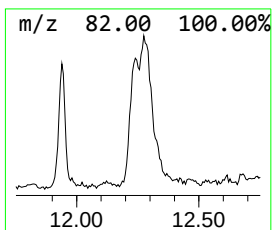
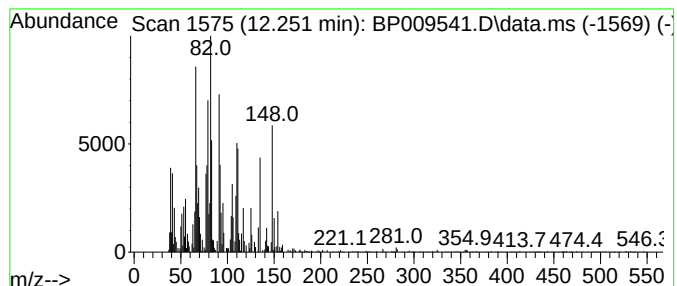
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Tricyclo[5.2.1.0(2,6)]dec-3... Concentration Rank 16

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 12.251 | 2.93 ng | 379053 | Naphthalene-d8   | 10.563 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Tricyclo[5.2.1.0(2,6)]dec-3-en-1... | 148 | C10H12O | 1000289-10-5 | 64   |
| 2    |    |   | 4,7-Methano-5H-inden-5-one, 3,3a... | 148 | C10H12O | 014888-58-5  | 60   |
| 3    |    |   | Bicyclo[2.2.1]hept-2-ene, 5-meth... | 106 | C8H10   | 000694-91-7  | 50   |
| 4    |    |   | Tetracyclo[3.3.1.1(3,7).0(4,6)]d... | 148 | C10H12O | 052750-53-5  | 49   |
| 5    |    |   | 4,7-Methano-1H-indenol, hexahydro-  | 150 | C10H14O | 037275-49-3  | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
Data File : BP009541.D  
Acq On : 24 Mar 2022 21:49  
Operator : CG/JU  
Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

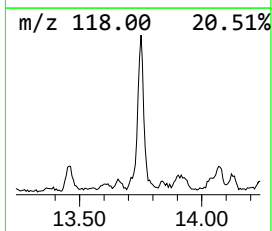
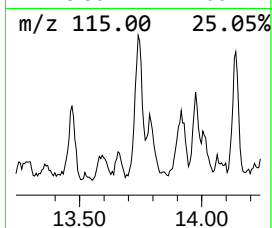
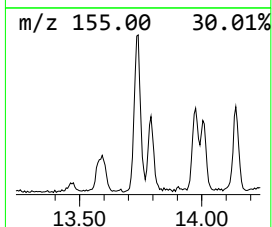
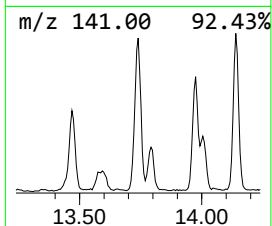
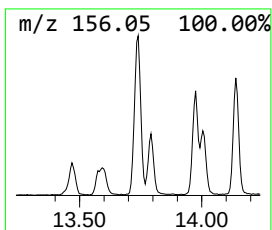
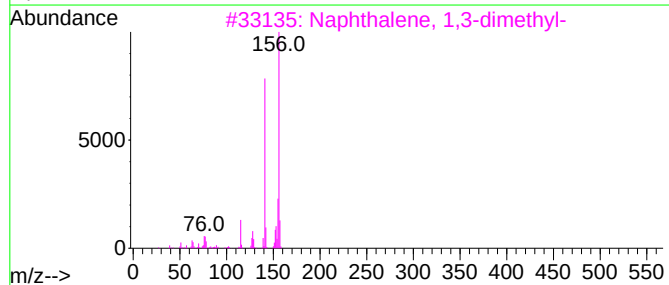
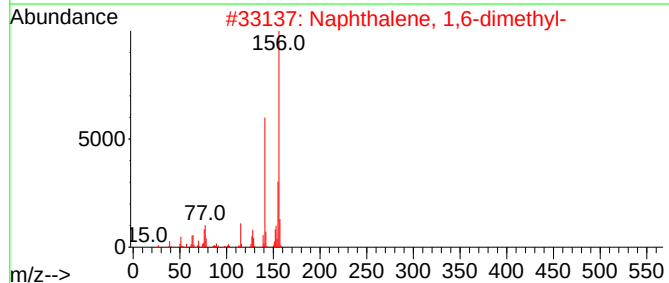
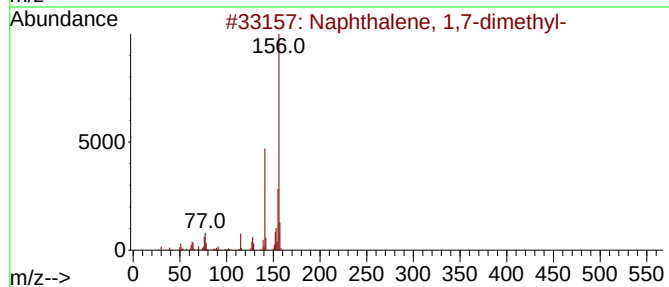
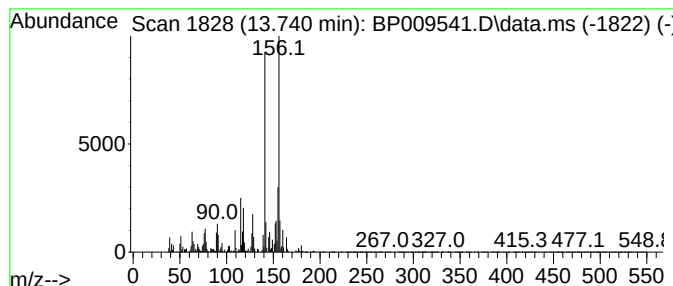
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 17 Naphthalene, 1,7-dimethyl- Concentration Rank 8

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 13.740 | 5.00 ng | 768004 | Acenaphthene-d10 | 14.416 |

| Hit# | of 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|------|----------------------------|-----|---------|-------------|------|
| 1    |      | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 96   |
| 2    |      | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |
| 3    |      | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 96   |
| 4    |      | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 95   |
| 5    |      | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 95   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
Data File : BP009541.D  
Acq On : 24 Mar 2022 21:49  
Operator : CG/JU  
Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

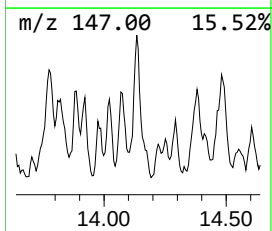
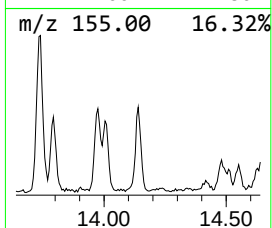
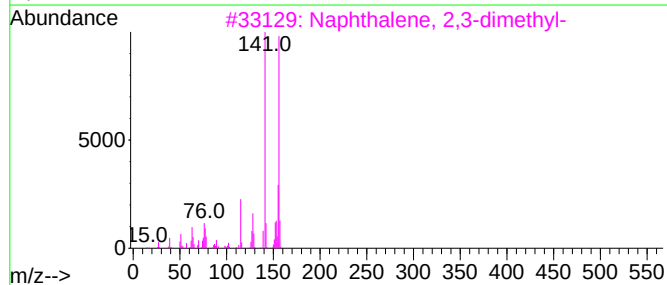
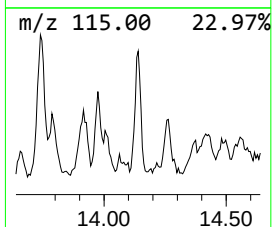
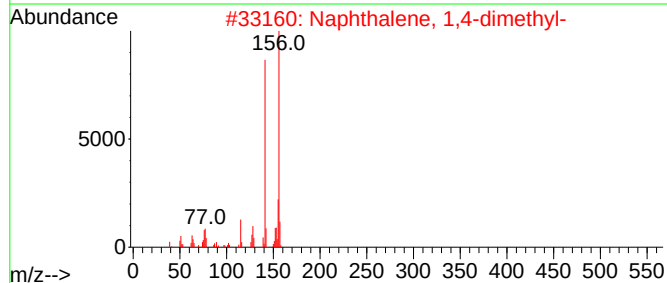
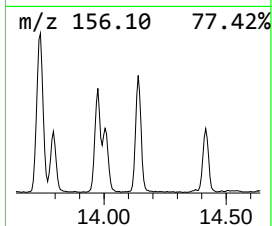
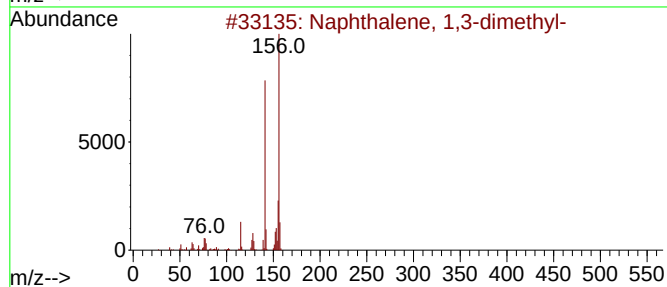
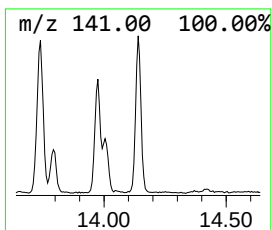
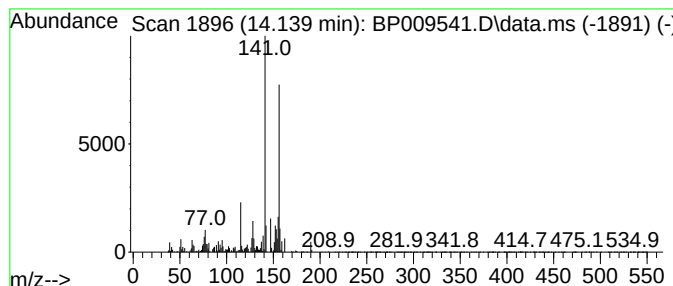
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Naphthalene, 1,3-dimethyl- Concentration Rank 13

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 14.140 | 3.62 ng | 556654 | Acenaphthene-d10 | 14.416 |

| Hit# | of 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|------|----------------------------|-----|---------|-------------|------|
| 1    |      | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 95   |
| 2    |      | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 95   |
| 3    |      | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 95   |
| 4    |      | Naphthalene, 1,2-dimethyl- | 156 | C12H12  | 000573-98-8 | 94   |
| 5    |      | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
Data File : BP009541.D  
Acq On : 24 Mar 2022 21:49  
Operator : CG/JU  
Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

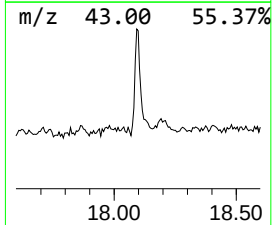
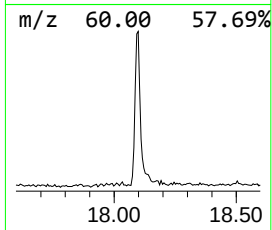
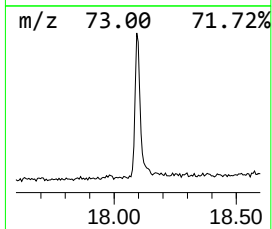
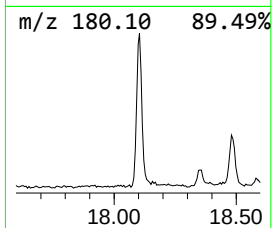
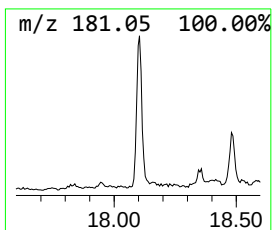
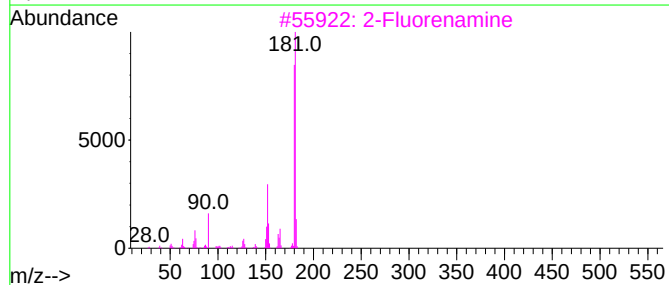
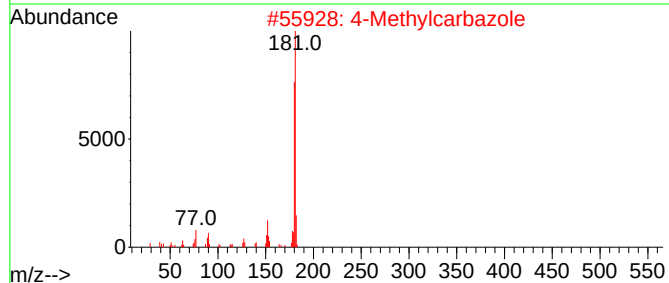
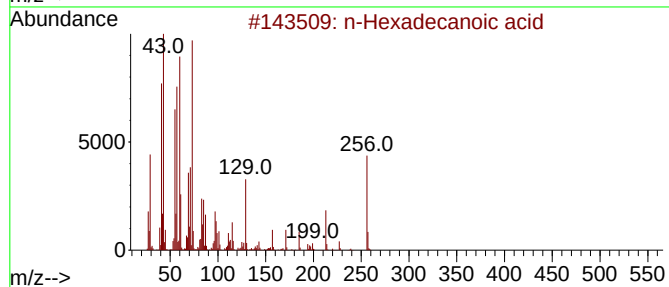
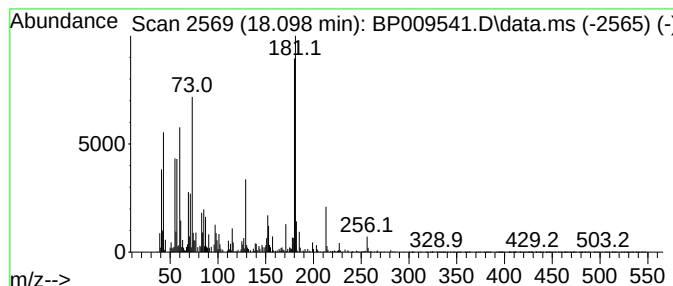
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 19 n-Hexadecanoic acid Concentration Rank 12

| R.T.      | EstConc                 | Area   | Relative to ISTD |          | R.T.        |      |
|-----------|-------------------------|--------|------------------|----------|-------------|------|
| 18.098    | 3.67 ng                 | 628357 | Phenanthrene-d10 |          | 17.180      |      |
| Hit# of 5 | Tentative ID            |        | MW               | MolForm  | CAS#        | Qual |
| 1         | n-Hexadecanoic acid     |        | 256              | C16H32O2 | 000057-10-3 | 91   |
| 2         | 4-Methylcarbazole       |        | 181              | C13H11N  | 003770-48-7 | 87   |
| 3         | 2-Fluorenamine          |        | 181              | C13H11N  | 000153-78-6 | 60   |
| 4         | 1-Aminofluorene         |        | 181              | C13H11N  | 006344-63-4 | 60   |
| 5         | 9H-Carbazole, 9-methyl- |        | 181              | C13H11N  | 001484-12-4 | 55   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
Data File : BP009541.D  
Acq On : 24 Mar 2022 21:49  
Operator : CG/JU  
Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

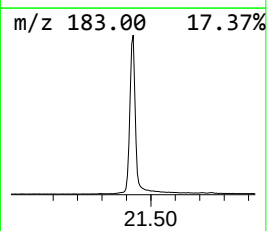
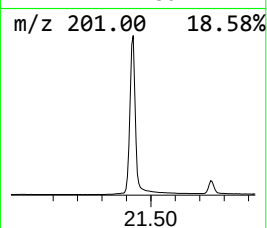
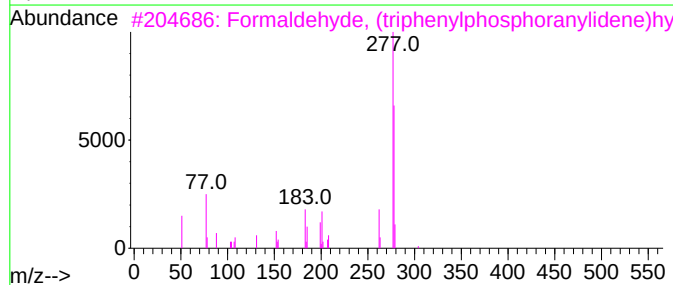
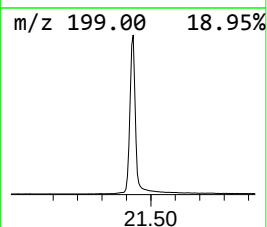
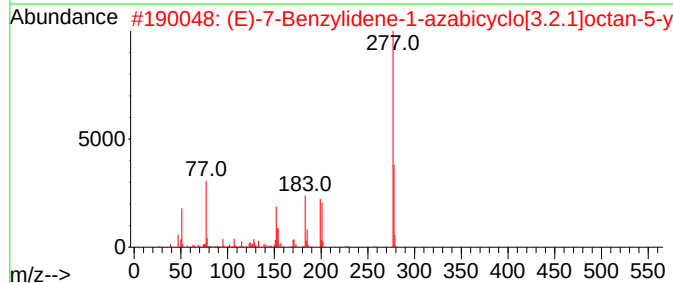
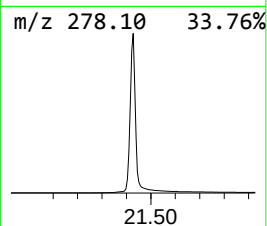
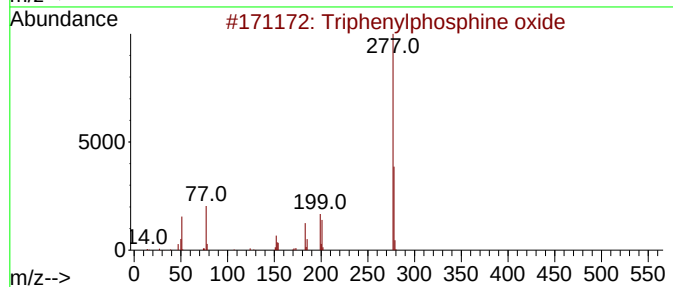
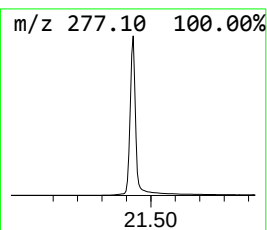
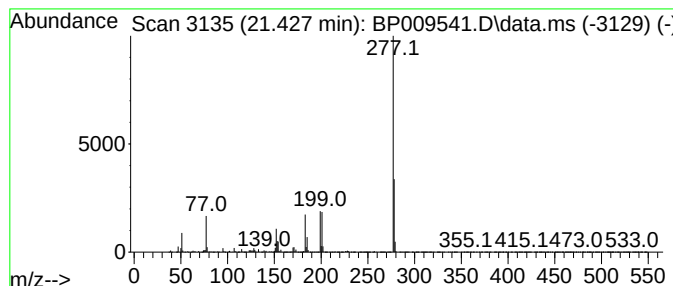
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 20 Triphenylphosphine oxide Concentration Rank 1

| R.T.   | EstConc   | Area     | Relative to ISTD | R.T.   |
|--------|-----------|----------|------------------|--------|
| 21.427 | 106.93 ng | 20714500 | Chrysene-d12     | 21.292 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm     | CAS#         | Qual |
|------|----|---|--------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Triphenylphosphine oxide             | 278 | C18H15OP    | 000791-28-6  | 94   |
| 2    |    |   | (E)-7-Benzylidene-1-azabicyclo[3...] | 293 | C15H19NO3S  | 1000483-83-4 | 91   |
| 3    |    |   | Formaldehyde, (triphenylphosphor...  | 304 | C19H17N2P   | 015990-54-2  | 47   |
| 4    |    |   | 5-Methyl-8-phenylfuro[3',2':5,6]...  | 277 | C16H11N3O2  | 1000460-07-9 | 28   |
| 5    |    |   | Thiophene-3-carbonitrile, 4-(4-m...  | 277 | C13H11NO2S2 | 1000268-75-0 | 9    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032422\  
Data File : BP009541.D  
Acq On : 24 Mar 2022 21:49  
Operator : CG/JU  
Sample : N2090-03  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-48

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.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... | 3.022  | 26.4    | ng    | 2096920  | 1                     | 7.769  | 1586570 | 20.0 |
| Di-sec-Butyl ether | 4.246  | 4.0     | ng    | 315863   | 1                     | 7.769  | 1586570 | 20.0 |
| Acetaldehyde, d... | 4.305  | 4.8     | ng    | 379964   | 1                     | 7.769  | 1586570 | 20.0 |
| Benzene, (1-met... | 6.269  | 8.6     | ng    | 680884   | 1                     | 7.769  | 1586570 | 20.0 |
| Indane             | 8.146  | 36.5    | ng    | 2897410  | 1                     | 7.769  | 1586570 | 20.0 |
| Benzene, 1,3-di... | 8.269  | 4.9     | ng    | 388640   | 1                     | 7.769  | 1586570 | 20.0 |
| Benzene, 1,2,4,... | 9.399  | 7.6     | ng    | 984851   | 2                     | 10.563 | 2583000 | 20.0 |
| Indan, 1-methyl-   | 9.816  | 2.6     | ng    | 332360   | 2                     | 10.563 | 2583000 | 20.0 |
| Benzene, 2-ethe... | 9.969  | 15.1    | ng    | 1953590  | 2                     | 10.563 | 2583000 | 20.0 |
| unknown10.616      | 10.616 | 2.5     | ng    | 327965   | 2                     | 10.563 | 2583000 | 20.0 |
| Naphthalene, 1,... | 11.146 | 2.8     | ng    | 366107   | 2                     | 10.563 | 2583000 | 20.0 |
| 1H-Indene, 2,3-... | 11.481 | 2.4     | ng    | 307793   | 2                     | 10.563 | 2583000 | 20.0 |
| Benzene, (3-met... | 11.681 | 3.5     | ng    | 448754   | 2                     | 10.563 | 2583000 | 20.0 |
| Naphthalene, 1,... | 12.122 | 3.0     | ng    | 383707   | 2                     | 10.563 | 2583000 | 20.0 |
| Tricyclo[5.2.1.... | 12.251 | 2.9     | ng    | 379053   | 2                     | 10.563 | 2583000 | 20.0 |
| Naphthalene, 1,... | 13.740 | 5.0     | ng    | 768004   | 3                     | 14.416 | 3071380 | 20.0 |
| Naphthalene, 1,... | 14.140 | 3.6     | ng    | 556654   | 3                     | 14.416 | 3071380 | 20.0 |
| n-Hexadecanoic ... | 18.098 | 3.7     | ng    | 628357   | 4                     | 17.180 | 3426750 | 20.0 |
| Triphenylphosph... | 21.427 | 106.9   | ng    | 20714500 | 5                     | 21.292 | 3874490 | 20.0 |