

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
 Data File : BP009065.D  
 Acq On : 08 Feb 2022 18:51  
 Operator : CG/JU  
 Sample : N1399-01  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 KTS1-GW-MH-02-8-13

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-BP020722.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP009065.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         | 5.393       | 403           | 409         | 424          | rBV      | 2309826        | 3875457       | 36.79%          | 5.524%        |
| 2         | 5.687       | 451           | 459         | 472          | rBV2     | 119331         | 288143        | 2.74%           | 0.411%        |
| 3         | 6.064       | 517           | 523         | 528          | rBV      | 263062         | 444824        | 4.22%           | 0.634%        |
| 4         | 6.358       | 566           | 573         | 578          | rBV      | 181784         | 352260        | 3.34%           | 0.502%        |
| 5         | 6.493       | 591           | 596         | 608          | rVB3     | 217403         | 484050        | 4.59%           | 0.690%        |
| 6         | 6.617       | 608           | 617         | 623          | rBV3     | 247219         | 482740        | 4.58%           | 0.688%        |
| 7         | 6.699       | 623           | 631         | 638          | rBV7     | 93815          | 297477        | 2.82%           | 0.424%        |
| 8         | 6.840       | 647           | 655         | 661          | rVB4     | 232094         | 548108        | 5.20%           | 0.781%        |
| 9         | 6.987       | 669           | 680         | 689          | rBV      | 1580836        | 3141914       | 29.82%          | 4.478%        |
| 10        | 7.105       | 693           | 700         | 704          | rBV3     | 257244         | 576046        | 5.47%           | 0.821%        |
| 11        | 7.187       | 710           | 714         | 718          | rVB      | 192314         | 274728        | 2.61%           | 0.392%        |
| 12        | 7.240       | 718           | 723         | 726          | rBV3     | 156961         | 303604        | 2.88%           | 0.433%        |
| 13        | 7.311       | 726           | 735         | 739          | rVB5     | 573859         | 1353368       | 12.85%          | 1.929%        |
| 14        | 7.393       | 739           | 749         | 756          | rVB4     | 627447         | 1660676       | 15.76%          | 2.367%        |
| 15        | 7.481       | 756           | 764         | 769          | rVB6     | 161814         | 358752        | 3.41%           | 0.511%        |
| 16        | 7.540       | 769           | 774         | 778          | rBV2     | 305481         | 560903        | 5.32%           | 0.799%        |
| 17        | 7.634       | 782           | 790         | 796          | rBV5     | 456920         | 1204705       | 11.44%          | 1.717%        |
| 18        | 7.728       | 800           | 806         | 812          | rBV2     | 1036830        | 2078305       | 19.73%          | 2.962%        |
| 19        | 7.805       | 813           | 819         | 827          | rBV2     | 919119         | 2362590       | 22.43%          | 3.367%        |
| 20        | 7.875       | 827           | 831         | 836          | rVB4     | 308782         | 444275        | 4.22%           | 0.633%        |
| 21        | 7.934       | 836           | 841         | 843          | rBV3     | 261958         | 446461        | 4.24%           | 0.636%        |
| 22        | 7.975       | 844           | 848         | 855          | rVB3     | 828806         | 1738719       | 16.50%          | 2.478%        |
| 23        | 8.040       | 855           | 859         | 862          | rBV3     | 402563         | 710169        | 6.74%           | 1.012%        |
| 24        | 8.087       | 862           | 867         | 870          | rBV4     | 584785         | 1114044       | 10.57%          | 1.588%        |
| 25        | 8.175       | 877           | 882         | 886          | rBV4     | 420594         | 756215        | 7.18%           | 1.078%        |
| 26        | 8.216       | 886           | 889         | 893          | rVB      | 318091         | 432278        | 4.10%           | 0.616%        |
| 27        | 8.287       | 893           | 901         | 906          | rBV5     | 269108         | 669270        | 6.35%           | 0.954%        |
| 28        | 8.363       | 910           | 914         | 917          | rBV3     | 240885         | 332149        | 3.15%           | 0.473%        |
| 29        | 8.440       | 924           | 927         | 929          | rBV      | 189088         | 267122        | 2.54%           | 0.381%        |
| 30        | 8.569       | 940           | 949         | 960          | rBV10    | 488376         | 2398940       | 22.77%          | 3.419%        |
| 31        | 8.681       | 962           | 968         | 973          | rBV2     | 688124         | 1605429       | 15.24%          | 2.288%        |
| 32        | 8.869       | 994           | 1000        | 1002         | rBV4     | 947603         | 1855740       | 17.62%          | 2.645%        |
| 33        | 8.969       | 1011          | 1017        | 1027         | rVB      | 2599741        | 5042449       | 47.87%          | 7.187%        |
| 34        | 9.169       | 1045          | 1051        | 1057         | rVB5     | 1081387        | 2549540       | 24.20%          | 3.634%        |
| 35        | 9.281       | 1065          | 1070        | 1075         | rVB5     | 711117         | 1378115       | 13.08%          | 1.964%        |
| 36        | 9.340       | 1076          | 1080        | 1084         | rBV4     | 602120         | 1068989       | 10.15%          | 1.524%        |

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Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

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-BP020722.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |          |         |         |
|----|--------|------|------|------|------|---------|----------|---------|---------|
| 37 | 9.475  | 1098 | 1103 | 1108 | rVB  | 1575064 | 2747851  | 26.08%  | 3.917%  |
| 38 | 9.658  | 1129 | 1134 | 1135 | rBV2 | 1296141 | 1933521  | 18.35%  | 2.756%  |
| 39 | 9.787  | 1151 | 1156 | 1166 | rBV5 | 1071820 | 3486122  | 33.09%  | 4.969%  |
| 40 | 10.487 | 1271 | 1275 | 1280 | rBV  | 1197536 | 2141143  | 20.32%  | 3.052%  |
| 41 | 11.116 | 1377 | 1382 | 1388 | rBV2 | 5018100 | 10534731 | 100.00% | 15.015% |
| 42 | 11.822 | 1498 | 1502 | 1510 | rVB  | 3217220 | 5858773  | 55.61%  | 8.351%  |

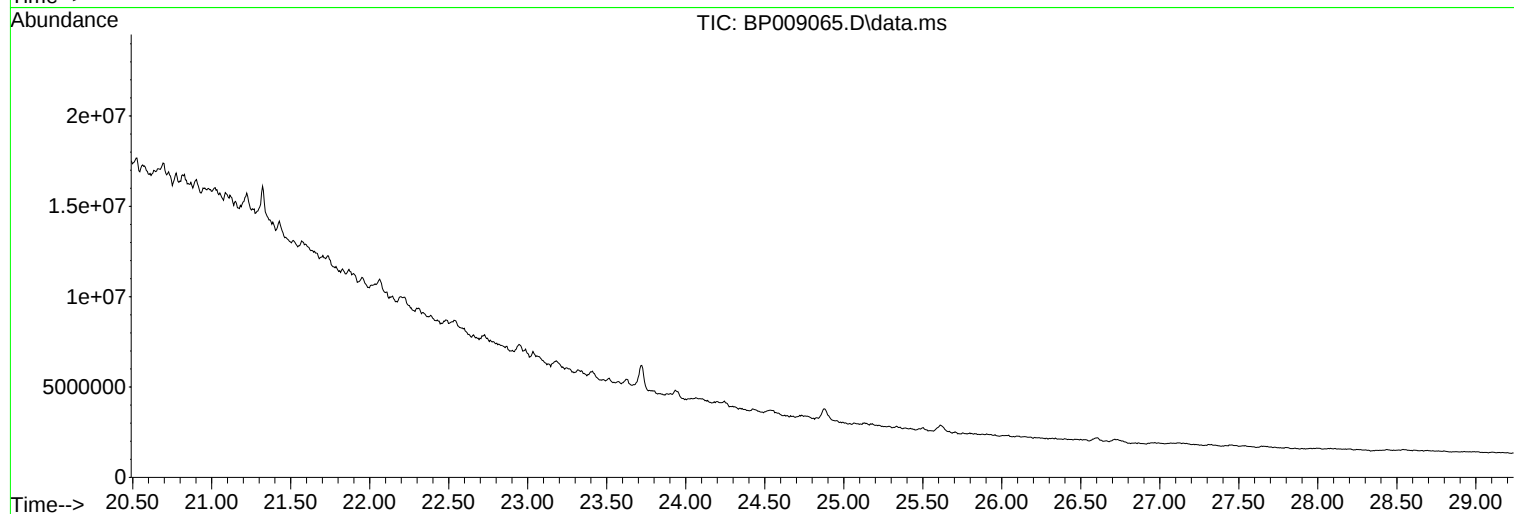
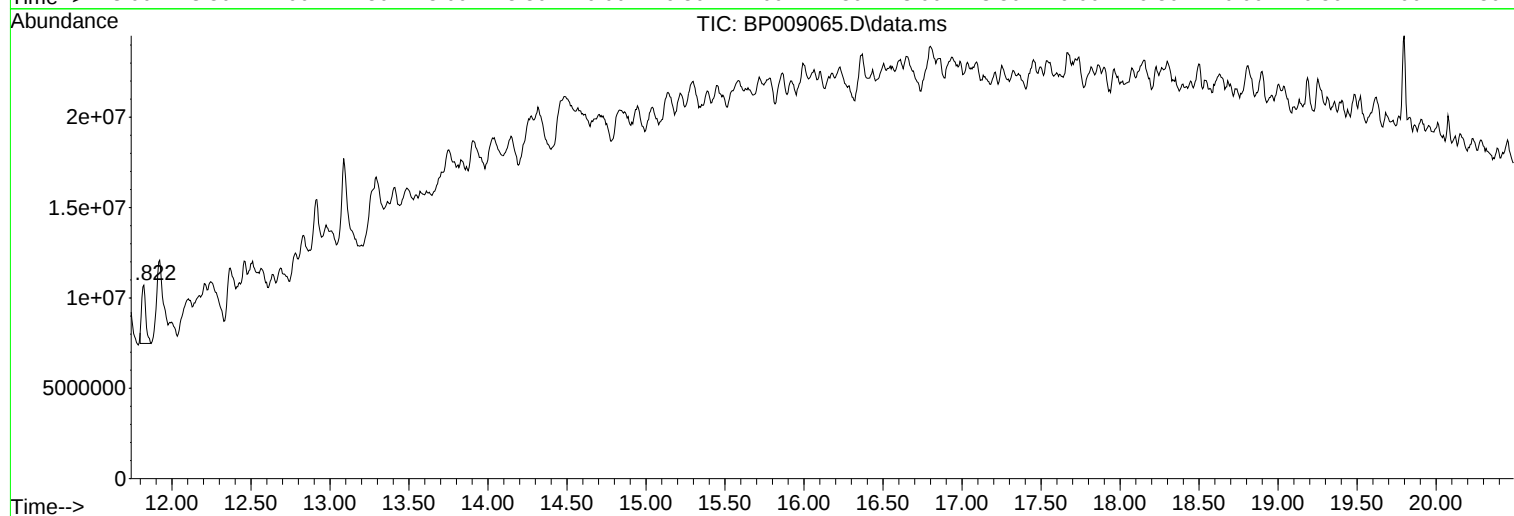
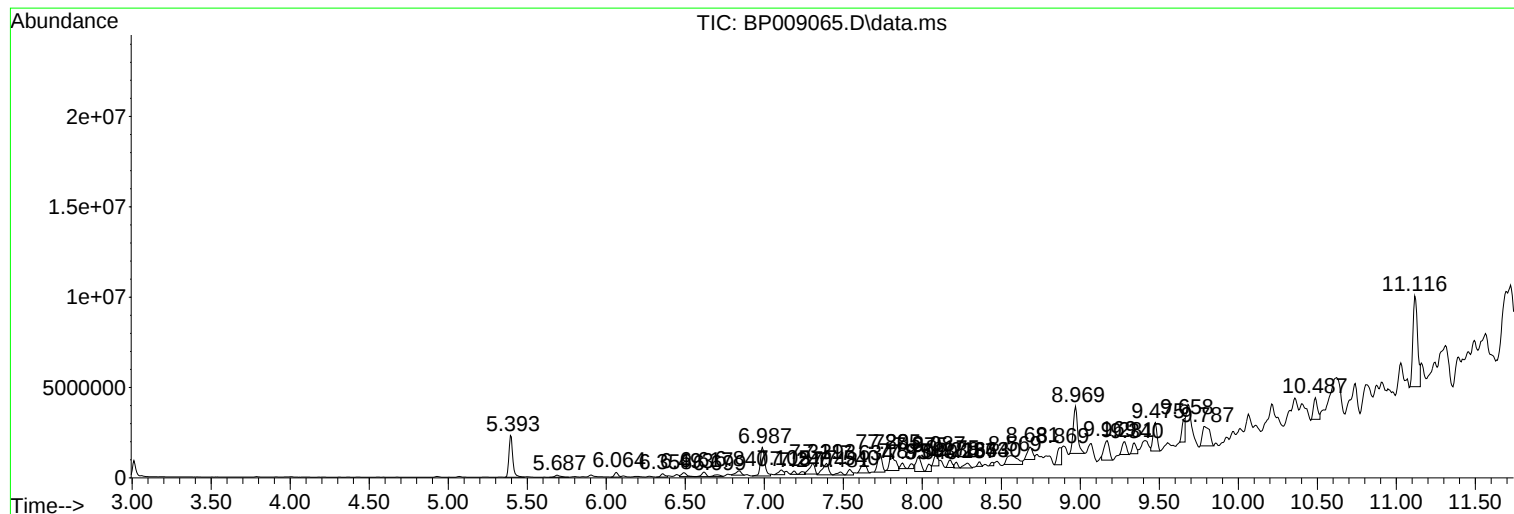
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Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
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Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP020722.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\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP020722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

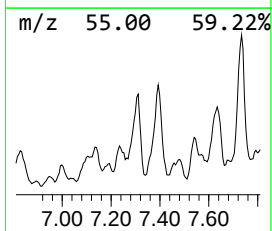
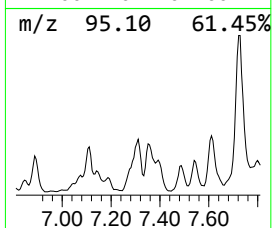
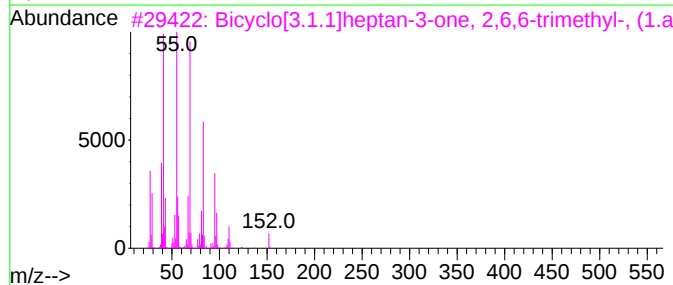
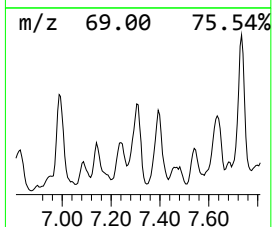
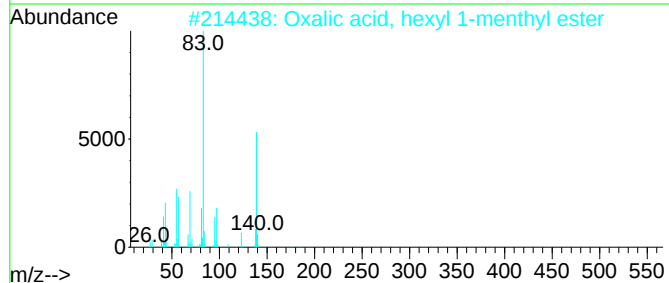
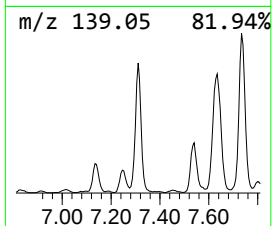
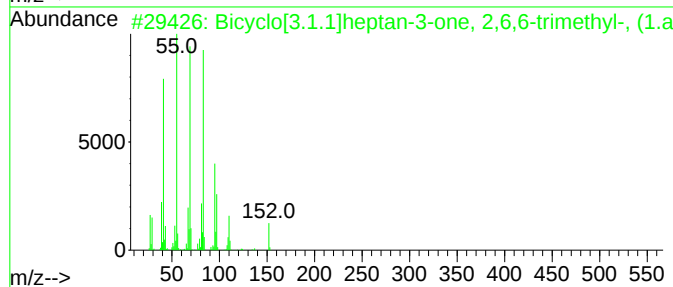
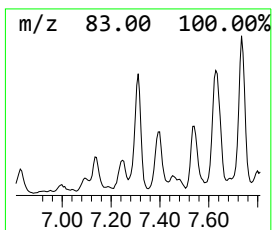
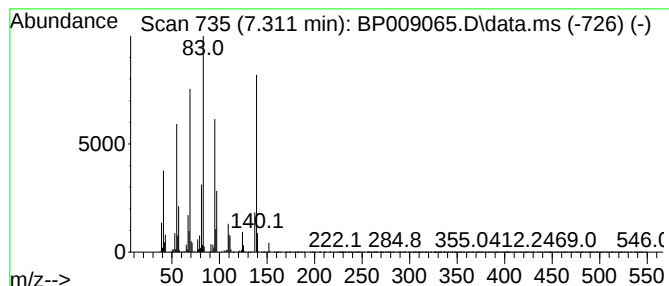
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Bicyclo[3.1.1]heptan-3-one,... Concentration Rank 15

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 7.311 | 11.46 ng | 1353370 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Bicyclo[3.1.1]heptan-3-one, 2,6,... | 152 | C10H16O   | 000547-60-4  | 58   |
| 2    |    |   | Oxalic acid, hexyl 1-menthyl ester  | 312 | C18H32O4  | 1000314-97-4 | 43   |
| 3    |    |   | Bicyclo[3.1.1]heptan-3-one, 2,6,... | 152 | C10H16O   | 015358-88-0  | 43   |
| 4    |    |   | Oxalic acid, heptadecyl 1-menthy... | 466 | C29H54O4  | 1000314-98-4 | 38   |
| 5    |    |   | Cyclohexanol, 5-methyl-2-(1-meth... | 358 | C20H38O3S | 026510-92-9  | 38   |



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ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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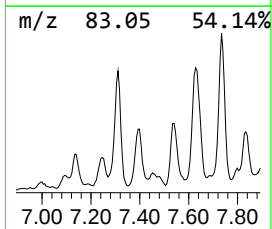
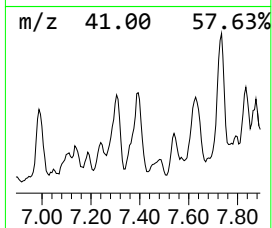
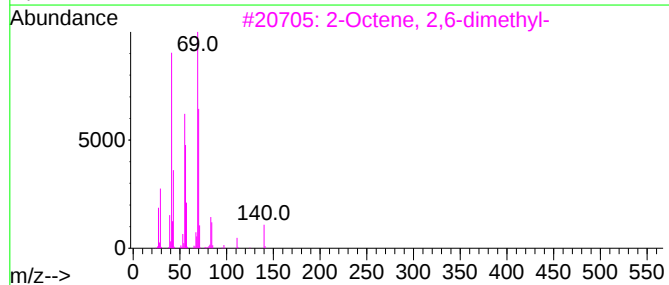
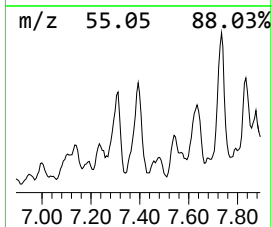
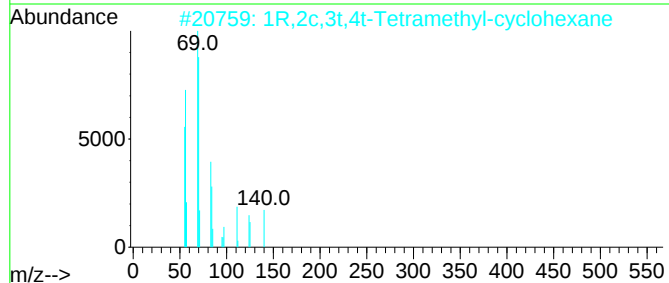
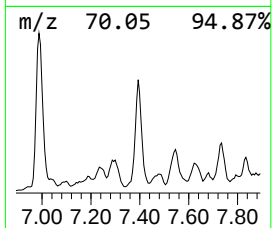
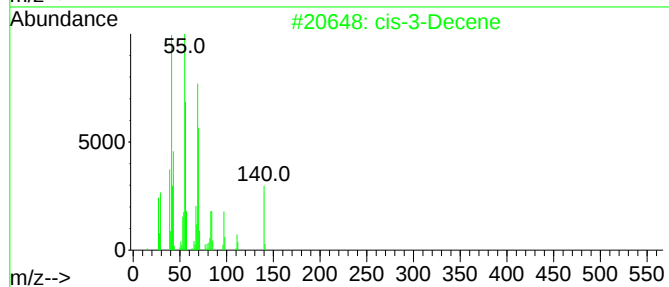
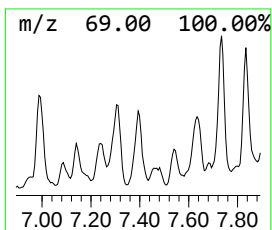
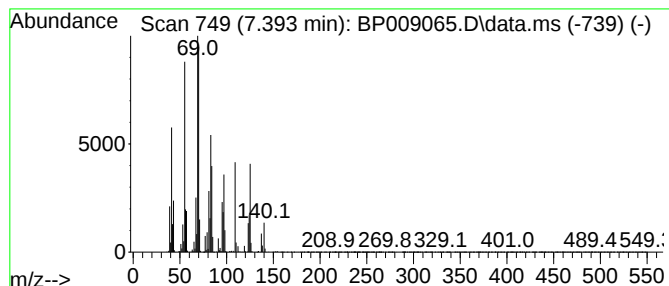
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP020722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 2 cis-3-Decene Concentration Rank 12

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 7.393 | 14.06 ng | 1660680 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | cis-3-Decene                        | 140 | C10H20  | 019398-86-8  | 53   |
| 2    |    |   | 1R,2c,3t,4t-Tetramethyl-cyclohexane | 140 | C10H20  | 1000144-07-3 | 47   |
| 3    |    |   | 2-Octene, 2,6-dimethyl-             | 140 | C10H20  | 004057-42-5  | 43   |
| 4    |    |   | Cyclopentane, 1,1,3,4-tetramethy... | 126 | C9H18   | 053907-60-1  | 35   |
| 5    |    |   | Cyclopropane, 1,1,2-trimethyl-3-... | 140 | C10H20  | 041977-43-9  | 25   |



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Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

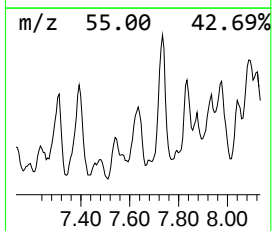
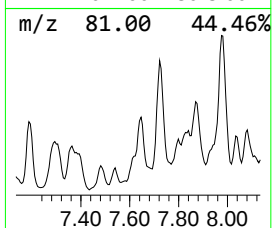
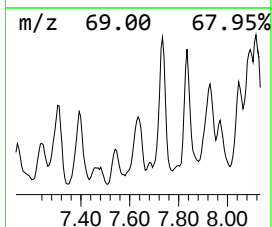
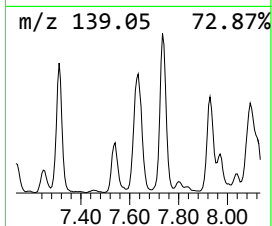
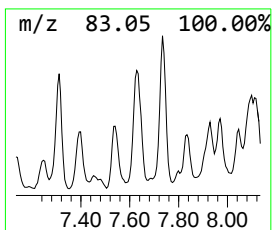
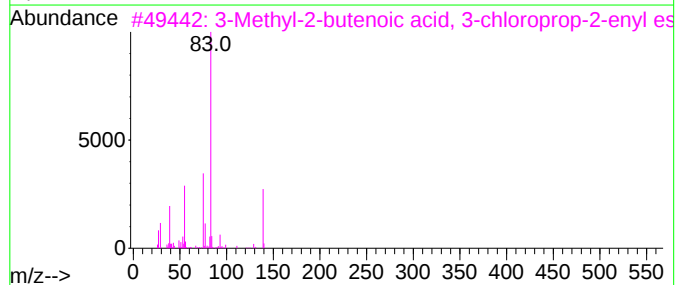
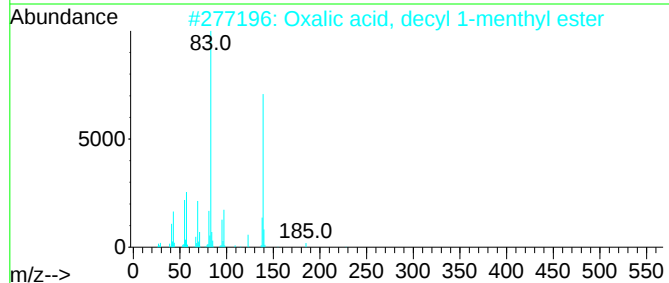
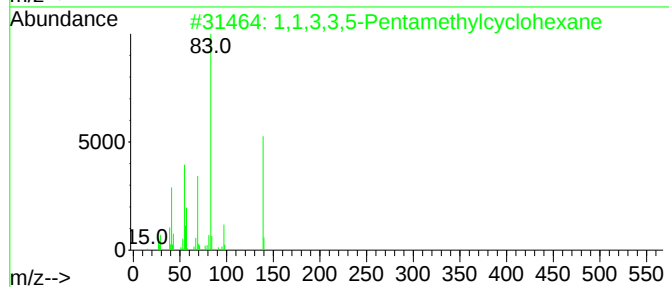
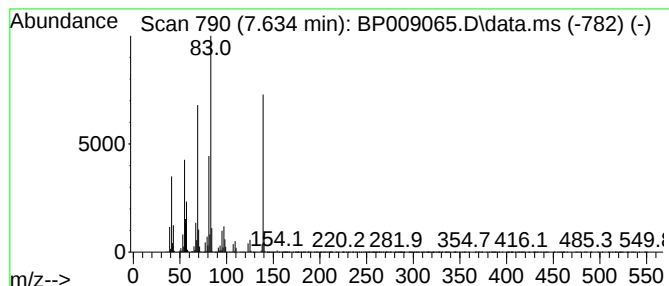
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP020722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 3 1,1,3,3,5-Pentamethylcycloh... Concentration Rank 16

| R.T.      | EstConc                             | Area    | Relative to ISTD       | R.T.         |      |
|-----------|-------------------------------------|---------|------------------------|--------------|------|
| 7.634     | 10.20 ng                            | 1204710 | 1,4-Dichlorobenzene-d4 | 7.805        |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm                | CAS#         | Qual |
| 1         | 1,1,3,3,5-Pentamethylcyclohexane    | 154     | C11H22                 | 070810-19-4  | 72   |
| 2         | Oxalic acid, decyl 1-menthyl ester  | 368     | C22H40O4               | 1000314-97-8 | 72   |
| 3         | 3-Methyl-2-butenic acid, 3-chlo...  | 174     | C8H11ClO2              | 1000299-31-8 | 59   |
| 4         | Oxalic acid, dodecyl 1-menthyl e... | 396     | C24H44O4               | 1000314-98-0 | 50   |
| 5         | Cyclohexanol, 5-methyl-2-(1-meth... | 358     | C20H38O3S              | 026510-92-9  | 50   |



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BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

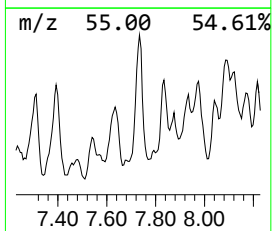
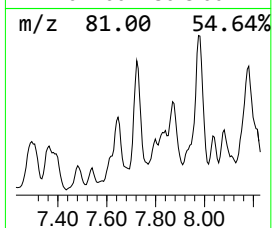
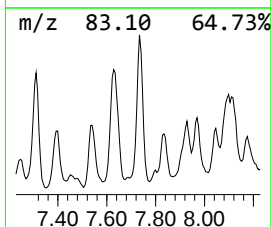
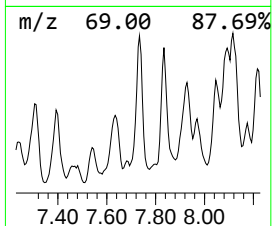
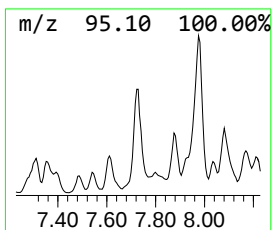
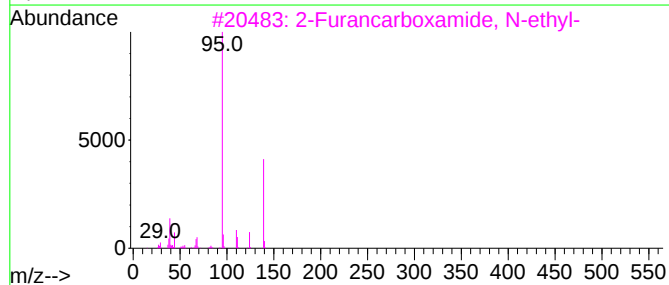
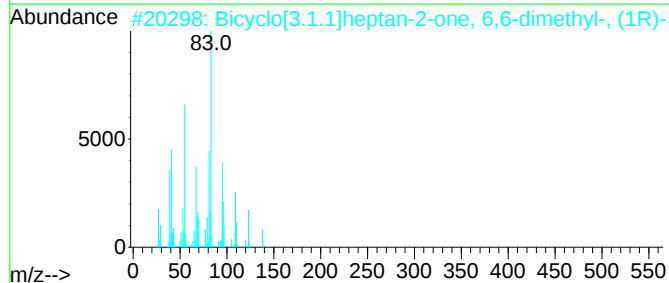
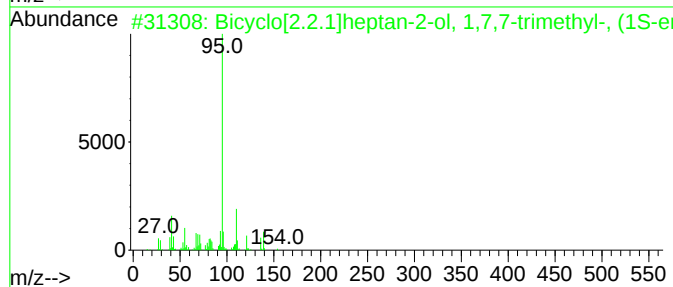
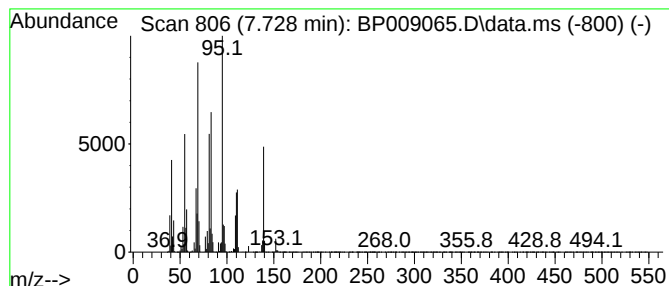
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 4 unknown7.728 Concentration Rank 9

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 7.728 | 17.59 ng | 2078310 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Bicyclo[2.2.1]heptan-2-ol, 1,7,7... | 154 | C10H18O  | 000464-45-9  | 38   |
| 2    |    |   | Bicyclo[3.1.1]heptan-2-one, 6,6-... | 138 | C9H14O   | 038651-65-9  | 38   |
| 3    |    |   | 2-Furancarboxamide, N-ethyl-        | 139 | C7H9NO2  | 1000407-23-8 | 35   |
| 4    |    |   | Cyclopentanecarboxylic acid, 3-m... | 278 | C18H30O2 | 1000152-25-8 | 27   |
| 5    |    |   | Cyclopentane, 1,2-dimethyl-3-met... | 110 | C8H14    | 1000150-99-3 | 27   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP020722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

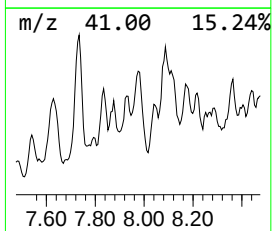
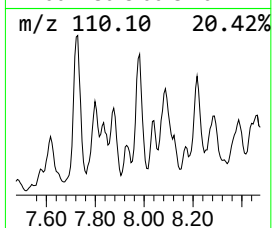
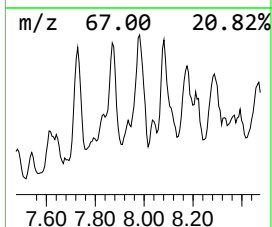
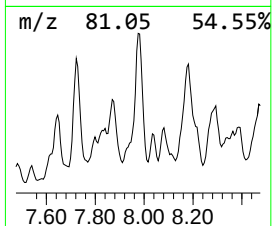
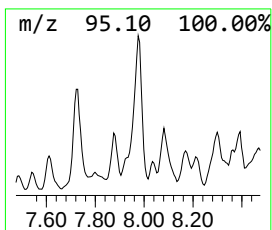
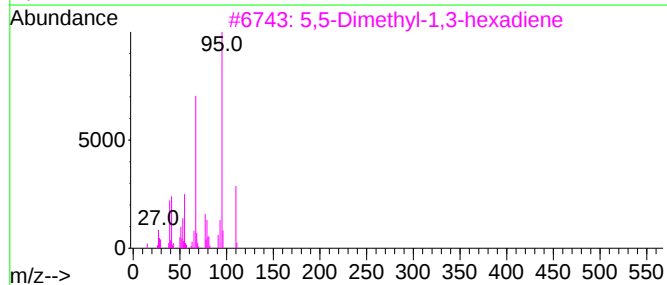
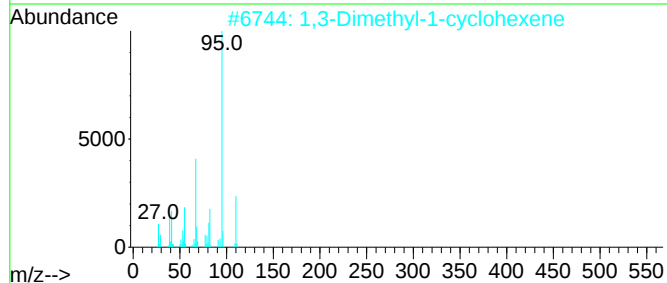
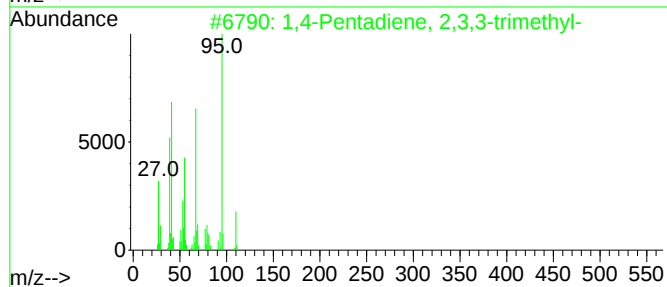
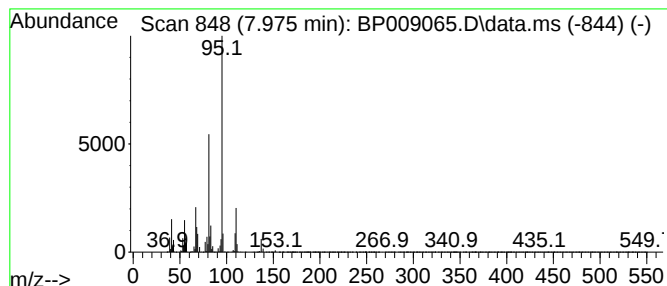
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 1,4-Pentadiene, 2,3,3-trime... Concentration Rank 11

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 7.975 | 14.72 ng | 1738720 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | 1,4-Pentadiene, 2,3,3-trimethyl-    | 110 | C8H14   | 000756-02-5  | 76   |
| 2    |    |   | 1,3-Dimethyl-1-cyclohexene          | 110 | C8H14   | 002808-76-6  | 64   |
| 3    |    |   | 5,5-Dimethyl-1,3-hexadiene          | 110 | C8H14   | 001515-79-3  | 59   |
| 4    |    |   | 1,4-Hexadiene, 2,3-dimethyl-        | 110 | C8H14   | 018669-52-8  | 53   |
| 5    |    |   | Bicyclo[3.1.0]hexane, 1,5-dimethyl- | 110 | C8H14   | 1010142-17-5 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

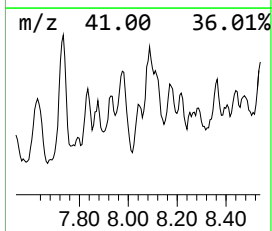
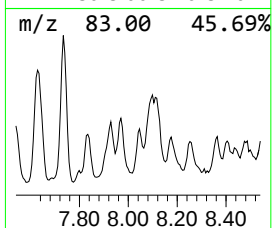
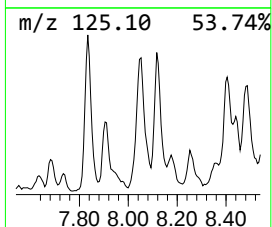
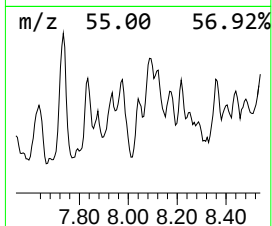
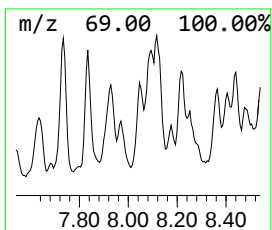
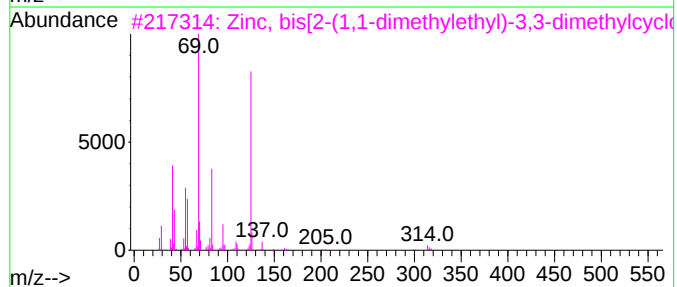
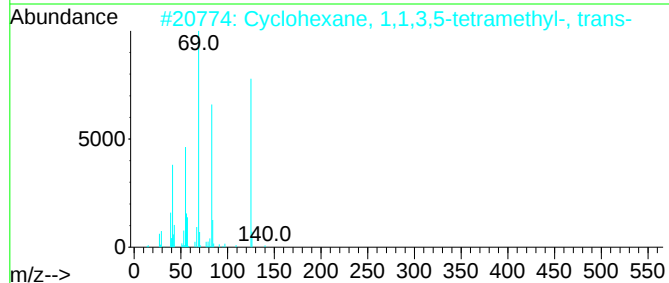
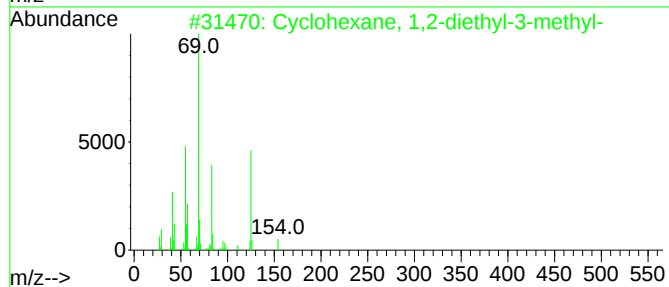
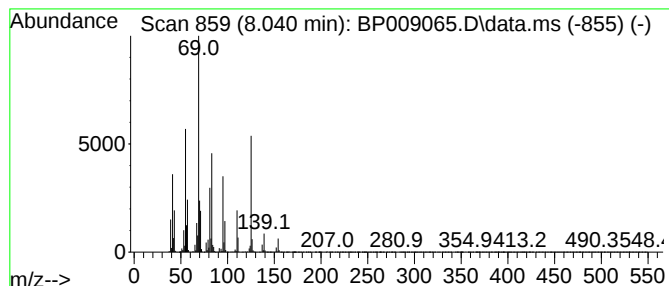
Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP020722.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 6 Cyclohexane, 1,2-diethyl-3-... Concentration Rank 20

| R.T.      | EstConc                             | Area   | Relative to ISTD       | R.T.         |      |
|-----------|-------------------------------------|--------|------------------------|--------------|------|
| 8.040     | 6.01 ng                             | 710169 | 1,4-Dichlorobenzene-d4 | 7.805        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm                | CAS#         | Qual |
| 1         | Cyclohexane, 1,2-diethyl-3-methyl-  | 154    | C11H22                 | 061141-80-8  | 58   |
| 2         | Cyclohexane, 1,1,3,5-tetramethyl... | 140    | C10H20                 | 050876-31-8  | 58   |
| 3         | Zinc, bis[2-(1,1-dimethylethyl)-... | 314    | C18H34Zn               | 074793-36-5  | 53   |
| 4         | Cyclohexane, 2,4-diethyl-1-methyl-  | 154    | C11H22                 | 061142-70-9  | 53   |
| 5         | (1,2,4-Trimethylcyclohexyl)metha... | 198    | C12H22O2               | 1000499-04-5 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP020722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

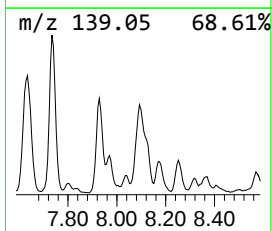
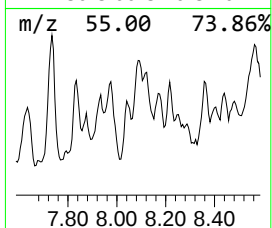
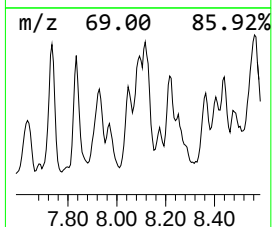
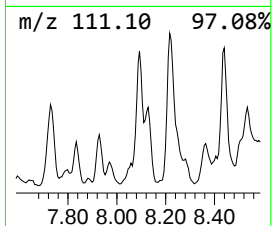
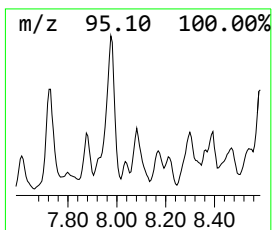
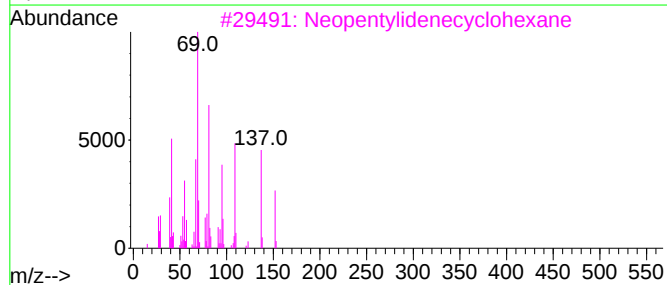
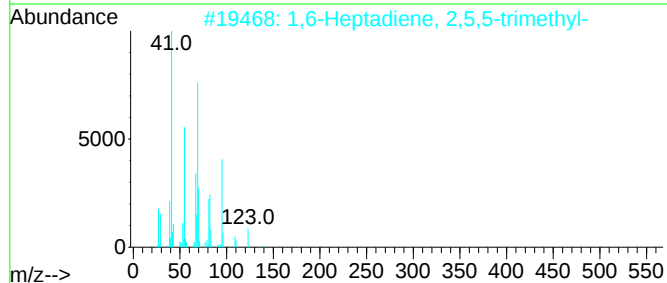
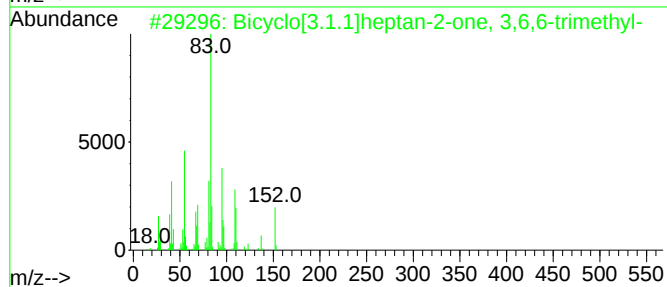
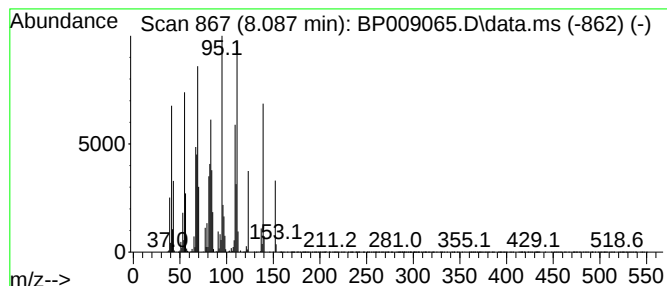
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 unknown8.087 Concentration Rank 18

| R.T.  | EstConc | Area    | Relative to ISTD       | R.T.  |
|-------|---------|---------|------------------------|-------|
| 8.087 | 9.43 ng | 1114040 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Bicyclo[3.1.1]heptan-2-one, 3,6,... | 152 | C10H16O | 016022-08-5 | 42   |
| 2    |    |   | 1,6-Heptadiene, 2,5,5-trimethyl-    | 138 | C10H18  | 062238-28-2 | 41   |
| 3    |    |   | Neopentylidenecyclohexane           | 152 | C11H20  | 039546-80-0 | 30   |
| 4    |    |   | 1-Cyclohexyl-2-methyl-prop-2-en-... | 152 | C10H16O | 025183-82-8 | 25   |
| 5    |    |   | 7-Butyl-3,4,5,6(2H)-tetrahydroaz... | 153 | C10H19N | 003338-06-5 | 18   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

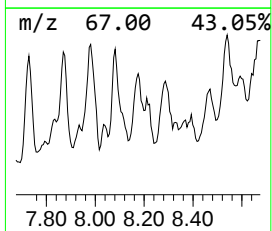
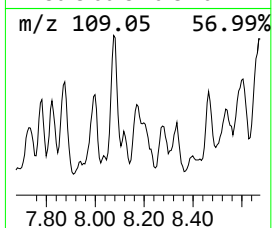
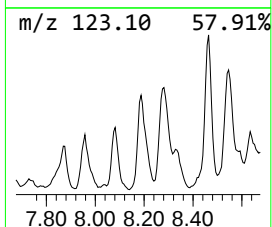
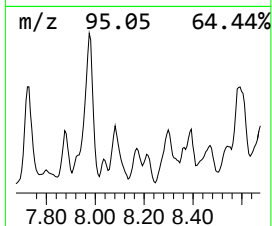
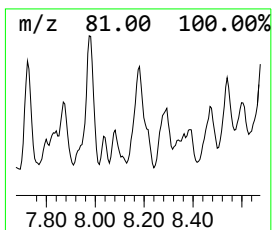
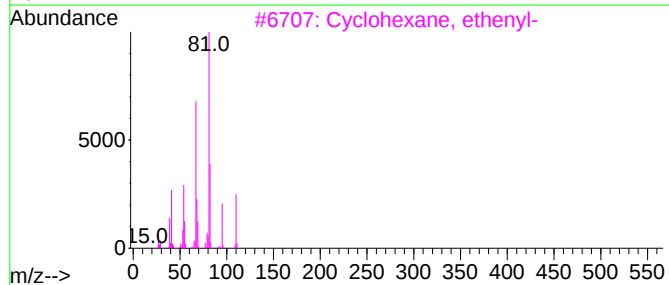
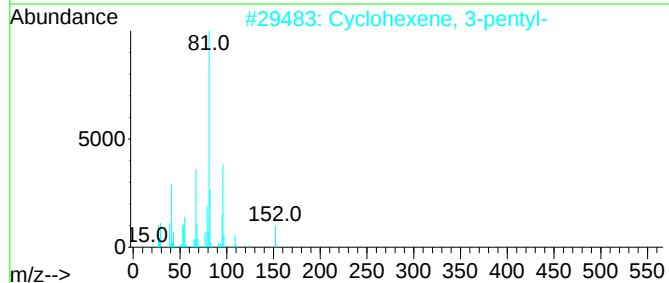
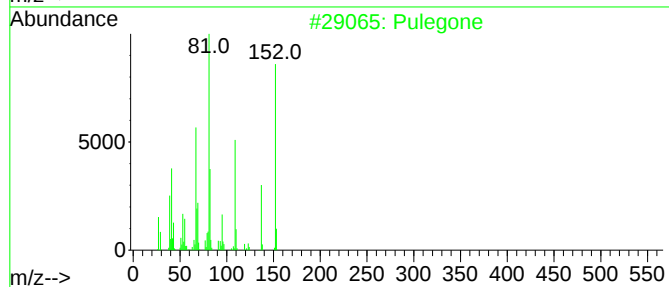
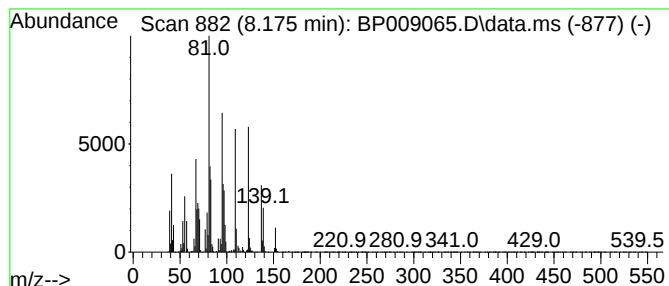
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BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP020722.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 8 Pulegone Concentration Rank 19

| R.T.      | EstConc                             | Area   | Relative to ISTD       |         |             | R.T.  |
|-----------|-------------------------------------|--------|------------------------|---------|-------------|-------|
| 8.175     | 6.40 ng                             | 756215 | 1,4-Dichlorobenzene-d4 |         |             | 7.805 |
| Hit# of 5 | Tentative ID                        |        | MW                     | MolForm | CAS#        | Qual  |
| 1         | Pulegone                            |        | 152                    | C10H16O | 000089-82-7 | 53    |
| 2         | Cyclohexene, 3-pentyl-              |        | 152                    | C11H20  | 015232-92-5 | 38    |
| 3         | Cyclohexane, ethenyl-               |        | 110                    | C8H14   | 000695-12-5 | 35    |
| 4         | Cyclohexanone, 5-methyl-2-(1-met... |        | 152                    | C10H16O | 015932-80-6 | 30    |
| 5         | 5-Eicosyne                          |        | 278                    | C20H38  | 074685-31-7 | 27    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP020722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

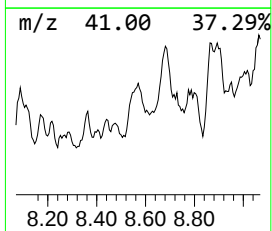
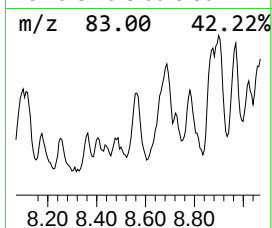
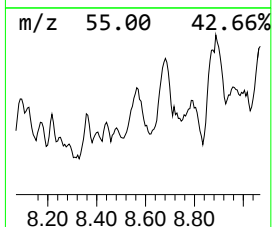
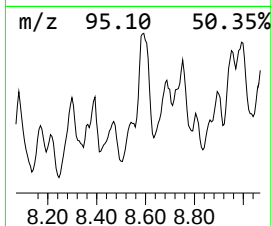
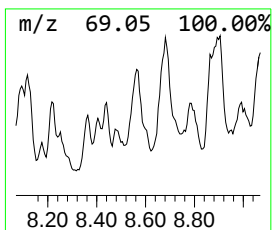
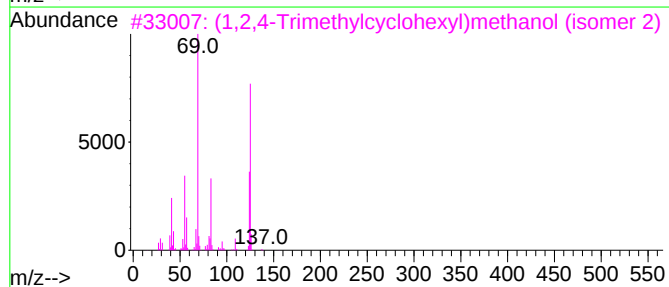
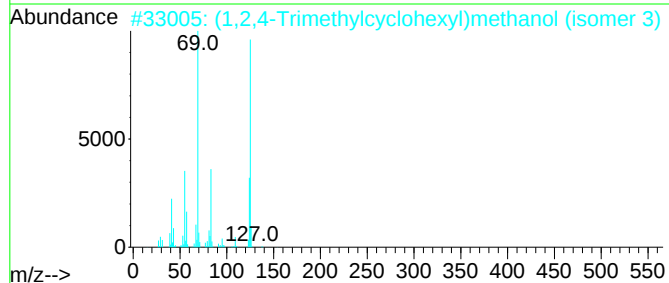
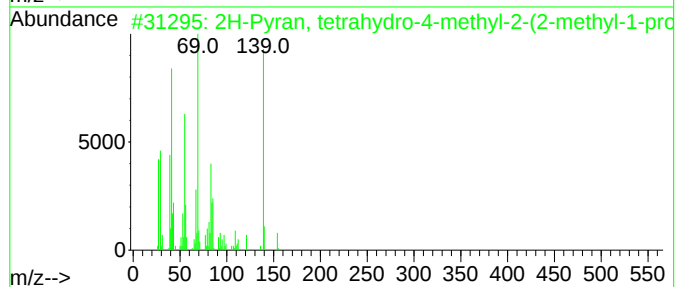
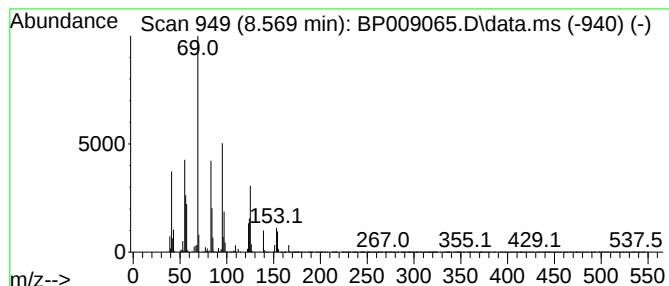
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TIC Integration Parameters: LSCINT.P

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Peak Number 9 unknown8.569 Concentration Rank 6

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 8.569 | 20.31 ng | 2398940 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 2H-Pyran, tetrahydro-4-methyl-2-... | 154 | C10H18O      | 016409-43-1  | 38      |      |      |
| 2    | (1,2,4-Trimethylcyclohexyl)metha... | 156 | C10H20O      | 1000499-04-0 | 38      |      |      |
| 3    | (1,2,4-Trimethylcyclohexyl)metha... | 156 | C10H20O      | 1000499-03-9 | 38      |      |      |
| 4    | 1-Isopropyl-1,4,5-trimethylcyclo... | 168 | C12H24       | 219783-06-9  | 37      |      |      |
| 5    | Cyclooctane, 1-methyl-3-propyl-     | 168 | C12H24       | 255885-37-1  | 37      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

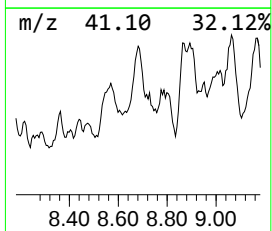
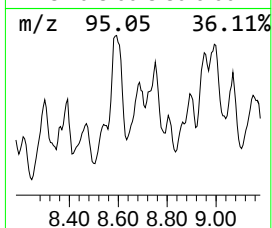
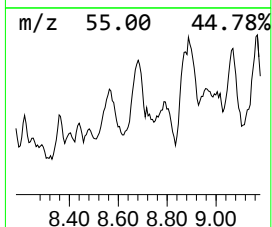
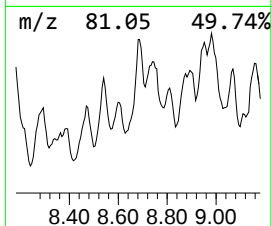
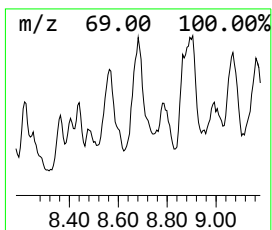
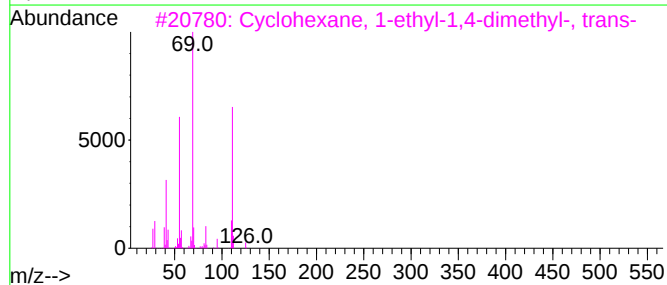
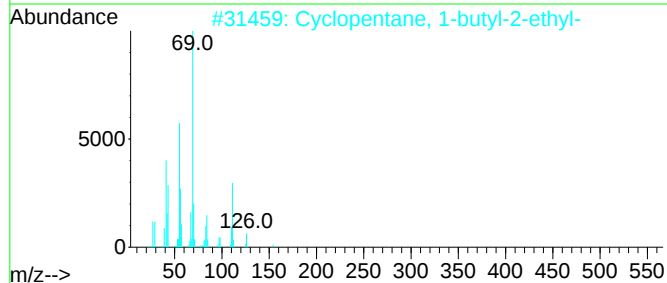
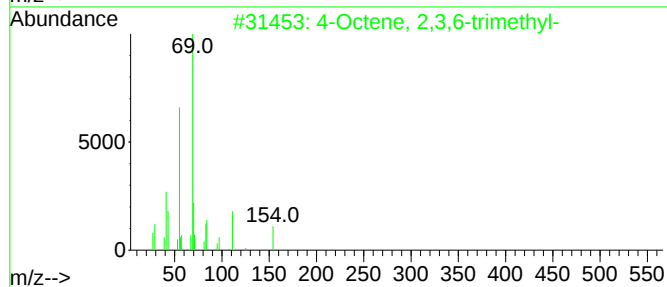
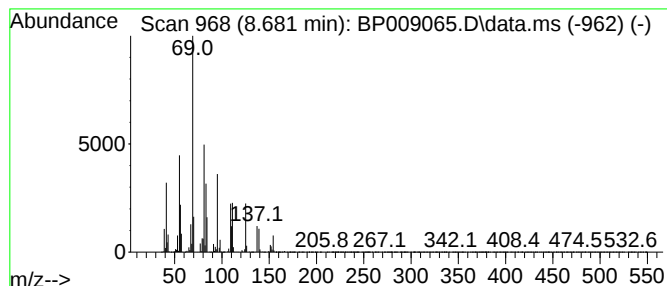
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 10 unknown8.681 Concentration Rank 13

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 8.681 | 13.59 ng | 1605430 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | 4-Octene, 2,3,6-trimethyl-          | 154 | C11H22  | 063830-65-9  | 43   |
| 2    |    |   | Cyclopentane, 1-butyl-2-ethyl-      | 154 | C11H22  | 072993-32-9  | 43   |
| 3    |    |   | Cyclohexane, 1-ethyl-1,4-dimethy... | 140 | C10H20  | 062238-32-8  | 43   |
| 4    |    |   | Cyclohexane, 1-ethyl-1,4-dimethy... | 140 | C10H20  | 062238-30-6  | 43   |
| 5    |    |   | (2,4-Dimethylcyclohexyl)methanol... | 142 | C9H18O  | 1000499-02-8 | 40   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

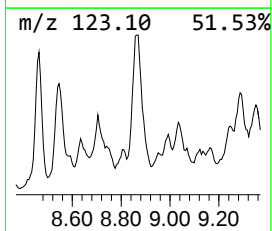
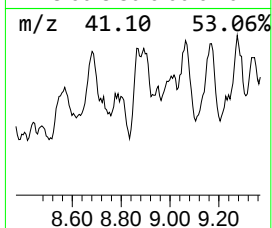
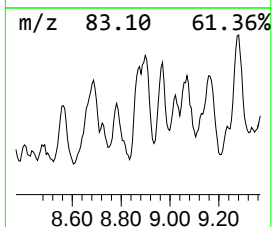
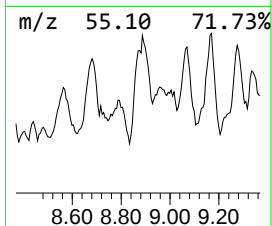
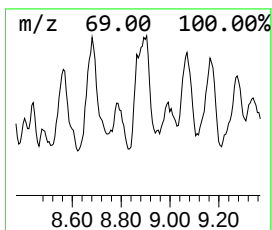
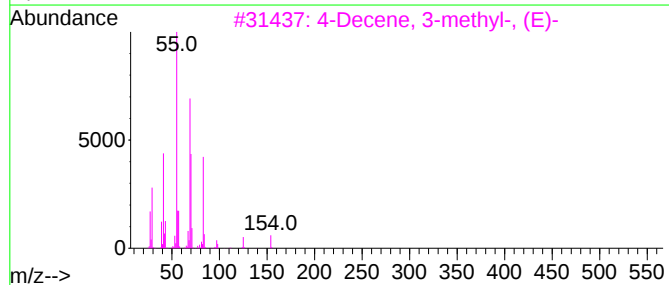
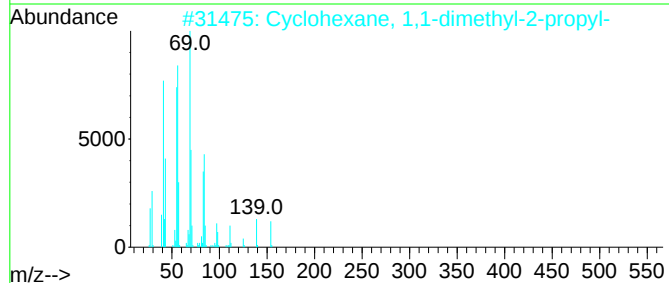
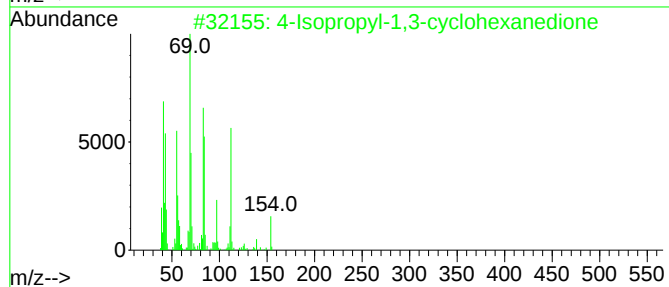
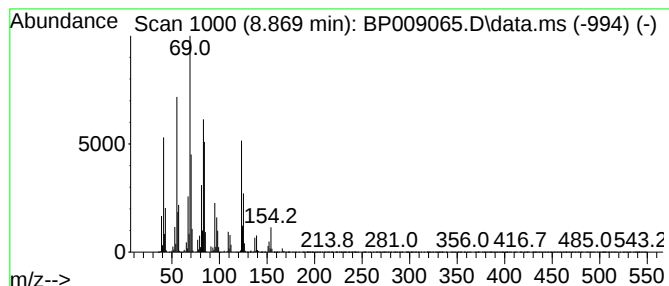
Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP020722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 11 4-Isopropyl-1,3-cyclohexane... Concentration Rank 10

| R.T.      | EstConc                             | Area    | Relative to ISTD       | R.T.        |      |
|-----------|-------------------------------------|---------|------------------------|-------------|------|
| 8.869     | 15.71 ng                            | 1855740 | 1,4-Dichlorobenzene-d4 | 7.805       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm                | CAS#        | Qual |
| 1         | 4-Isopropyl-1,3-cyclohexanedione    | 154     | C9H14O2                | 062831-62-3 | 64   |
| 2         | Cyclohexane, 1,1-dimethyl-2-propyl- | 154     | C11H22                 | 081983-71-3 | 49   |
| 3         | 4-Decene, 3-methyl-, (E)-           | 154     | C11H22                 | 062338-47-0 | 46   |
| 4         | 5-Undecene, (E)-                    | 154     | C11H22                 | 000764-97-6 | 43   |
| 5         | Cyclohexane, (1-ethylpropyl)-       | 154     | C11H22                 | 026321-98-2 | 42   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

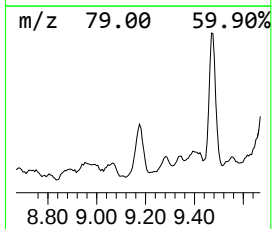
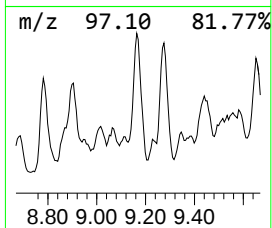
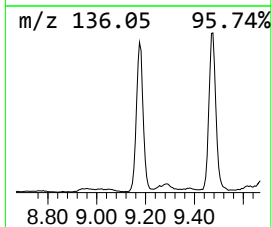
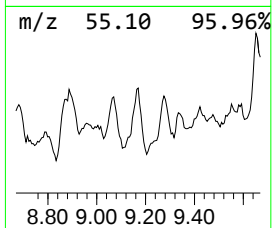
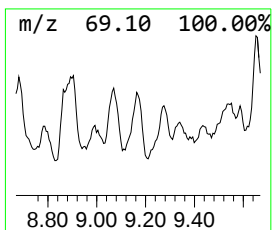
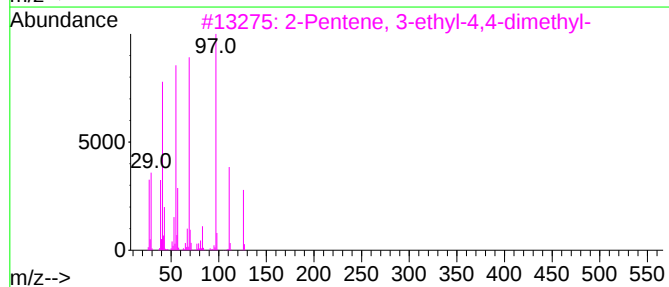
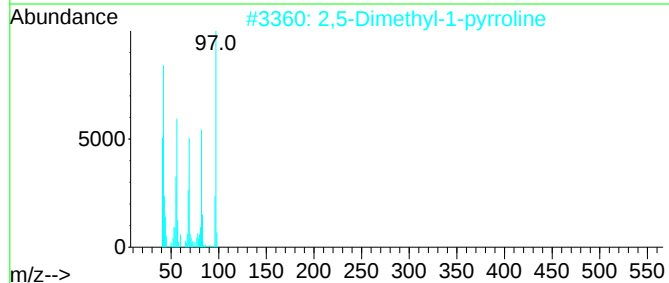
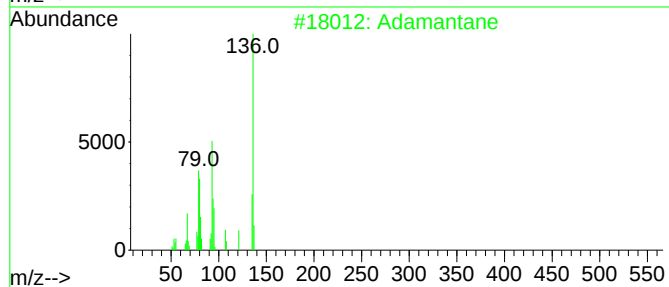
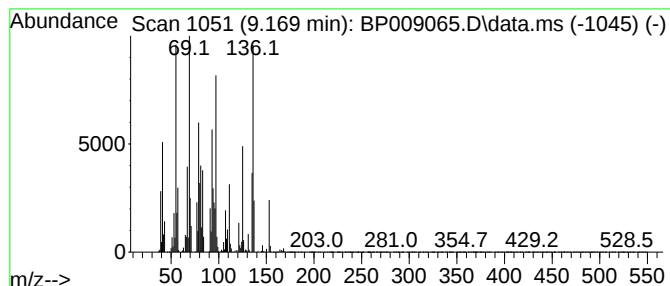
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 12 Adamantane Concentration Rank 5

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 9.169 | 21.58 ng | 2549540 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Adamantane                          | 136 | C10H16   | 000281-23-2 | 53   |
| 2    |    |   | 2,5-Dimethyl-1-pyrroline            | 97  | C6H11N   | 000822-64-0 | 35   |
| 3    |    |   | 2-Pentene, 3-ethyl-4,4-dimethyl-    | 126 | C9H18    | 053907-59-8 | 27   |
| 4    |    |   | 6-Methyl-1,5-diazabicyclo[3.1.0]... | 98  | C5H10N2  | 100463-00-1 | 27   |
| 5    |    |   | 2-Butenoic acid, 2-methyl-, 3-me... | 168 | C10H16O2 | 083783-86-2 | 14   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

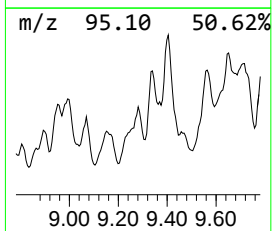
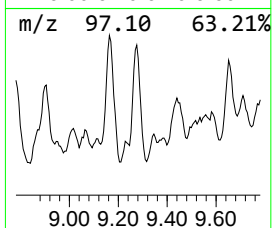
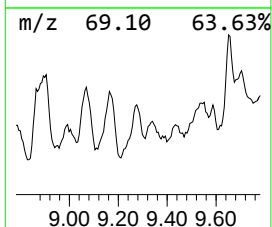
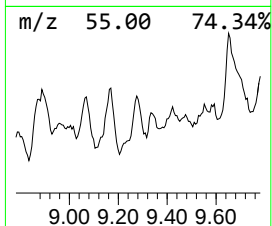
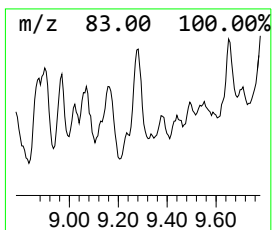
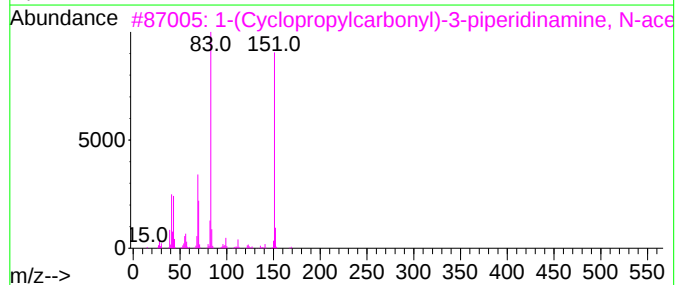
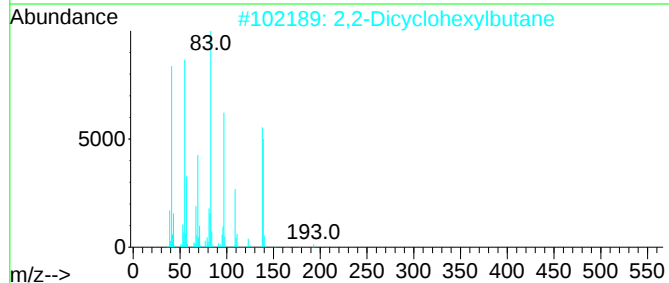
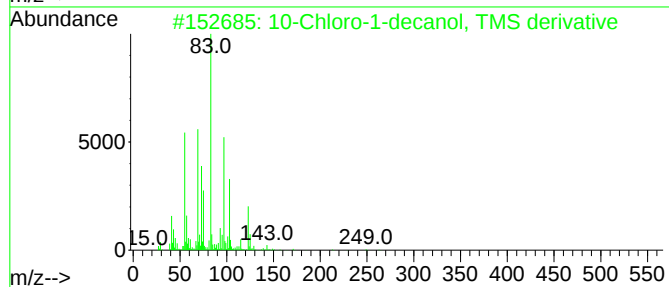
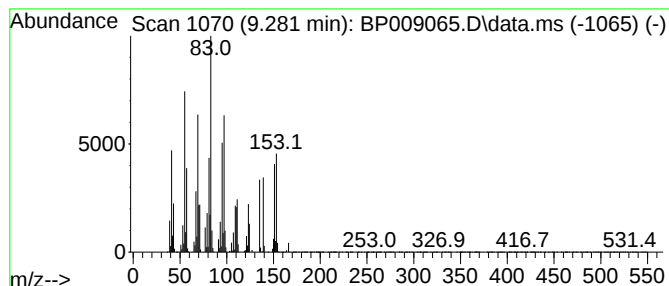
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 13 unknown9.281 Concentration Rank 14

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.   |
|-------|----------|---------|------------------|--------|
| 9.281 | 12.87 ng | 1378120 | Naphthalene-d8   | 10.610 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 10-Chloro-1-decanol, TMS derivative | 264 | C13H29ClOSi  | 034714-01-7  | 27      |      |      |
| 2    | 2,2-Dicyclohexylbutane              | 222 | C16H30       | 054890-02-7  | 27      |      |      |
| 3    | 1-(Cyclopropylcarbonyl)-3-piperi... | 210 | C11H18N2O2   | 1457693-07-0 | 22      |      |      |
| 4    | 2-Isopropyl-5-methylcyclohexylme... | 170 | C11H22O      | 1000223-18-0 | 22      |      |      |
| 5    | Benzene, 2-chloro-1,4-bis(cycloh... | 364 | C20H25ClO4   | 294874-04-7  | 22      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP020722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

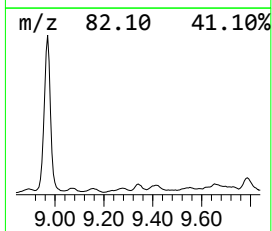
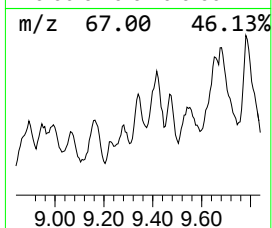
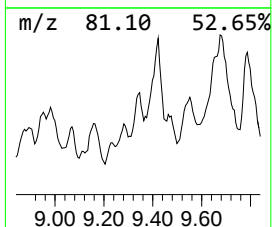
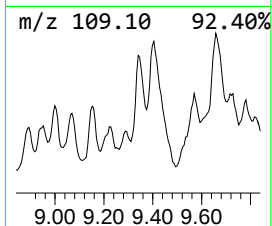
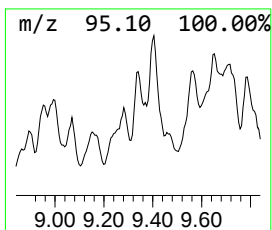
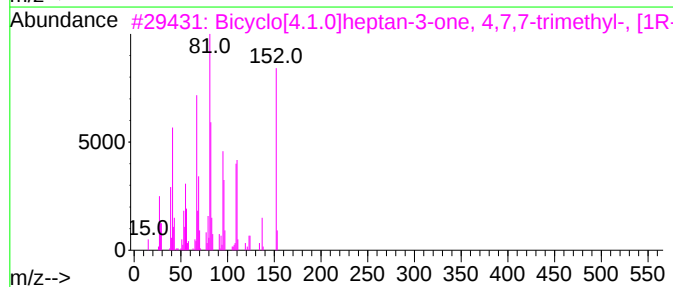
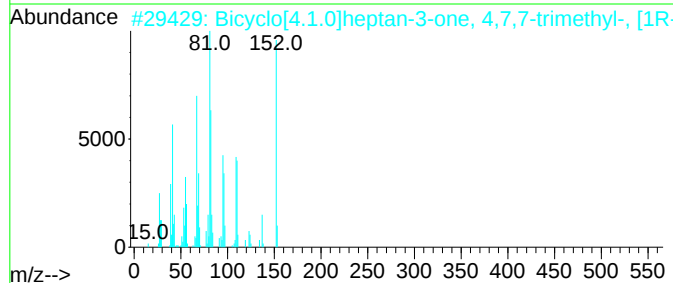
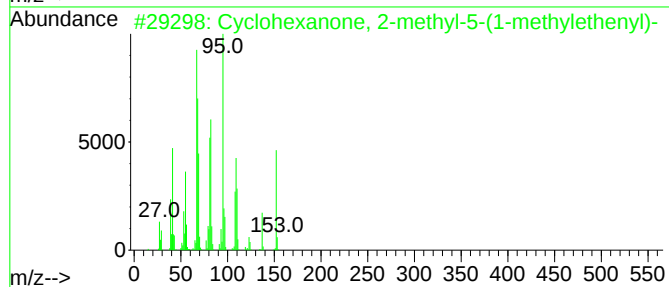
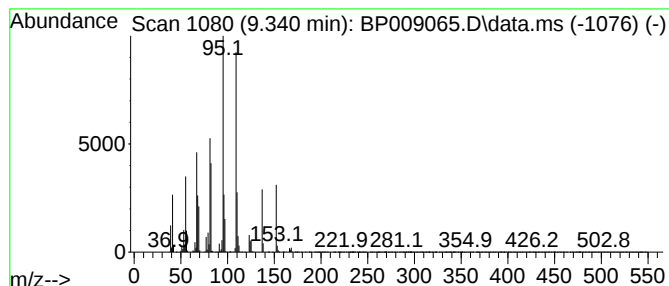
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TIC Integration Parameters: LSCINT.P

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Peak Number 14 Cyclohexanone, 2-methyl-5-(... Concentration Rank 17

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.   |
|-------|---------|---------|------------------|--------|
| 9.340 | 9.99 ng | 1068990 | Naphthalene-d8   | 10.610 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Cyclohexanone, 2-methyl-5-(1-met... | 152 | C10H16O | 007764-50-3 | 68   |
| 2    |      | Bicyclo[4.1.0]heptan-3-one, 4,7,... | 152 | C10H16O | 004176-01-6 | 49   |
| 3    |      | Bicyclo[4.1.0]heptan-3-one, 4,7,... | 152 | C10H16O | 004176-04-9 | 49   |
| 4    |      | 2-Dodecyne                          | 166 | C12H22  | 000629-49-2 | 43   |
| 5    |      | Bicyclo[2.2.1]heptane, 2-butyl-     | 152 | C11H20  | 061177-16-0 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

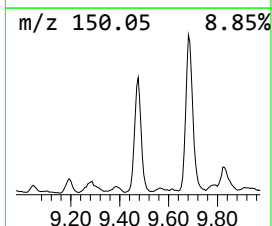
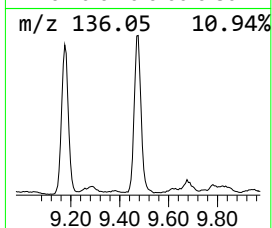
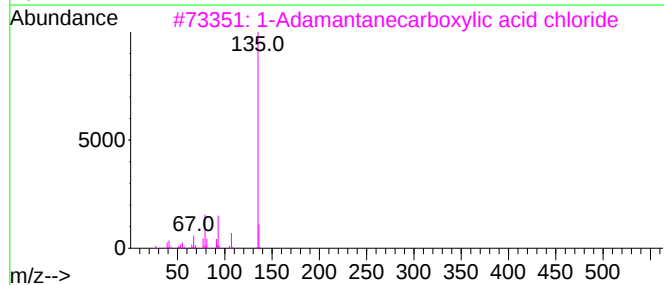
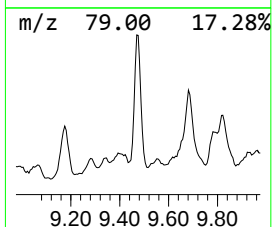
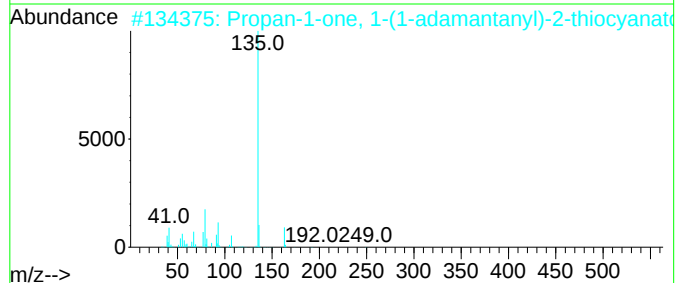
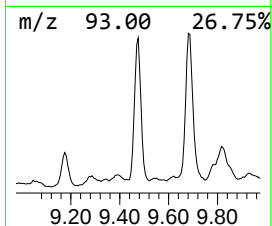
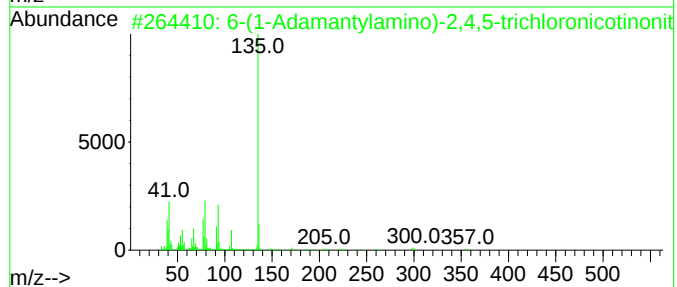
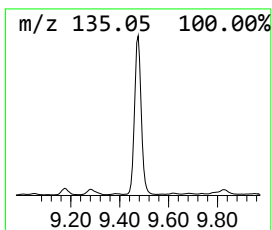
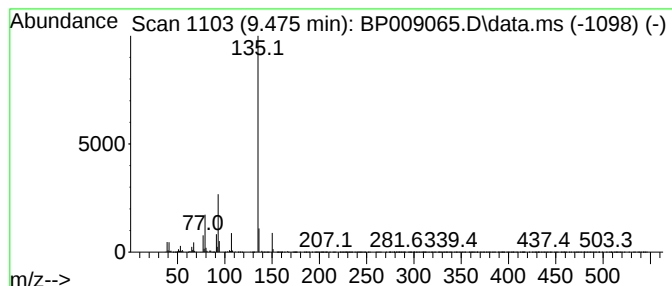
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 15 6-(1-Adamantylamino)-2,4,5-... Concentration Rank 4

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.   |
|-------|----------|---------|------------------|--------|
| 9.475 | 25.67 ng | 2747850 | Naphthalene-d8   | 10.610 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | 6-(1-Adamantylamino)-2,4,5-trich... | 355 | C16H16Cl3N3 | 339165-88-7  | 72   |
| 2    |    |   | Propan-1-one, 1-(1-adamantanyl)-... | 249 | C14H19NOS   | 313231-18-4  | 72   |
| 3    |    |   | 1-Adamantanecarboxylic acid chlo... | 198 | C11H15ClO   | 002094-72-6  | 72   |
| 4    |    |   | 1-Adamantanecarboxylic acid, 2-m... | 270 | C18H22O2    | 1010293-75-3 | 72   |
| 5    |    |   | 1-isocyanatocyclopropyl tricyclo... | 233 | C14H19NO2   | 1000396-40-9 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP020722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

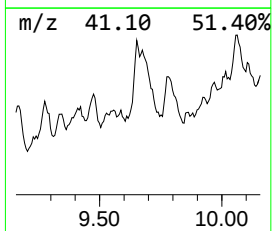
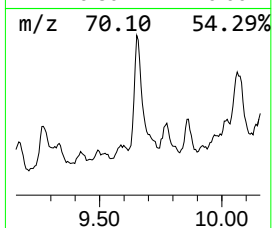
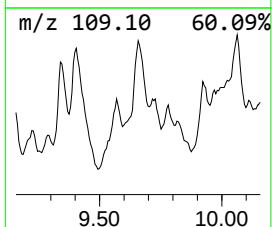
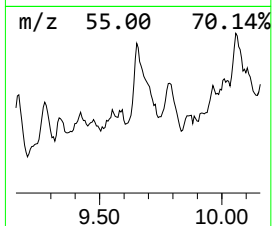
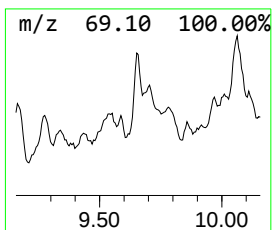
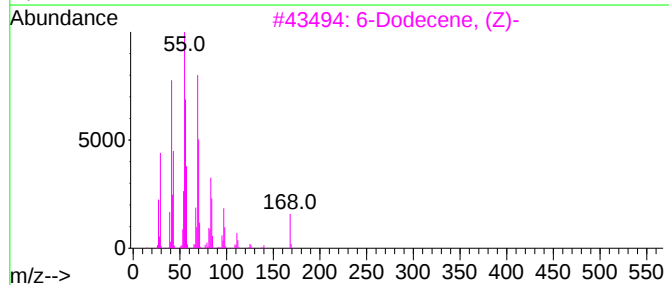
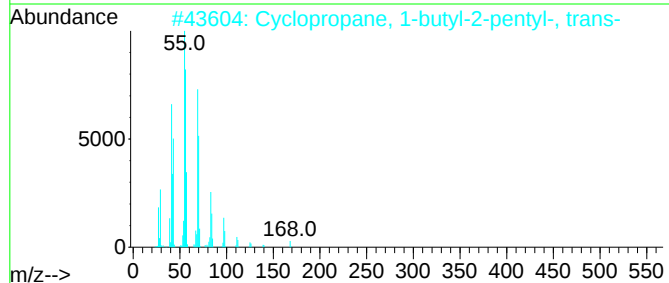
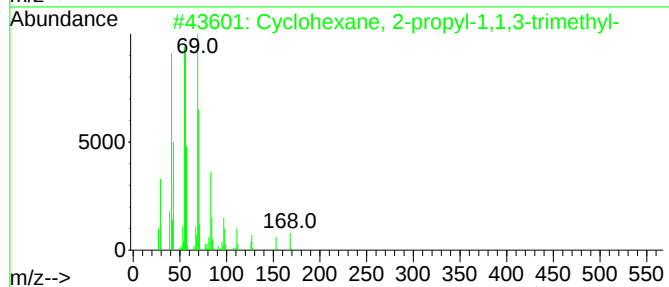
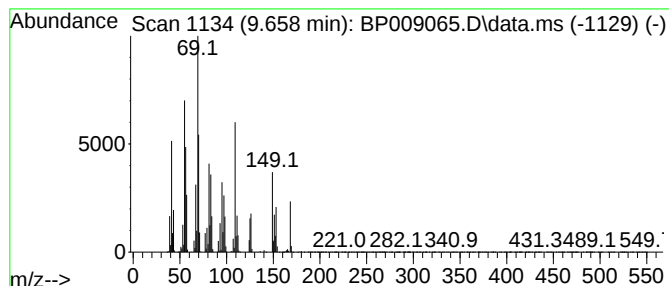
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Cyclohexane, 2-propyl-1,1,3... Concentration Rank 8

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.   |
|-------|----------|---------|------------------|--------|
| 9.658 | 18.06 ng | 1933520 | Naphthalene-d8   | 10.610 |

| Hit# | of 5 | Tentative ID                                          | MW  | MolForm  | CAS#        | Qual |
|------|------|-------------------------------------------------------|-----|----------|-------------|------|
| 1    |      | Cyclohexane, 2-propyl-1,1,3-trimethyl-                | 168 | C12H24   | 081983-70-2 | 70   |
| 2    |      | Cyclopropane, 1-butyl-2-pentyl-, trans-               | 168 | C12H24   | 074663-87-9 | 64   |
| 3    |      | 6-Dodecene, (Z)-                                      | 168 | C12H24   | 007206-29-3 | 64   |
| 4    |      | (R)-(+)-3-Aminoquinuclidine, N-allyl-                 | 168 | C9H16N2O | 134869-61-7 | 46   |
| 5    |      | 1.alpha.,2.beta.,3.alpha.,4.beta.-dicyclohexylmethane | 126 | C9H18    | 002532-67-4 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

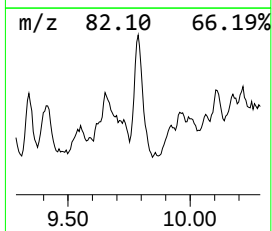
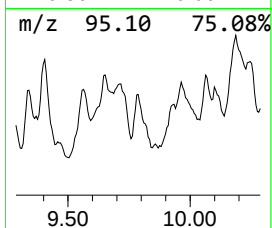
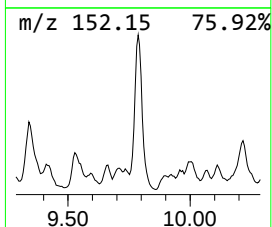
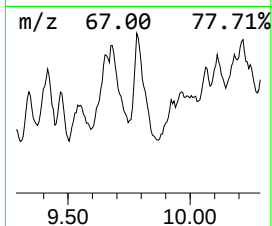
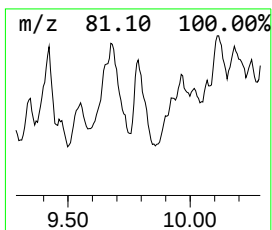
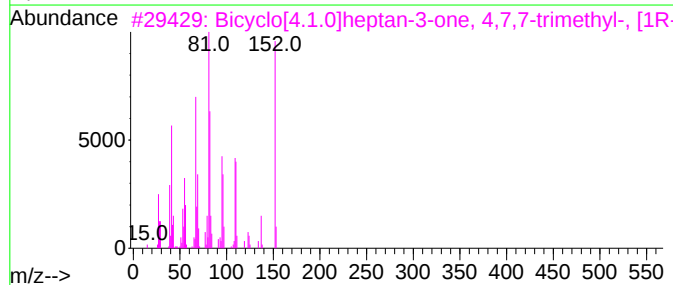
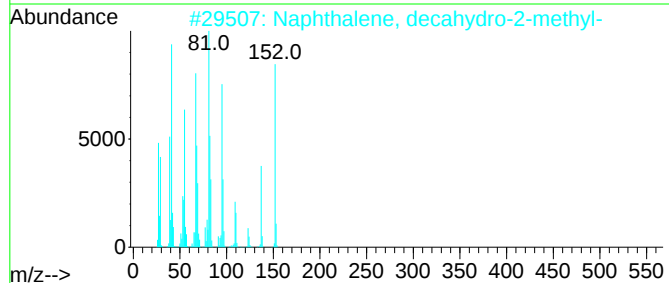
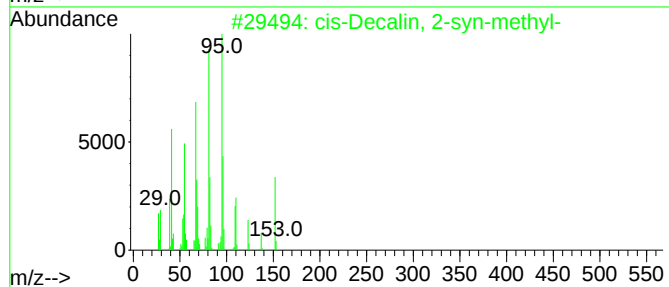
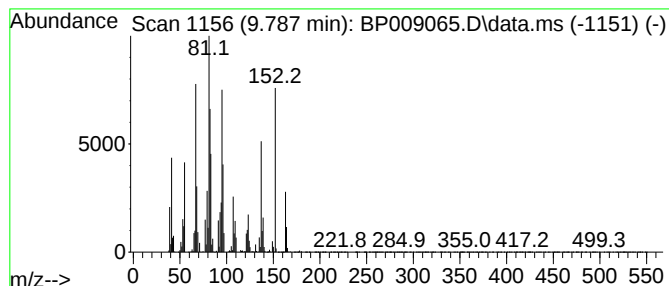
Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP020722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 17 cis-Decalin, 2-syn-methyl- Concentration Rank 3

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|--------------|------|
| 9.787     | 32.56 ng                            | 3486120 | Naphthalene-d8   | 10.610       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#         | Qual |
| 1         | cis-Decalin, 2-syn-methyl-          | 152     | C11H20           | 1000155-85-6 | 76   |
| 2         | Naphthalene, decahydro-2-methyl-    | 152     | C11H20           | 002958-76-1  | 72   |
| 3         | Bicyclo[4.1.0]heptan-3-one, 4,7,... | 152     | C10H16O          | 004176-01-6  | 53   |
| 4         | 7-Octadecyne, 2-methyl-             | 264     | C19H36           | 035354-38-2  | 52   |
| 5         | 9-Methylbicyclo[3.3.1]nonane        | 138     | C10H18           | 025107-01-1  | 46   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

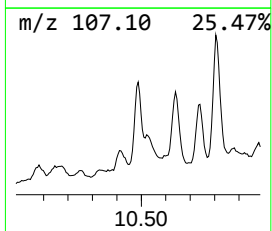
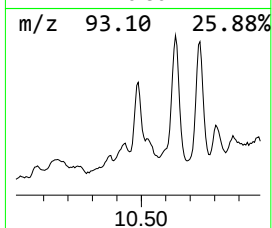
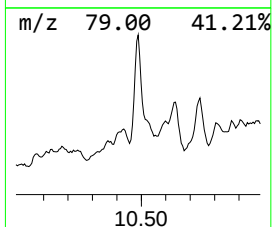
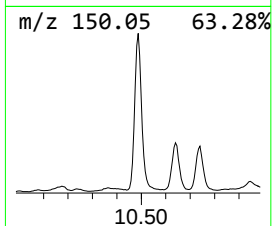
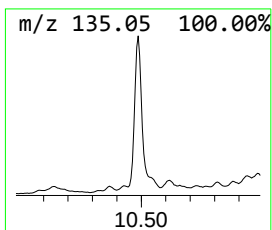
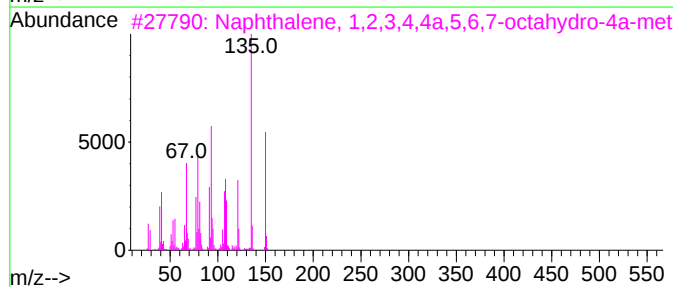
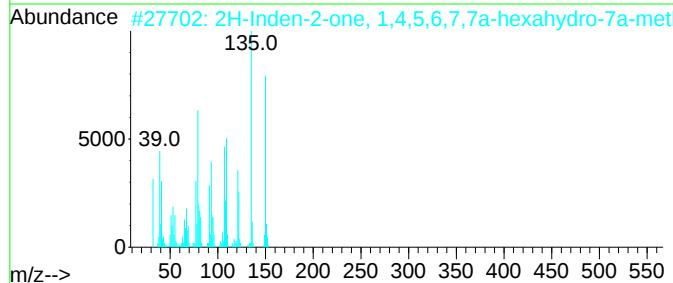
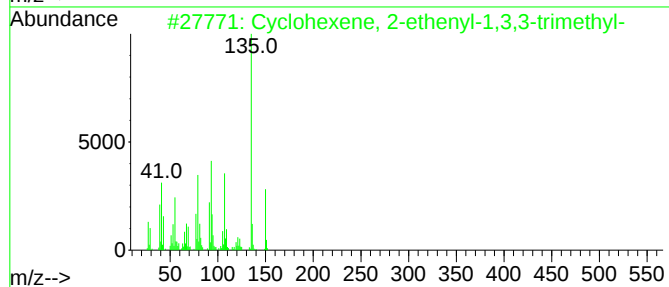
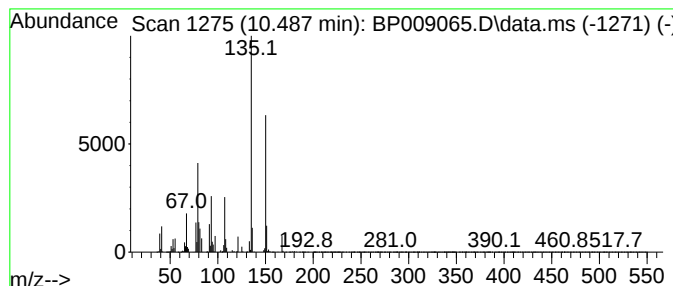
Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP020722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 18 Cyclohexene, 2-ethenyl-1,3,... Concentration Rank 7

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.    |              |      |
|-----------|-------------------------------------|---------|------------------|---------|--------------|------|
| 10.487    | 20.00 ng                            | 2141140 | Naphthalene-d8   | 10.610  |              |      |
| Hit# of 5 | Tentative ID                        |         | MW               | MolForm | CAS#         | Qual |
| 1         | Cyclohexene, 2-ethenyl-1,3,3-tri... |         | 150              | C11H18  | 005293-90-3  | 80   |
| 2         | 2H-Inden-2-one, 1,4,5,6,7,7a-hex... |         | 150              | C10H14O | 054725-16-5  | 74   |
| 3         | Naphthalene, 1,2,3,4,4a,5,6,7-oc... |         | 150              | C11H18  | 013943-77-6  | 72   |
| 4         | 3,9-Epoxy-p-mentha-1,8(10)-diene    |         | 150              | C10H14O | 1000111-14-8 | 64   |
| 5         | Phenol, 2-(1,1-dimethylethyl)-      |         | 150              | C10H14O | 000088-18-6  | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP020722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

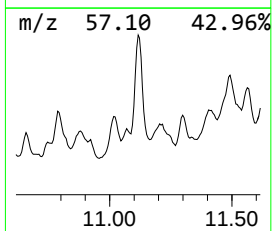
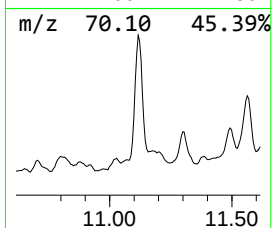
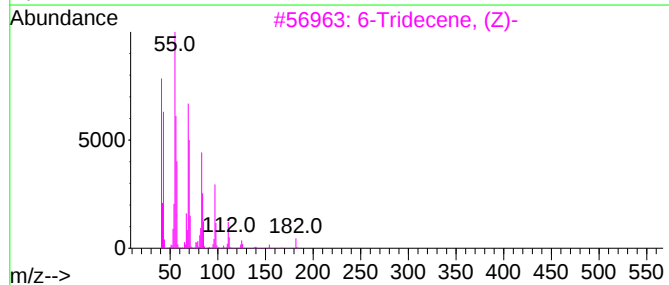
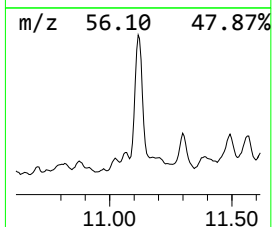
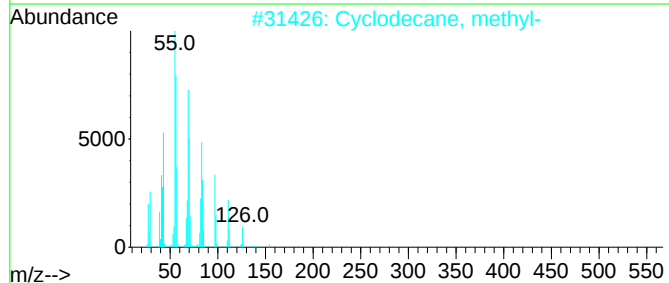
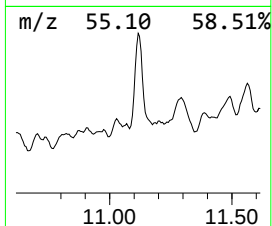
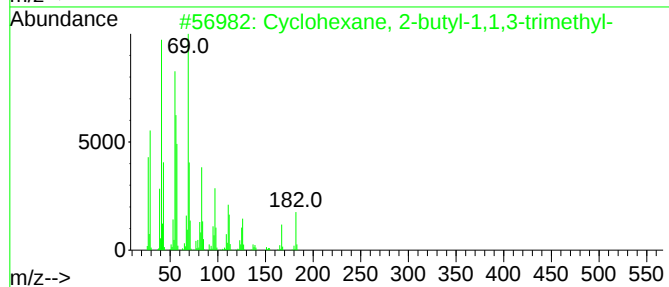
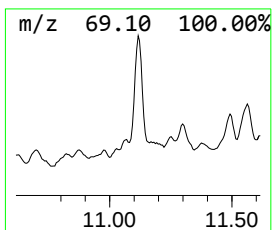
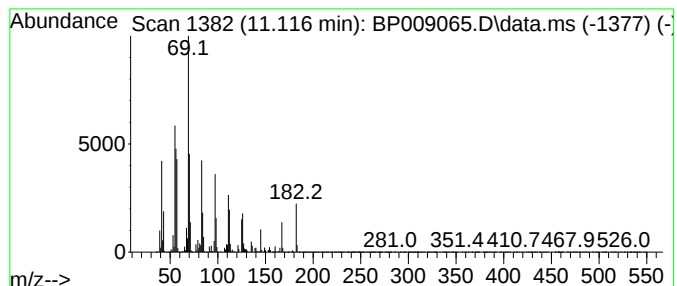
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Peak Number 19 Cyclohexane, 2-butyl-1,1,3-... Concentration Rank 1

| R.T.   | EstConc  | Area     | Relative to ISTD | R.T.   |
|--------|----------|----------|------------------|--------|
| 11.116 | 98.40 ng | 10534700 | Naphthalene-d8   | 10.610 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Cyclohexane, 2-butyl-1,1,3-trime... | 182 | C13H26  | 054676-39-0 | 96   |
| 2    |      | Cyclodecane, methyl-                | 154 | C11H22  | 013151-43-4 | 55   |
| 3    |      | 6-Tridecene, (Z)-                   | 182 | C13H26  | 006508-77-6 | 53   |
| 4    |      | 5-Dodecene, (E)-                    | 168 | C12H24  | 007206-16-8 | 50   |
| 5    |      | 6-Tridecene                         | 182 | C13H26  | 024949-38-0 | 49   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

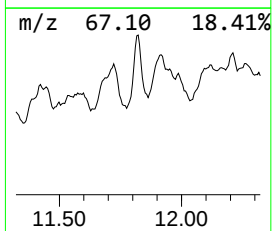
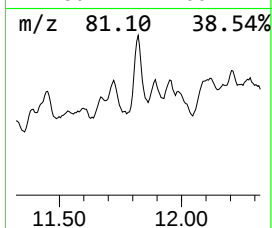
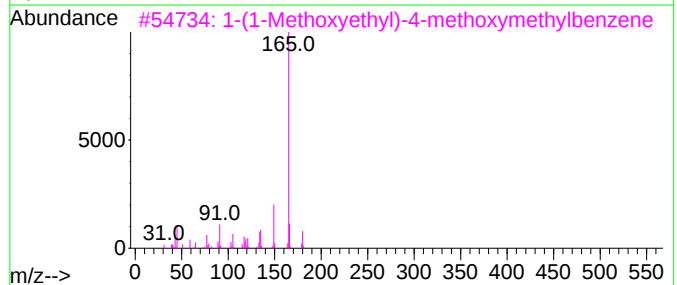
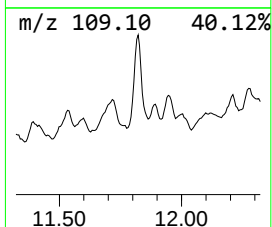
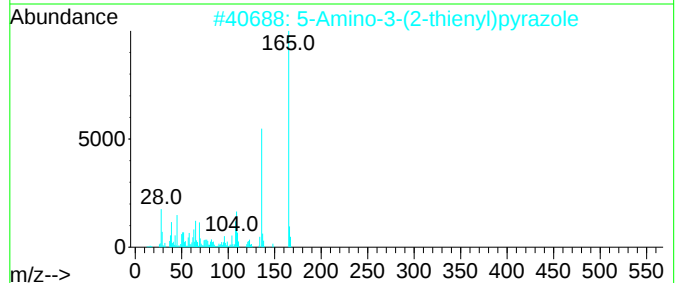
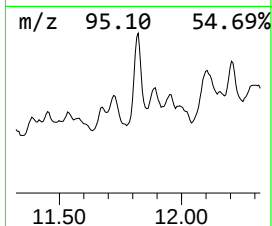
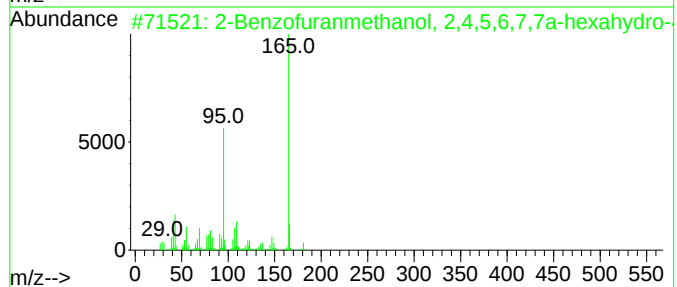
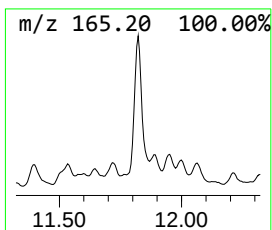
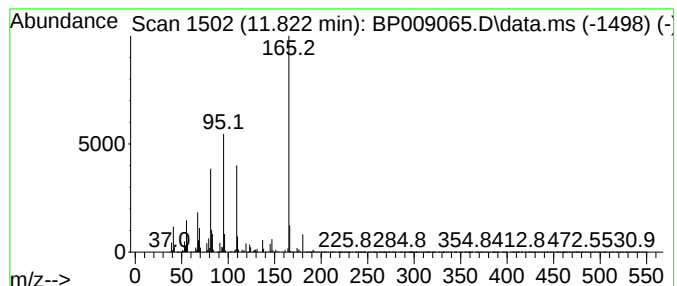
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 20 2-Benzofuranmethanol, 2,4,5... Concentration Rank 2

| R.T.      | EstConc                             | Area    | Relative to ISTD |          | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|----------|-------------|------|
| 11.822    | 54.73 ng                            | 5858770 | Naphthalene-d8   |          | 10.610      |      |
| Hit# of 5 | Tentative ID                        |         | MW               | MolForm  | CAS#        | Qual |
| 1         | 2-Benzofuranmethanol, 2,4,5,6,7,... |         | 196              | C12H20O2 | 077384-15-7 | 59   |
| 2         | 5-Amino-3-(2-thienyl)pyrazole       |         | 165              | C7H7N3S  | 096799-03-0 | 43   |
| 3         | 1-(1-Methoxyethyl)-4-methoxymeth... |         | 180              | C11H16O2 | 061145-46-8 | 38   |
| 4         | 2-Methyl-5-benzothiazolol           |         | 165              | C8H7NOS  | 068867-14-1 | 35   |
| 5         | 3-Methoxy-7-methyl-7H-pirazolo[4... |         | 165              | C6H7N5O  | 037526-58-2 | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020822\  
Data File : BP009065.D  
Acq On : 08 Feb 2022 18:51  
Operator : CG/JU  
Sample : N1399-01  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
KTS1-GW-MH-02-8-13

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP020722.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 |
| Bicyclo[3.1.1]h... | 7.311  | 11.5    | ng    | 1353370  | 1                     | 7.805  | 2362590 | 20.0 |
| cis-3-Decene       | 7.393  | 14.1    | ng    | 1660680  | 1                     | 7.805  | 2362590 | 20.0 |
| 1,1,3,3,5-Penta... | 7.634  | 10.2    | ng    | 1204710  | 1                     | 7.805  | 2362590 | 20.0 |
| unknown7.728       | 7.728  | 17.6    | ng    | 2078310  | 1                     | 7.805  | 2362590 | 20.0 |
| 1,4-Pentadiene,... | 7.975  | 14.7    | ng    | 1738720  | 1                     | 7.805  | 2362590 | 20.0 |
| Cyclohexane, 1,... | 8.040  | 6.0     | ng    | 710169   | 1                     | 7.805  | 2362590 | 20.0 |
| unknown8.087       | 8.087  | 9.4     | ng    | 1114040  | 1                     | 7.805  | 2362590 | 20.0 |
| Pulegone           | 8.175  | 6.4     | ng    | 756215   | 1                     | 7.805  | 2362590 | 20.0 |
| unknown8.569       | 8.569  | 20.3    | ng    | 2398940  | 1                     | 7.805  | 2362590 | 20.0 |
| unknown8.681       | 8.681  | 13.6    | ng    | 1605430  | 1                     | 7.805  | 2362590 | 20.0 |
| 4-Isopropyl-1,3... | 8.869  | 15.7    | ng    | 1855740  | 1                     | 7.805  | 2362590 | 20.0 |
| Adamantane         | 9.169  | 21.6    | ng    | 2549540  | 1                     | 7.805  | 2362590 | 20.0 |
| unknown9.281       | 9.281  | 12.9    | ng    | 1378120  | 2                     | 10.610 | 2141140 | 20.0 |
| Cyclohexanone, ... | 9.340  | 10.0    | ng    | 1068990  | 2                     | 10.610 | 2141140 | 20.0 |
| 6-(1-Adamantyla... | 9.475  | 25.7    | ng    | 2747850  | 2                     | 10.610 | 2141140 | 20.0 |
| Cyclohexane, 2-... | 9.658  | 18.1    | ng    | 1933520  | 2                     | 10.610 | 2141140 | 20.0 |
| cis-Decalin, 2-... | 9.787  | 32.6    | ng    | 3486120  | 2                     | 10.610 | 2141140 | 20.0 |
| Cyclohexene, 2-... | 10.487 | 20.0    | ng    | 2141140  | 2                     | 10.610 | 2141140 | 20.0 |
| Cyclohexane, 2-... | 11.116 | 98.4    | ng    | 10534700 | 2                     | 10.610 | 2141140 | 20.0 |
| 2-Benzofuranmet... | 11.822 | 54.7    | ng    | 5858770  | 2                     | 10.610 | 2141140 | 20.0 |