

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
 Data File : BP013585.D  
 Acq On : 24 Jan 2023 23:18  
 Operator : CG/JU  
 Sample : 01190-12  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 WS-1

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-BP012323.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP013585.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         | 3.223       | 3             | 6           | 12           | rVB      | 6868235        | 7787009       | 20.59%          | 2.554%        |
| 2         | 3.528       | 53            | 58          | 65           | rVB      | 496439         | 619210        | 1.64%           | 0.203%        |
| 3         | 5.217       | 336           | 345         | 351          | rBV      | 1183192        | 1635110       | 4.32%           | 0.536%        |
| 4         | 5.687       | 419           | 425         | 433          | rVV      | 3950936        | 5689512       | 15.05%          | 1.866%        |
| 5         | 7.299       | 691           | 699         | 708          | rVB      | 19576429       | 37814748      | 100.00%         | 12.402%       |
| 6         | 7.769       | 766           | 779         | 782          | rVV      | 395929         | 788084        | 2.08%           | 0.258%        |
| 7         | 7.799       | 782           | 784         | 793          | rVB      | 408792         | 572676        | 1.51%           | 0.188%        |
| 8         | 8.087       | 821           | 833         | 837          | rBV4     | 194027         | 524499        | 1.39%           | 0.172%        |
| 9         | 8.158       | 837           | 845         | 852          | rVV      | 1386889        | 2428071       | 6.42%           | 0.796%        |
| 10        | 8.393       | 880           | 885         | 891          | rBV      | 2564731        | 4169243       | 11.03%          | 1.367%        |
| 11        | 8.446       | 891           | 894         | 903          | rVB6     | 201147         | 435454        | 1.15%           | 0.143%        |
| 12        | 8.728       | 937           | 942         | 950          | rVB3     | 512965         | 1091142       | 2.89%           | 0.358%        |
| 13        | 8.805       | 950           | 955         | 957          | rBV2     | 236989         | 387422        | 1.02%           | 0.127%        |
| 14        | 8.846       | 957           | 962         | 968          | rVB2     | 777866         | 1346728       | 3.56%           | 0.442%        |
| 15        | 8.940       | 968           | 978         | 980          | rBV8     | 336144         | 918631        | 2.43%           | 0.301%        |
| 16        | 8.993       | 980           | 987         | 992          | rVV      | 6966864        | 11445122      | 30.27%          | 3.754%        |
| 17        | 9.052       | 992           | 997         | 1005         | rVB5     | 360206         | 791461        | 2.09%           | 0.260%        |
| 18        | 9.275       | 1023          | 1035        | 1039         | rBV3     | 1072574        | 3014724       | 7.97%           | 0.989%        |
| 19        | 9.334       | 1039          | 1045        | 1050         | rVV      | 4331324        | 7618570       | 20.15%          | 2.499%        |
| 20        | 9.393       | 1051          | 1055        | 1059         | rVV2     | 1158724        | 1915425       | 5.07%           | 0.628%        |
| 21        | 9.434       | 1059          | 1062        | 1067         | rVB4     | 501395         | 853119        | 2.26%           | 0.280%        |
| 22        | 9.528       | 1074          | 1078        | 1086         | rVB4     | 537857         | 1168305       | 3.09%           | 0.383%        |
| 23        | 9.605       | 1086          | 1091        | 1092         | rBV2     | 293645         | 408946        | 1.08%           | 0.134%        |
| 24        | 9.640       | 1093          | 1097        | 1104         | rVB4     | 664412         | 1364734       | 3.61%           | 0.448%        |
| 25        | 9.710       | 1104          | 1109        | 1113         | rBV4     | 398950         | 863580        | 2.28%           | 0.283%        |
| 26        | 9.810       | 1121          | 1126        | 1130         | rBV4     | 360610         | 600943        | 1.59%           | 0.197%        |
| 27        | 9.863       | 1130          | 1135        | 1139         | rBV3     | 500678         | 893715        | 2.36%           | 0.293%        |
| 28        | 9.905       | 1139          | 1142        | 1145         | rBV2     | 330752         | 427448        | 1.13%           | 0.140%        |
| 29        | 10.063      | 1159          | 1169        | 1174         | rBV3     | 783621         | 2198694       | 5.81%           | 0.721%        |
| 30        | 10.163      | 1181          | 1186        | 1193         | rBV4     | 871732         | 2278722       | 6.03%           | 0.747%        |
| 31        | 10.234      | 1194          | 1198        | 1200         | rBV3     | 404545         | 646386        | 1.71%           | 0.212%        |
| 32        | 10.952      | 1316          | 1320        | 1324         | rBV2     | 2028194        | 3789309       | 10.02%          | 1.243%        |
| 33        | 10.999      | 1324          | 1328        | 1336         | rVB3     | 2222670        | 4713314       | 12.46%          | 1.546%        |
| 34        | 11.146      | 1349          | 1353        | 1357         | rVB2     | 1843681        | 2740501       | 7.25%           | 0.899%        |
| 35        | 11.205      | 1359          | 1363        | 1367         | rBV3     | 936649         | 1658439       | 4.39%           | 0.544%        |
| 36        | 11.487      | 1406          | 1411        | 1417         | rVB3     | 2925611        | 5046579       | 13.35%          | 1.655%        |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
 Data File : BP013585.D  
 Acq On : 24 Jan 2023 23:18  
 Operator : CG/JU  
 Sample : 01190-12  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 WS-1

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-BP012323.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |          |          |        |        |
|----|--------|------|------|------|------|----------|----------|--------|--------|
| 37 | 11.657 | 1437 | 1440 | 1444 | rBV4 | 1188055  | 2311777  | 6.11%  | 0.758% |
| 38 | 11.757 | 1452 | 1457 | 1462 | rBV5 | 1429319  | 3310451  | 8.75%  | 1.086% |
| 39 | 11.922 | 1479 | 1485 | 1490 | rVB6 | 1969739  | 5007587  | 13.24% | 1.642% |
| 40 | 11.975 | 1490 | 1494 | 1496 | rBV2 | 1241467  | 1700069  | 4.50%  | 0.558% |
| 41 | 12.010 | 1497 | 1500 | 1502 | rVV  | 1649489  | 2264433  | 5.99%  | 0.743% |
| 42 | 12.046 | 1502 | 1506 | 1510 | rVV2 | 5112281  | 8563495  | 22.65% | 2.809% |
| 43 | 12.093 | 1510 | 1514 | 1520 | rVB3 | 3181195  | 5981000  | 15.82% | 1.962% |
| 44 | 12.181 | 1524 | 1529 | 1535 | rVB  | 3943157  | 7023348  | 18.57% | 2.303% |
| 45 | 12.269 | 1540 | 1544 | 1548 | rBV2 | 2907463  | 4382260  | 11.59% | 1.437% |
| 46 | 12.575 | 1591 | 1596 | 1609 | rVB2 | 7224397  | 19046164 | 50.37% | 6.246% |
| 47 | 12.716 | 1615 | 1620 | 1630 | rBV5 | 3806395  | 9993930  | 26.43% | 3.278% |
| 48 | 12.804 | 1631 | 1635 | 1639 | rBV  | 2084814  | 3760591  | 9.94%  | 1.233% |
| 49 | 13.028 | 1669 | 1673 | 1678 | rVB3 | 2599750  | 4382801  | 11.59% | 1.437% |
| 50 | 13.257 | 1708 | 1712 | 1717 | rVB2 | 4177960  | 6224781  | 16.46% | 2.042% |
| 51 | 13.428 | 1736 | 1741 | 1747 | rVB2 | 8780161  | 15910759 | 42.08% | 5.218% |
| 52 | 13.599 | 1766 | 1770 | 1773 | rBV  | 5802233  | 9617161  | 25.43% | 3.154% |
| 53 | 14.234 | 1874 | 1878 | 1883 | rBV2 | 9746150  | 17342163 | 45.86% | 5.688% |
| 54 | 14.551 | 1928 | 1932 | 1938 | rBV5 | 4043714  | 9784879  | 25.88% | 3.209% |
| 55 | 14.640 | 1943 | 1947 | 1952 | rVB  | 9371885  | 15463253 | 40.89% | 5.071% |
| 56 | 19.822 | 2824 | 2828 | 2836 | rVB3 | 6385663  | 11529483 | 30.49% | 3.781% |
| 57 | 20.086 | 2870 | 2873 | 2877 | rVB  | 10233195 | 12184610 | 32.22% | 3.996% |
| 58 | 21.628 | 3131 | 3135 | 3140 | rVB  | 2733776  | 4000747  | 10.58% | 1.312% |
| 59 | 24.210 | 3566 | 3574 | 3590 | rVB  | 1940487  | 4478971  | 11.84% | 1.469% |

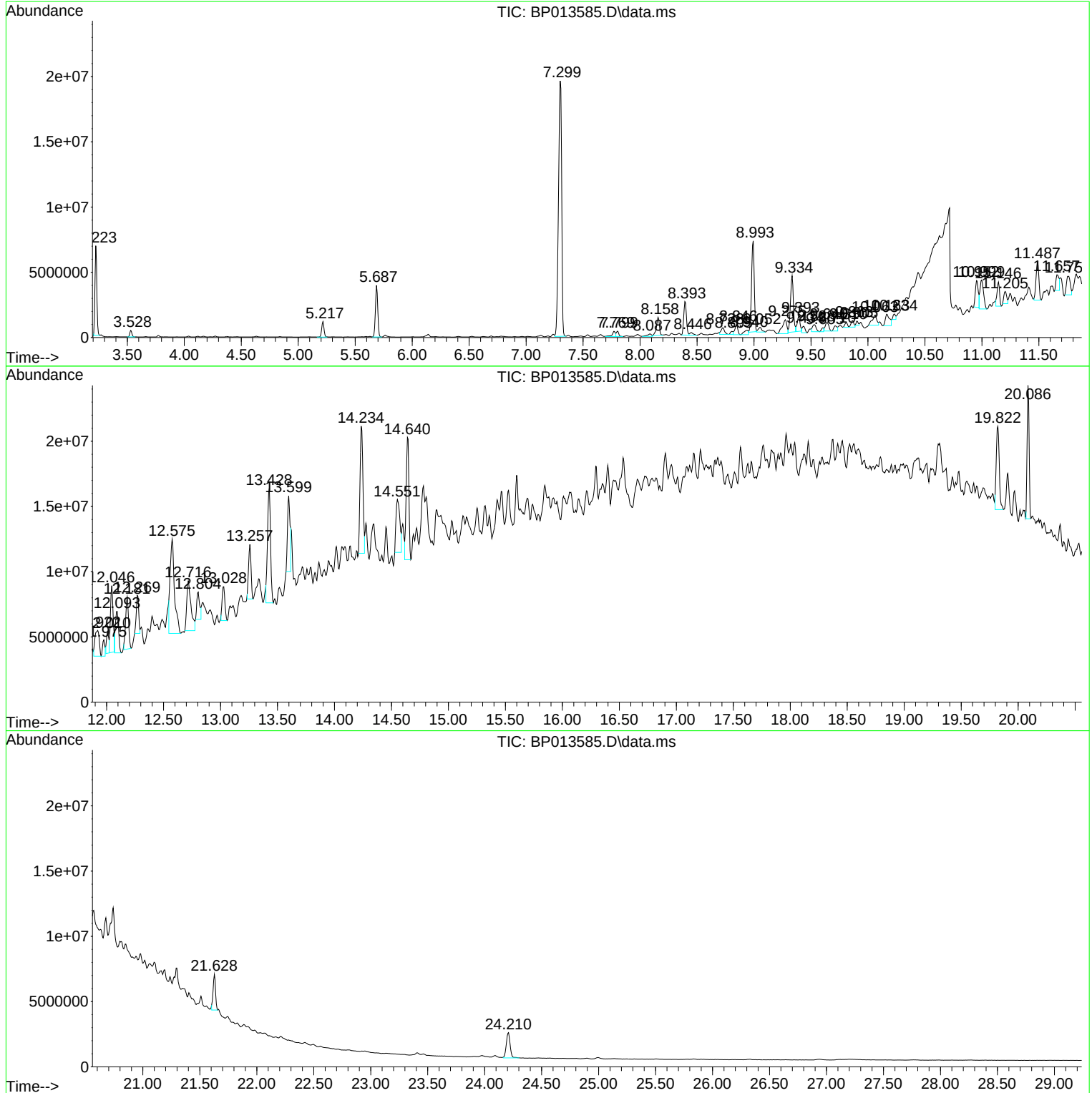
Sum of corrected areas: 304910288

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.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\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

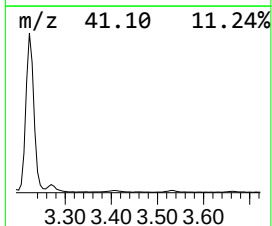
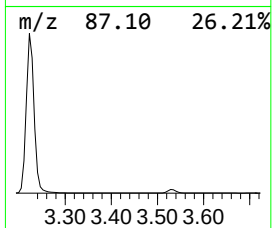
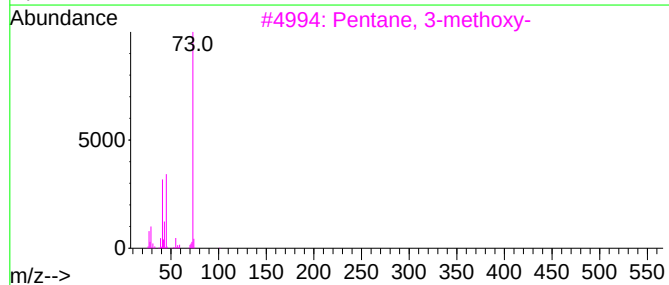
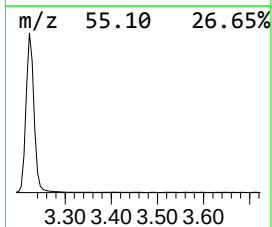
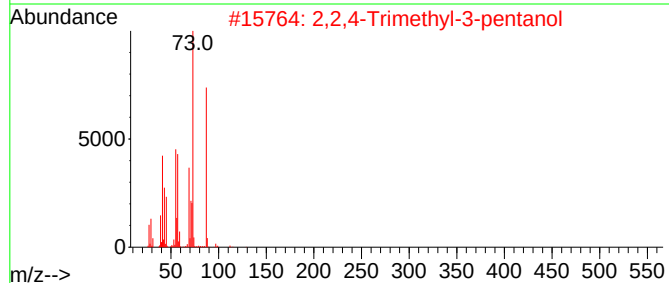
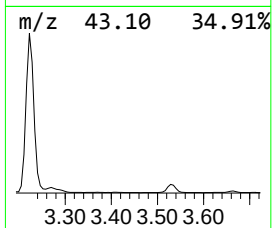
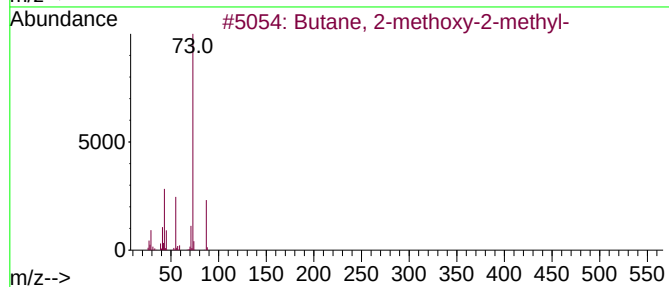
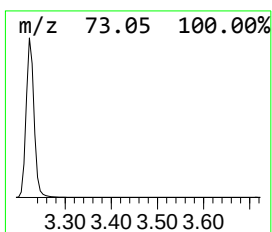
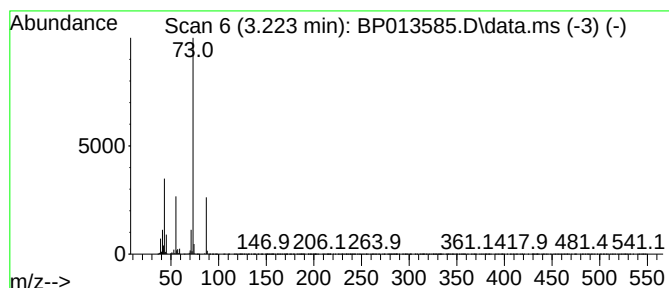
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

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

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 3.223 | 64.14 ng | 7787010 | 1,4-Dichlorobenzene-d4 | 8.158 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 78   |
| 2    |    |   | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 40   |
| 3    |    |   | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 17   |
| 4    |    |   | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 12   |
| 5    |    |   | Acetamide, N-ethyl-         | 87  | C4H9NO  | 000625-50-3 | 9    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

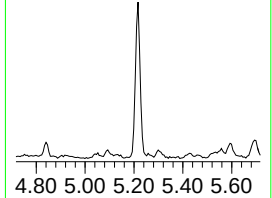
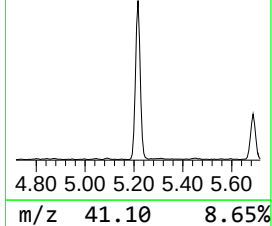
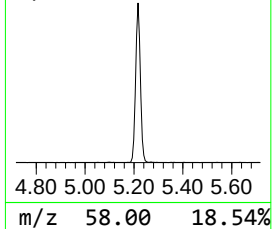
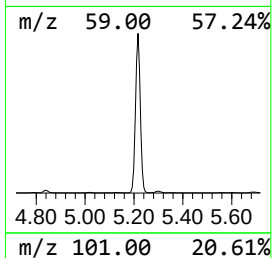
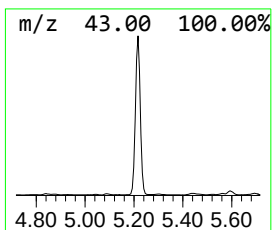
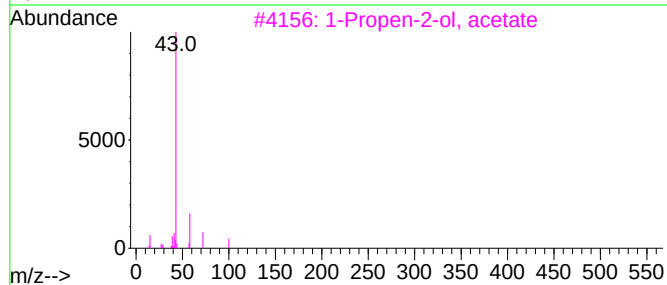
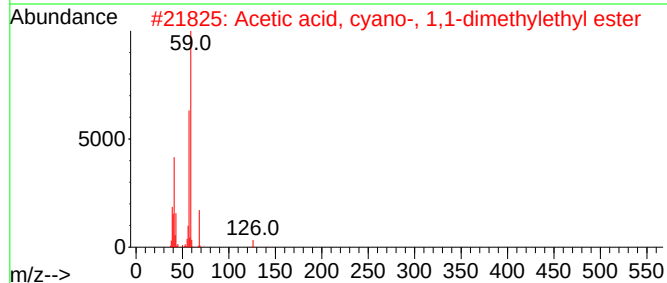
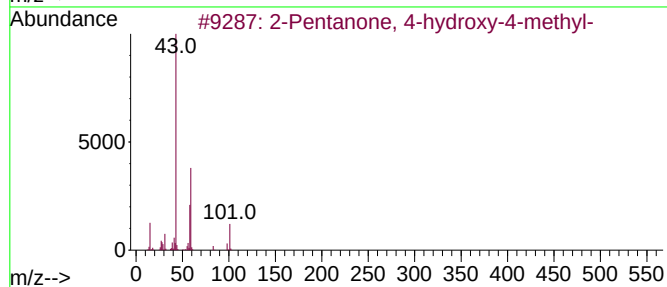
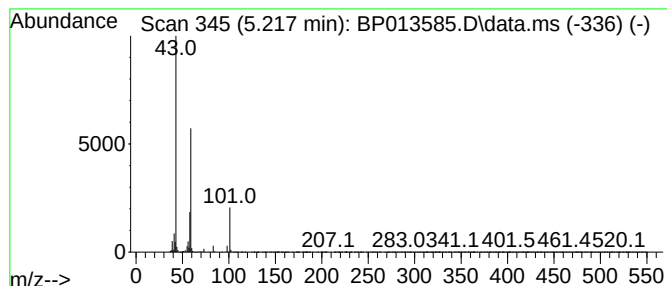
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 16

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 5.217 | 13.47 ng | 1635110 | 1,4-Dichlorobenzene-d4 | 8.158 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 50   |
| 2    |    |   | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 23   |
| 3    |    |   | 1-Propen-2-ol, acetate              | 100 | C5H8O2   | 000108-22-5 | 10   |
| 4    |    |   | Morpholine, 4-methyl-               | 101 | C5H11NO  | 000109-02-4 | 9    |
| 5    |    |   | Acetone                             | 58  | C3H6O    | 000067-64-1 | 9    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

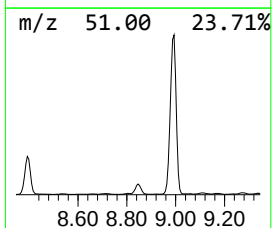
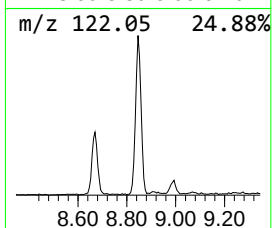
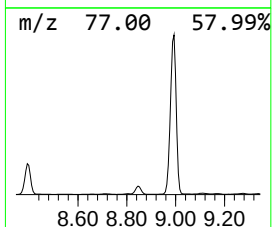
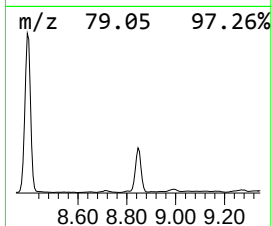
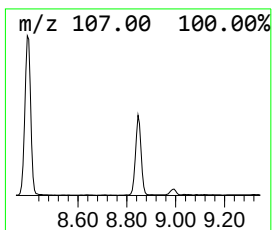
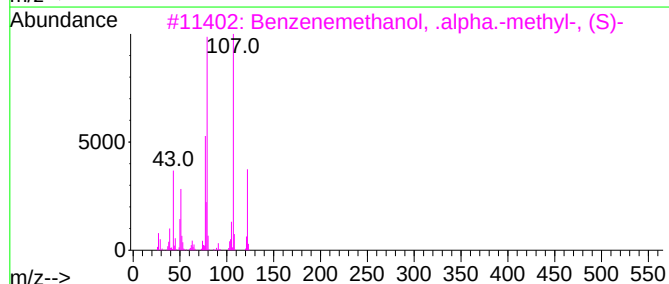
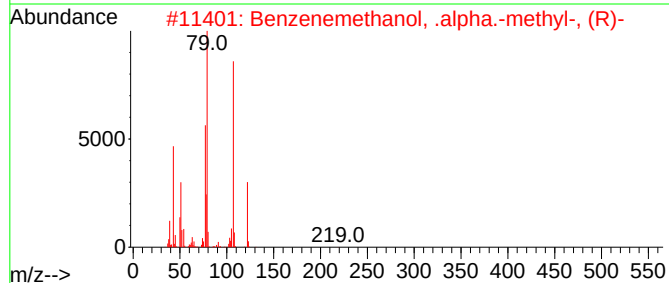
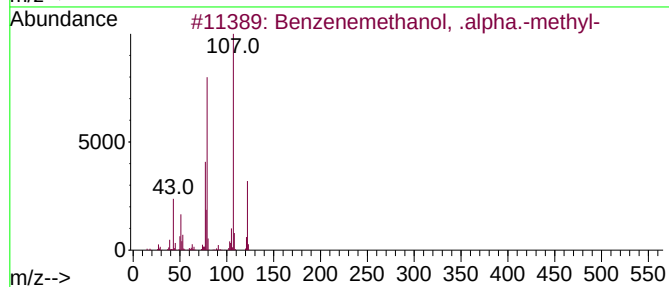
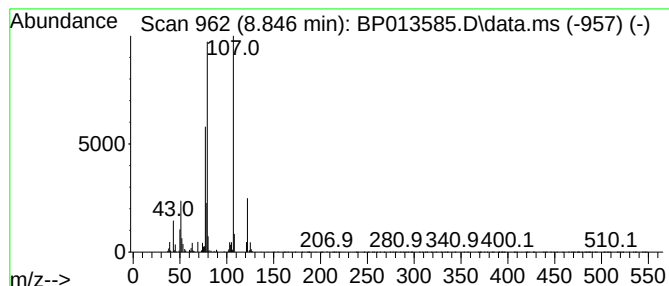
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 3 Benzenemethanol, .alpha.-me... Concentration Rank 20

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 8.846 | 11.09 ng | 1346730 | 1,4-Dichlorobenzene-d4 | 8.158 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzenemethanol, .alpha.-methyl-    | 122 | C8H10O  | 000098-85-1 | 94   |
| 2    |    |   | Benzenemethanol, .alpha.-methyl-... | 122 | C8H10O  | 001517-69-7 | 94   |
| 3    |    |   | Benzenemethanol, .alpha.-methyl-... | 122 | C8H10O  | 001445-91-6 | 91   |
| 4    |    |   | Benzenepropanenitrile, .beta.-hy... | 147 | C9H9NO  | 017190-29-3 | 72   |
| 5    |    |   | S-(-)-1-Phenylpropanol              | 136 | C9H12O  | 000613-87-6 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

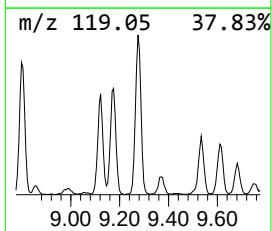
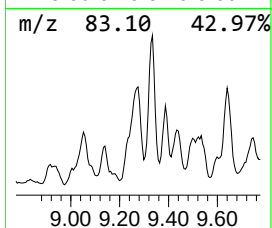
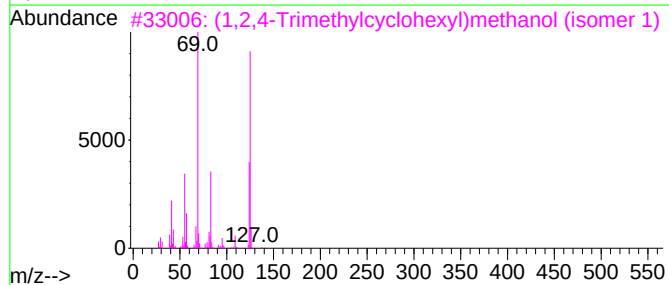
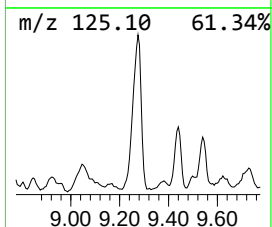
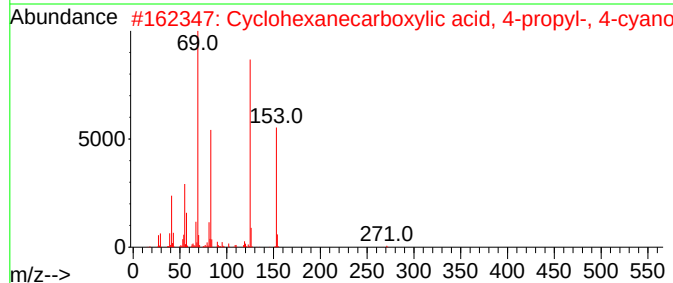
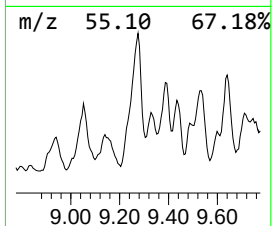
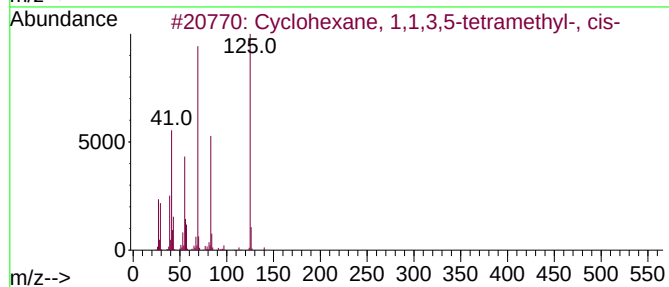
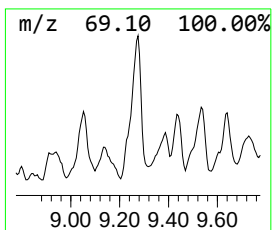
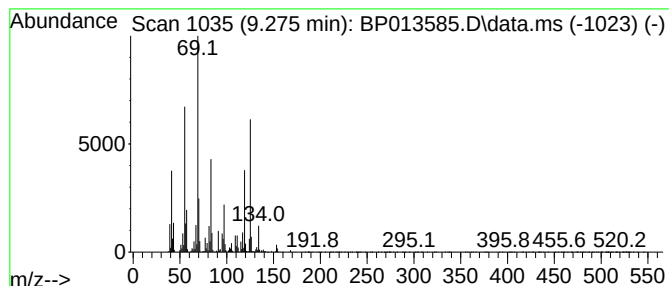
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 Cyclohexane, 1,1,3,5-tetram... Concentration Rank 8

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 9.275 | 24.83 ng | 3014720 | 1,4-Dichlorobenzene-d4 | 8.158 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Cyclohexane, 1,1,3,5-tetramethyl... | 140 | C10H20    | 050876-32-9  | 50   |
| 2    |    |   | Cyclohexanecarboxylic acid, 4-pr... | 271 | C17H21NO2 | 062439-33-2  | 50   |
| 3    |    |   | (1,2,4-Trimethylcyclohexyl)metha... | 156 | C10H20O   | 1000499-03-8 | 50   |
| 4    |    |   | Cyclohexane, 1,1,4,4-tetramethyl-   | 140 | C10H20    | 002223-52-1  | 50   |
| 5    |    |   | (1,2,4-Trimethylcyclohexyl)metha... | 156 | C10H20O   | 1000499-04-1 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

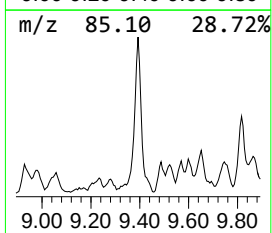
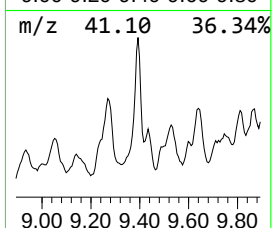
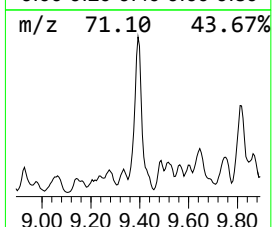
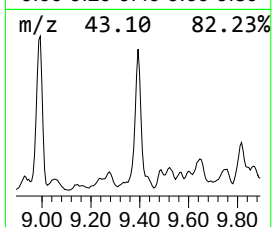
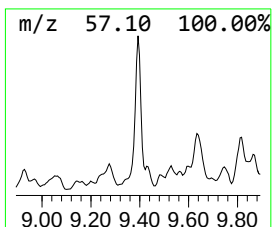
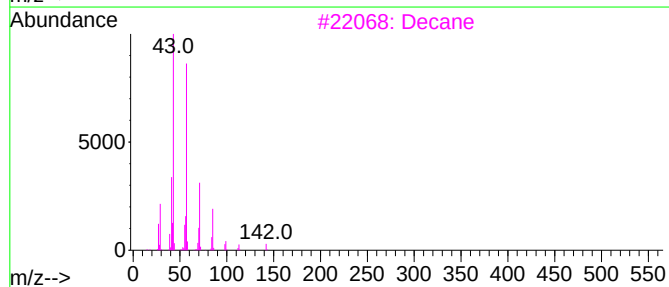
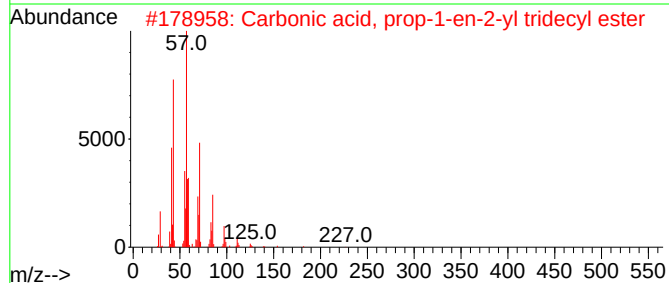
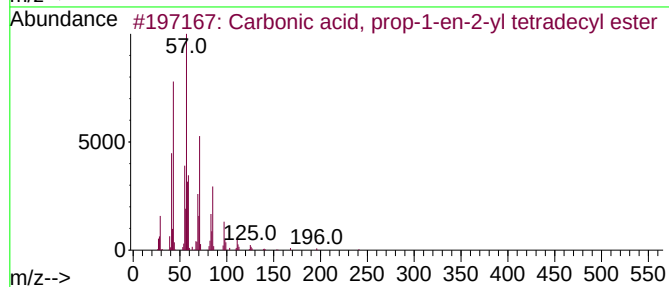
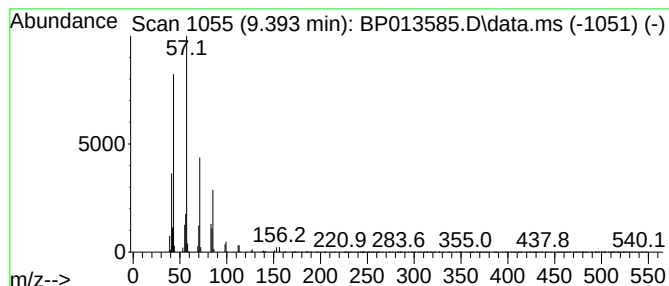
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 5 Carbonic acid, prop-1-en-2-... Concentration Rank 14

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 9.393 | 15.78 ng | 1915430 | 1,4-Dichlorobenzene-d4 | 8.158 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Carbonic acid, prop-1-en-2-yl te... | 298 | C18H34O3 | 1000382-90-9 | 90   |
| 2    |    |   | Carbonic acid, prop-1-en-2-yl tr... | 284 | C17H32O3 | 1000382-90-8 | 90   |
| 3    |    |   | Decane                              | 142 | C10H22   | 000124-18-5  | 86   |
| 4    |    |   | Undecane                            | 156 | C11H24   | 001120-21-4  | 81   |
| 5    |    |   | Hexadecane                          | 226 | C16H34   | 000544-76-3  | 80   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

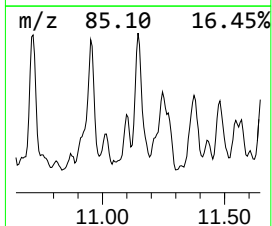
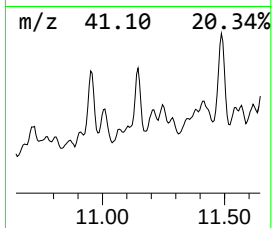
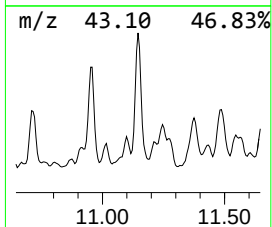
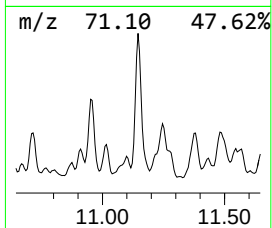
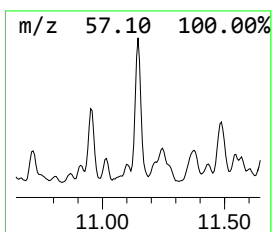
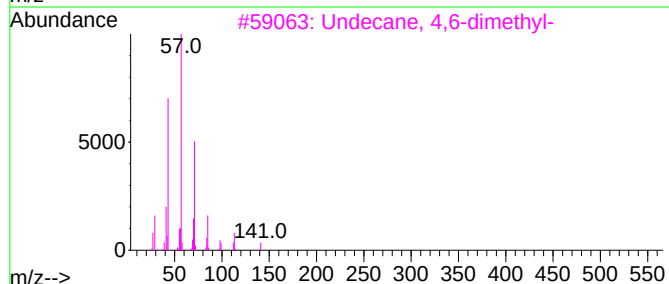
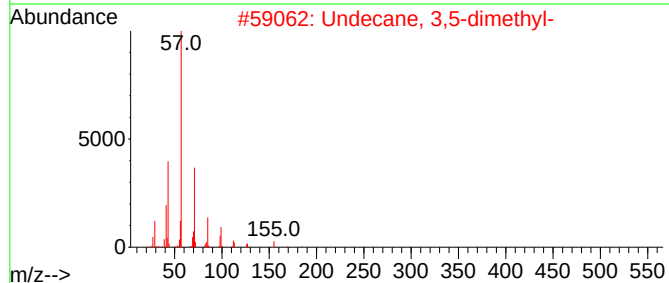
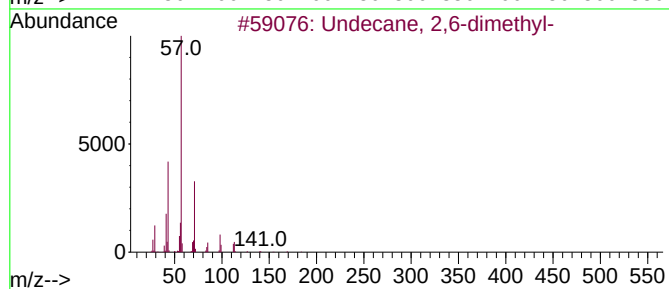
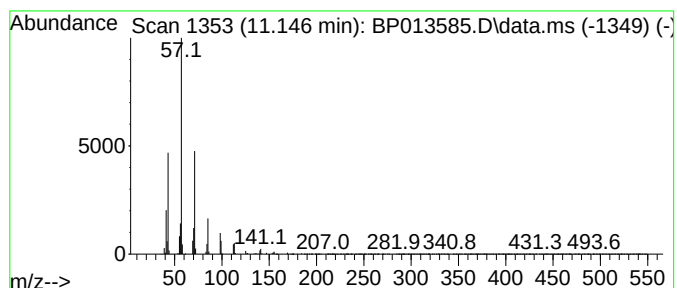
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*

Peak Number 6 Undecane, 2,6-dimethyl- Concentration Rank 19

| R.T.      | EstConc                        | Area    | Relative to ISTD | R.T.        |      |
|-----------|--------------------------------|---------|------------------|-------------|------|
| 11.146    | 11.63 ng                       | 2740500 | Naphthalene-d8   | 10.993      |      |
| Hit# of 5 | Tentative ID                   | MW      | MolForm          | CAS#        | Qual |
| 1         | Undecane, 2,6-dimethyl-        | 184     | C13H28           | 017301-23-4 | 90   |
| 2         | Undecane, 3,5-dimethyl-        | 184     | C13H28           | 017312-81-1 | 90   |
| 3         | Undecane, 4,6-dimethyl-        | 184     | C13H28           | 017312-82-2 | 87   |
| 4         | Dodecane, 2,6,10-trimethyl-    | 212     | C15H32           | 003891-98-3 | 78   |
| 5         | Pentadecane, 2,6,10-trimethyl- | 254     | C18H38           | 003892-00-0 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

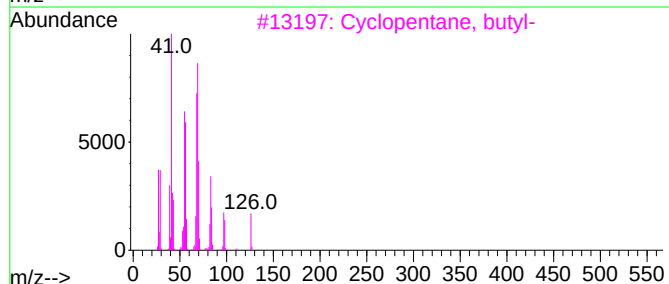
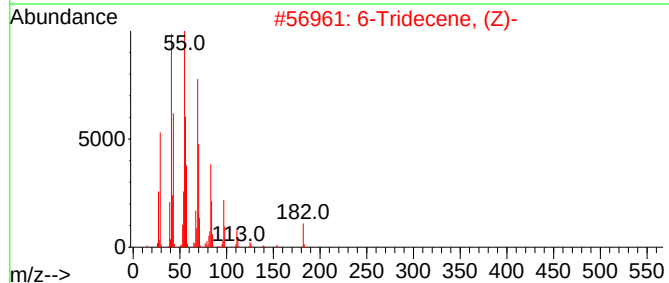
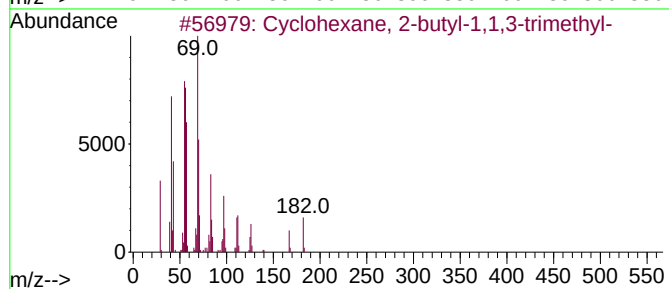
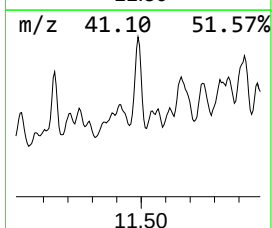
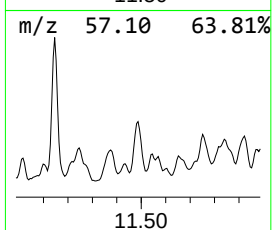
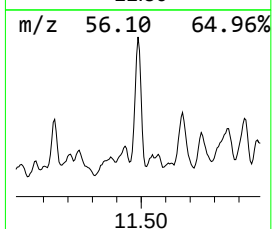
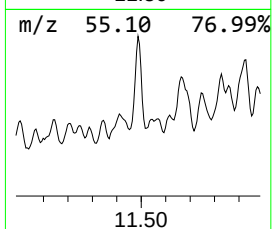
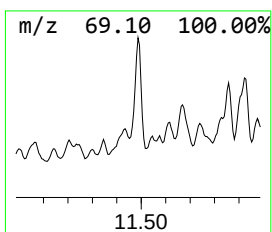
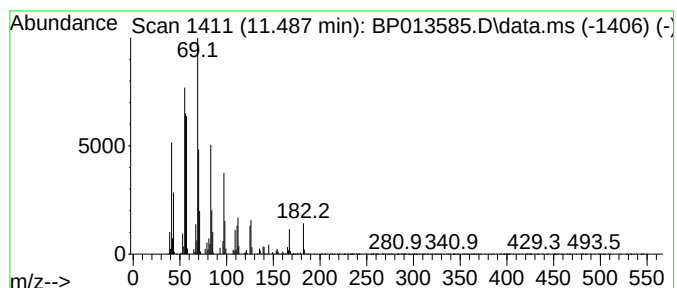
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*

Peak Number 7 Cyclohexane, 2-butyl-1,1,3-... Concentration Rank 9

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 11.487    | 21.41 ng                            | 5046580 | Naphthalene-d8   | 10.993      |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | Cyclohexane, 2-butyl-1,1,3-trime... | 182     | C13H26           | 054676-39-0 | 94   |
| 2         | 6-Tridecene, (Z)-                   | 182     | C13H26           | 006508-77-6 | 90   |
| 3         | Cyclopentane, butyl-                | 126     | C9H18            | 002040-95-1 | 64   |
| 4         | 5-Dodecene, (E)-                    | 168     | C12H24           | 007206-16-8 | 58   |
| 5         | 1.alpha.,2.beta.,3.alpha.,4.beta... | 126     | C9H18            | 002532-67-4 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

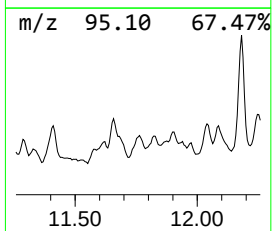
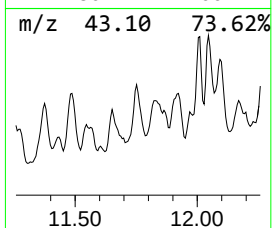
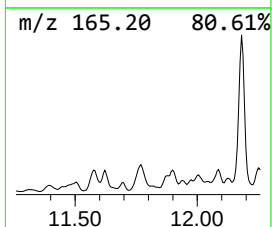
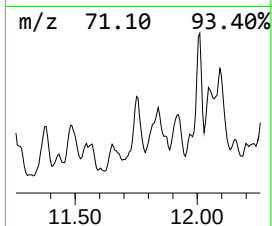
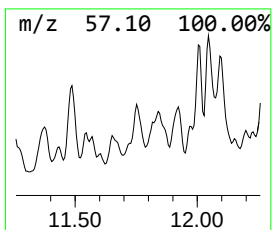
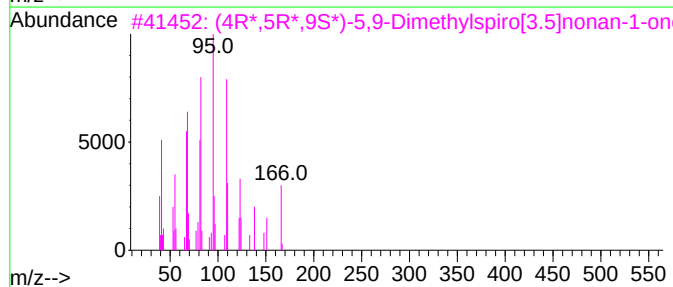
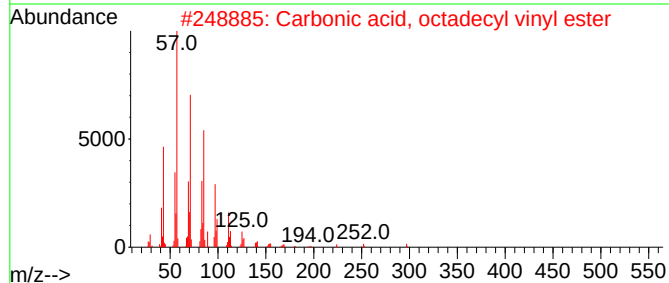
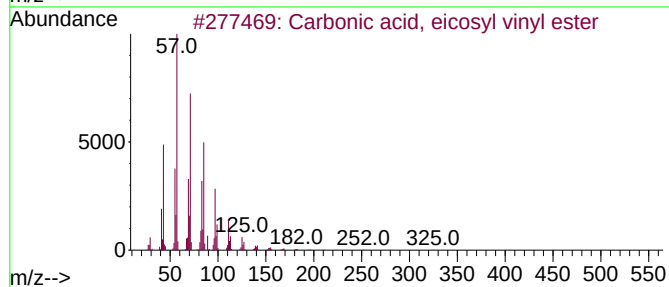
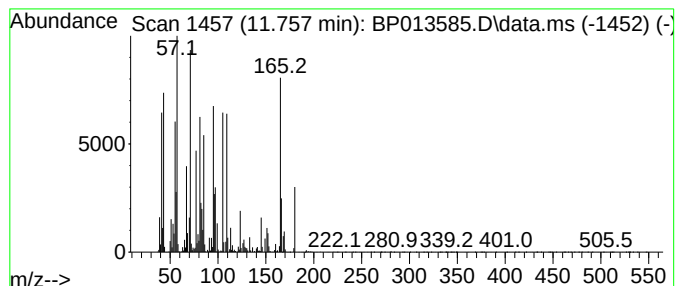
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 8 unknown11.757 Concentration Rank 15

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 11.757 | 14.05 ng | 3310450 | Naphthalene-d8   | 10.993 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Carbonic acid, eicosyl vinyl ester  | 368 | C23H44O3  | 1000382-54-3 | 25   |
| 2    |    |   | Carbonic acid, octadecyl vinyl e... | 340 | C21H40O3  | 1000382-54-4 | 25   |
| 3    |    |   | (4R*,5R*,9S*)-5,9-Dimethylspiro[... | 166 | C11H18O   | 063088-19-7  | 25   |
| 4    |    |   | Sulfurous acid, pentadecyl 2-pro... | 334 | C18H38O3S | 1000309-12-6 | 22   |
| 5    |    |   | Nonadecane                          | 268 | C19H40    | 000629-92-5  | 18   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

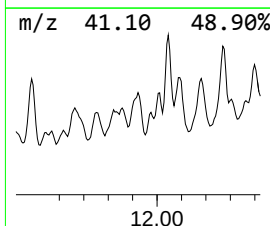
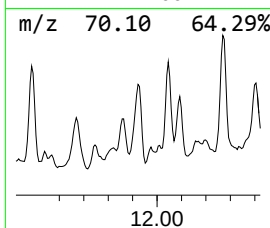
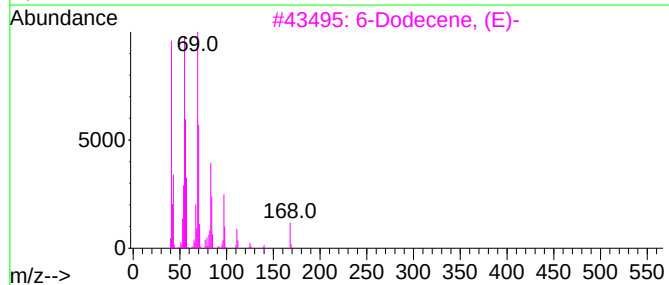
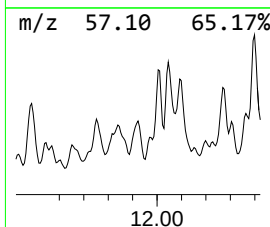
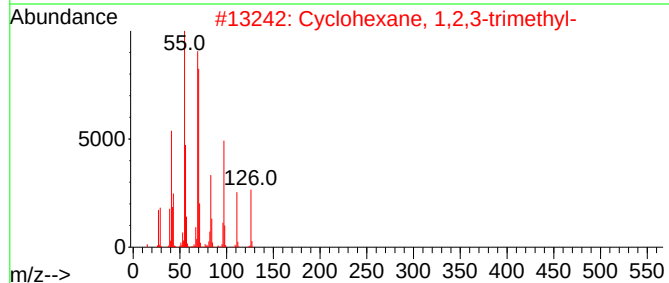
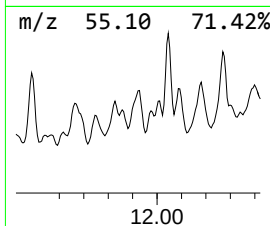
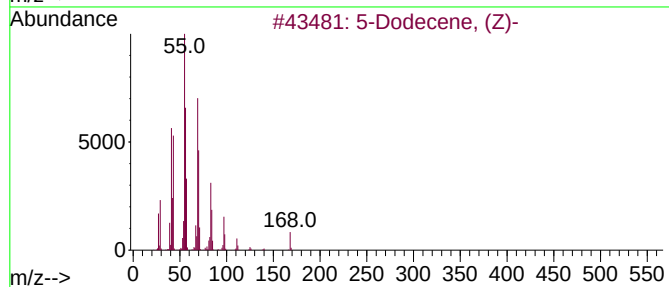
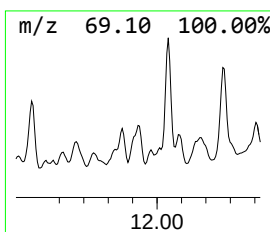
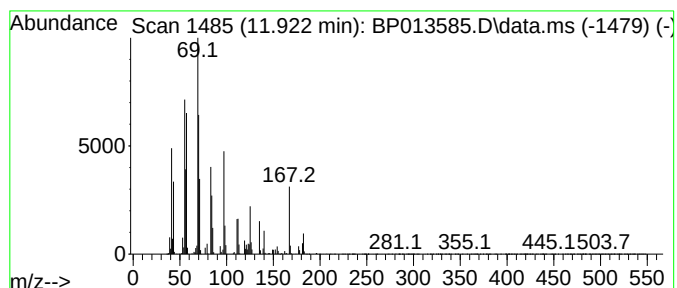
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*

Peak Number 9 5-Dodecene, (Z)- Concentration Rank 10

| R.T.      | EstConc                       | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------|---------|------------------|-------------|------|
| 11.922    | 21.25 ng                      | 5007590 | Naphthalene-d8   | 10.993      |      |
| Hit# of 5 | Tentative ID                  | MW      | MolForm          | CAS#        | Qual |
| 1         | 5-Dodecene, (Z)-              | 168     | C12H24           | 007206-28-2 | 76   |
| 2         | Cyclohexane, 1,2,3-trimethyl- | 126     | C9H18            | 001678-97-3 | 76   |
| 3         | 6-Dodecene, (E)-              | 168     | C12H24           | 007206-17-9 | 68   |
| 4         | 4-Decene                      | 140     | C10H20           | 019689-18-0 | 62   |
| 5         | 2-Tridecene, (E)-             | 182     | C13H26           | 041446-58-6 | 60   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

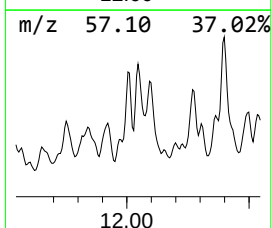
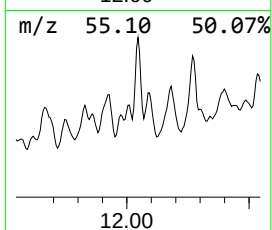
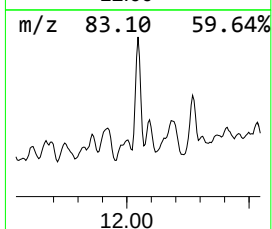
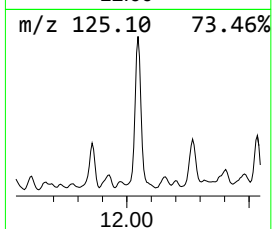
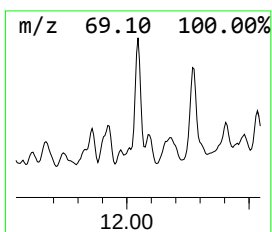
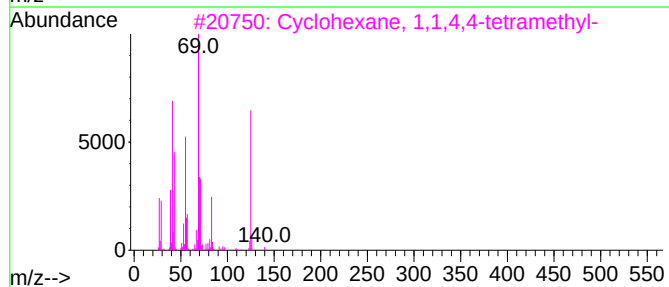
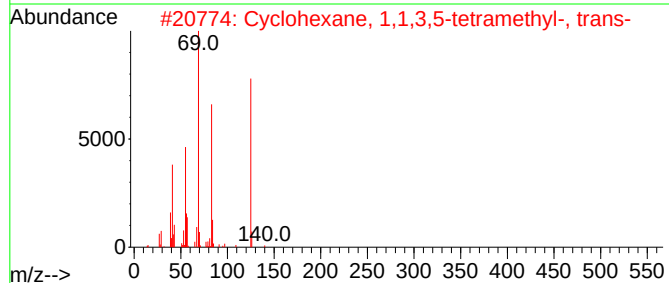
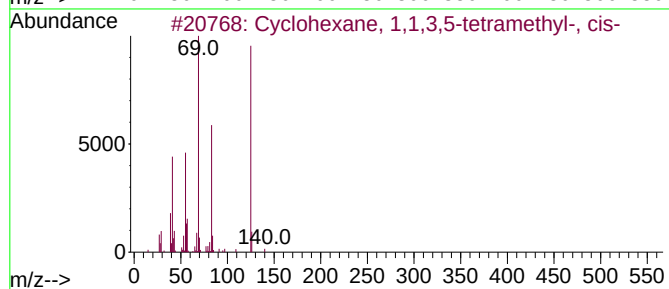
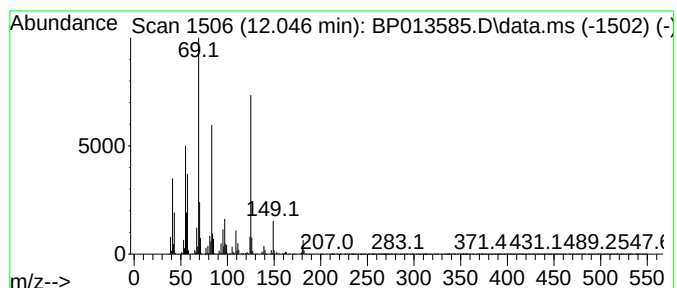
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*

Peak Number 10 Cyclohexane, 1,1,4,4-tetram... Concentration Rank 5

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|--------------|------|
| 12.046    | 36.34 ng                            | 8563500 | Naphthalene-d8   | 10.993       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#         | Qual |
| 1         | Cyclohexane, 1,1,3,5-tetramethyl... | 140     | C10H20           | 050876-32-9  | 74   |
| 2         | Cyclohexane, 1,1,3,5-tetramethyl... | 140     | C10H20           | 050876-31-8  | 59   |
| 3         | Cyclohexane, 1,1,4,4-tetramethyl-   | 140     | C10H20           | 002223-52-1  | 58   |
| 4         | Cyclooctane, 1-methyl-3-propyl-     | 168     | C12H24           | 255885-37-1  | 53   |
| 5         | (1,2,4-Trimethylcyclohexyl)metha... | 156     | C10H20O          | 1000499-04-1 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

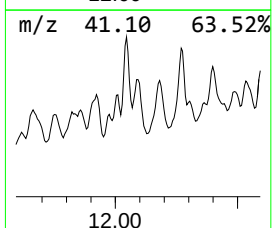
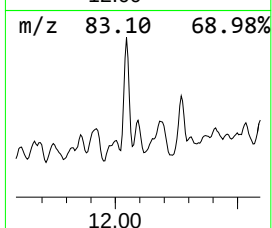
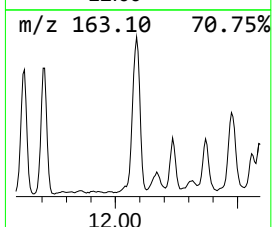
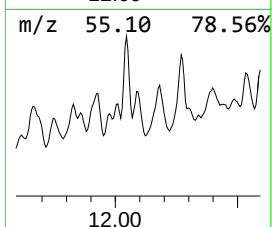
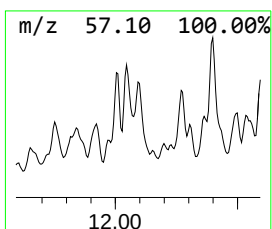
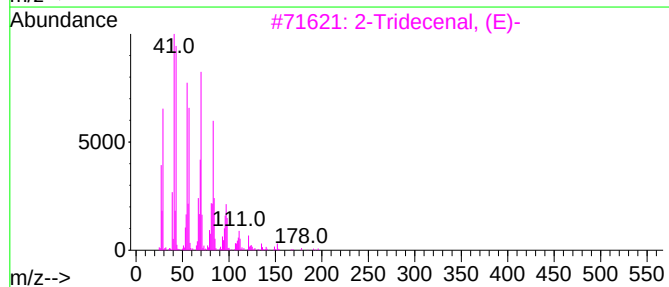
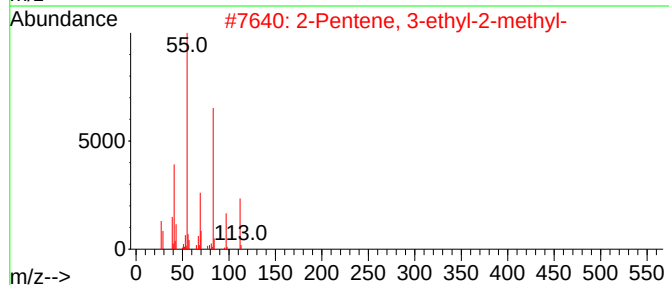
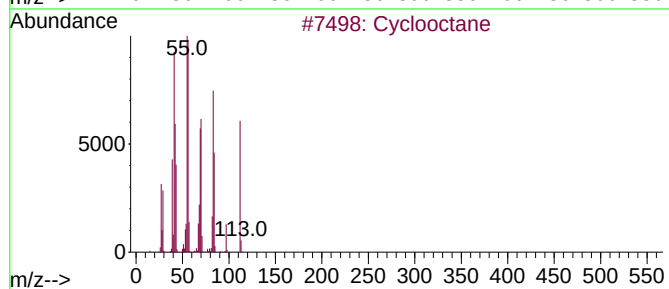
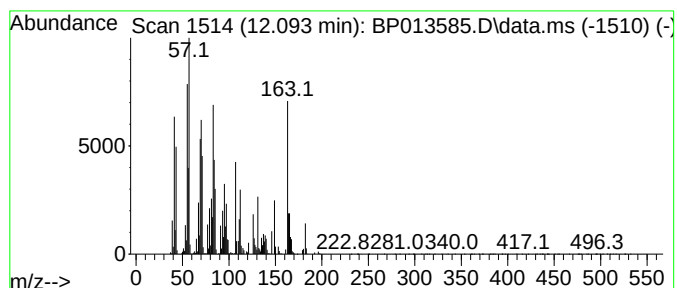
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*

Peak Number 11 unknown12.093 Concentration Rank 7

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 12.093    | 25.38 ng                            | 5981000 | Naphthalene-d8   | 10.993      |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | Cyclooctane                         | 112     | C8H16            | 000292-64-8 | 27   |
| 2         | 2-Pentene, 3-ethyl-2-methyl-        | 112     | C8H16            | 019780-67-7 | 18   |
| 3         | 2-Tridecenal, (E)-                  | 196     | C13H24O          | 007069-41-2 | 15   |
| 4         | Cyclopentane, 1-hexyl-3-methyl-     | 168     | C12H24           | 061142-68-5 | 11   |
| 5         | (2,2,6-Trimethyl-bicyclo[4.1.0]h... | 168     | C11H20O          | 078996-11-9 | 11   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

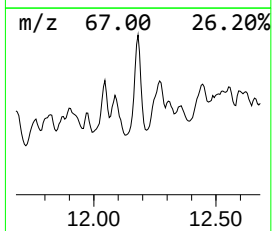
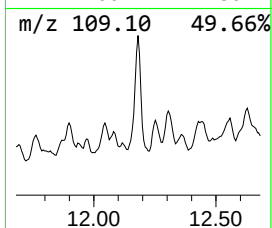
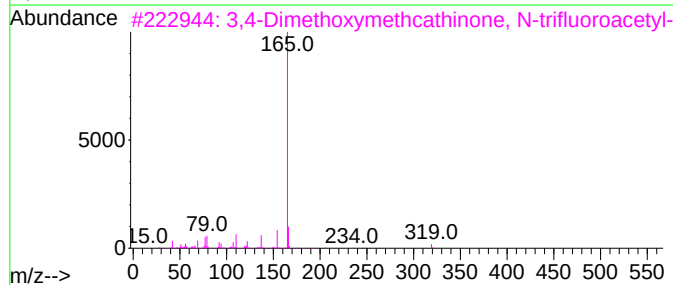
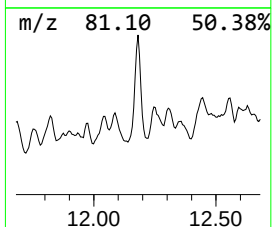
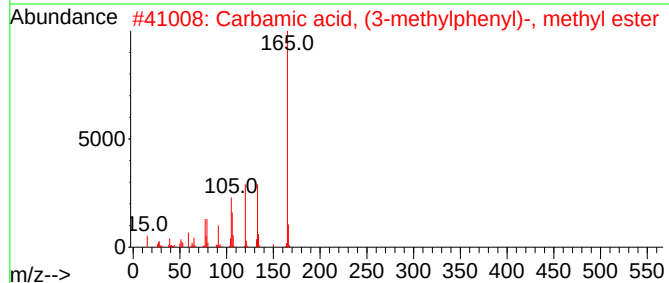
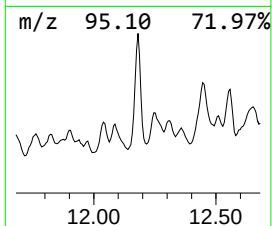
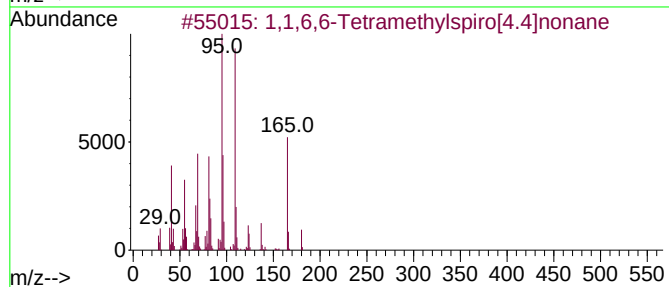
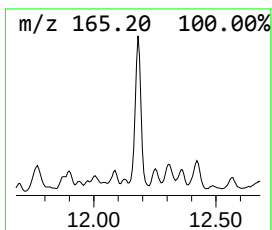
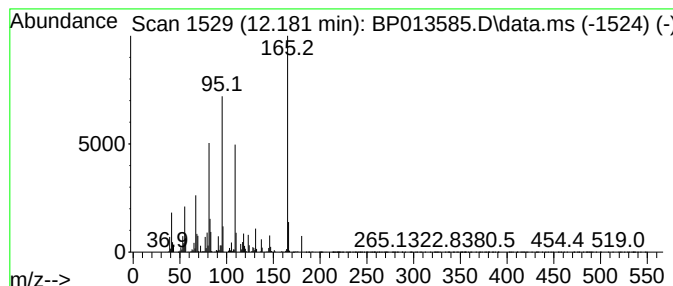
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*

Peak Number 12 unknown12.181 Concentration Rank 6

| R.T.   | EstConc                             | Area            | Relative to ISTD | R.T.      |
|--------|-------------------------------------|-----------------|------------------|-----------|
| 12.181 | 29.80 ng                            | 7023350         | Naphthalene-d8   | 10.993    |
| Hit#   | of 5                                | Tentative ID    | MW MolForm       | CAS# Qual |
| 1      | 1,1,6,6-Tetramethylspiro[4.4]nonane | 180 C13H24      | 074054-92-5      | 38        |
| 2      | Carbamic acid, (3-methylphenyl)-... | 165 C9H11NO2    | 039076-18-1      | 30        |
| 3      | 3,4-Dimethoxymethcathinone, N-tr... | 319 C14H16F3NO4 | 1000475-76-9     | 27        |
| 4      | cis,cis- and cis,trans-1,9-dimet... | 166 C12H22      | 1000111-72-5     | 25        |
| 5      | trans,trans-1,7-Dimethylspiro[4...  | 166 C12H22      | 1000111-72-6     | 20        |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

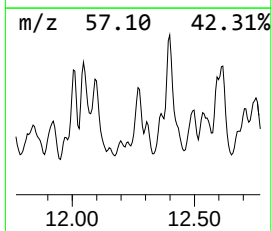
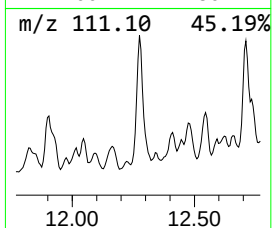
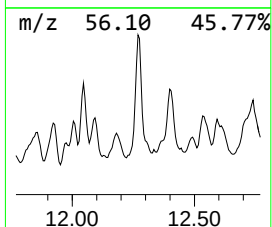
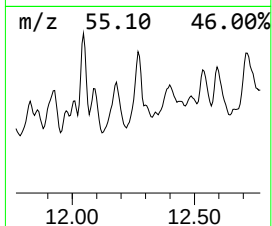
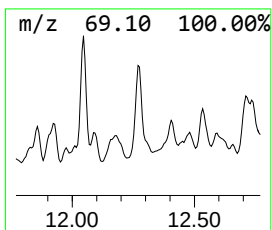
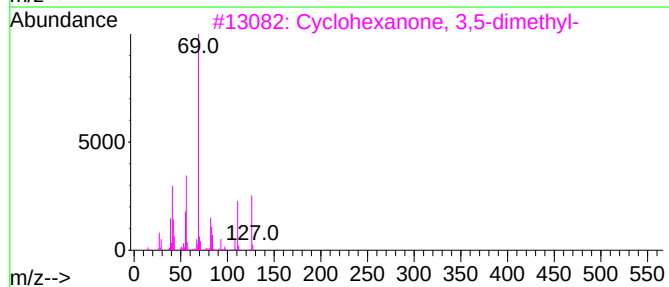
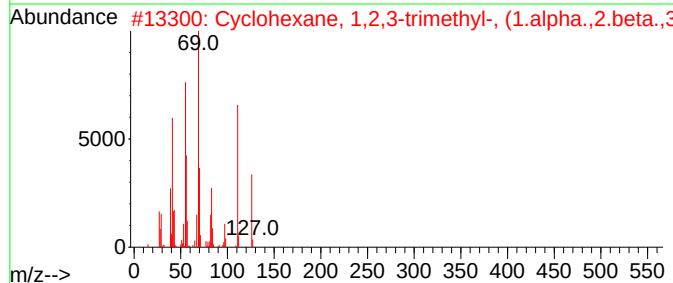
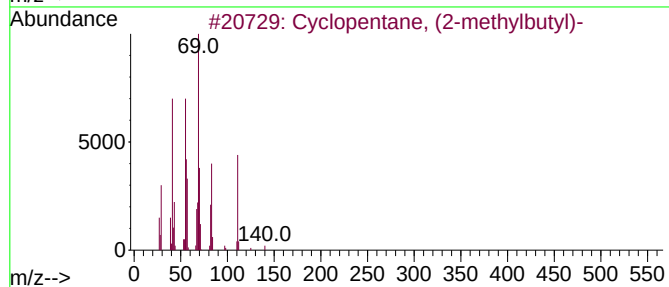
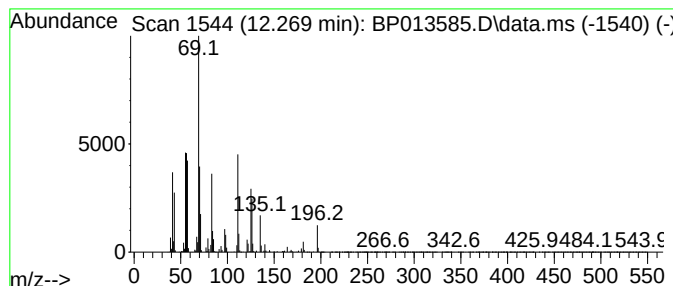
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 13 Cyclopentane, (2-methylbutyl)- Concentration Rank 12

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 12.269 | 18.60 ng | 4382260 | Naphthalene-d8   | 10.993 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclopentane, (2-methylbutyl)-      | 140 | C10H20  | 053366-38-4 | 64   |
| 2    |    |   | Cyclohexane, 1,2,3-trimethyl-, (... | 126 | C9H18   | 001678-81-5 | 53   |
| 3    |    |   | Cyclohexanone, 3,5-dimethyl-        | 126 | C8H14O  | 002320-30-1 | 47   |
| 4    |    |   | trans-3,5-Dimethylcyclohexanone     | 126 | C8H14O  | 007214-49-5 | 47   |
| 5    |    |   | 3,4-Dimethyl cyclohexanone          | 126 | C8H14O  | 005465-09-8 | 43   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

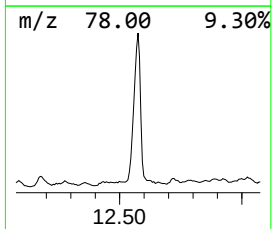
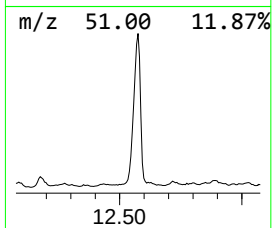
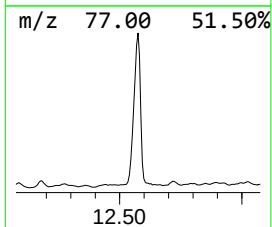
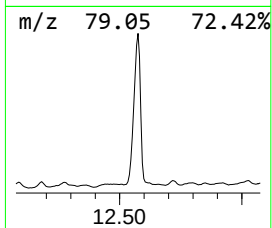
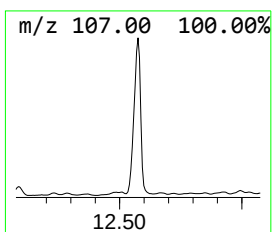
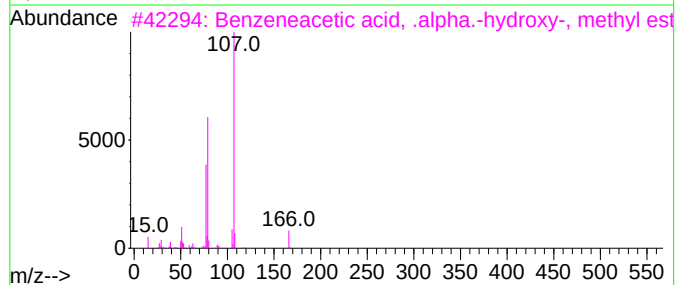
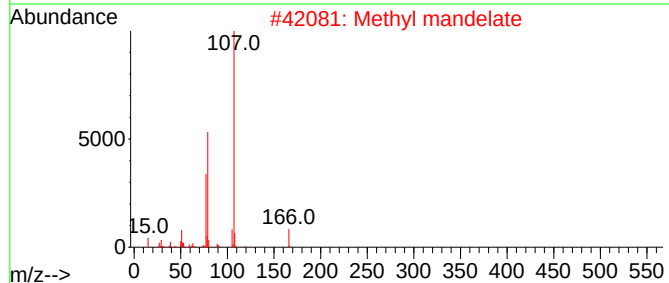
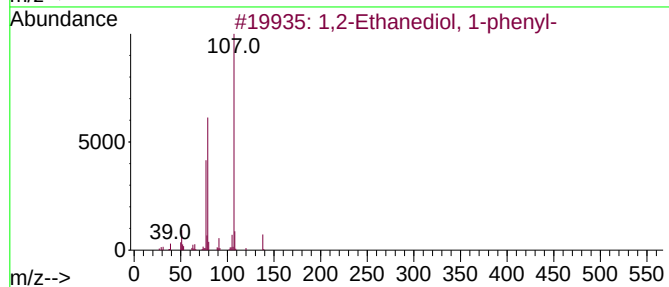
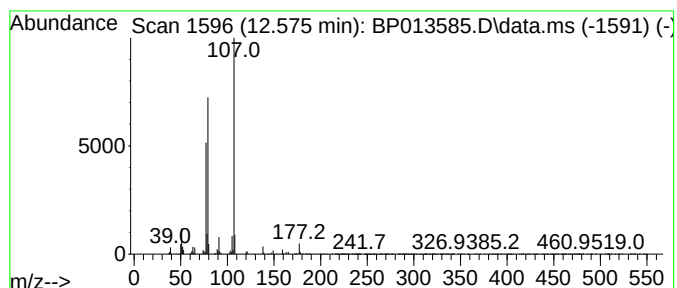
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*

Peak Number 14 1,2-Ethanediol, 1-phenyl- Concentration Rank 1

| R.T.      | EstConc                             | Area     | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|----------|------------------|-------------|------|
| 12.575    | 80.82 ng                            | 19046200 | Naphthalene-d8   | 10.993      |      |
| Hit# of 5 | Tentative ID                        | MW       | MolForm          | CAS#        | Qual |
| 1         | 1,2-Ethanediol, 1-phenyl-           | 138      | C8H10O2          | 000093-56-1 | 94   |
| 2         | Methyl mandelate                    | 166      | C9H10O3          | 004358-87-6 | 87   |
| 3         | Benzeneacetic acid, .alpha.-hydr... | 166      | C9H10O3          | 021210-43-5 | 86   |
| 4         | Benzeneacetic acid, .alpha.-hydr... | 166      | C9H10O3          | 020698-91-3 | 80   |
| 5         | 2-Methyl-1-phenylbut-3-en-1-ol      | 162      | C11H14O          | 025201-44-9 | 78   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

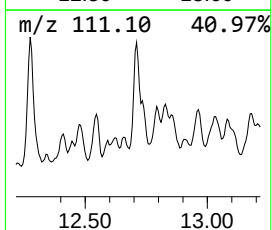
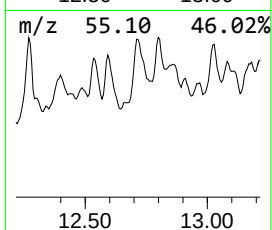
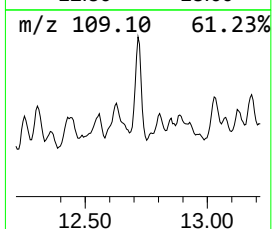
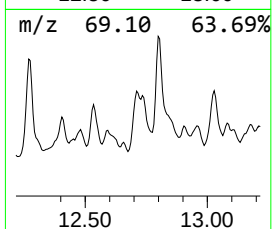
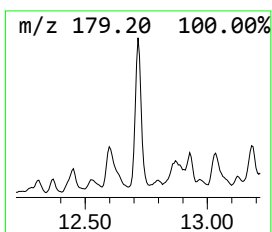
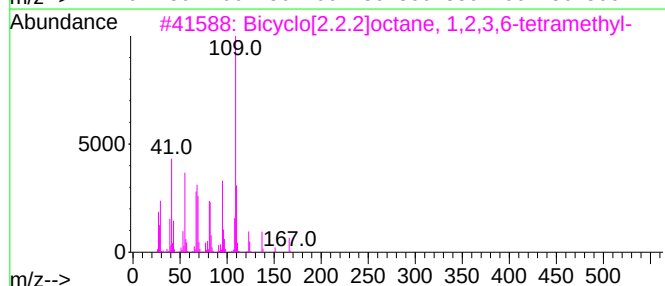
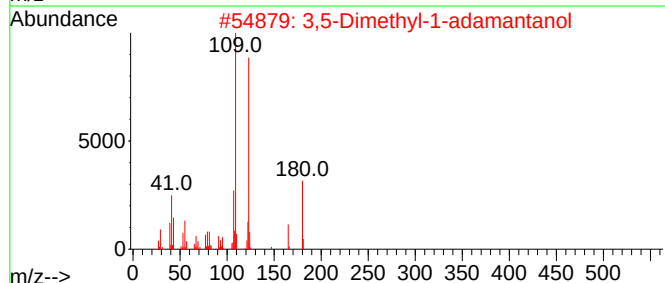
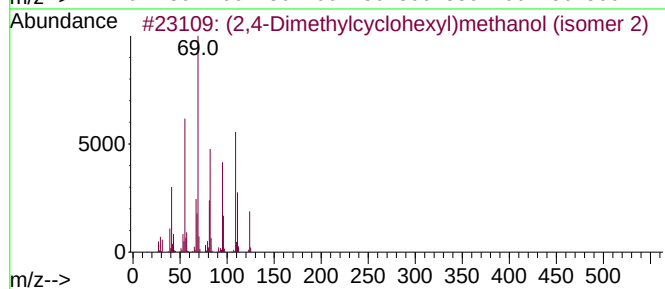
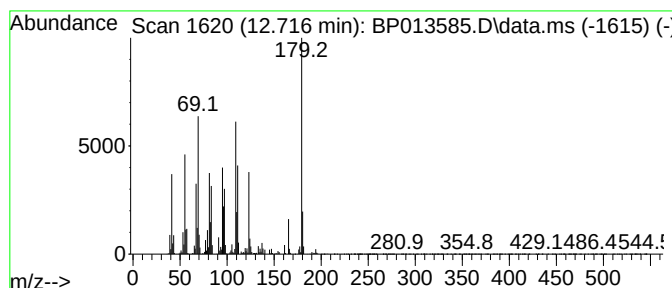
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 15 unknown12.716 Concentration Rank 4

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 12.716 | 42.41 ng | 9993930 | Naphthalene-d8   | 10.993 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | (2,4-Dimethylcyclohexyl)methanol... | 142 | C9H18O  | 1000499-02-7 | 25   |
| 2    |    |   | 3,5-Dimethyl-1-adamantanol          | 180 | C12H20O | 000707-37-9  | 14   |
| 3    |    |   | Bicyclo[2.2.2]octane, 1,2,3,6-te... | 166 | C12H22  | 062338-45-8  | 14   |
| 4    |    |   | Cyclopentaneethanol, .beta.,2,3-... | 156 | C10H20O | 021721-18-6  | 14   |
| 5    |    |   | Naphthalene, decahydro-1,6-dimet... | 208 | C15H28  | 029788-41-8  | 12   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

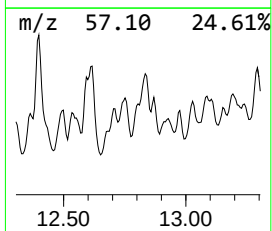
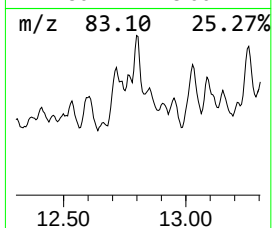
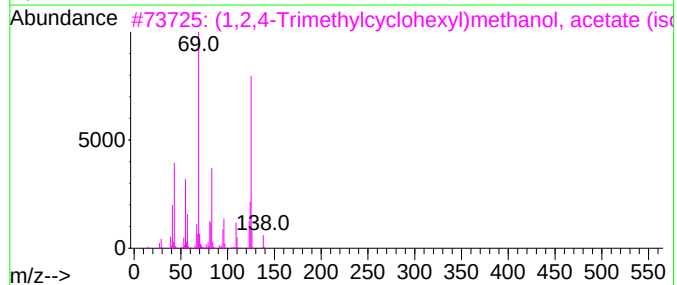
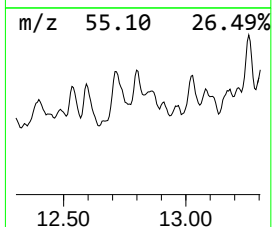
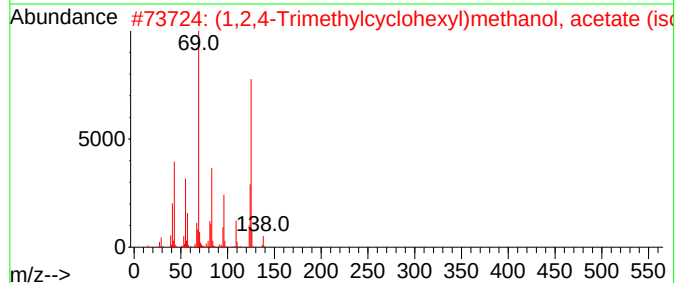
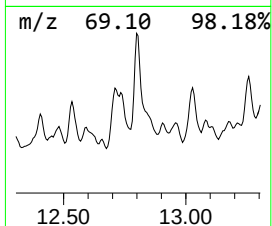
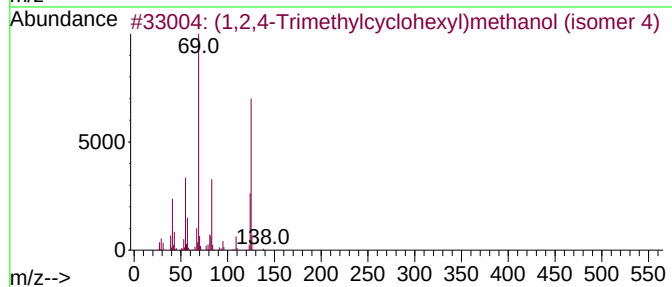
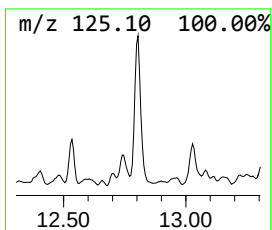
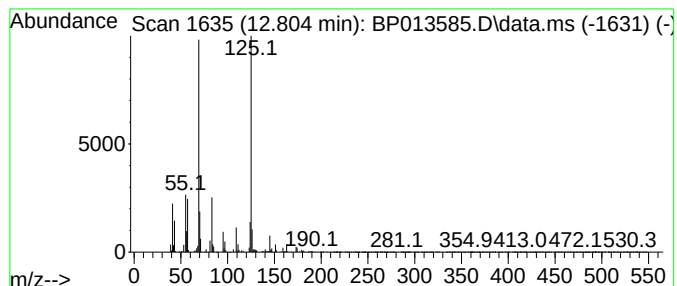
\*\*\*\*\*

Peak Number 16 (1,2,4-Trimethylcyclohexyl)... Concentration Rank 13

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 12.804 | 15.96 ng | 3760590 | Naphthalene-d8   | 10.993 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|------|-------------------------------------|-----|----------|--------------|------|
| 1    |      | (1,2,4-Trimethylcyclohexyl)metha... | 156 | C10H20O  | 1000499-04-1 | 72   |
| 2    |      | (1,2,4-Trimethylcyclohexyl)metha... | 198 | C12H22O2 | 1000499-04-5 | 72   |
| 3    |      | (1,2,4-Trimethylcyclohexyl)metha... | 198 | C12H22O2 | 1010499-04-6 | 64   |
| 4    |      | (1,2,4-Trimethylcyclohexyl)metha... | 156 | C10H20O  | 1000499-03-8 | 56   |
| 5    |      | Zinc, bis[2-(1,1-dimethylethyl)-... | 314 | C18H34Zn | 074793-36-5  | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

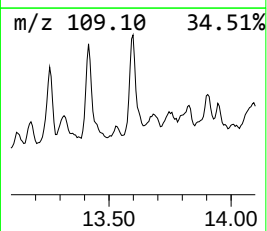
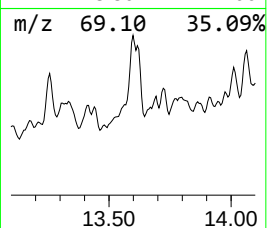
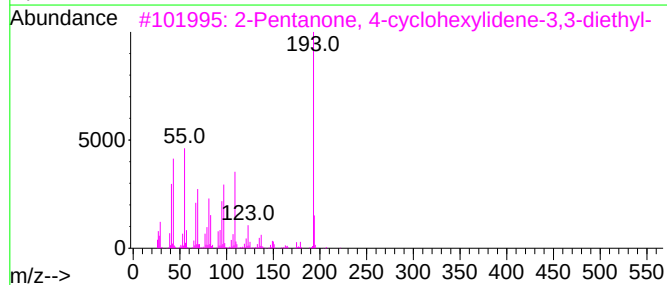
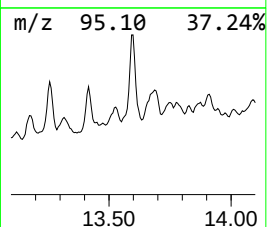
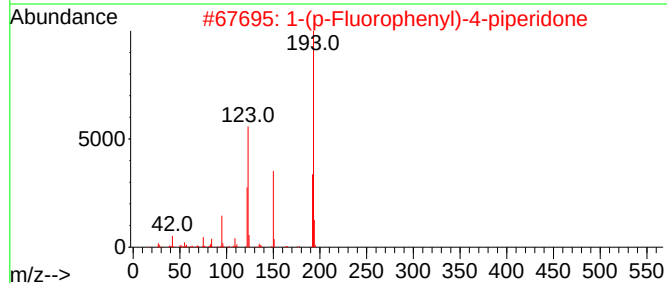
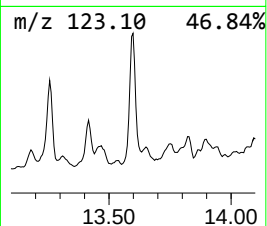
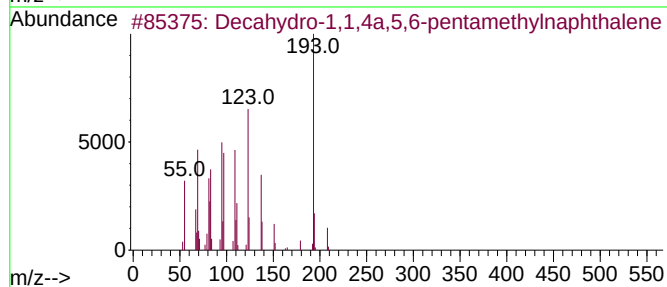
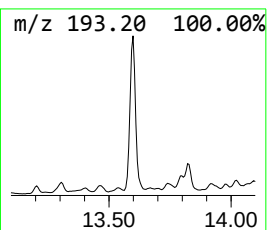
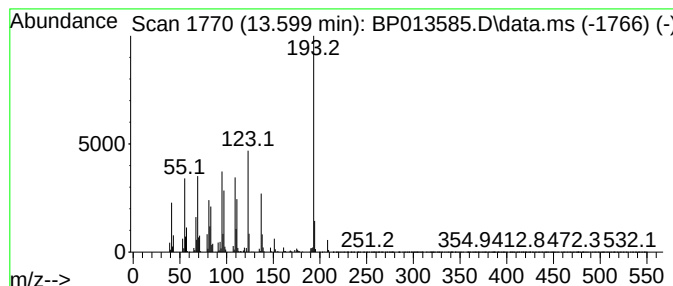
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*

Peak Number 17 Decahydro-1,1,4a,5,6-pentam... Concentration Rank 18

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|--------------|------|
| 13.599    | 12.44 ng                            | 9617160 | Acenaphthene-d10 | 14.804       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#         | Qual |
| 1         | Decahydro-1,1,4a,5,6-pentamethyl... | 208     | C15H28           | 080655-44-3  | 55   |
| 2         | 1-(p-Fluorophenyl)-4-piperidone     | 193     | C11H12FNO        | 1000238-56-7 | 49   |
| 3         | 2-Pentanone, 4-cyclohexylidene-3... | 222     | C15H26O          | 313253-65-5  | 40   |
| 4         | Borinic acid, diethyl-, 3,3,5-tr... | 208     | C13H25BO         | 057387-76-5  | 37   |
| 5         | N-(4-Fluorophenyl)succinamic acid   | 211     | C10H10FNO3       | 199461-14-8  | 37   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

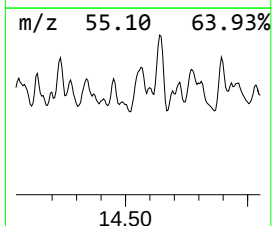
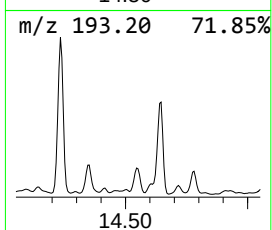
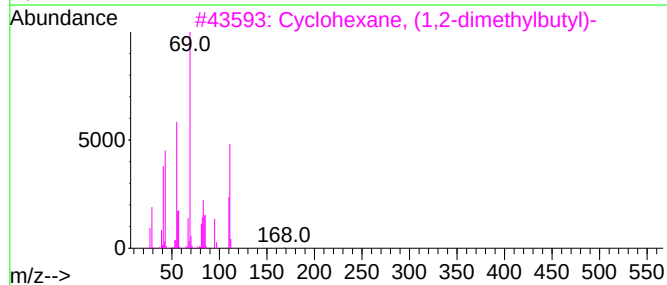
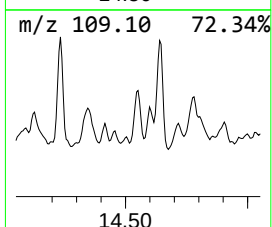
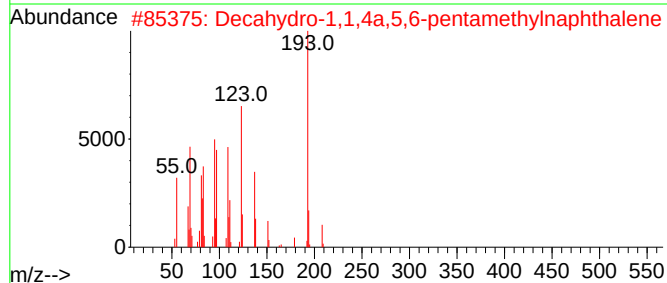
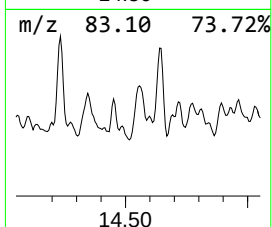
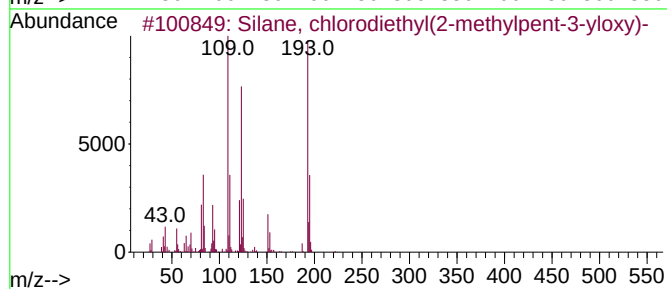
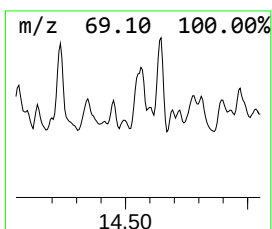
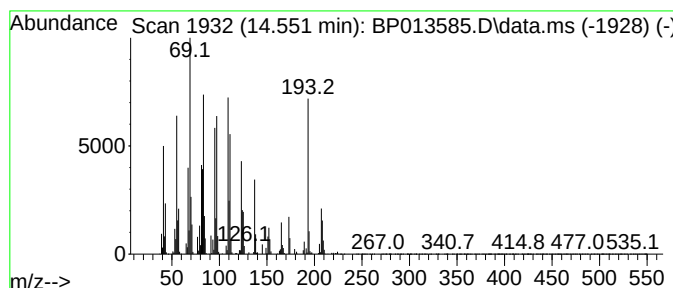
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 18 unknown14.551 Concentration Rank 17

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 14.551 | 12.66 ng | 9784880 | Acenaphthene-d10 | 14.804 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Silane, chlorodiethyl(2-methylpe... | 222 | C10H23ClOSi | 1000363-79-1 | 45   |
| 2    |    |   | Decahydro-1,1,4a,5,6-pentamethyl... | 208 | C15H28      | 080655-44-3  | 41   |
| 3    |    |   | Cyclohexane, (1,2-dimethylbutyl)-   | 168 | C12H24      | 061142-37-8  | 38   |
| 4    |    |   | Cyclopentane, 1,1,3-trimethyl-      | 112 | C8H16       | 004516-69-2  | 38   |
| 5    |    |   | 1-Cyclohexyl-1-(4-methylcyclohex... | 208 | C15H28      | 002027-12-5  | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

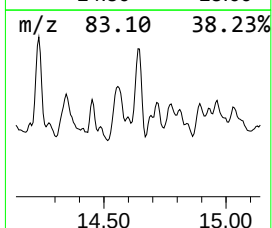
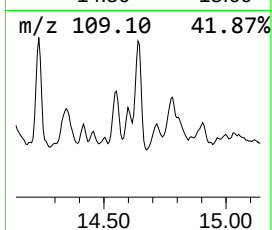
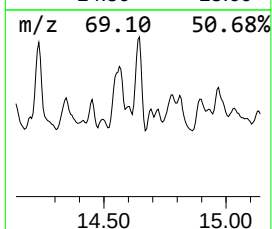
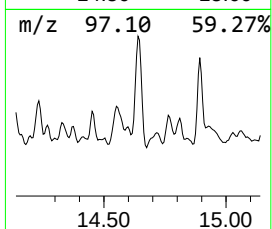
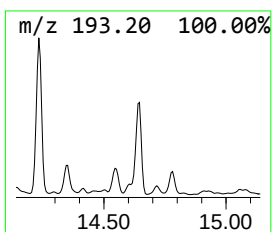
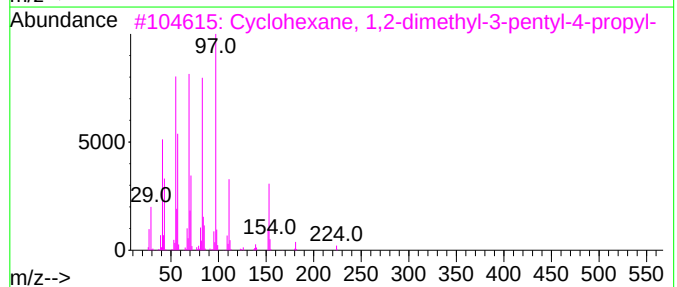
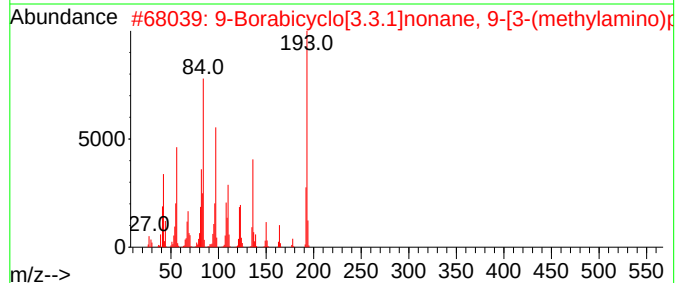
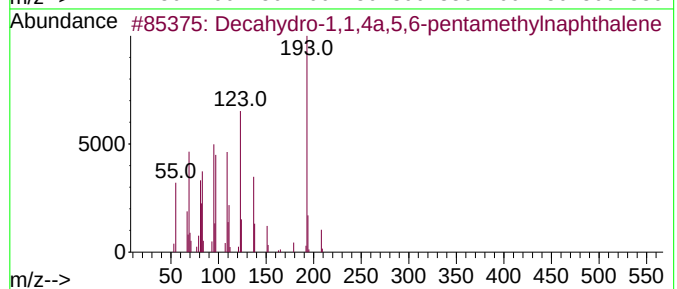
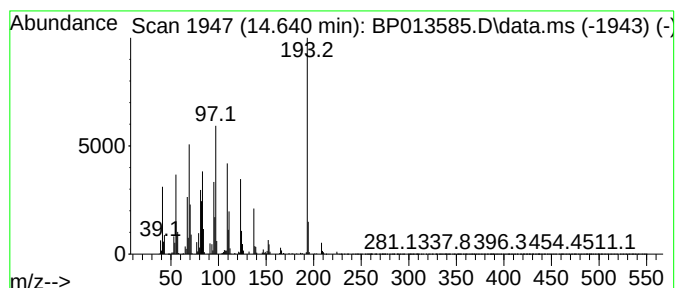
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*

Peak Number 19 unknown14.640 Concentration Rank 11

| R.T.   | EstConc  | Area                                | Relative to ISTD | R.T.            |
|--------|----------|-------------------------------------|------------------|-----------------|
| 14.640 | 20.00 ng | 15463300                            | Acenaphthene-d10 | 14.804          |
| Hit#   | of 5     | Tentative ID                        | MW MolForm       | CAS# Qual       |
| 1      |          | Decahydro-1,1,4a,5,6-pentamethyl... | 208 C15H28       | 080655-44-3 49  |
| 2      |          | 9-Borabicyclo[3.3.1]nonane, 9-[3... | 193 C12H24BN     | 1010162-56-0 38 |
| 3      |          | Cyclohexane, 1,2-dimethyl-3-pent... | 224 C16H32       | 062376-17-4 38  |
| 4      |          | 2-Anthracenamine                    | 193 C14H11N      | 000613-13-8 35  |
| 5      |          | [1,1'-Biphenyl]-4-acetonitrile      | 193 C14H11N      | 031603-77-7 18  |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

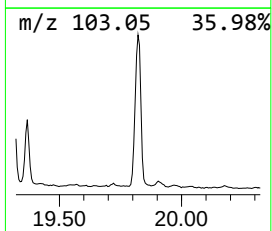
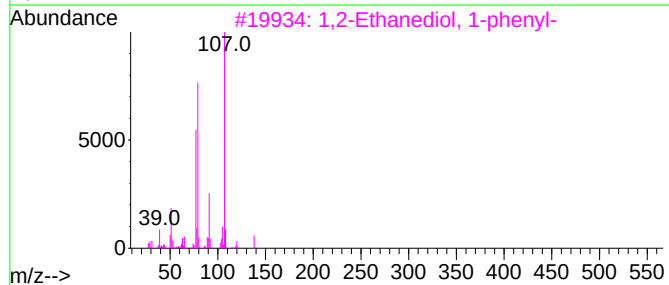
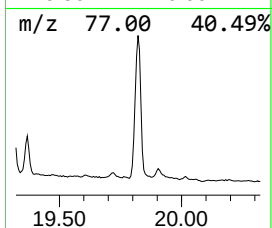
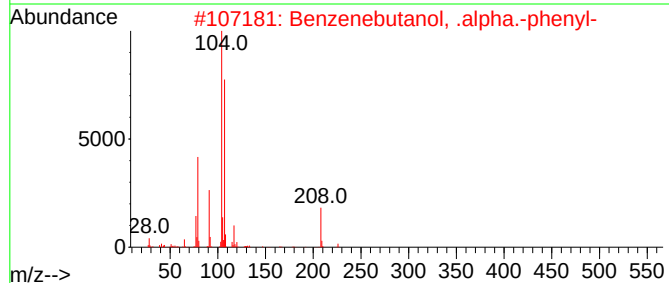
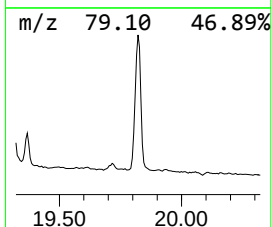
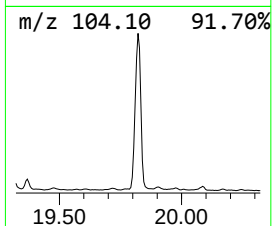
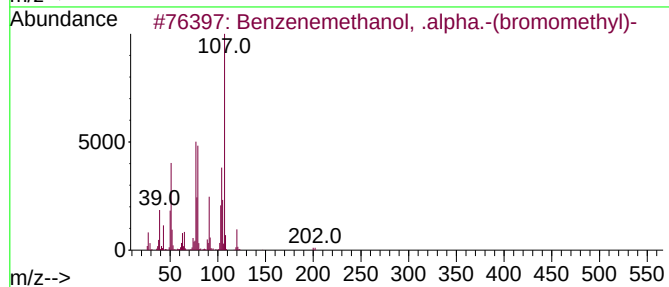
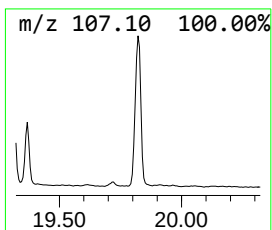
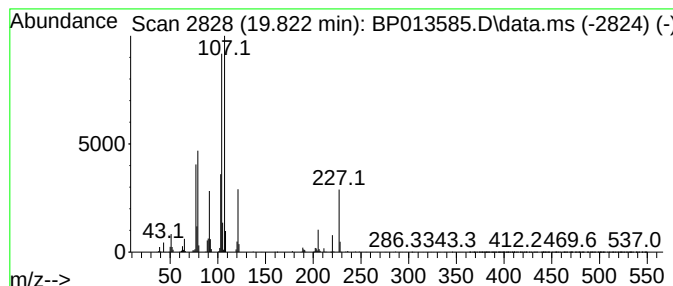
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 20 Benzenemethanol, .alpha.-(b... Concentration Rank 3

| R.T.   | EstConc  | Area     | Relative to ISTD | R.T.   |
|--------|----------|----------|------------------|--------|
| 19.822 | 57.64 ng | 11529500 | Chrysene-d12     | 21.628 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Benzenemethanol, .alpha.-(bromom... | 200 | C8H9BrO | 002425-28-7 | 59   |
| 2    |      | Benzenebutanol, .alpha.-phenyl-     | 226 | C16H18O | 030078-89-8 | 53   |
| 3    |      | 1,2-Ethanediol, 1-phenyl-           | 138 | C8H10O2 | 000093-56-1 | 43   |
| 4    |      | Mandelamide                         | 151 | C8H9NO2 | 004410-31-5 | 43   |
| 5    |      | Benzenemethanol, .alpha.-pentyl-    | 178 | C12H18O | 004471-05-0 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012423\  
Data File : BP013585.D  
Acq On : 24 Jan 2023 23:18  
Operator : CG/JU  
Sample : 01190-12  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
WS-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP012323.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...  | 3.223  | 64.1    | ng    | 7787010  | 1                     | 8.158  | 2428070  | 20.0 |
| 2-Pentanone, 4-...  | 5.217  | 13.5    | ng    | 1635110  | 1                     | 8.158  | 2428070  | 20.0 |
| Benzenemethanol...  | 8.846  | 11.1    | ng    | 1346730  | 1                     | 8.158  | 2428070  | 20.0 |
| Cyclohexane, 1,...  | 9.275  | 24.8    | ng    | 3014720  | 1                     | 8.158  | 2428070  | 20.0 |
| Carbonic acid, ...  | 9.393  | 15.8    | ng    | 1915430  | 1                     | 8.158  | 2428070  | 20.0 |
| Undecane, 2,6-d...  | 11.146 | 11.6    | ng    | 2740500  | 2                     | 10.993 | 4713310  | 20.0 |
| Cyclohexane, 2-...  | 11.487 | 21.4    | ng    | 5046580  | 2                     | 10.993 | 4713310  | 20.0 |
| unknown11.757       | 11.757 | 14.1    | ng    | 3310450  | 2                     | 10.993 | 4713310  | 20.0 |
| 5-Dodecene, (Z)-    | 11.922 | 21.3    | ng    | 5007590  | 2                     | 10.993 | 4713310  | 20.0 |
| Cyclohexane, 1,...  | 12.046 | 36.3    | ng    | 8563500  | 2                     | 10.993 | 4713310  | 20.0 |
| unknown12.093       | 12.093 | 25.4    | ng    | 5981000  | 2                     | 10.993 | 4713310  | 20.0 |
| unknown12.181       | 12.181 | 29.8    | ng    | 7023350  | 2                     | 10.993 | 4713310  | 20.0 |
| Cyclopentane, (...) | 12.269 | 18.6    | ng    | 4382260  | 2                     | 10.993 | 4713310  | 20.0 |
| 1,2-Ethanediol,...  | 12.575 | 80.8    | ng    | 19046200 | 2                     | 10.993 | 4713310  | 20.0 |
| unknown12.716       | 12.716 | 42.4    | ng    | 9993930  | 2                     | 10.993 | 4713310  | 20.0 |
| (1,2,4-Trimethy...  | 12.804 | 16.0    | ng    | 3760590  | 2                     | 10.993 | 4713310  | 20.0 |
| Decahydro-1,1,4...  | 13.599 | 12.4    | ng    | 9617160  | 3                     | 14.804 | 15463300 | 20.0 |
| unknown14.551       | 14.551 | 12.7    | ng    | 9784880  | 3                     | 14.804 | 15463300 | 20.0 |
| unknown14.640       | 14.640 | 20.0    | ng    | 15463300 | 3                     | 14.804 | 15463300 | 20.0 |
| Benzenemethanol...  | 19.822 | 57.6    | ng    | 11529500 | 5                     | 21.628 | 4000750  | 20.0 |