

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM041625\  
 Data File : BM049958.D  
 Acq On : 16 Apr 2025 18:42  
 Operator : RC/JU  
 Sample : Q1808-03  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 LAW-25-0060

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\_M\Methods\8270-BM040825.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM049958.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         | 4.887       | 317           | 323         | 333          | rBV      | 2136189        | 2878414       | 20.92%          | 3.458%        |
| 2         | 5.357       | 396           | 403         | 421          | rBV      | 4438729        | 6471837       | 47.04%          | 7.775%        |
| 3         | 5.963       | 500           | 506         | 518          | rVB      | 514186         | 744679        | 5.41%           | 0.895%        |
| 4         | 6.951       | 666           | 674         | 689          | rBV      | 3966345        | 6610264       | 48.05%          | 7.941%        |
| 5         | 7.769       | 805           | 813         | 821          | rBV      | 1450156        | 2293738       | 16.67%          | 2.755%        |
| 6         | 8.922       | 1002          | 1009        | 1022         | rBV      | 2433830        | 4203268       | 30.55%          | 5.049%        |
| 7         | 10.563      | 1280          | 1288        | 1300         | rBV      | 1692094        | 2847795       | 20.70%          | 3.421%        |
| 8         | 13.033      | 1700          | 1708        | 1728         | rBV      | 7443595        | 11033570      | 80.20%          | 13.255%       |
| 9         | 14.416      | 1934          | 1943        | 1951         | rBV      | 2593226        | 3963670       | 28.81%          | 4.762%        |
| 10        | 14.939      | 2028          | 2032        | 2041         | rBV3     | 74645          | 158390        | 1.15%           | 0.190%        |
| 11        | 15.204      | 2072          | 2077        | 2085         | rBV3     | 200857         | 353394        | 2.57%           | 0.425%        |
| 12        | 15.768      | 2168          | 2173        | 2180         | rBV      | 1165226        | 1736835       | 12.62%          | 2.086%        |
| 13        | 15.904      | 2190          | 2196        | 2203         | rBV      | 6083591        | 8765104       | 63.71%          | 10.530%       |
| 14        | 16.157      | 2235          | 2239        | 2244         | rVB      | 711849         | 1028450       | 7.48%           | 1.235%        |
| 15        | 17.157      | 2405          | 2409        | 2415         | rBV2     | 3057781        | 4558097       | 33.13%          | 5.476%        |
| 16        | 18.068      | 2560          | 2564        | 2568         | rBV2     | 2105304        | 3077294       | 22.37%          | 3.697%        |
| 17        | 19.786      | 2852          | 2856        | 2865         | rVB      | 10987191       | 13757910      | 100.00%         | 16.527%       |
| 18        | 21.403      | 3127          | 3131        | 3136         | rVB      | 2918748        | 4051932       | 29.45%          | 4.868%        |
| 19        | 24.403      | 3634          | 3641        | 3652         | rVB      | 1899167        | 4708432       | 34.22%          | 5.656%        |

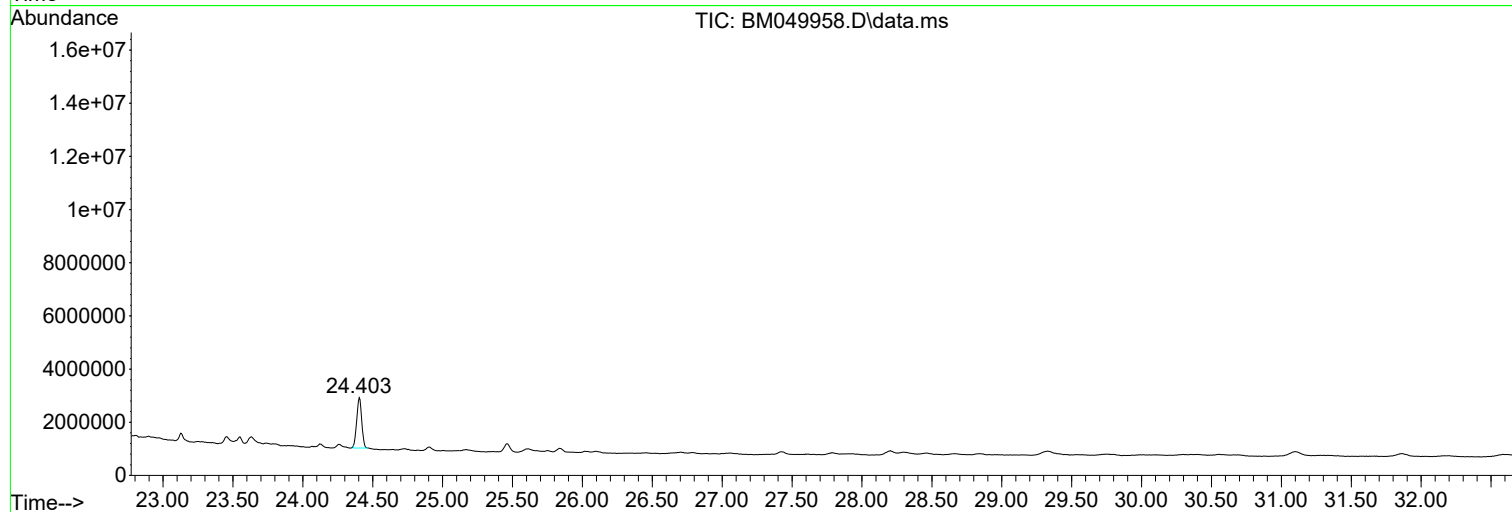
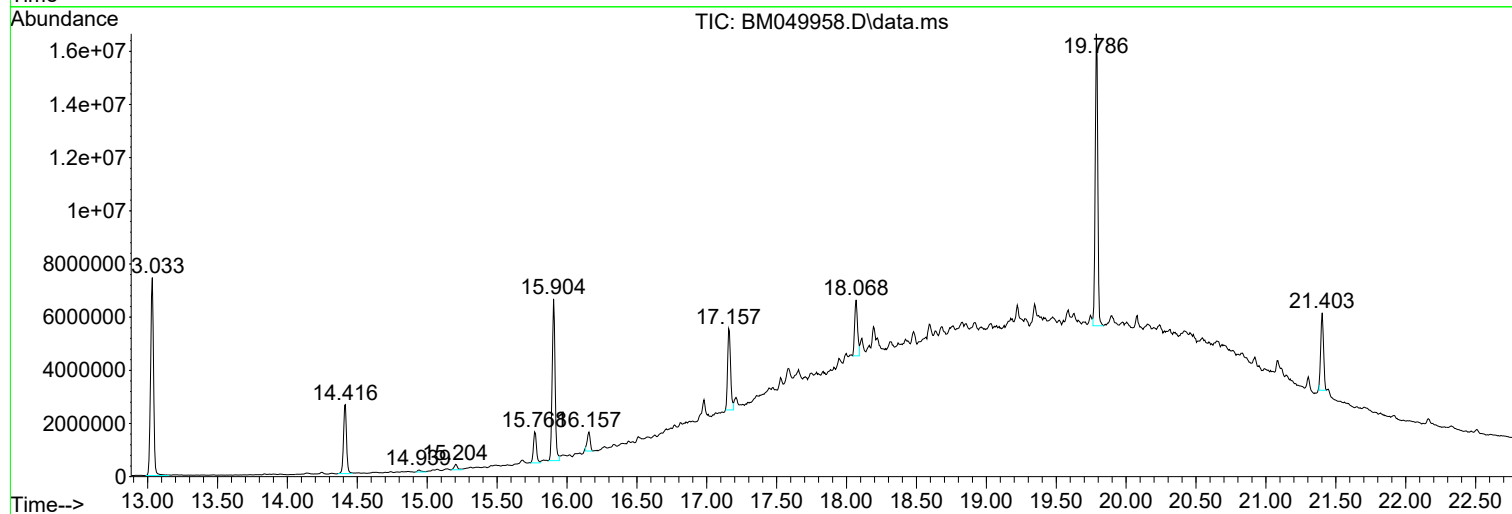
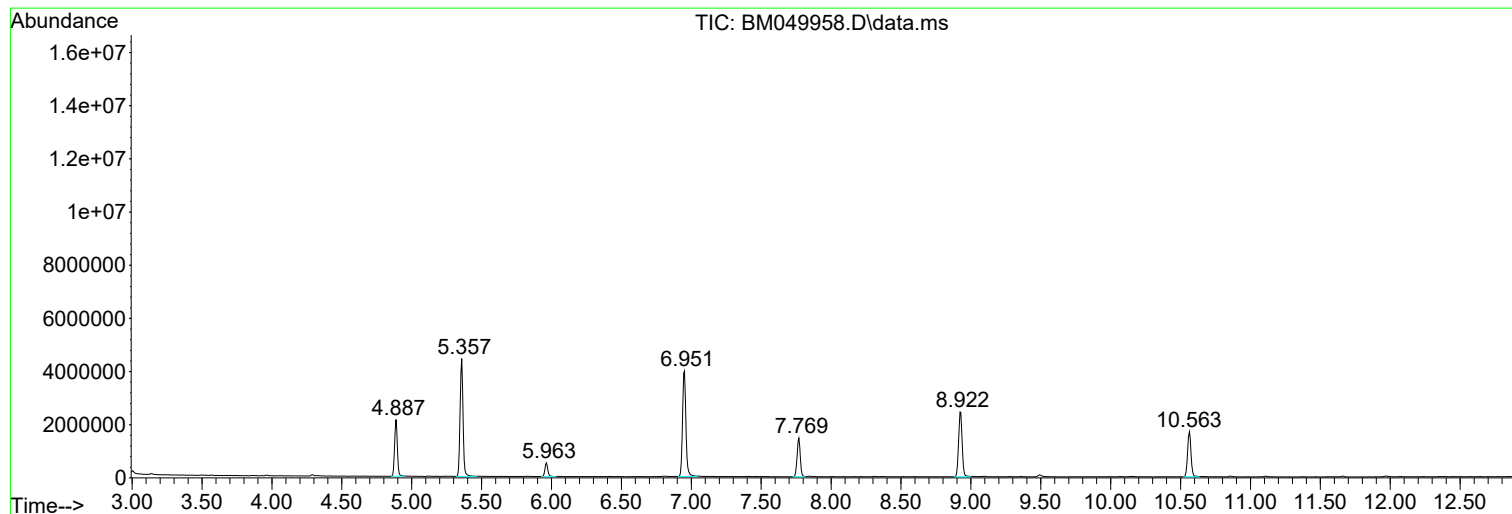
Sum of corrected areas: 83243073

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM041625\  
Data File : BM049958.D  
Acq On : 16 Apr 2025 18:42  
Operator : RC/JU  
Sample : Q1808-03  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
LAW-25-0060

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM040825.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\_M\Data\BM041625\  
Data File : BM049958.D  
Acq On : 16 Apr 2025 18:42  
Operator : RC/JU  
Sample : Q1808-03  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
LAW-25-0060

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM040825.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

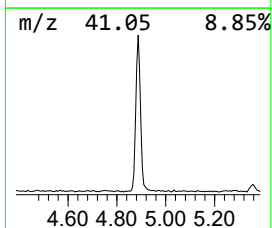
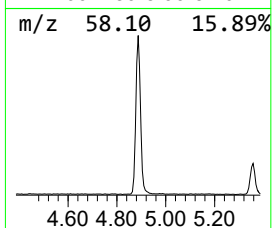
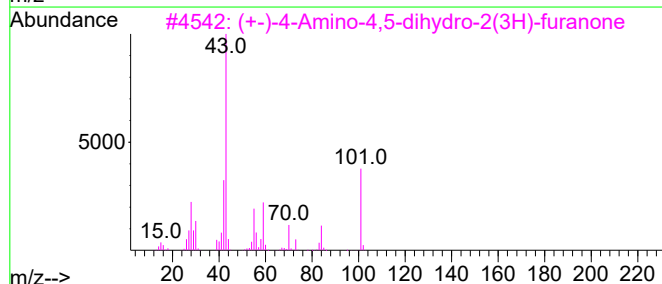
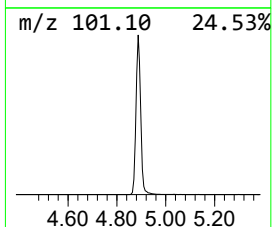
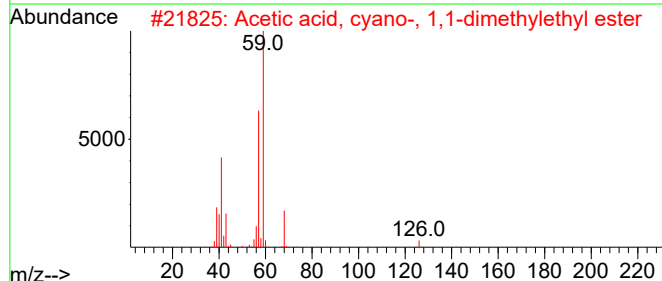
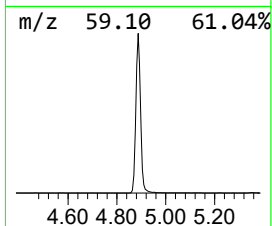
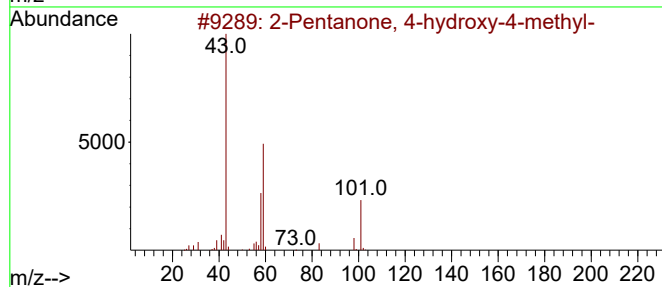
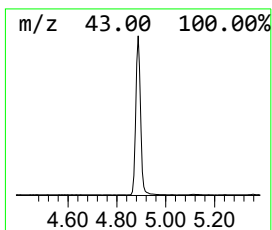
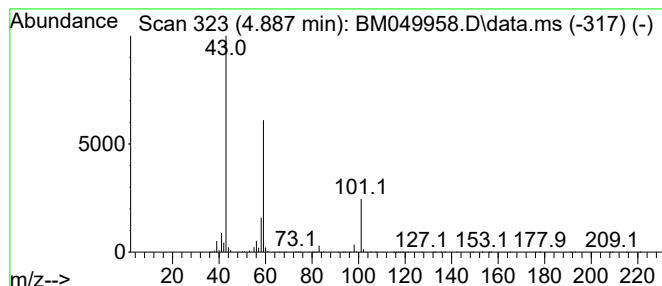
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 4.887 | 25.10 ng | 2878410 | 1,4-Dichlorobenzene-d4 | 7.769 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2      | 000123-42-2 | 64      |      |      |
| 2    | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2     | 001116-98-9 | 17      |      |      |
| 3    | (+)-4-Amino-4,5-dihydro-2(3H)-f...  | 101 | C4H7NO2      | 016504-58-8 | 9       |      |      |
| 4    | Morpholine, 4-methyl-               | 101 | C5H11NO      | 000109-02-4 | 9       |      |      |
| 5    | 5-Hexen-2-one                       | 98  | C6H10O       | 000109-49-9 | 9       |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM041625\  
Data File : BM049958.D  
Acq On : 16 Apr 2025 18:42  
Operator : RC/JU  
Sample : Q1808-03  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
LAW-25-0060

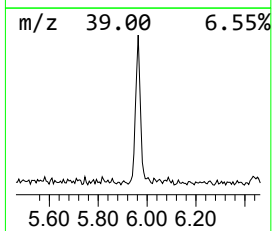
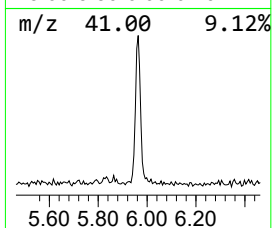
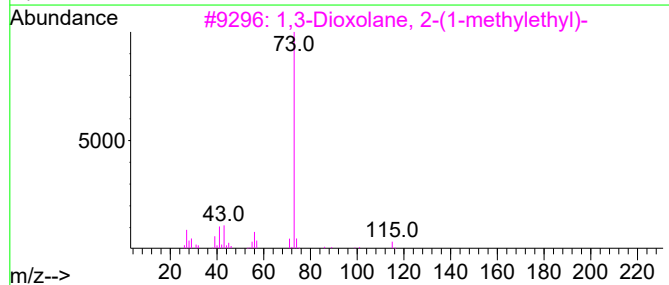
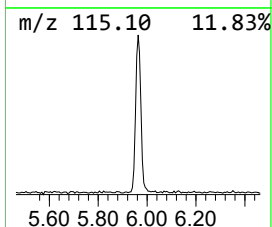
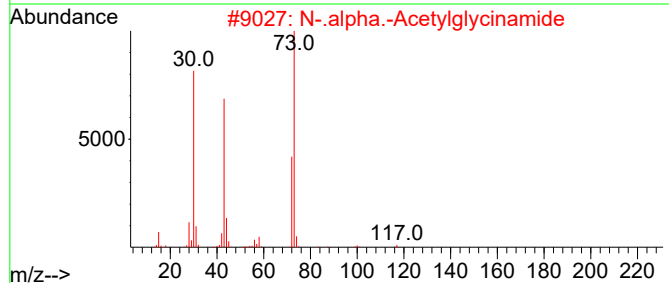
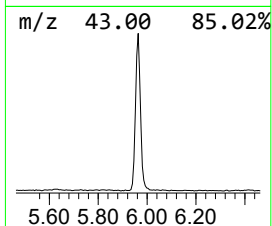
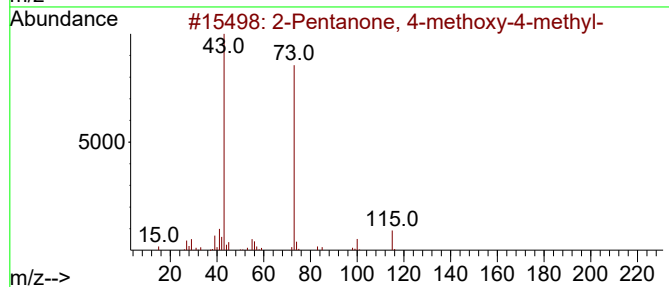
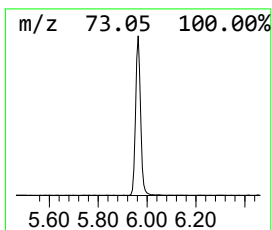
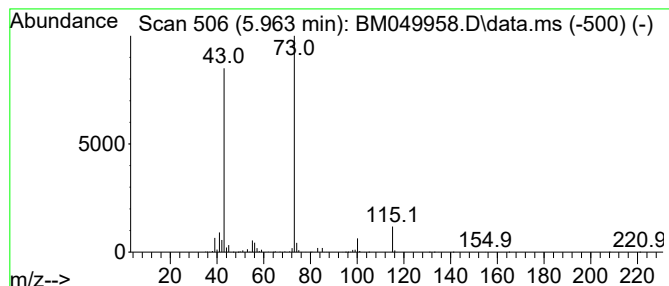
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM040825.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 2-Pentanone, 4-methoxy-4-me... Concentration Rank 4

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 5.963 | 6.49 ng | 744679 | 1,4-Dichlorobenzene-d4 | 7.769 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-----------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-methoxy-4-methyl-  | 130 | C7H14O2  | 000107-70-0 | 83   |
| 2    |    |   | N-.alpha.-Acetylglycinamide       | 116 | C4H8N2O2 | 002620-63-5 | 25   |
| 3    |    |   | 1,3-Dioxolane, 2-(1-methylethyl)- | 116 | C6H12O2  | 000822-83-3 | 10   |
| 4    |    |   | N-Formylmorpholine                | 115 | C5H9NO2  | 004394-85-8 | 9    |
| 5    |    |   | 2-Pentanone, 3-methyl-            | 100 | C6H12O   | 000565-61-7 | 9    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM041625\  
Data File : BM049958.D  
Acq On : 16 Apr 2025 18:42  
Operator : RC/JU  
Sample : Q1808-03  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
LAW-25-0060

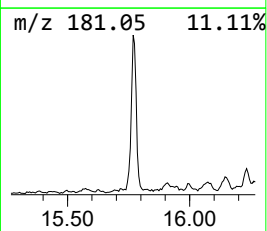
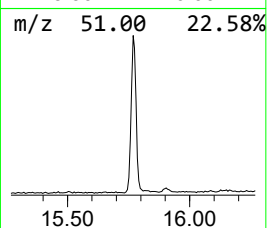
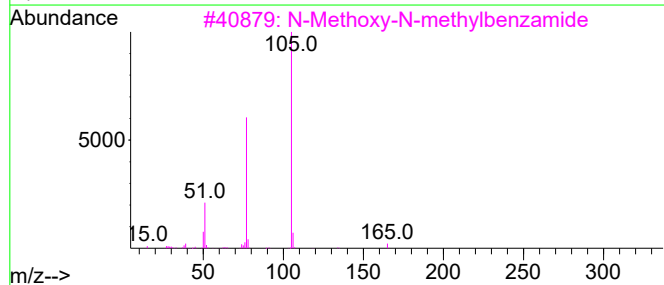
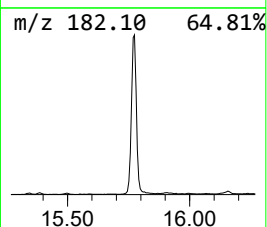
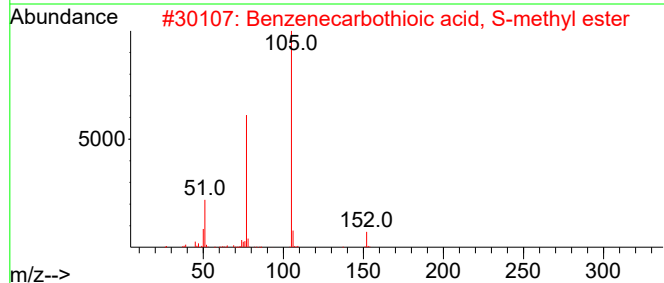
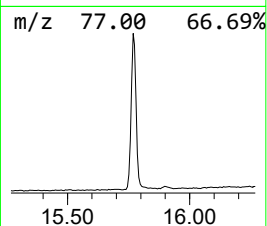
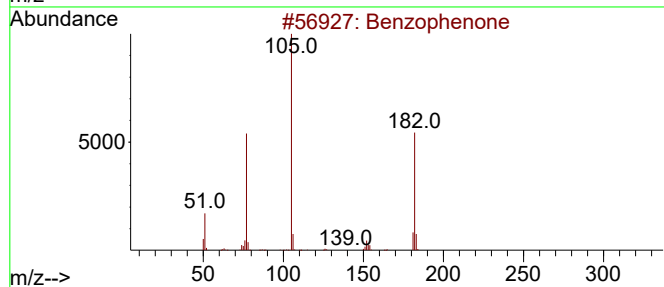
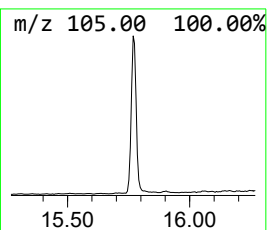
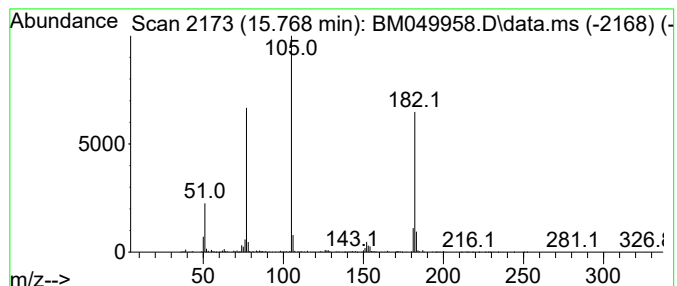
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM040825.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 Benzophenone Concentration Rank 3

| R.T.   | EstConc | Area    | Relative to ISTD | R.T.   |
|--------|---------|---------|------------------|--------|
| 15.769 | 8.76 ng | 1736840 | Acenaphthene-d10 | 14.416 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Benzophenone                        | 182 | C13H10O  | 000119-61-9 | 96   |
| 2    |    |   | Benzenecarbothioic acid, S-methy... | 152 | C8H8OS   | 005925-68-8 | 47   |
| 3    |    |   | N-Methoxy-N-methylbenzamide         | 165 | C9H11NO2 | 006919-61-5 | 46   |
| 4    |    |   | Benzenecarbothioic acid             | 138 | C7H6OS   | 000098-91-9 | 43   |
| 5    |    |   | 1,2-Propanedione, 1-phenyl-         | 148 | C9H8O2   | 000579-07-7 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM041625\  
Data File : BM049958.D  
Acq On : 16 Apr 2025 18:42  
Operator : RC/JU  
Sample : Q1808-03  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
LAW-25-0060

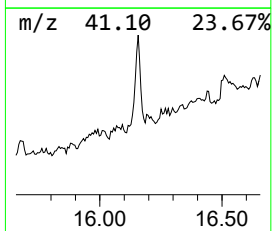
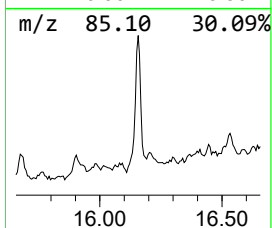
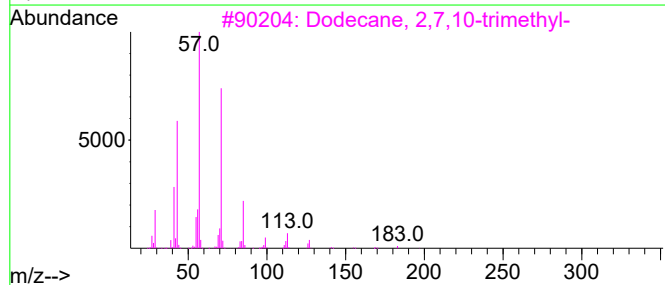
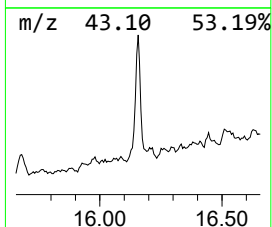
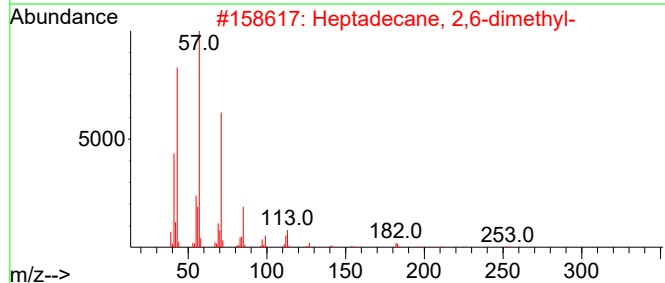
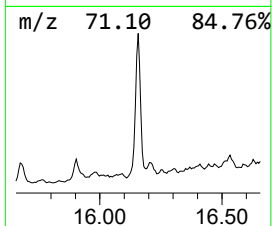
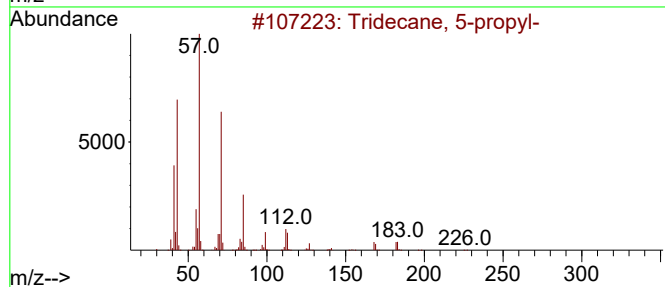
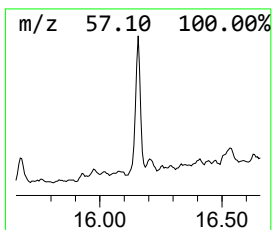
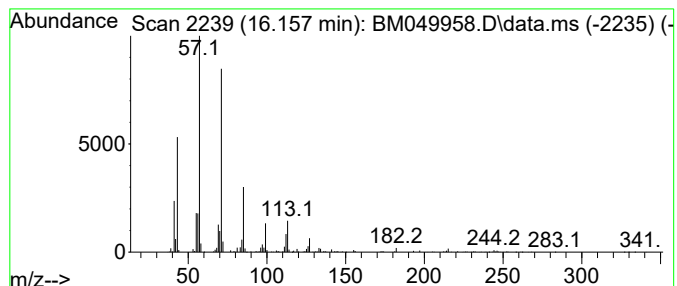
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM040825.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 Tridecane, 5-propyl- Concentration Rank 5

| R.T.   | EstConc | Area    | Relative to ISTD | R.T.   |
|--------|---------|---------|------------------|--------|
| 16.157 | 4.51 ng | 1028450 | Phenanthrene-d10 | 17.157 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Tridecane, 5-propyl-                | 226 | C16H34    | 055045-11-9  | 91   |
| 2    |    |   | Heptadecane, 2,6-dimethyl-          | 268 | C19H40    | 054105-67-8  | 87   |
| 3    |    |   | Dodecane, 2,7,10-trimethyl-         | 212 | C15H32    | 074645-98-0  | 86   |
| 4    |    |   | Sulfurous acid, 2-ethylhexyl non... | 320 | C17H36O3S | 1010309-19-2 | 83   |
| 5    |    |   | Sulfurous acid, decyl 2-ethylhex... | 334 | C18H38O3S | 1000309-19-3 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM041625\  
Data File : BM049958.D  
Acq On : 16 Apr 2025 18:42  
Operator : RC/JU  
Sample : Q1808-03  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
LAW-25-0060

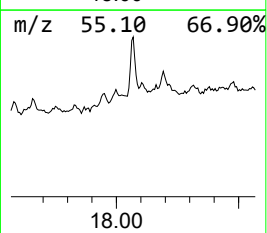
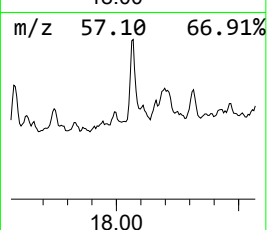
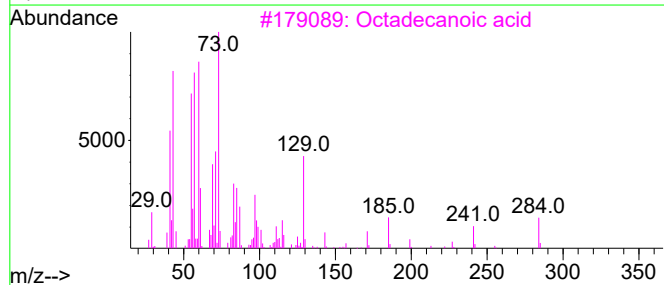
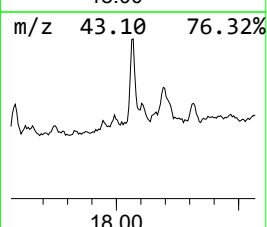
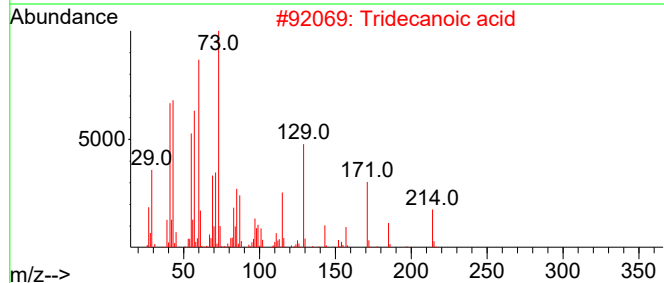
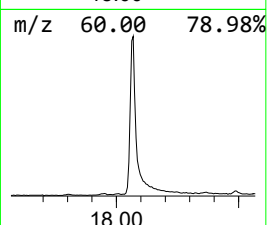
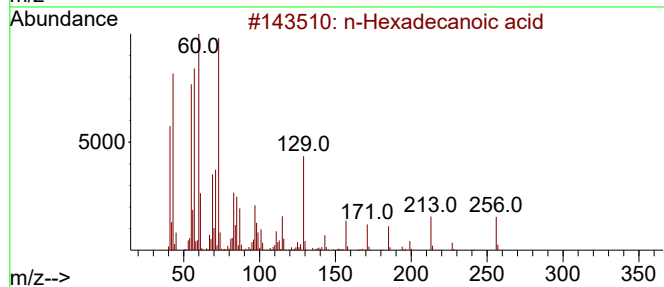
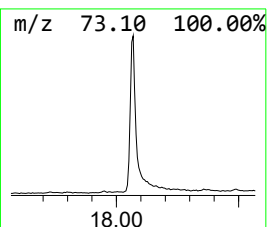
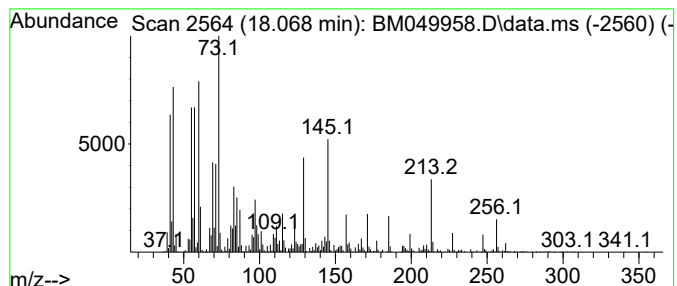
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM040825.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 n-Hexadecanoic acid Concentration Rank 2

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 18.068 | 13.50 ng | 3077290 | Phenanthrene-d10 | 17.157 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|---------------------|-----|----------|-------------|------|
| 1    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 99   |
| 2    |    |   | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 92   |
| 3    |    |   | Octadecanoic acid   | 284 | C18H36O2 | 000057-11-4 | 81   |
| 4    |    |   | Undecanoic acid     | 186 | C11H22O2 | 000112-37-8 | 68   |
| 5    |    |   | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 68   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM041625\  
Data File : BM049958.D  
Acq On : 16 Apr 2025 18:42  
Operator : RC/JU  
Sample : Q1808-03  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
LAW-25-0060

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM040825.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
|                    |        |         |       |          | #                     | RT     | Resp    | Conc |
| 2-Pentanone, 4-... | 4.887  | 25.1    | ng    | 2878410  | 1                     | 7.769  | 2293740 | 20.0 |
| 2-Pentanone, 4-... | 5.963  | 6.5     | ng    | 744679   | 1                     | 7.769  | 2293740 | 20.0 |
| Benzophenone       | 15.769 | 8.8     | ng    | 1736840  | 3                     | 14.416 | 3963670 | 20.0 |
| Tridecane, 5-pr... | 16.157 | 4.5     | ng    | 1028450  | 4                     | 17.157 | 4558100 | 20.0 |
| n-Hexadecanoic ... | 18.068 | 13.5    | ng    | 3077290  | 4                     | 17.157 | 4558100 | 20.0 |