

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082823\  
 Data File : BP016836.D  
 Acq On : 29 Aug 2023 17:21  
 Operator : MA/JU  
 Sample : 04182-05  
 Misc :  
 ALS Vial : 54 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SVOC-GPC-BLANK

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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\SFAM-EPA-BP082423.MA.M

Title : SVOA CALIBRATION

Signal : TIC: BP016836.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         | 1.811       | 274           | 285         | 289          | rBV2     | 16330          | 56744         | 1.17%           | 0.273%        |
| 2         | 2.070       | 326           | 329         | 330          | rBV      | 45743          | 42069         | 0.87%           | 0.202%        |
| 3         | 3.005       | 483           | 488         | 498          | rBV      | 105321         | 199148        | 4.11%           | 0.958%        |
| 4         | 3.087       | 498           | 502         | 511          | rVB      | 67458          | 102147        | 2.11%           | 0.492%        |
| 5         | 4.128       | 676           | 679         | 685          | rBV2     | 18063          | 25918         | 0.54%           | 0.125%        |
| 6         | 8.028       | 1335          | 1342        | 1350         | rBV      | 1121143        | 1764791       | 36.45%          | 8.494%        |
| 7         | 10.857      | 1815          | 1823        | 1831         | rBV      | 1646481        | 2773511       | 57.29%          | 13.349%       |
| 8         | 13.522      | 2272          | 2276        | 2282         | rBV3     | 13471          | 24537         | 0.51%           | 0.118%        |
| 9         | 14.681      | 2466          | 2473        | 2489         | rBV      | 2732316        | 3999405       | 82.61%          | 19.249%       |
| 10        | 17.445      | 2937          | 2943        | 2965         | rBV      | 3115351        | 4841260       | 100.00%         | 23.300%       |
| 11        | 21.545      | 3635          | 3640        | 3649         | rBV      | 2748285        | 3736531       | 77.18%          | 17.983%       |
| 12        | 24.110      | 4070          | 4076        | 4095         | rVB      | 1412515        | 3211592       | 66.34%          | 15.457%       |

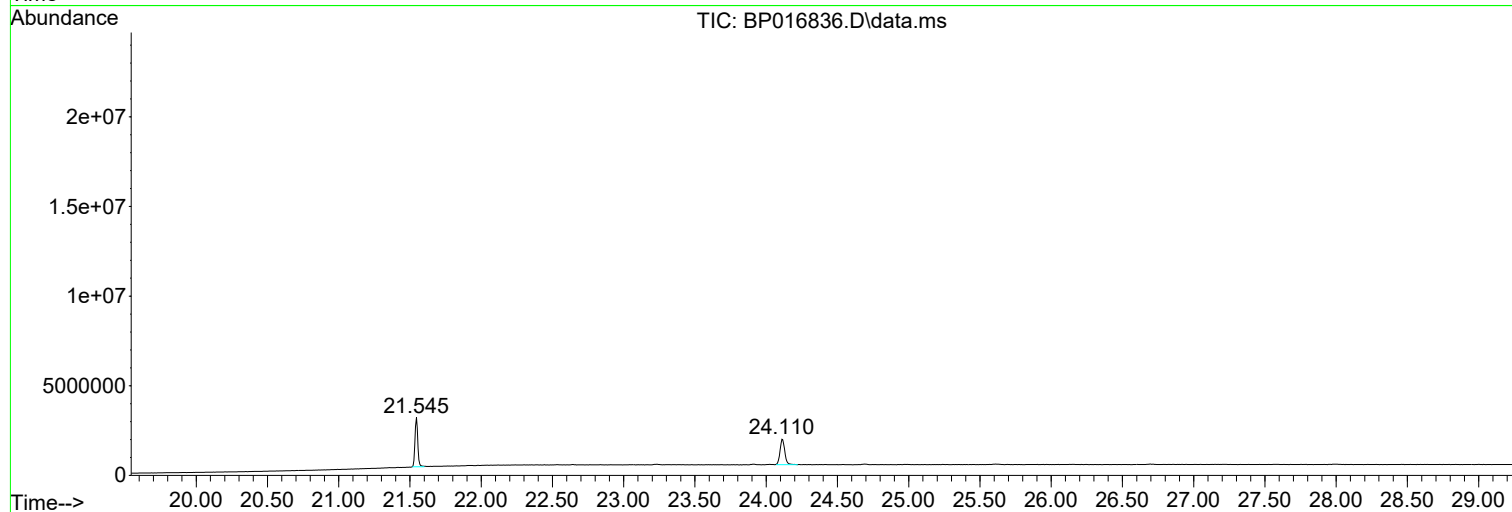
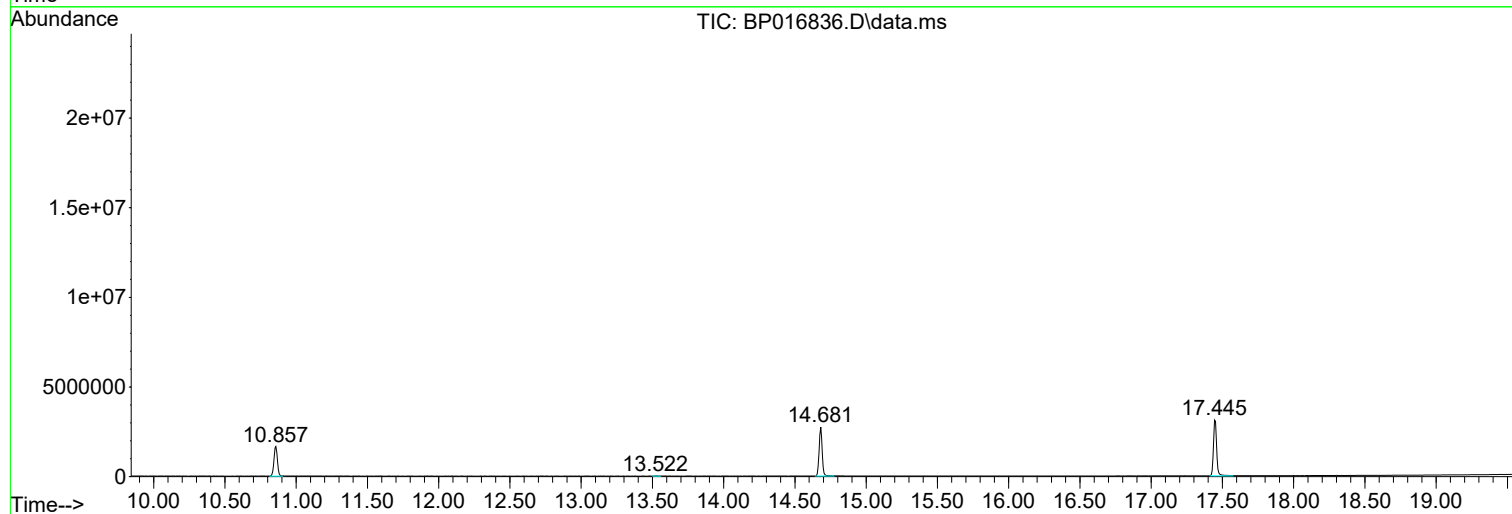
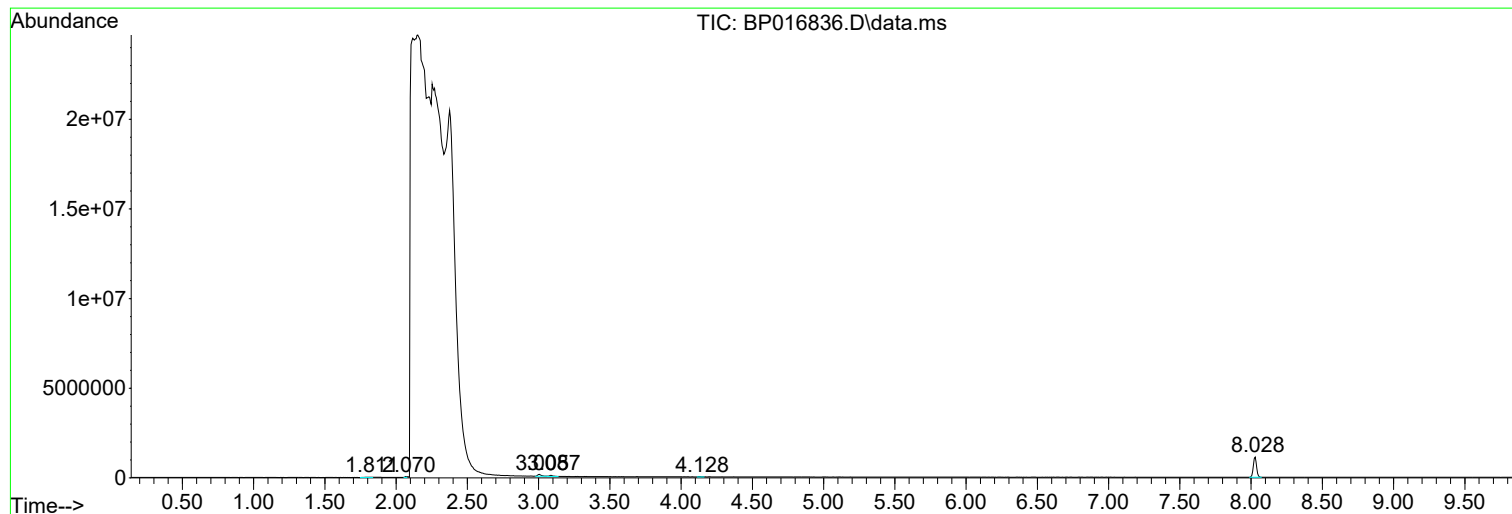
Sum of corrected areas: 20777653

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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP082423.MA.M  
Quant Title : SVOA CALIBRATION

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



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Quant Title : SVOA CALIBRATION

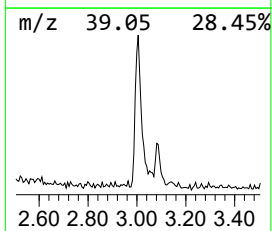
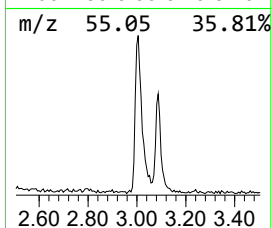
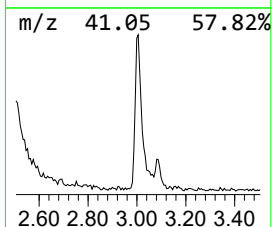
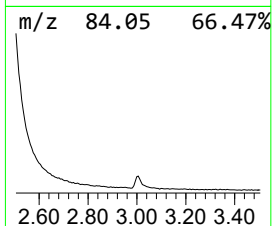
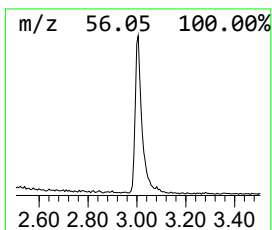
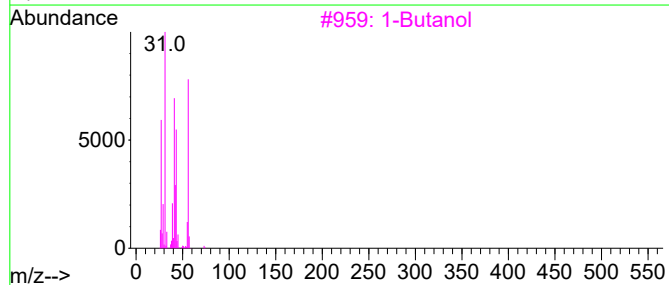
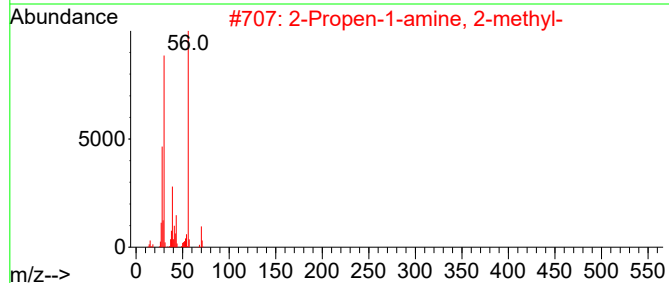
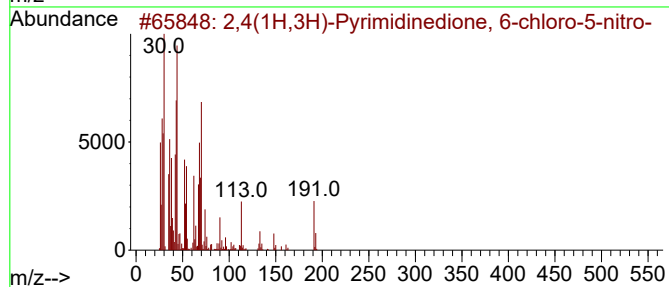
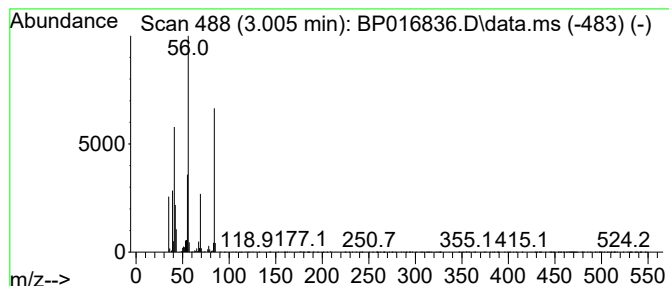
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 unknown-01 Concentration Rank 1

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 3.005 | 2.26 ng/ul | 199148 | 1,4-Dichlorobenzene-d4 | 8.028 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2,4(1H,3H)-Pyrimidinedione, 6-ch... | 191 | C4H2ClN3O4   | 006630-30-4 | 12      |      |      |
| 2    | 2-Propen-1-amine, 2-methyl-         | 71  | C4H9N        | 002878-14-0 | 10      |      |      |
| 3    | 1-Butanol                           | 74  | C4H10O       | 000071-36-3 | 5       |      |      |
| 4    | 1-Butanol                           | 74  | C4H10O       | 000071-36-3 | 5       |      |      |
| 5    | Methan-d3-ol                        | 35  | CHD3O        | 001849-29-2 | 4       |      |      |



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| TIC Top Hit name | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                  |       |         |       |          | #                     | RT    | Resp    | Conc |
| unknown-01       | 3.005 | 2.3     | ng/ul | 199148   | 1                     | 8.028 | 1764790 | 20.0 |