

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091520\  
 Data File : BG046625.D  
 Acq On : 15 Sep 2020 15:26  
 Operator : CG/JU  
 Sample : L3993-14  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PAV-B41-091120-GW

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\_G\METHODS\8270-BG082520.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| 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         | 3.118       | 3             | 5           | 27           | rVB      | 153621         | 256376        | 5.77%           | 1.331%        |
| 2         | 5.509       | 405           | 412         | 430          | rBV      | 422853         | 709368        | 15.95%          | 3.684%        |
| 3         | 7.096       | 673           | 682         | 694          | rBV      | 274965         | 556382        | 12.51%          | 2.889%        |
| 4         | 7.918       | 816           | 822         | 831          | rBV      | 225379         | 374189        | 8.42%           | 1.943%        |
| 5         | 9.076       | 1012          | 1019        | 1029         | rBV      | 696912         | 1250582       | 28.12%          | 6.494%        |
| 6         | 10.715      | 1291          | 1298        | 1314         | rBV      | 340009         | 638569        | 14.36%          | 3.316%        |
| 7         | 13.177      | 1709          | 1717        | 1727         | rBV      | 2157467        | 3467025       | 77.97%          | 18.005%       |
| 8         | 13.999      | 1852          | 1857        | 1869         | rBV      | 168356         | 269780        | 6.07%           | 1.401%        |
| 9         | 14.551      | 1944          | 1951        | 1961         | rBV2     | 674583         | 1057738       | 23.79%          | 5.493%        |
| 10        | 16.038      | 2197          | 2204        | 2219         | rBV      | 1525641        | 2430933       | 54.67%          | 12.624%       |
| 11        | 17.295      | 2411          | 2418        | 2429         | rBV      | 755622         | 1154276       | 25.96%          | 5.994%        |
| 12        | 18.200      | 2567          | 2572        | 2577         | rBV      | 39398          | 67152         | 1.51%           | 0.349%        |
| 13        | 19.910      | 2857          | 2863        | 2874         | rBV      | 3418520        | 4446541       | 100.00%         | 23.091%       |
| 14        | 21.214      | 3080          | 3085        | 3093         | rBV2     | 90516          | 169316        | 3.81%           | 0.879%        |
| 15        | 21.543      | 3134          | 3141        | 3148         | rBV2     | 693477         | 1169857       | 26.31%          | 6.075%        |
| 16        | 24.634      | 3657          | 3667        | 3684         | rBV2     | 420950         | 1238196       | 27.85%          | 6.430%        |

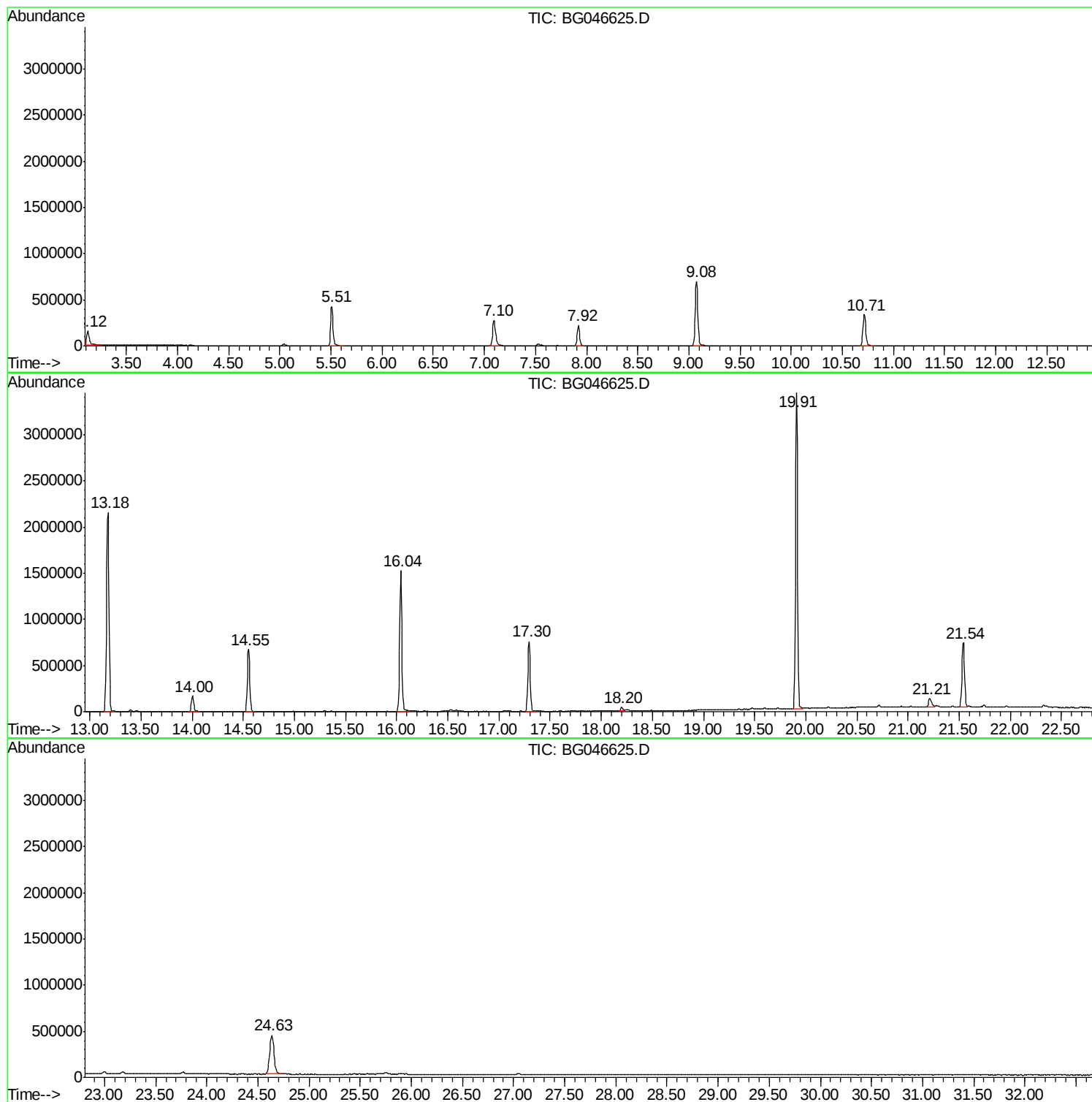
Sum of corrected areas: 19256280

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TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P



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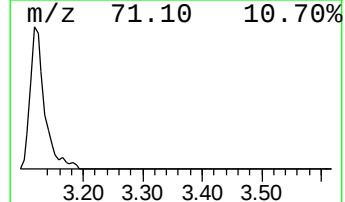
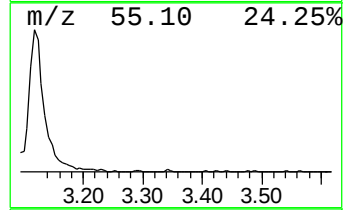
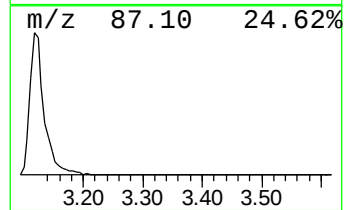
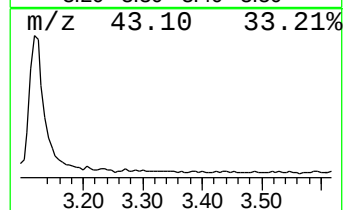
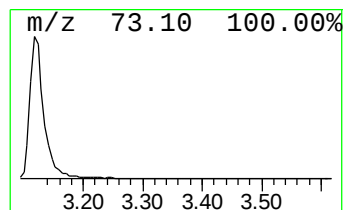
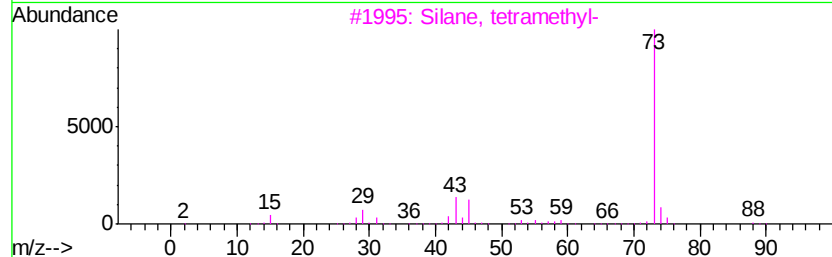
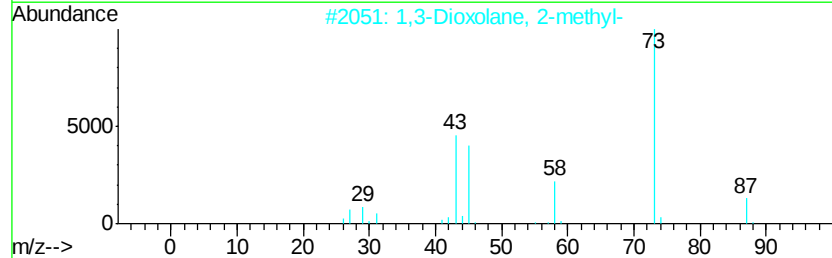
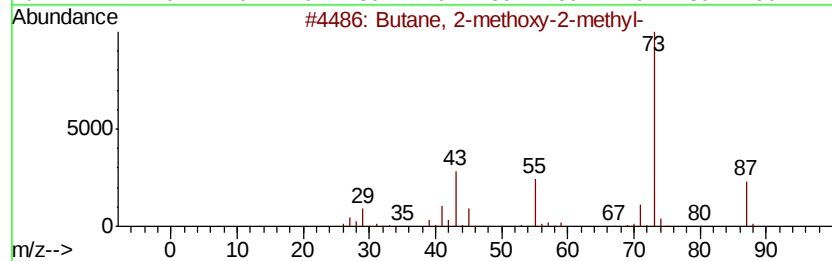
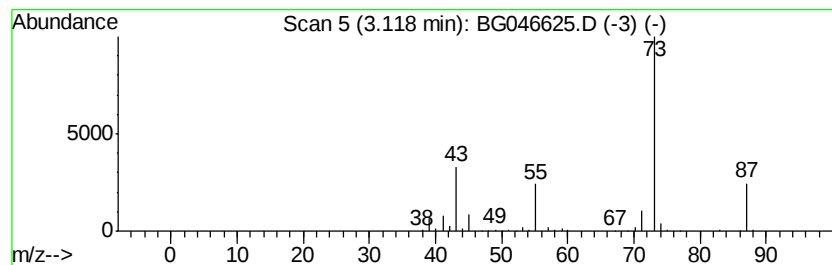
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082520.M  
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TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 3.12 | 13.70 ng | 256376 | 1,4-Dichlorobenzene-d4 | 7.92 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl-         | 102 | C6H14O  | 000994-05-8  | 78   |
| 2    |    | 1,3-Dioxolane, 2-methyl-            | 88  | C4H8O2  | 000497-26-7  | 32   |
| 3    |    | Silane, tetramethyl-                | 88  | C4H12Si | 000075-76-3  | 12   |
| 4    |    | N-Ethylformamide                    | 73  | C3H7NO  | 000627-45-2  | 9    |
| 5    |    | 1-(2-Methoxyethoxy)-2-methyl-2-p... | 162 | C8H18O3 | 1000364-24-4 | 9    |



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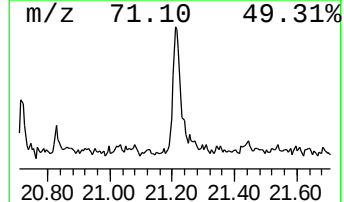
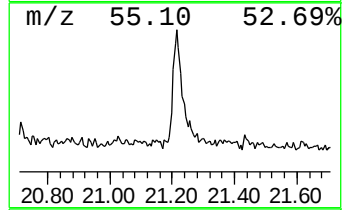
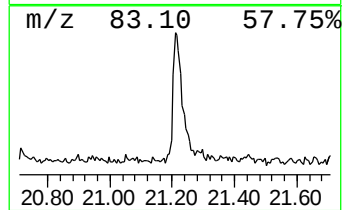
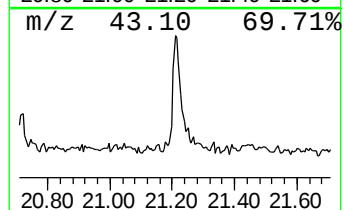
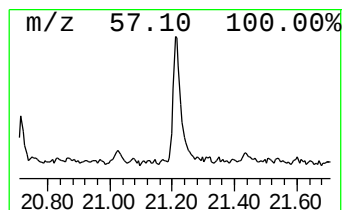
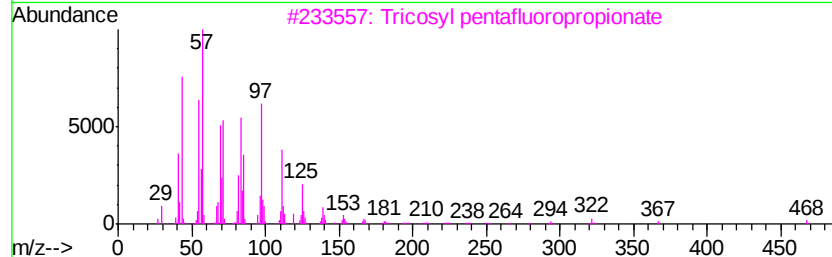
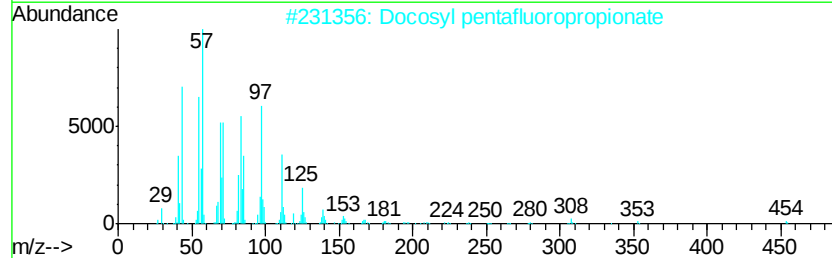
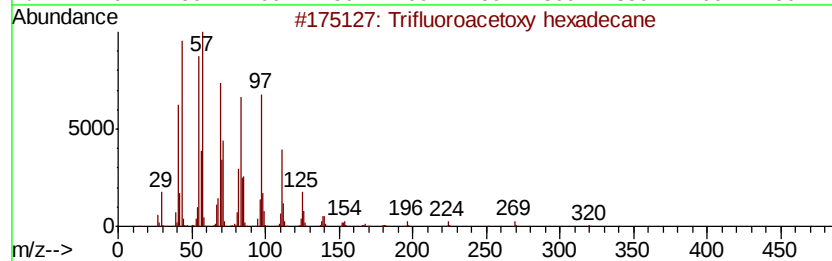
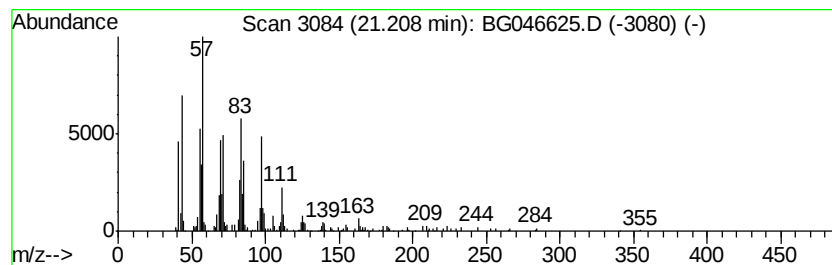
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TIC Integration Parameters: LSCINT.P

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Peak Number 2 Trifluoroacetoxy hexadecane Concentration Rank 2

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 21.21 | 2.89 ng | 169316 | Chrysene-d12     | 21.54 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Trifluoroacetoxy hexadecane         | 338 | C18H33F3O2 | 006222-03-3  | 93   |
| 2    |    | Docosyl pentafluoropropionate       | 472 | C25H45F5O2 | 1000351-80-9 | 91   |
| 3    |    | Tricosyl pentafluoropropionate      | 486 | C26H47F5O2 | 1000351-81-0 | 91   |
| 4    |    | Pentafluoropropionic acid, hepta... | 402 | C20H35F5O2 | 959218-78-1  | 91   |
| 5    |    | Octadecyl trifluoroacetate          | 366 | C20H37F3O2 | 079392-43-1  | 91   |



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| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| Butane, 2-methoxy... | 3.12  | 13.7    | ng    | 256376   | 1                     | 7.92  | 374189  | 20.0 |
| Trifluoroacetoxy ... | 21.21 | 2.9     | ng    | 169316   | 5                     | 21.54 | 1169860 | 20.0 |