

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081919\  
 Data File : BG042339.D  
 Acq On : 19 Aug 2019 23:05  
 Operator : HP/JU  
 Sample : K4331-05  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 0625-COMP-B

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

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-BG081919.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         | 3.123       | 3             | 6           | 21           | rVB      | 482502         | 747344        | 20.76%          | 3.740%        |
| 2         | 5.567       | 415           | 422         | 435          | rBV      | 717067         | 1182759       | 32.86%          | 5.919%        |
| 3         | 7.177       | 687           | 696         | 712          | rBV      | 728859         | 1344619       | 37.36%          | 6.729%        |
| 4         | 7.541       | 750           | 758         | 773          | rBV      | 790963         | 1373223       | 38.15%          | 6.872%        |
| 5         | 8.006       | 829           | 837         | 850          | rBV      | 195011         | 334916        | 9.30%           | 1.676%        |
| 6         | 8.335       | 885           | 893         | 904          | rBV      | 620242         | 1102352       | 30.63%          | 5.517%        |
| 7         | 9.175       | 1027          | 1036        | 1052         | rBV      | 418971         | 802260        | 22.29%          | 4.015%        |
| 8         | 10.832      | 1310          | 1318        | 1337         | rBV      | 239480         | 473375        | 13.15%          | 2.369%        |
| 9         | 13.282      | 1728          | 1735        | 1749         | rBV      | 1382900        | 2278769       | 63.31%          | 11.404%       |
| 10        | 14.110      | 1869          | 1876        | 1893         | rBV      | 170651         | 273896        | 7.61%           | 1.371%        |
| 11        | 14.663      | 1963          | 1970        | 1983         | rBV      | 479040         | 758707        | 21.08%          | 3.797%        |
| 12        | 16.155      | 2218          | 2224        | 2244         | rBV      | 1119868        | 1824062       | 50.68%          | 9.129%        |
| 13        | 17.418      | 2432          | 2439        | 2450         | rBV2     | 619329         | 967791        | 26.89%          | 4.843%        |
| 14        | 17.536      | 2455          | 2459        | 2466         | rVB3     | 30215          | 44736         | 1.24%           | 0.224%        |
| 15        | 18.305      | 2586          | 2590        | 2599         | rBV4     | 23167          | 45043         | 1.25%           | 0.225%        |
| 16        | 20.027      | 2877          | 2883        | 2896         | rBV      | 2677103        | 3599474       | 100.00%         | 18.014%       |
| 17        | 21.343      | 3102          | 3107        | 3126         | rBV3     | 52323          | 135249        | 3.76%           | 0.677%        |
| 18        | 21.690      | 3160          | 3166        | 3180         | rVB2     | 697038         | 1160583       | 32.24%          | 5.808%        |
| 19        | 22.500      | 3299          | 3304        | 3309         | rBV2     | 30027          | 55663         | 1.55%           | 0.279%        |
| 20        | 23.205      | 3419          | 3424        | 3430         | rVB2     | 36577          | 67114         | 1.86%           | 0.336%        |
| 21        | 24.016      | 3558          | 3562        | 3569         | rVB4     | 36768          | 82665         | 2.30%           | 0.414%        |
| 22        | 24.933      | 3708          | 3718        | 3736         | rVB3     | 433157         | 1327119       | 36.87%          | 6.642%        |

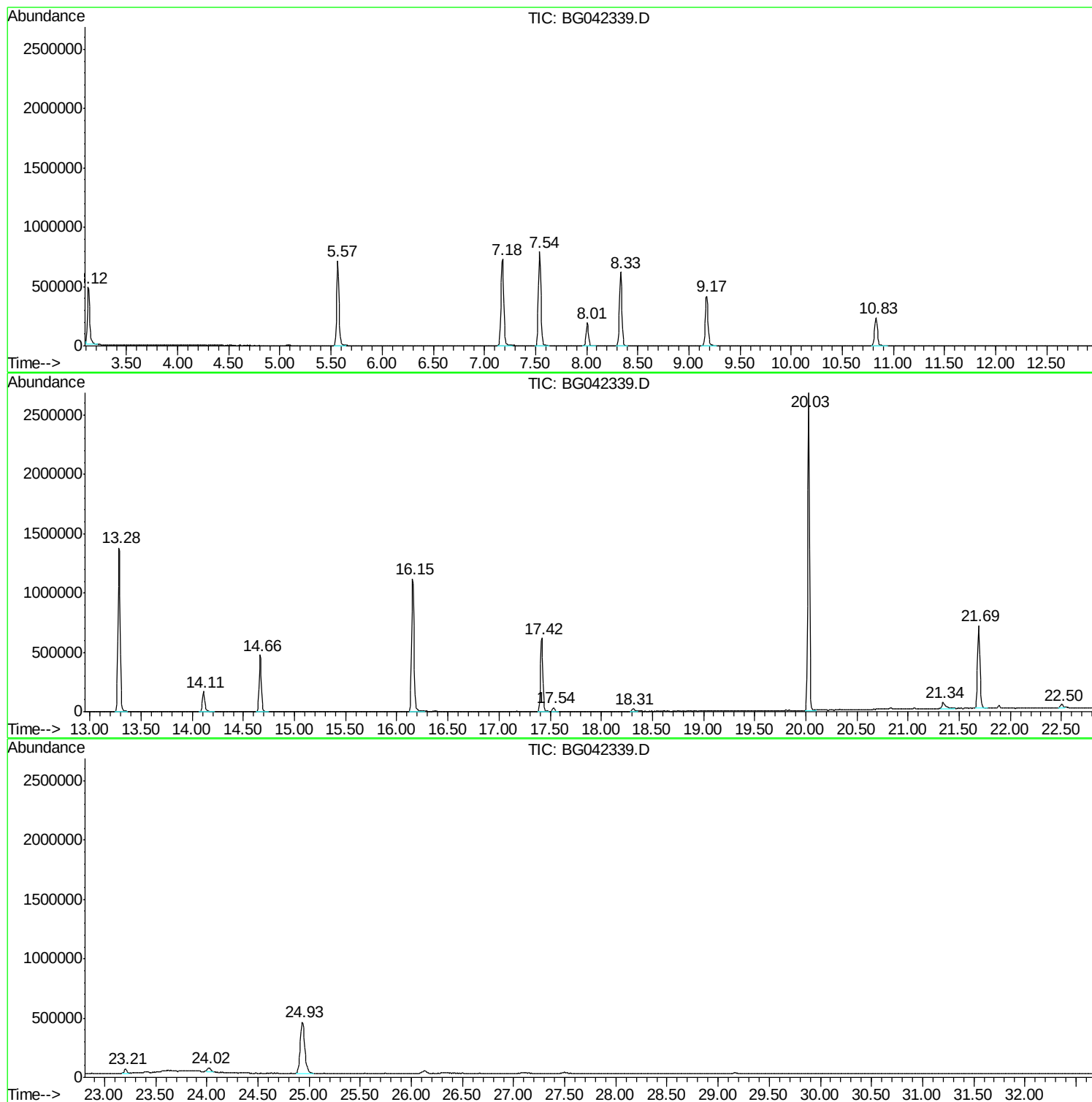
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BNA\_G  
ClientSampleId :  
0625-COMP-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG081919.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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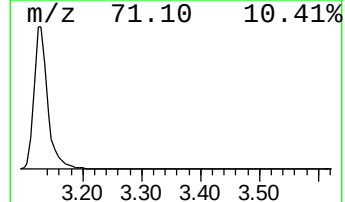
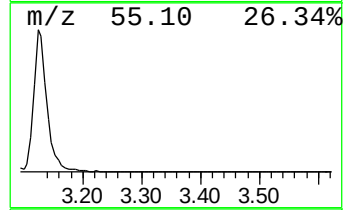
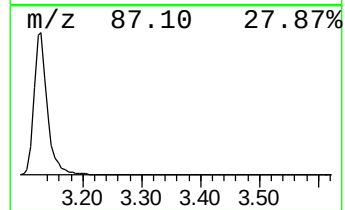
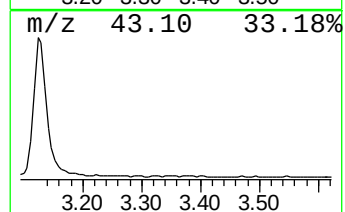
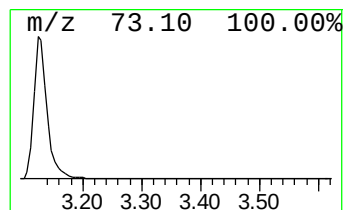
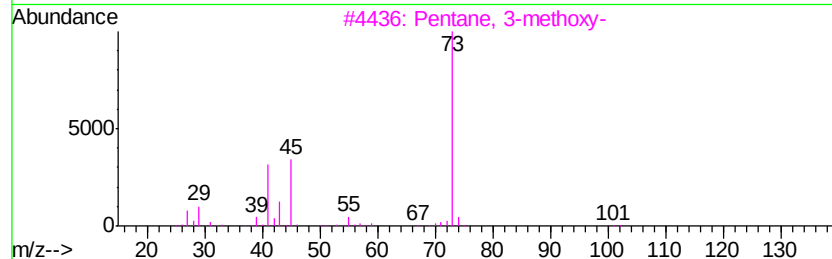
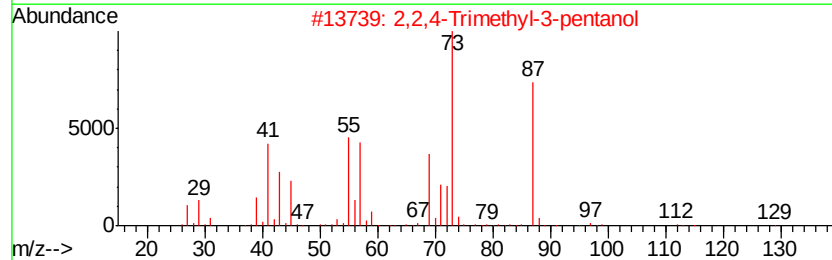
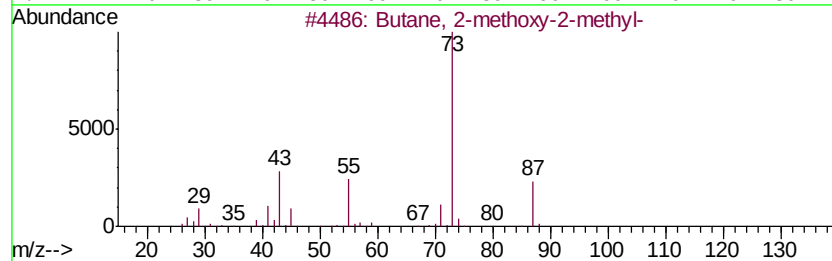
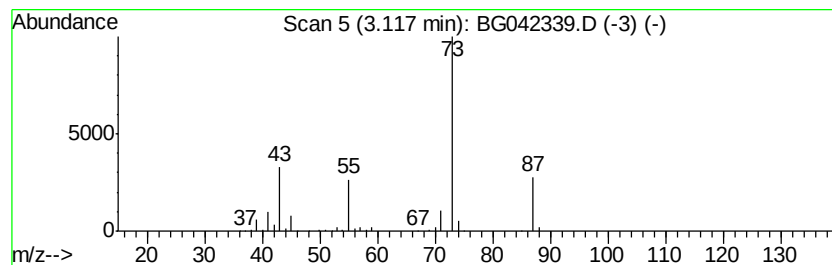
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG081919.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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 3

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 3.12 | 44.63 ng | 747344 | 1,4-Dichlorobenzene-d4 | 8.01 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 86   |
| 2    |    | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 39   |
| 3    |    | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 23   |
| 4    |    | N-Ethylformamide            | 73  | C3H7NO  | 000627-45-2 | 9    |
| 5    |    | Acetamide, N-ethyl-         | 87  | C4H9NO  | 000625-50-3 | 9    |



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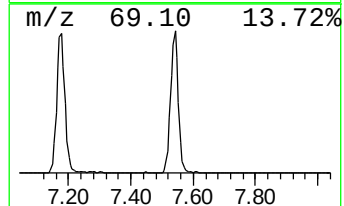
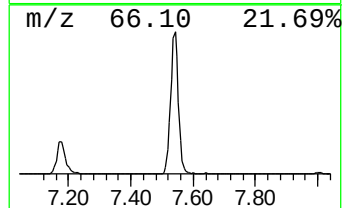
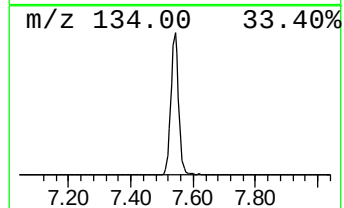
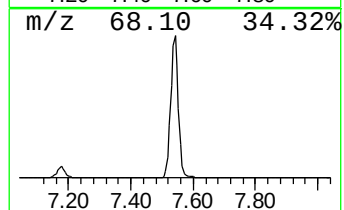
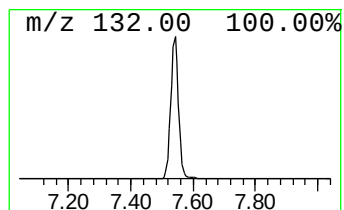
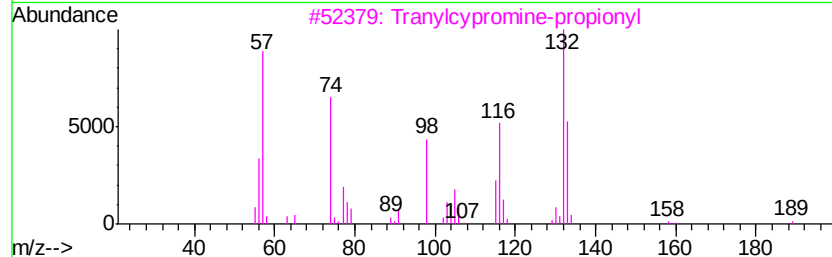
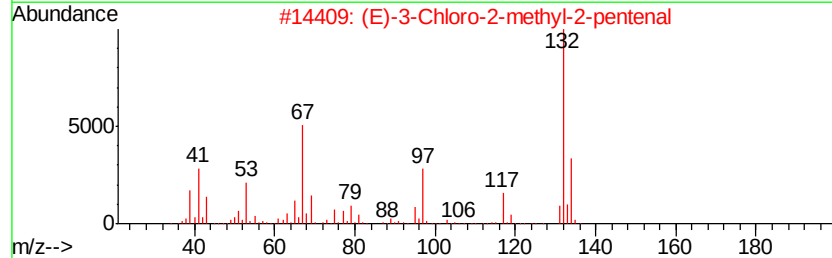
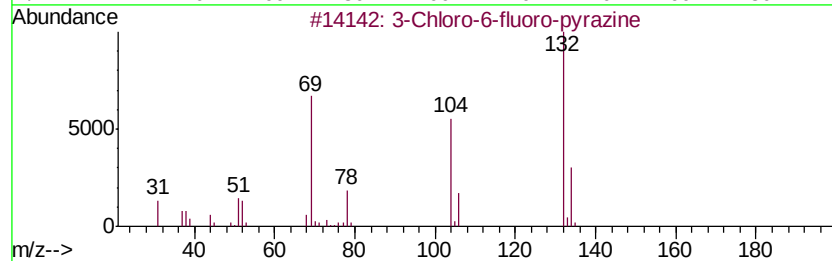
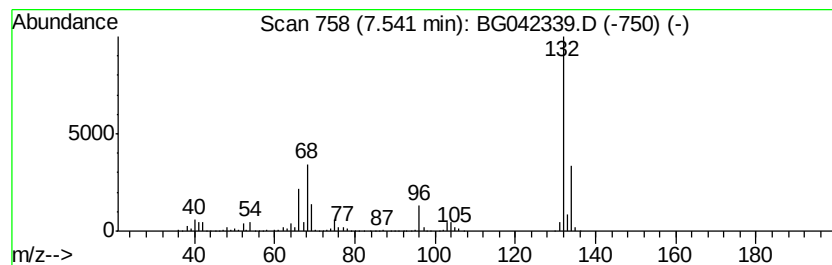
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Peak Number 2 unknown7.54 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 7.54 | 82.00 ng | 1373220 | 1,4-Dichlorobenzene-d4 | 8.01 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 38   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 25   |
| 3    |    | Tranvlcypropine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 16   |
| 4    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 16   |
| 5    |    | 1,3-Cyclopentanedione, 2-chloro- | 132 | C5H5ClO2  | 014203-19-1  | 9    |



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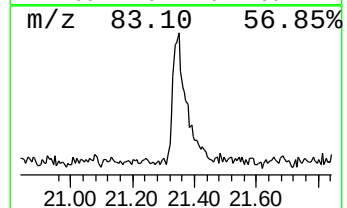
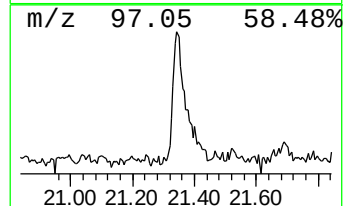
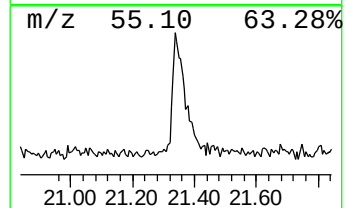
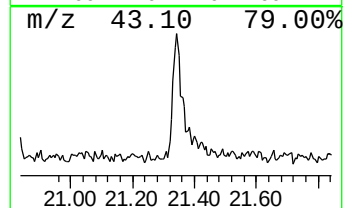
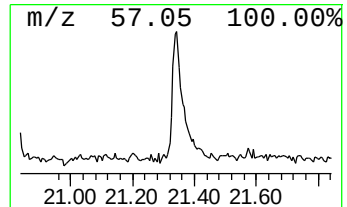
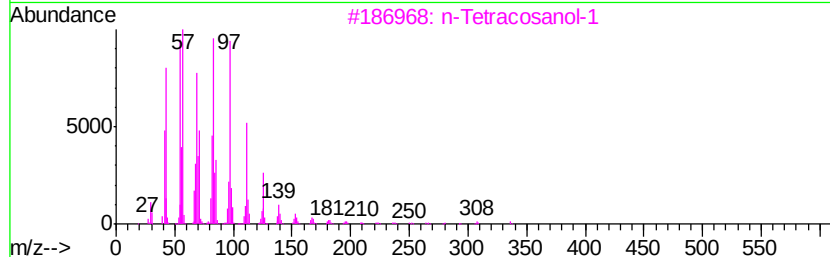
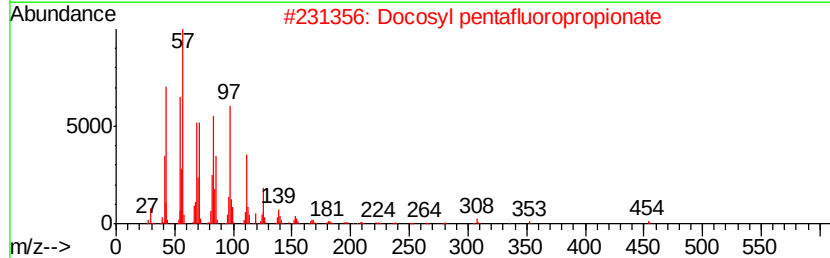
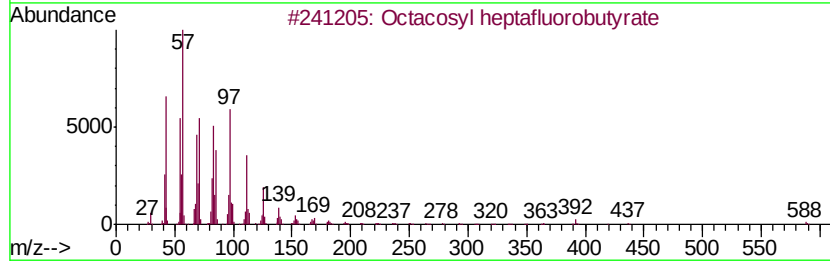
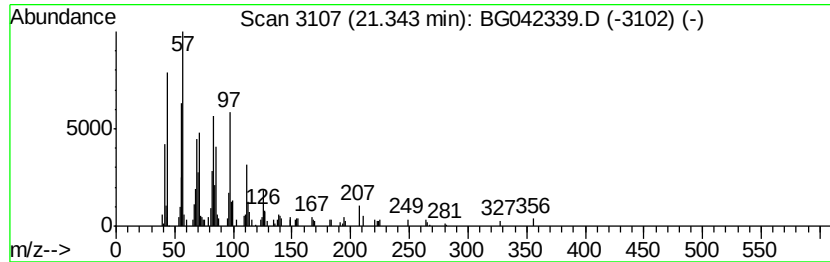
Instrument :  
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ClientSampleId :  
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Peak Number 4 Octacosyl heptafluorobutyrate Concentration Rank 4

| R.T.    | EstConc | Area                                | Relative to ISTD |            | R.T.         |      |
|---------|---------|-------------------------------------|------------------|------------|--------------|------|
| 21.34   | 2.33 ng | 135249                              | Chrysene-d12     |            | 21.69        |      |
| Hit# of | 5       | Tentative ID                        | MW               | MolForm    | CAS#         | Qual |
| 1       |         | Octacosyl heptafluorobutyrate       | 606              | C32H57F7O2 | 1000351-83-6 | 91   |
| 2       |         | Docosyl pentafluoropropionate       | 472              | C25H45F5O2 | 1000351-80-9 | 91   |
| 3       |         | n-Tetracosanol-1                    | 354              | C24H50O    | 000506-51-4  | 91   |
| 4       |         | Heptafluorobutyric acid, n-octad... | 466              | C22H37F7O2 | 000400-57-7  | 90   |
| 5       |         | Heneicosyl heptafluorobutyrate      | 508              | C25H43F7O2 | 1000351-83-8 | 90   |



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BNA\_G  
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| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| Butane, 2-methoxy... | 3.12  | 44.6    | ng    | 747344   | 1                     | 8.01  | 334916  | 20.0 |
| unknown7.54          | 7.54  | 82.0    | ng    | 1373220  | 1                     | 8.01  | 334916  | 20.0 |
| Octacosyl heptafl... | 21.34 | 2.3     | ng    | 135249   | 5                     | 21.69 | 1160580 | 20.0 |