

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101019\  
 Data File : BP000619.D  
 Acq On : 10 Oct 2019 16:28  
 Operator : JU  
 Sample : K5292-06  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 20190855-AMENDED-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\_P\METHODS\8270-BP100119.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         | 4.993       | 482           | 494         | 497          | rBV      | 4853570        | 11981067      | 100.00%         | 27.976%       |
| 2         | 6.387       | 728           | 731         | 741          | rBV3     | 625636         | 657803        | 5.49%           | 1.536%        |
| 3         | 6.522       | 751           | 754         | 761          | rVB      | 214204         | 179544        | 1.50%           | 0.419%        |
| 4         | 6.752       | 789           | 793         | 796          | rBV      | 1885798        | 1444806       | 12.06%          | 3.374%        |
| 5         | 6.905       | 816           | 819         | 823          | rBV      | 3265491        | 2943381       | 24.57%          | 6.873%        |
| 6         | 7.311       | 884           | 888         | 891          | rBV      | 2776523        | 2155455       | 17.99%          | 5.033%        |
| 7         | 8.034       | 1007          | 1011        | 1014         | rBV      | 2451565        | 1892470       | 15.80%          | 4.419%        |
| 8         | 9.116       | 1191          | 1195        | 1198         | rBV      | 5245264        | 4126696       | 34.44%          | 9.636%        |
| 9         | 9.510       | 1259          | 1262        | 1265         | rBV      | 997802         | 772807        | 6.45%           | 1.805%        |
| 10        | 9.799       | 1307          | 1311        | 1314         | rVB      | 2861470        | 2355341       | 19.66%          | 5.500%        |
| 11        | 10.734      | 1465          | 1470        | 1475         | rVB4     | 89220          | 141034        | 1.18%           | 0.329%        |
| 12        | 11.157      | 1538          | 1542        | 1547         | rBV2     | 310214         | 353369        | 2.95%           | 0.825%        |
| 13        | 11.293      | 1561          | 1565        | 1572         | rBV      | 3035155        | 2525229       | 21.08%          | 5.897%        |
| 14        | 12.899      | 1834          | 1838        | 1841         | rBV      | 5504221        | 4645224       | 38.77%          | 10.847%       |
| 15        | 13.822      | 1992          | 1995        | 2009         | rVB2     | 205058         | 330459        | 2.76%           | 0.772%        |
| 16        | 13.963      | 2015          | 2019        | 2023         | rBV      | 2439407        | 2542001       | 21.22%          | 5.936%        |
| 17        | 14.457      | 2100          | 2103        | 2106         | rBV      | 192966         | 171066        | 1.43%           | 0.399%        |
| 18        | 14.763      | 2152          | 2155        | 2162         | rVB      | 218049         | 247692        | 2.07%           | 0.578%        |
| 19        | 15.016      | 2195          | 2198        | 2207         | rVB      | 70525          | 136171        | 1.14%           | 0.318%        |
| 20        | 15.104      | 2210          | 2213        | 2221         | rVB2     | 244792         | 279263        | 2.33%           | 0.652%        |
| 21        | 15.440      | 2265          | 2270        | 2275         | rBV      | 1923695        | 2395787       | 20.00%          | 5.594%        |
| 22        | 15.487      | 2275          | 2278        | 2291         | rVB      | 218501         | 333603        | 2.78%           | 0.779%        |
| 23        | 15.928      | 2349          | 2353        | 2358         | rBV      | 142773         | 215530        | 1.80%           | 0.503%        |

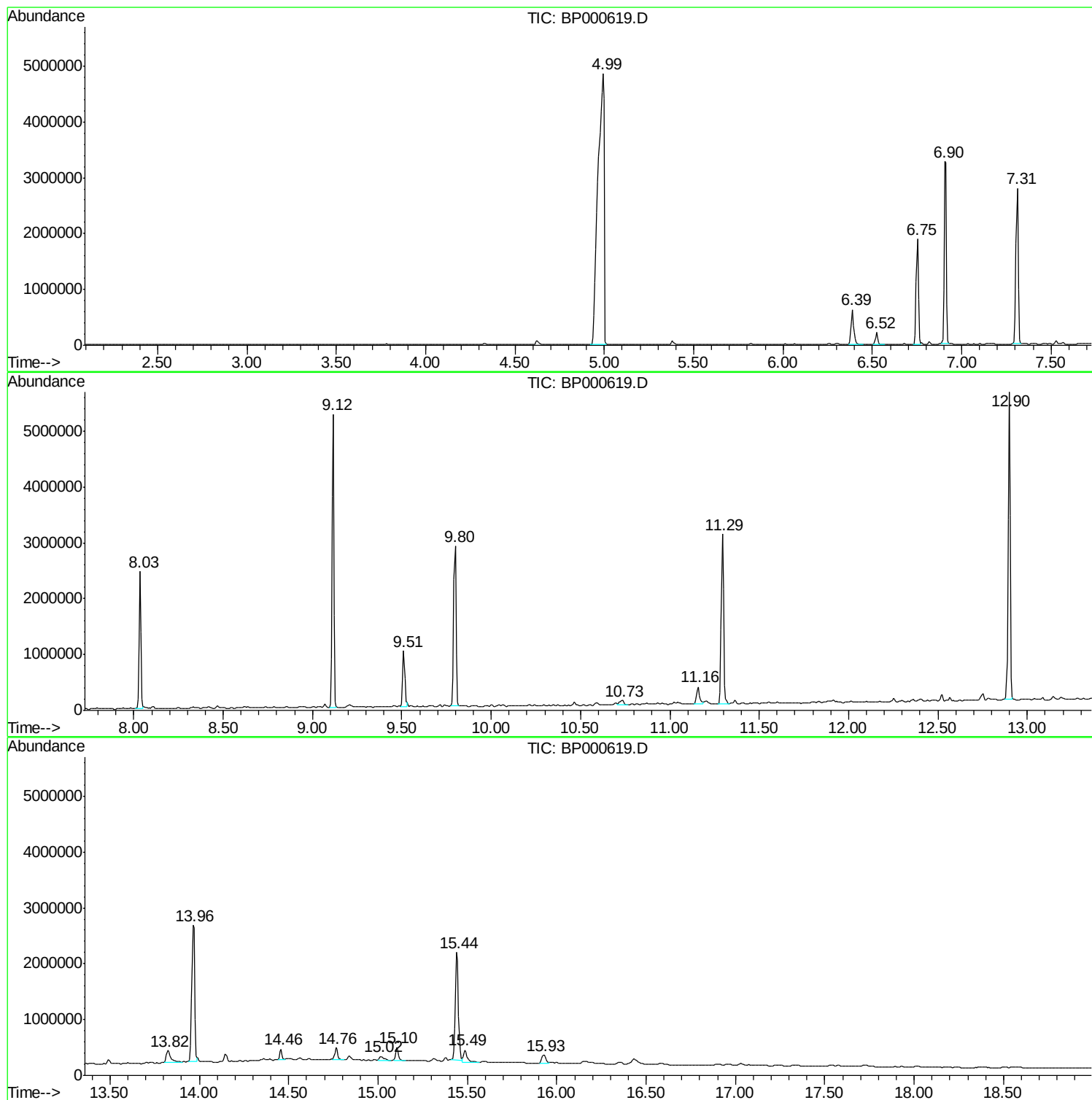
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ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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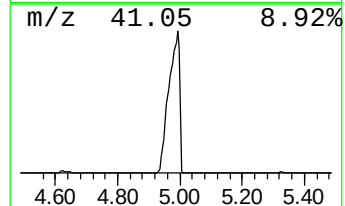
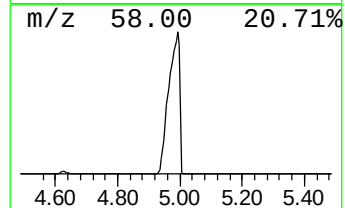
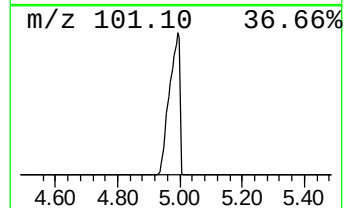
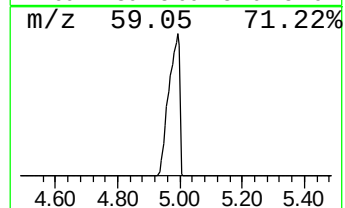
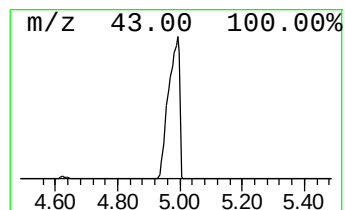
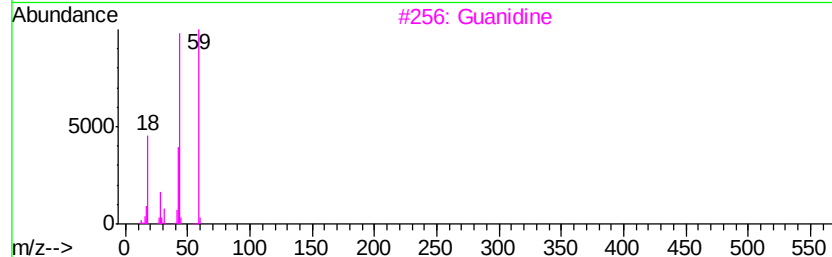
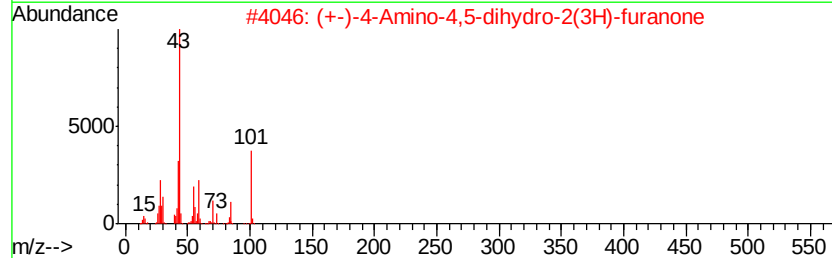
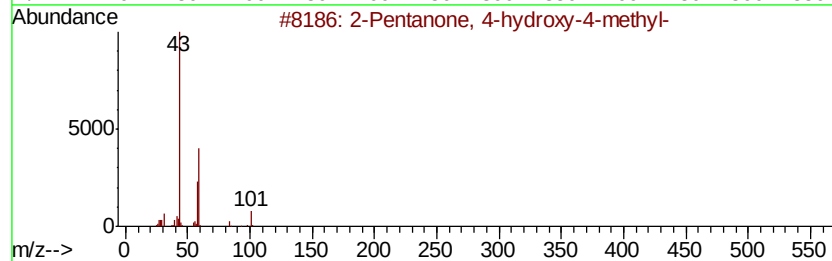
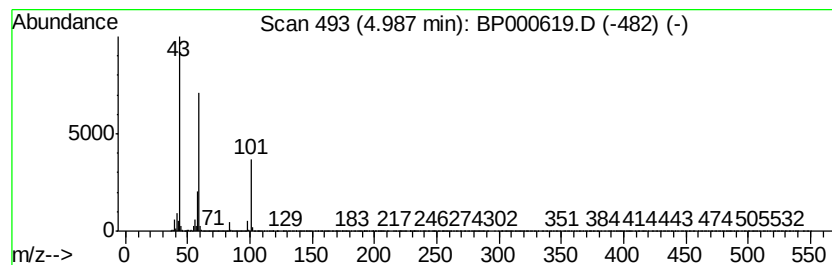
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ClientSampleId :  
20190855-AMENDED-COMP-B

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

| R.T.      | EstConc                             | Area     | Relative to ISTD       | R.T.        |      |
|-----------|-------------------------------------|----------|------------------------|-------------|------|
| 4.99      | 165.85 ng                           | 11981100 | 1,4-Dichlorobenzene-d4 | 6.75        |      |
| Hit# of 5 | Tentative ID                        | MW       | MolForm                | CAS#        | Qual |
| 1         | 2-Pentanone, 4-hydroxy-4-methyl-    | 116      | C6H12O2                | 000123-42-2 | 45   |
| 2         | (+)-4-Amino-4,5-dihydro-2(3H)-f...  | 101      | C4H7NO2                | 016504-58-8 | 43   |
| 3         | Guanidine                           | 59       | CH5N3                  | 000113-00-8 | 38   |
| 4         | Acetone                             | 58       | C3H6O                  | 000067-64-1 | 12   |
| 5         | Acetic acid, cyano-, 1,1-dimethy... | 141      | C7H11NO2               | 001116-98-9 | 12   |



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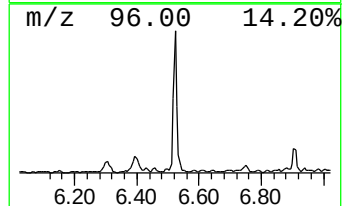
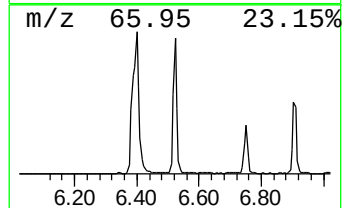
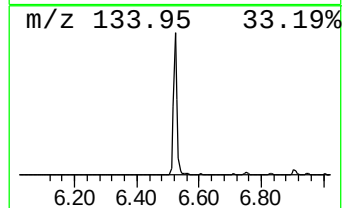
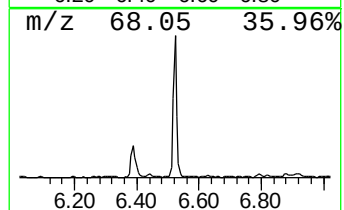
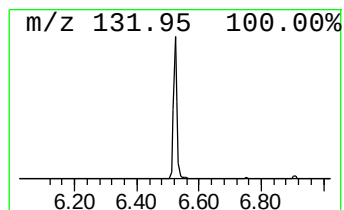
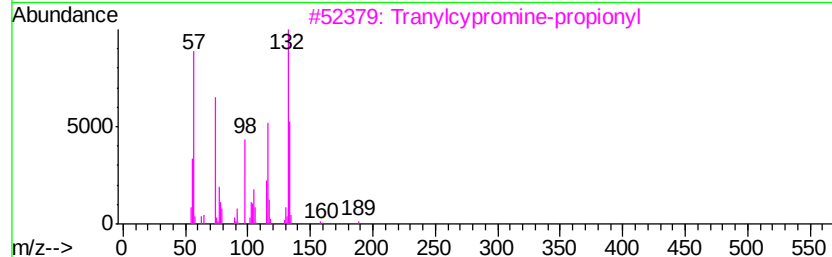
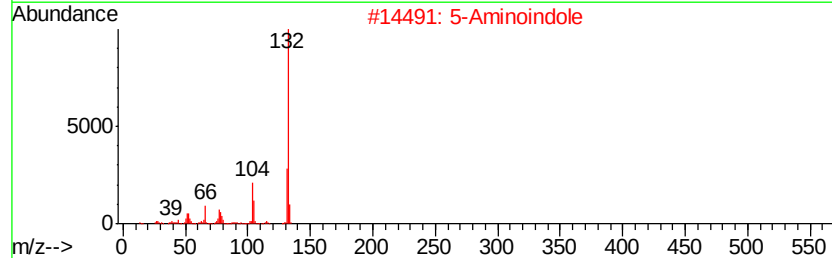
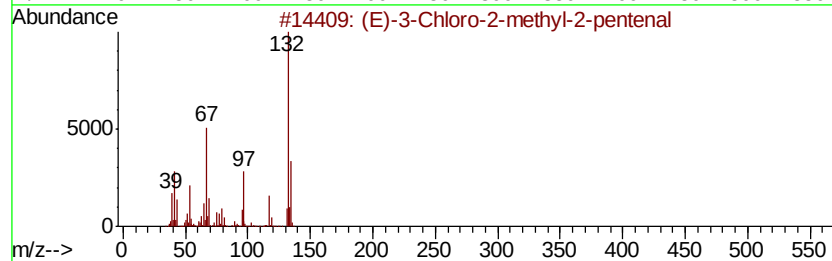
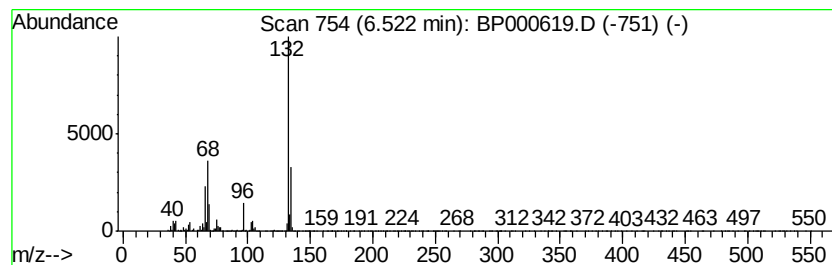
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 unknown6.52 Concentration Rank 5

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.52 | 2.49 ng | 179544 | 1,4-Dichlorobenzene-d4 | 6.75 |

| Hit# | of | Tentative ID                     | MW  | MolForm  | CAS#         | Qual |
|------|----|----------------------------------|-----|----------|--------------|------|
| 1    | 5  | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO  | 031357-76-3  | 35   |
| 2    |    | 5-Aminoindole                    | 132 | C8H8N2   | 005192-03-0  | 16   |
| 3    |    | Tranvlcypropine-propionyl        | 189 | C12H15NO | 1000123-86-3 | 16   |
| 4    |    | Amphetaminil                     | 250 | C17H18N2 | 017590-01-1  | 12   |
| 5    |    | Imidazole, 4,5-dicyano-1-methyl- | 132 | C6H4N4   | 019485-35-9  | 9    |



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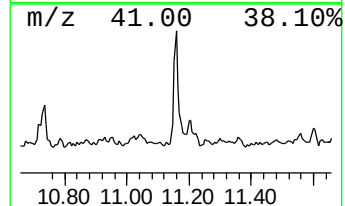
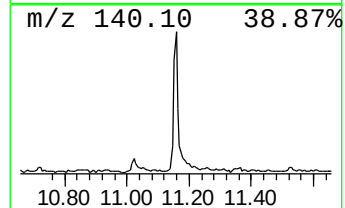
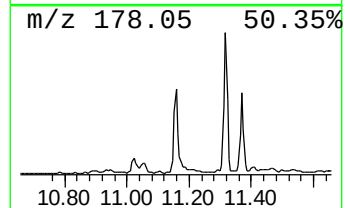
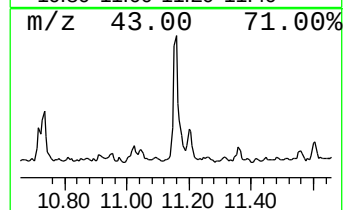
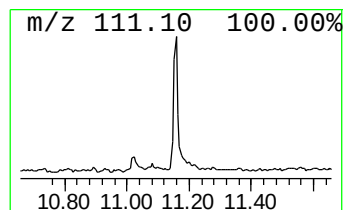
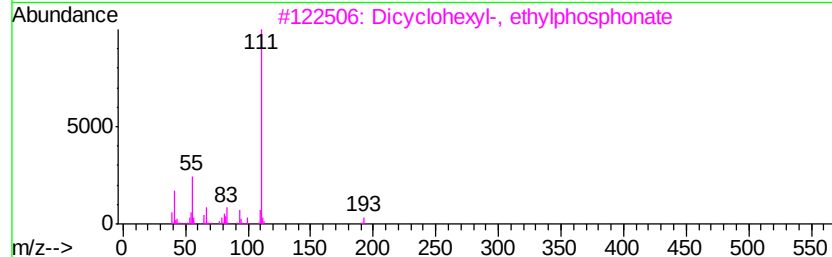
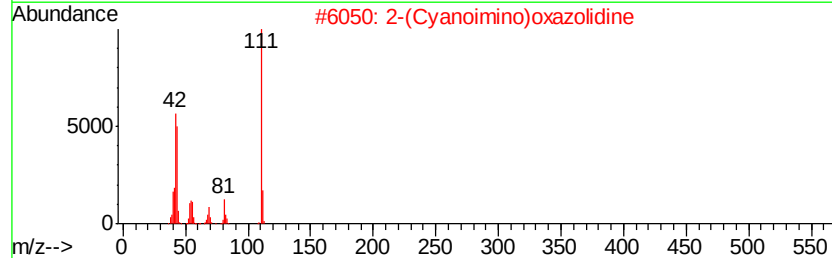
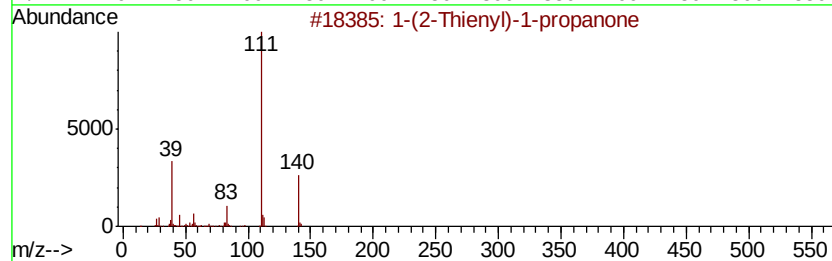
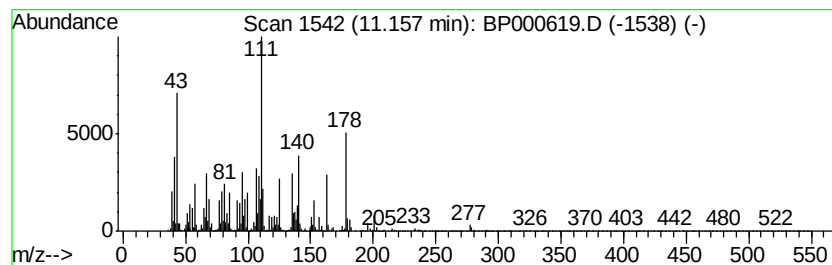
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 unknown11.16 Concentration Rank 3

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.16 | 2.80 ng | 353369 | Phenanthrene-d10 | 11.29 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | 1-(2-Thienyl)-1-propanone           | 140 | C7H8OS    | 013679-75-9 | 27   |
| 2    |    | 2-(Cyanoimino)oxazolidine           | 111 | C4H5N3O   | 050411-06-8 | 18   |
| 3    |    | Dicyclohexyl-, ethylphosphonate     | 274 | C14H27O3P | 169739-47-3 | 14   |
| 4    |    | 2-(4-Nitrobutyryl)-cyclopentanone   | 199 | C9H13NO4  | 079630-91-4 | 14   |
| 5    |    | 2-Thiophenecarboxylic acid, phen... | 204 | C11H8O2S  | 000881-89-0 | 14   |



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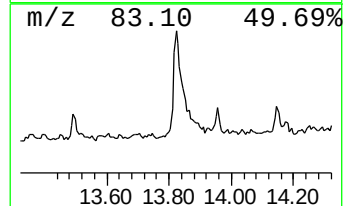
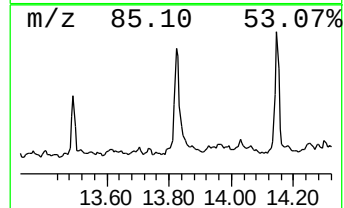
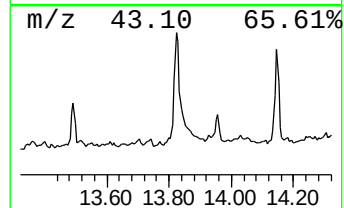
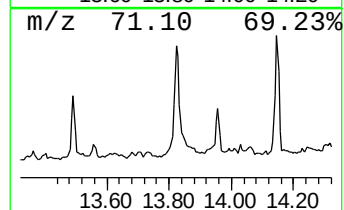
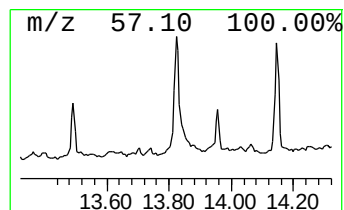
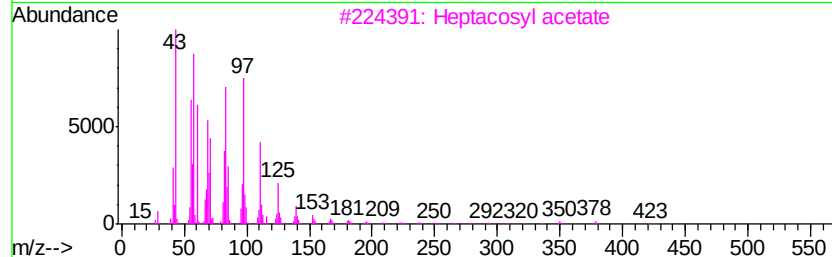
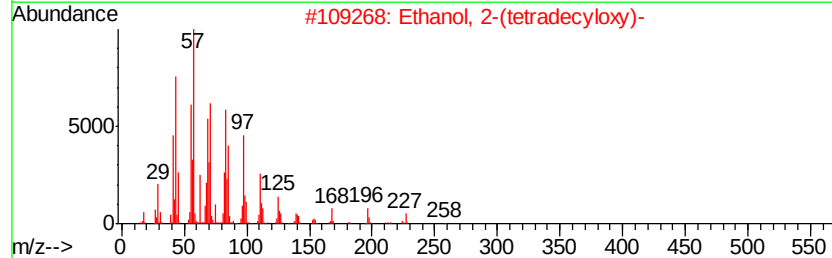
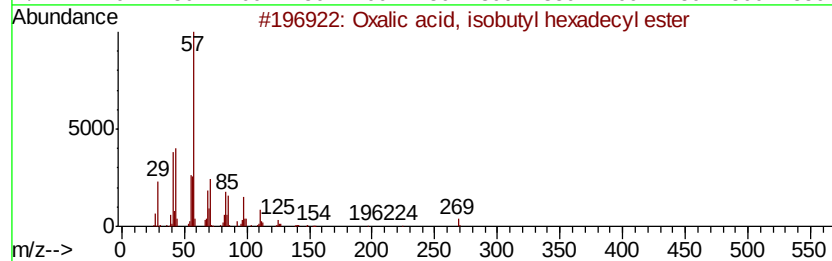
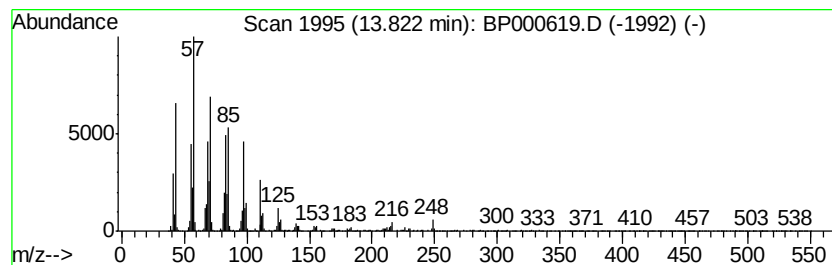
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Peak Number 5 Oxalic acid, isobutyl hexad... Concentration Rank 4

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.82 | 2.60 ng | 330459 | Chrysene-d12     | 13.97 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Oxalic acid, isobutyl hexadecyl ... | 370 | C22H42O4 | 1000309-38-1 | 91   |
| 2    |    | Ethanol, 2-(tetradecyloxy)-         | 258 | C16H34O2 | 002136-70-1  | 91   |
| 3    |    | Heptacosyl acetate                  | 438 | C29H58O2 | 1000351-78-2 | 90   |
| 4    |    | Octacosyl acetate                   | 452 | C30H60O2 | 018206-97-8  | 87   |
| 5    |    | 1-Decanol, 2-hexyl-                 | 242 | C16H34O  | 002425-77-6  | 87   |



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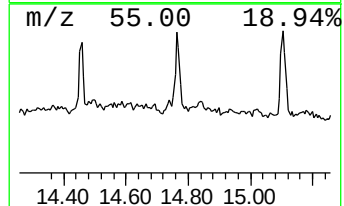
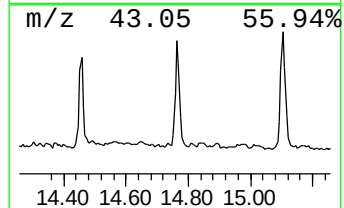
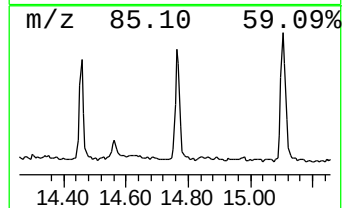
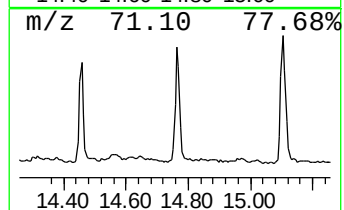
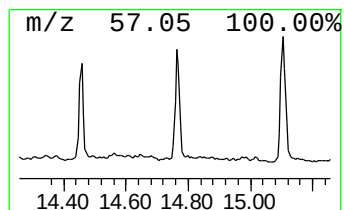
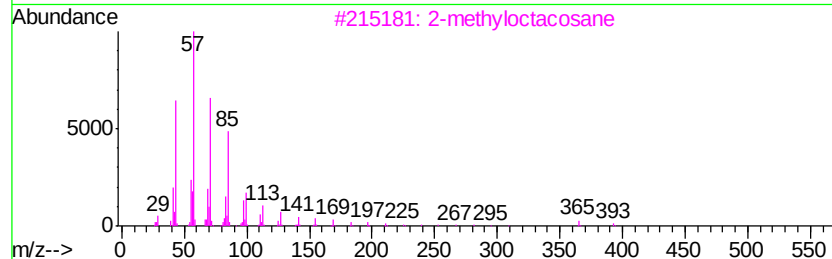
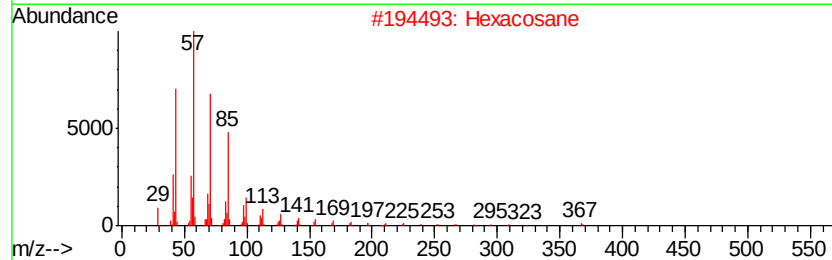
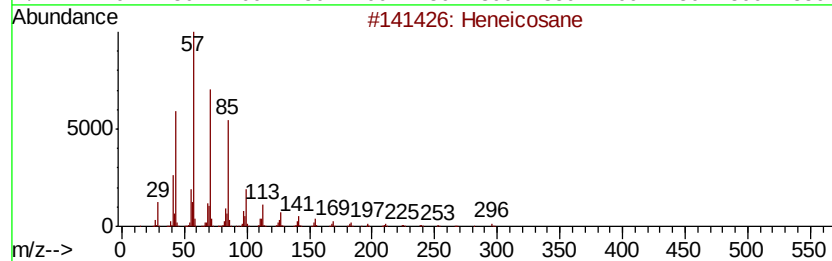
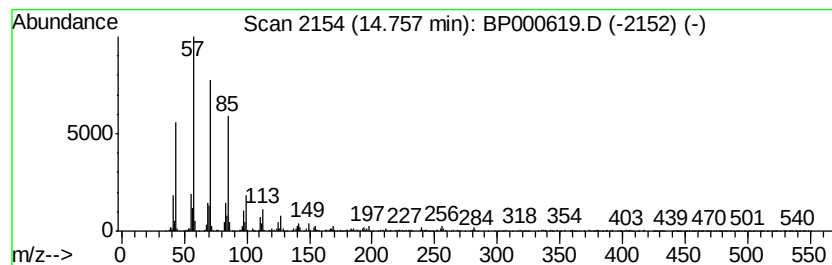
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 Heneicosane Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.76 | 2.07 ng | 247692 | Perylene-d12     | 15.44 |

| Hit# of | 5                     | Tentative ID | MW  | MolForm | CAS#         | Qual |
|---------|-----------------------|--------------|-----|---------|--------------|------|
| 1       | Heneicosane           |              | 296 | C21H44  | 000629-94-7  | 91   |
| 2       | Hexacosane            |              | 366 | C26H54  | 000630-01-3  | 91   |
| 3       | 2-methyloctacosane    |              | 408 | C29H60  | 1000376-72-8 | 91   |
| 4       | Heptadecane, 9-octyl- |              | 352 | C25H52  | 007225-64-1  | 91   |
| 5       | Octacosane            |              | 394 | C28H58  | 000630-02-4  | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101019\  
Data File : BP000619.D  
Acq On : 10 Oct 2019 16:28  
Operator : JU  
Sample : K5292-06  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
20190855-AMENDED-COMP-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

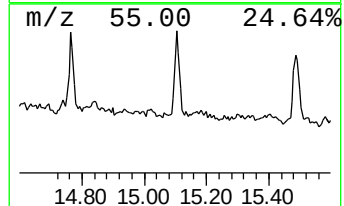
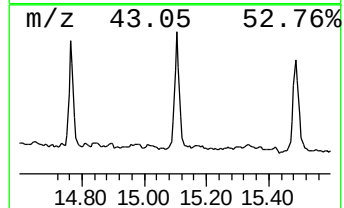
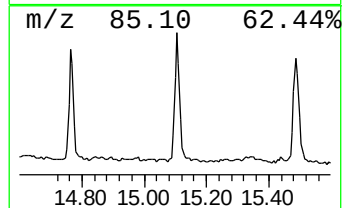
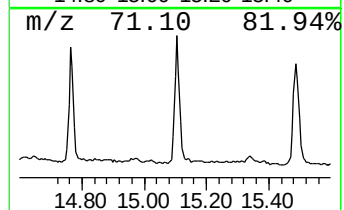
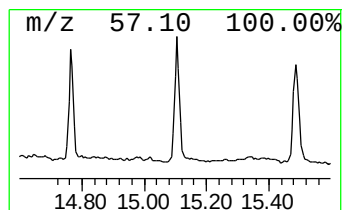
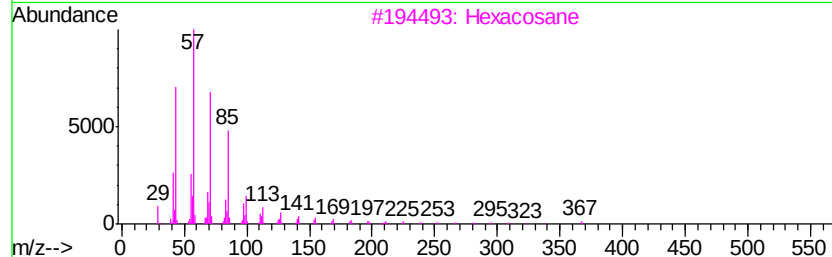
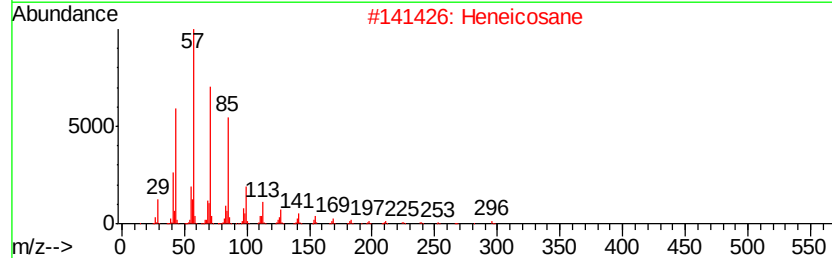
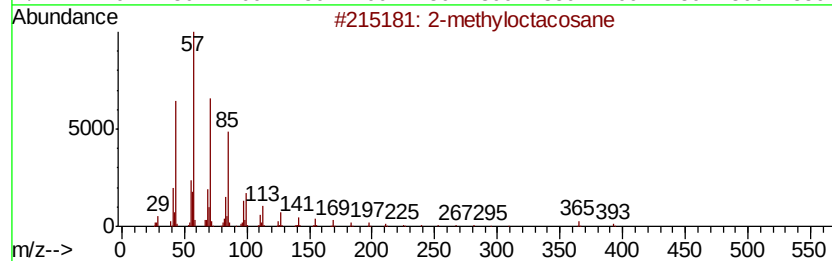
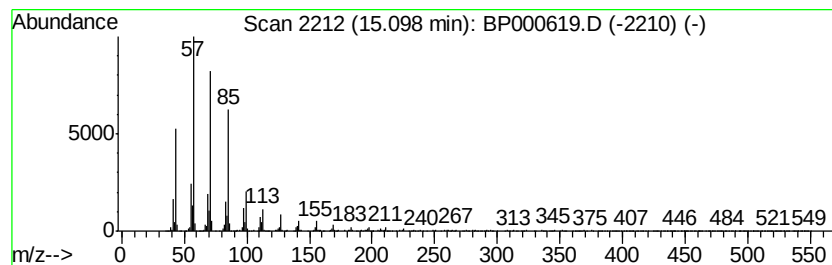
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 2-methyloctacosane Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.10 | 2.33 ng | 279263 | Perylene-d12     | 15.44 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------|-----|---------|--------------|------|
| 1    | 5  | 2-methyloctacosane      | 408 | C29H60  | 1000376-72-8 | 91   |
| 2    |    | Heneicosane             | 296 | C21H44  | 000629-94-7  | 91   |
| 3    |    | Hexacosane              | 366 | C26H54  | 000630-01-3  | 90   |
| 4    |    | Heneicosane, 11-pentyl- | 366 | C26H54  | 014739-72-1  | 90   |
| 5    |    | Pentadecane, 8-hexyl-   | 296 | C21H44  | 013475-75-7  | 87   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA\_P\DATA\BP101019\  
Data File : BP000619.D  
Acq On : 10 Oct 2019 16:28  
Operator : JU  
Sample : K5292-06  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
20190855-AMENDED-COMP-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\8270-BP100119.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| 2-Pentanone, 4-hy... | 4.99  | 165.8   | ng    | 11981100 | 1                     | 6.75  | 1444810 | 20.0 |
| unknown6.52          | 6.52  | 2.5     | ng    | 179544   | 1                     | 6.75  | 1444810 | 20.0 |
| unknown11.16         | 11.16 | 2.8     | ng    | 353369   | 4                     | 11.29 | 2525230 | 20.0 |
| Oxalic acid, isob... | 13.82 | 2.6     | ng    | 330459   | 5                     | 13.97 | 2542000 | 20.0 |
| Heneicosane          | 14.76 | 2.1     | ng    | 247692   | 6                     | 15.44 | 2395790 | 20.0 |
| 2-methyloctacosane   | 15.10 | 2.3     | ng    | 279263   | 6                     | 15.44 | 2395790 | 20.0 |