

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM012117\  
 Data File : BM008771.D  
 Acq On : 21 Jan 2017 14:28  
 Operator : UM/SJ  
 Sample : I1323-01  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DA9P2

## Integration Parameters: LSCINT.P

Integrator: RTE  
 Smoothing : OFF  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0

Filtering: 5  
 Min Area: 1 % 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:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010617.M  
 Title : SVOA 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         | 2.928       | 52            | 58          | 85           | rVB2     | 74550158       | 123926390     | 100.00%         | 59.040%       |
| 2         | 10.369      | 1315          | 1323        | 1326         | rBV2     | 720919         | 1314393       | 1.06%           | 0.626%        |
| 3         | 12.922      | 1742          | 1757        | 1771         | rBV      | 785046         | 3469862       | 2.80%           | 1.653%        |
| 4         | 13.992      | 1933          | 1939        | 1945         | rBV      | 1324562        | 1787811       | 1.44%           | 0.852%        |
| 5         | 14.280      | 1981          | 1988        | 1995         | rVB      | 1313493        | 1917480       | 1.55%           | 0.914%        |
| 6         | 14.586      | 2033          | 2040        | 2045         | rBV      | 931075         | 1328126       | 1.07%           | 0.633%        |
| 7         | 15.398      | 2172          | 2178        | 2184         | rVB      | 1110173        | 1396890       | 1.13%           | 0.665%        |
| 8         | 15.574      | 2201          | 2208        | 2213         | rBV      | 1633831        | 2382182       | 1.92%           | 1.135%        |
| 9         | 15.633      | 2213          | 2218        | 2222         | rVV      | 912840         | 1302285       | 1.05%           | 0.620%        |
| 10        | 16.627      | 2375          | 2387        | 2391         | rBV      | 1148661        | 2005551       | 1.62%           | 0.955%        |
| 11        | 16.863      | 2413          | 2427        | 2431         | rBV      | 336960         | 1276382       | 1.03%           | 0.608%        |
| 12        | 16.968      | 2441          | 2445        | 2453         | rVB2     | 1618366        | 2316169       | 1.87%           | 1.103%        |
| 13        | 17.074      | 2453          | 2463        | 2468         | rBV      | 2205350        | 3609309       | 2.91%           | 1.720%        |
| 14        | 17.327      | 2499          | 2506        | 2514         | rBV      | 1366361        | 1907059       | 1.54%           | 0.909%        |
| 15        | 17.427      | 2517          | 2523        | 2526         | rVV      | 2153415        | 3105512       | 2.51%           | 1.479%        |
| 16        | 17.457      | 2526          | 2528        | 2539         | rVB      | 1035861        | 1421806       | 1.15%           | 0.677%        |
| 17        | 19.709      | 2904          | 2911        | 2914         | rBV      | 2759086        | 4036891       | 3.26%           | 1.923%        |
| 18        | 19.739      | 2914          | 2916        | 2922         | rVB      | 1704028        | 1917585       | 1.55%           | 0.914%        |
| 19        | 20.627      | 3062          | 3067        | 3072         | rBV      | 3430961        | 3653369       | 2.95%           | 1.740%        |
| 20        | 20.992      | 3121          | 3129        | 3134         | rBV      | 1911281        | 3732154       | 3.01%           | 1.778%        |
| 21        | 21.056      | 3134          | 3140        | 3149         | rVB      | 1050293        | 2169748       | 1.75%           | 1.034%        |
| 22        | 21.392      | 3193          | 3197        | 3202         | rBV      | 3305313        | 3418759       | 2.76%           | 1.629%        |
| 23        | 21.486      | 3206          | 3213        | 3217         | rVV2     | 2481665        | 5056429       | 4.08%           | 2.409%        |
| 24        | 21.527      | 3217          | 3220        | 3225         | rVB      | 2045268        | 2481969       | 2.00%           | 1.182%        |
| 25        | 21.668      | 3237          | 3244        | 3249         | rBV2     | 1615473        | 3446012       | 2.78%           | 1.642%        |
| 26        | 21.727      | 3249          | 3254        | 3262         | rVB      | 1022065        | 2052716       | 1.66%           | 0.978%        |
| 27        | 21.845      | 3267          | 3274        | 3282         | rBV      | 900570         | 1818255       | 1.47%           | 0.866%        |
| 28        | 23.139      | 3486          | 3494        | 3498         | rBV      | 2100798        | 3877387       | 3.13%           | 1.847%        |
| 29        | 23.186      | 3498          | 3502        | 3509         | rVB      | 1281660        | 2098764       | 1.69%           | 1.000%        |
| 30        | 23.703      | 3584          | 3590        | 3595         | rBV      | 1830941        | 3680545       | 2.97%           | 1.753%        |
| 31        | 23.750      | 3595          | 3598        | 3606         | rVB      | 1075440        | 1816271       | 1.47%           | 0.865%        |
| 32        | 23.856      | 3609          | 3616        | 3623         | rVB2     | 1118844        | 2225415       | 1.80%           | 1.060%        |
| 33        | 24.509      | 3716          | 3727        | 3729         | rBV      | 903153         | 2811844       | 2.27%           | 1.340%        |
| 34        | 26.280      | 4018          | 4028        | 4042         | rVB3     | 1612018        | 5142349       | 4.15%           | 2.450%        |

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Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
DA9P2

Integration Parameters: LSCINT.P  
Integrator: RTE  
Smoothing : OFF Filtering: 5  
Sampling : 1 Min Area: 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:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010617.M  
Title : SVOA CALIBRATION

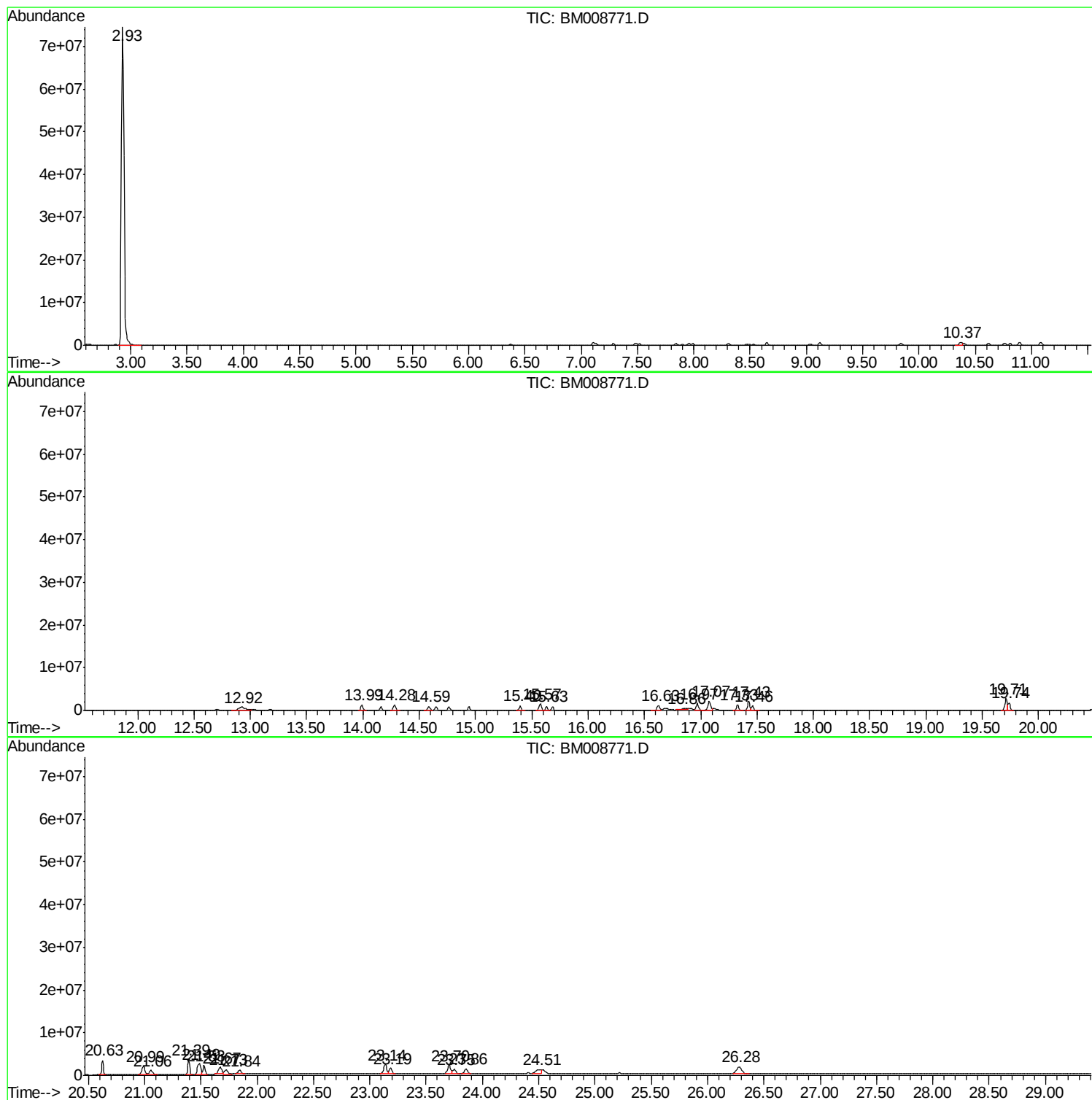
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Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010617.M  
Quant Title : SVOA CALIBRATION

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



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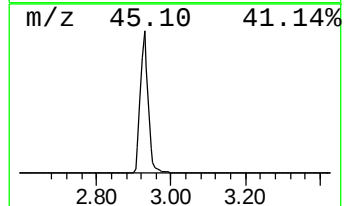
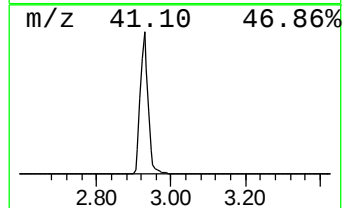
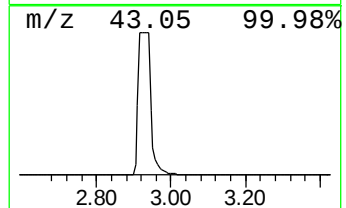
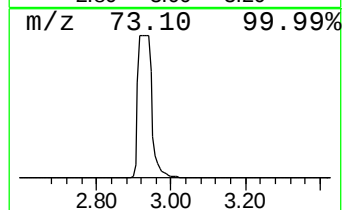
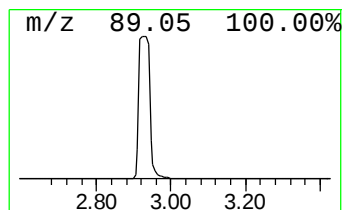
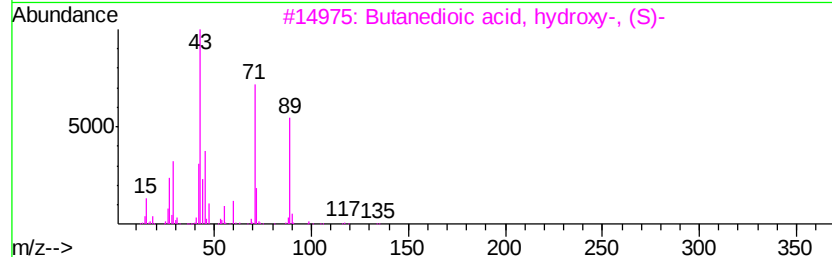
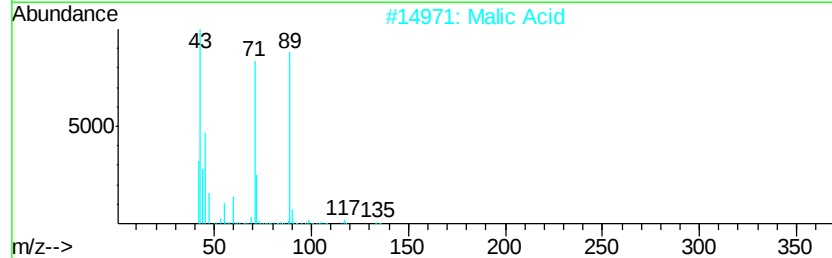
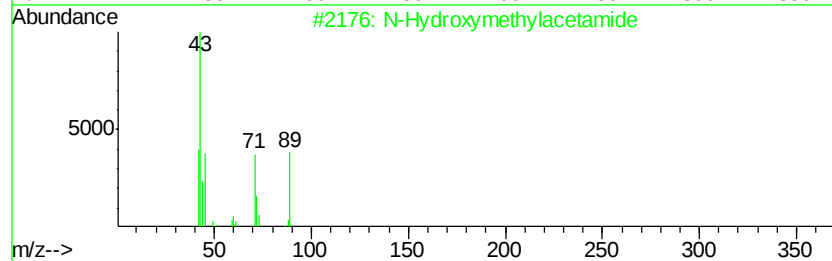
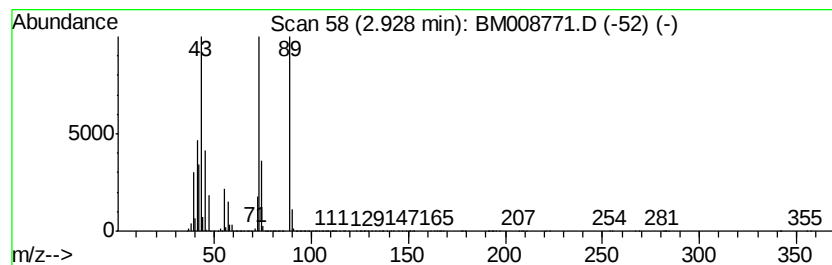
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Quant Title : SVOA CALIBRATION

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

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Peak Number 1 N-Hydroxymethylacetamide Concentration Rank 1

| R.T. | EstConc       | Area      | Relative to ISTD       | R.T. |
|------|---------------|-----------|------------------------|------|
| 2.93 | 1885.68 ng/ul | 123926000 | 1,4-Dichlorobenzene-d4 | 7.96 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | N-Hydroxymethylacetamide         | 89  | C3H7NO2 | 000625-51-4 | 64   |
| 2    |    | Malic Acid                       | 134 | C4H6O5  | 006915-15-7 | 56   |
| 3    |    | Butanedioic acid, hydroxy-, (S)- | 134 | C4H6O5  | 000097-67-6 | 56   |
| 4    |    | Malic Acid                       | 134 | C4H6O5  | 006915-15-7 | 56   |
| 5    |    | Thiazole, tetrahydro-            | 89  | C3H7NS  | 000504-78-9 | 9    |



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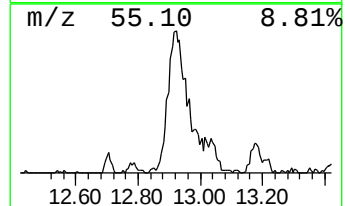
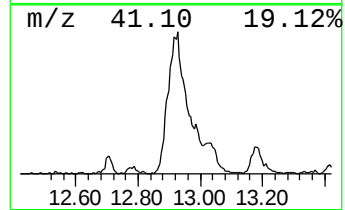
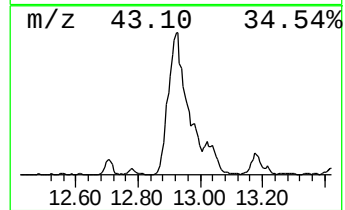
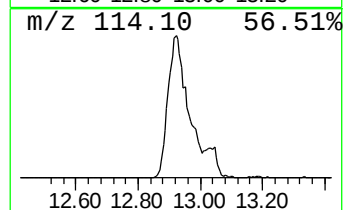
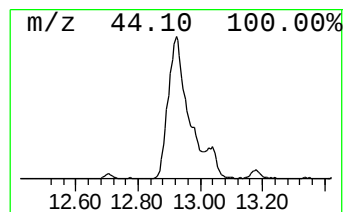
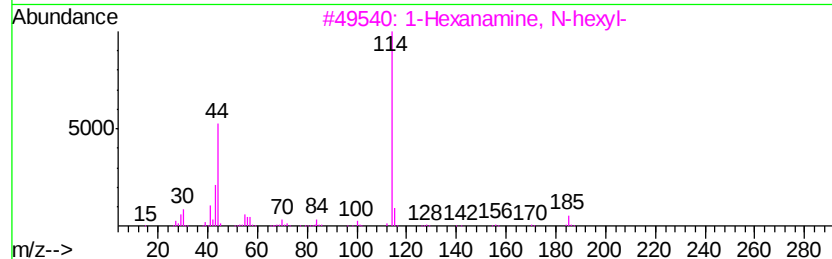
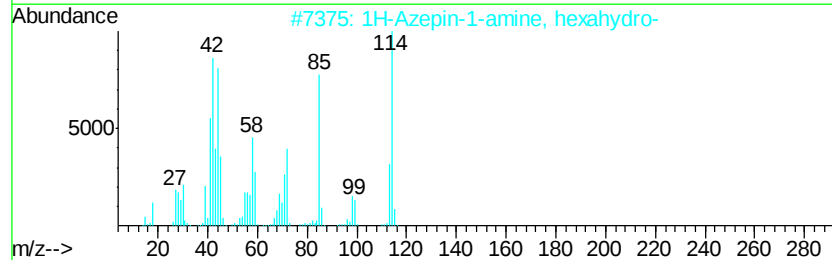
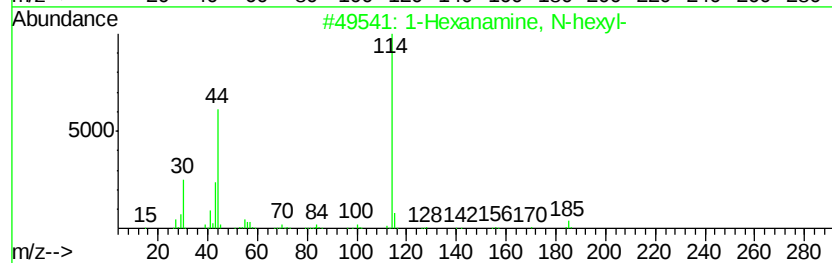
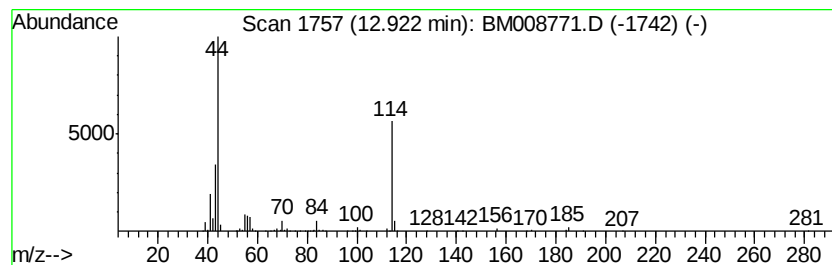
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TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 1-Hexanamine, N-hexyl- Concentration Rank 2

| R.T.    | EstConc     | Area                          | Relative to ISTD | R.T.            |
|---------|-------------|-------------------------------|------------------|-----------------|
| 12.92   | 52.25 ng/ul | 3469860                       | Acenaphthene-d10 | 14.59           |
| Hit# of | 5           | Tentative ID                  | MW MolForm       | CAS# Qual       |
| 1       |             | 1-Hexanamine, N-hexyl-        | 185 C12H27N      | 000143-16-8 78  |
| 2       |             | 1H-Azepin-1-amine, hexahydro- | 114 C6H14N2      | 005906-35-4 50  |
| 3       |             | 1-Hexanamine, N-hexyl-        | 185 C12H27N      | 000143-16-8 47  |
| 4       |             | Cyclobutanol                  | 72 C4H8O         | 002919-23-5 43  |
| 5       |             | Butanoic acid, 2-hexylimino-  | 185 C10H19NO2    | 1000258-67-1 43 |



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Instrument :  
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Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010617.M  
Quant Title : SVOA CALIBRATION

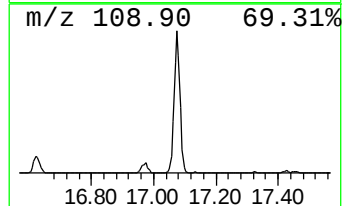
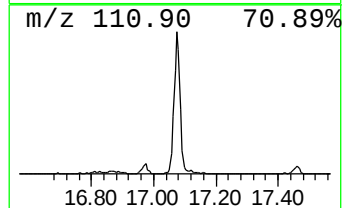
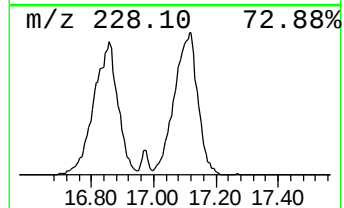
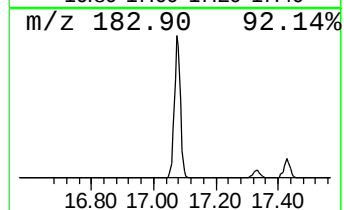
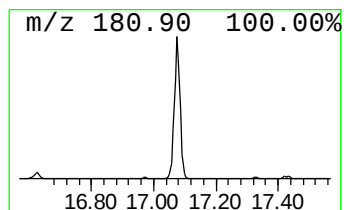
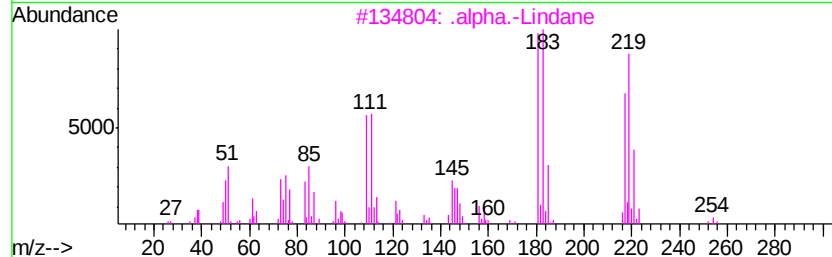
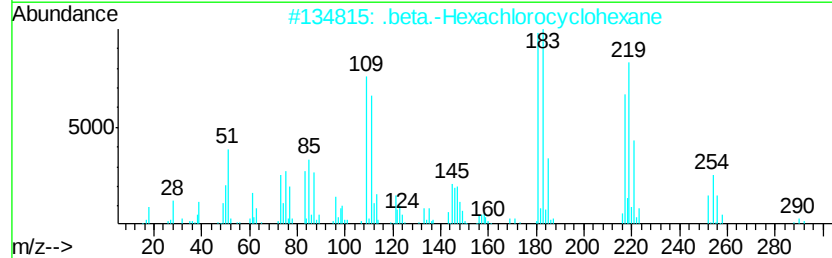
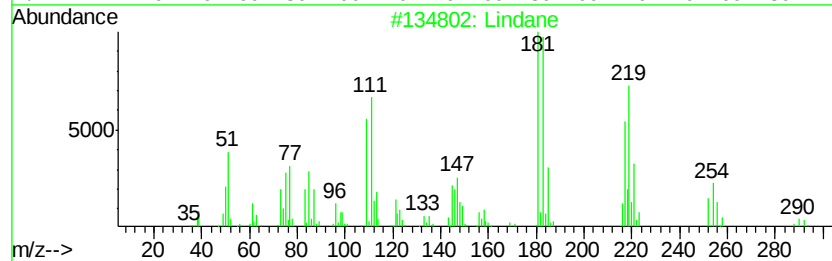
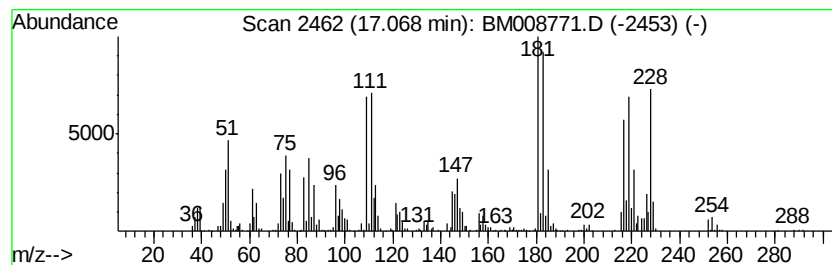
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 Lindane Concentration Rank 3

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 17.07 | 37.85 ng/ul | 3609310 | Phenanthrene-d10 | 17.33 |

| Hit# of | 5                            | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|------------------------------|--------------|-----|---------|-------------|------|
| 1       | Lindane                      |              | 288 | C6H6Cl6 | 000058-89-9 | 93   |
| 2       | .beta.-Hexachlorocyclohexane |              | 288 | C6H6Cl6 | 000319-85-7 | 91   |
| 3       | .alpha.-Lindane              |              | 288 | C6H6Cl6 | 000319-84-6 | 90   |
| 4       | .alpha.-Lindane              |              | 288 | C6H6Cl6 | 000319-84-6 | 90   |
| 5       | Lindane                      |              | 288 | C6H6Cl6 | 000058-89-9 | 86   |



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ALS Vial : 3 Sample Multiplier: 1

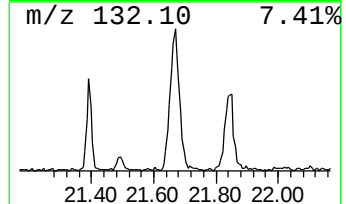
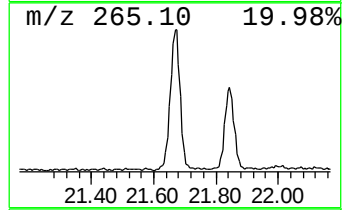
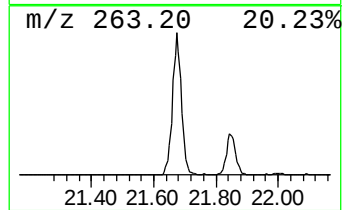
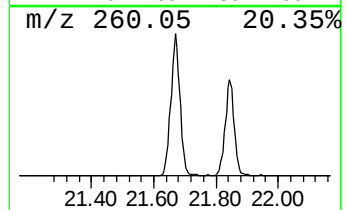
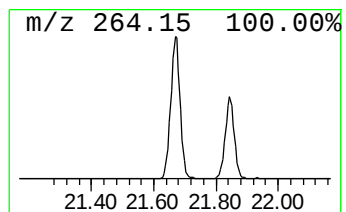
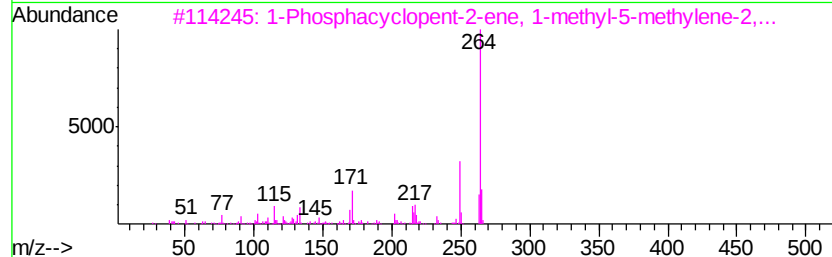
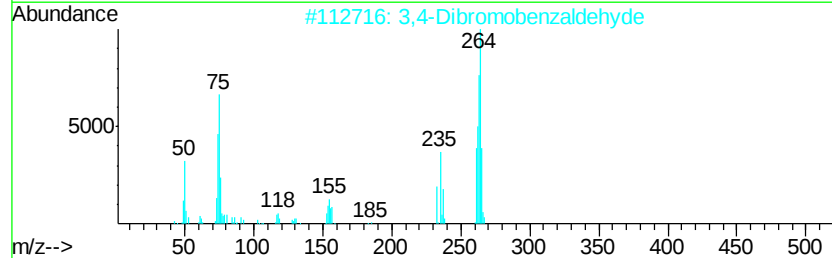
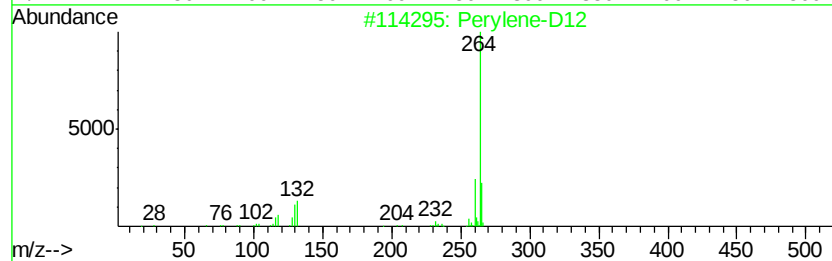
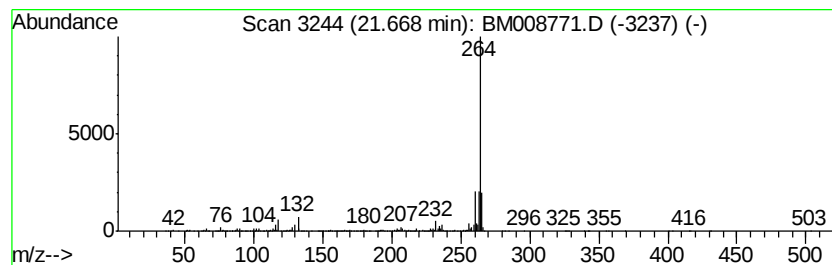
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TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 3,4-Dibromobenzaldehyde Concentration Rank 6

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.      |
|---------|-------------------------------------|--------------|------------------|-----------|
| 21.67   | 13.63 ng/ul                         | 3446010      | Chrysene-d12     | 21.49     |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Perylene-D12                        | 264 C20D12   | 001520-96-3      | 91        |
| 2       | 3,4-Dibromobenzaldehyde             | 262 C7H4Br2O | 074003-55-7      | 64        |
| 3       | 1-Phosphacyclopent-2-ene, 1-meth... | 264 C18H17P  | 1000161-67-7     | 59        |
| 4       | Pyrazolo[1,5-a]pyrimidin-7-amine... | 264 C14H12N6 | 1000276-36-9     | 52        |
| 5       | 5,12-Dioxa-chrysene-6,11-dione      | 264 C16H8O4  | 002288-98-4      | 50        |



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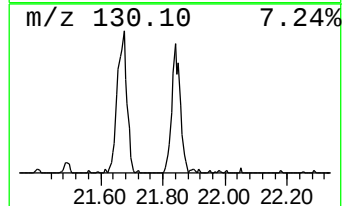
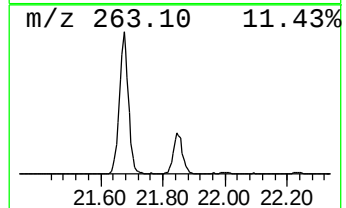
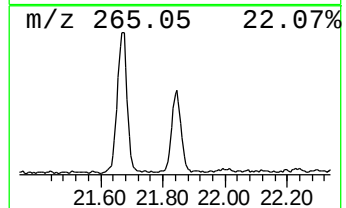
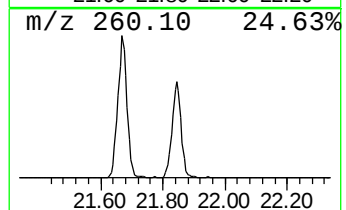
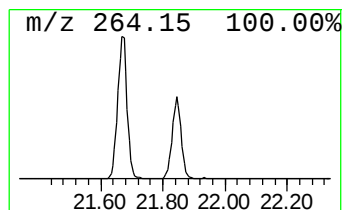
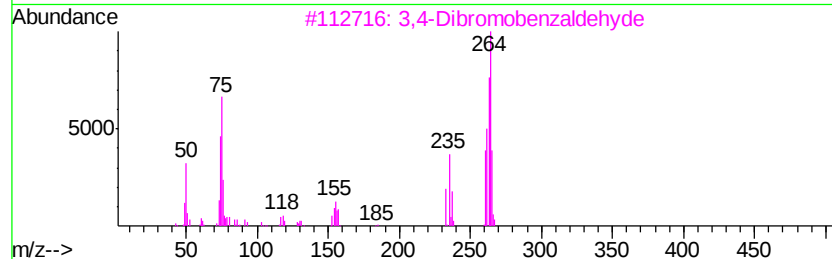
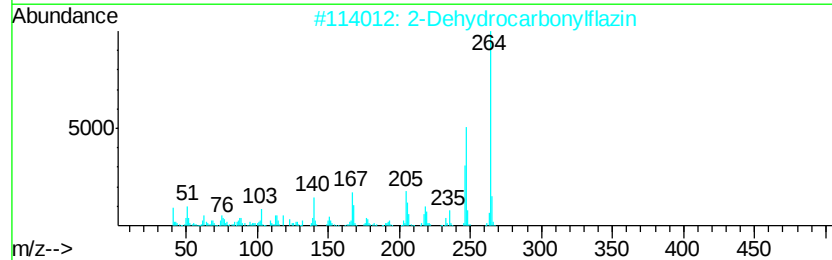
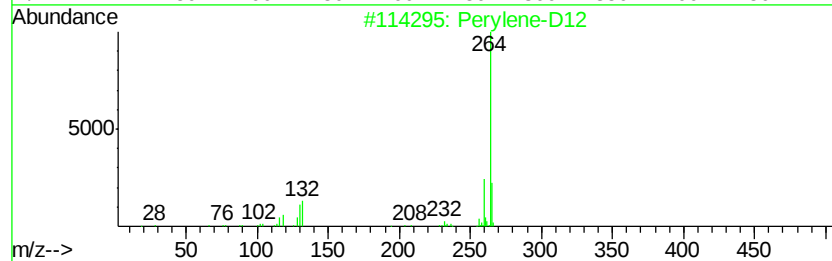
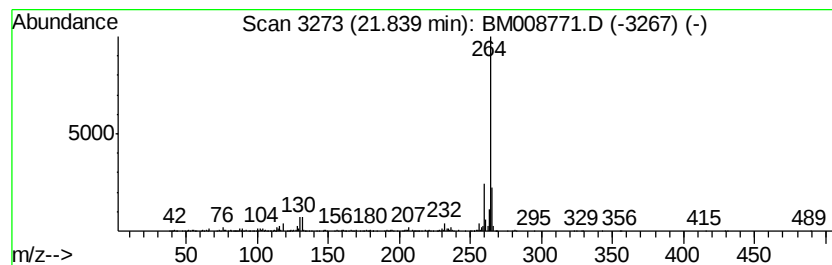
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 2-Dehydrocarbonylfazin Concentration Rank 10

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 21.84 | 7.19 ng/ul | 1818260 | Chrysene-d12     | 21.49 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | Perylene-D12                        | 264 | C20D12      | 001520-96-3  | 94   |
| 2    |    | 2-Dehydrocarbonylfazin              | 264 | C16H12N2O2  | 029700-20-7  | 64   |
| 3    |    | 3,4-Dibromobenzaldehyde             | 262 | C7H4Br2O    | 074003-55-7  | 64   |
| 4    |    | Benzamide, N-[(1,5-dimethyl-1H-p... | 264 | C14H14F2N2O | 1000337-83-3 | 53   |
| 5    |    | Indole, 5-(4-nitrostyryl)-          | 264 | C16H12N2O2  | 1000151-40-1 | 53   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM012117\  
Data File : BM008771.D  
Acq On : 21 Jan 2017 14:28  
Operator : UM/SJ  
Sample : I1323-01  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
DA9P2

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010617.M  
Quant Title : SVOA 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 |
| N-Hydroxymethylac... | 2.93  | 1885.7  | ng/ul | 123926000 | 1                     | 7.96  | 1314390 | 20.0 |
| 1-Hexanamine, N-h... | 12.92 | 52.3    | ng/ul | 3469860   | 3                     | 14.59 | 1328130 | 20.0 |
| Lindane              | 17.07 | 37.9    | ng/ul | 3609310   | 4                     | 17.33 | 1907060 | 20.0 |
| 3,4-Dibromobenzal... | 21.67 | 13.6    | ng/ul | 3446010   | 5                     | 21.49 | 5056430 | 20.0 |
| 2-Dehydrocarbonyl... | 21.84 | 7.2     | ng/ul | 1818260   | 5                     | 21.49 | 5056430 | 20.0 |