

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080318\  
 Data File : BM016174.D  
 Acq On : 03 Aug 2018 11:39  
 Operator : SJ/JU  
 Sample : J4178-03  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C0AB9

## 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:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM071318MA.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         | 8.075       | 803           | 814         | 821          | rBV      | 4103058        | 6158844       | 2.45%           | 0.621%        |
| 2         | 8.228       | 828           | 840         | 850          | rBV      | 4840923        | 7479364       | 2.97%           | 0.754%        |
| 3         | 8.563       | 885           | 897         | 910          | rBV      | 17087426       | 26969111      | 10.72%          | 2.719%        |
| 4         | 10.869      | 1282          | 1289        | 1296         | rBV      | 2811227        | 4601157       | 1.83%           | 0.464%        |
| 5         | 14.222      | 1855          | 1859        | 1863         | rVB      | 2619488        | 3282104       | 1.30%           | 0.331%        |
| 6         | 14.533      | 1906          | 1912        | 1917         | rBV4     | 1363975        | 2619570       | 1.04%           | 0.264%        |
| 7         | 14.627      | 1923          | 1928        | 1933         | rVB      | 3589890        | 4977022       | 1.98%           | 0.502%        |
| 8         | 14.763      | 1945          | 1951        | 1958         | rBV5     | 1074589        | 2854843       | 1.13%           | 0.288%        |
| 9         | 15.516      | 2075          | 2079        | 2085         | rVB2     | 1788989        | 2872920       | 1.14%           | 0.290%        |
| 10        | 15.586      | 2086          | 2091        | 2097         | rVV2     | 6794096        | 10049085      | 3.99%           | 1.013%        |
| 11        | 16.527      | 2245          | 2251        | 2257         | rBV3     | 15006209       | 28598446      | 11.36%          | 2.883%        |
| 12        | 16.627      | 2265          | 2268        | 2272         | rBV      | 3945385        | 6247634       | 2.48%           | 0.630%        |
| 13        | 16.751      | 2285          | 2289        | 2292         | rBV3     | 3961091        | 6433873       | 2.56%           | 0.649%        |
| 14        | 19.062      | 2679          | 2682        | 2686         | rVB      | 12711942       | 16019737      | 6.37%           | 1.615%        |
| 15        | 19.668      | 2781          | 2785        | 2805         | rVB3     | 11605496       | 31067004      | 12.35%          | 3.132%        |
| 16        | 19.827      | 2807          | 2812        | 2818         | rBV      | 14957677       | 22934395      | 9.11%           | 2.312%        |
| 17        | 19.945      | 2828          | 2832        | 2844         | rVB      | 8232887        | 18516811      | 7.36%           | 1.867%        |
| 18        | 20.051      | 2845          | 2850        | 2861         | rVB      | 16630883       | 36529488      | 14.52%          | 3.683%        |
| 19        | 20.168      | 2862          | 2870        | 2877         | rBV4     | 47915166       | 194187025     | 77.17%          | 19.577%       |
| 20        | 20.380      | 2898          | 2906        | 2930         | rVB      | 10880752       | 37598710      | 14.94%          | 3.791%        |
| 21        | 20.603      | 2933          | 2944        | 2952         | rBV5     | 49365778       | 209656185     | 83.31%          | 21.137%       |
| 22        | 20.798      | 2973          | 2977        | 2987         | rBV      | 15482095       | 29829281      | 11.85%          | 3.007%        |
| 23        | 21.003      | 3005          | 3012        | 3024         | rBV3     | 48697971       | 251644076     | 100.00%         | 25.370%       |
| 24        | 21.568      | 3105          | 3108        | 3112         | rVB      | 2644385        | 2886890       | 1.15%           | 0.291%        |
| 25        | 22.203      | 3209          | 3216        | 3220         | rBV      | 20803082       | 27880723      | 11.08%          | 2.811%        |

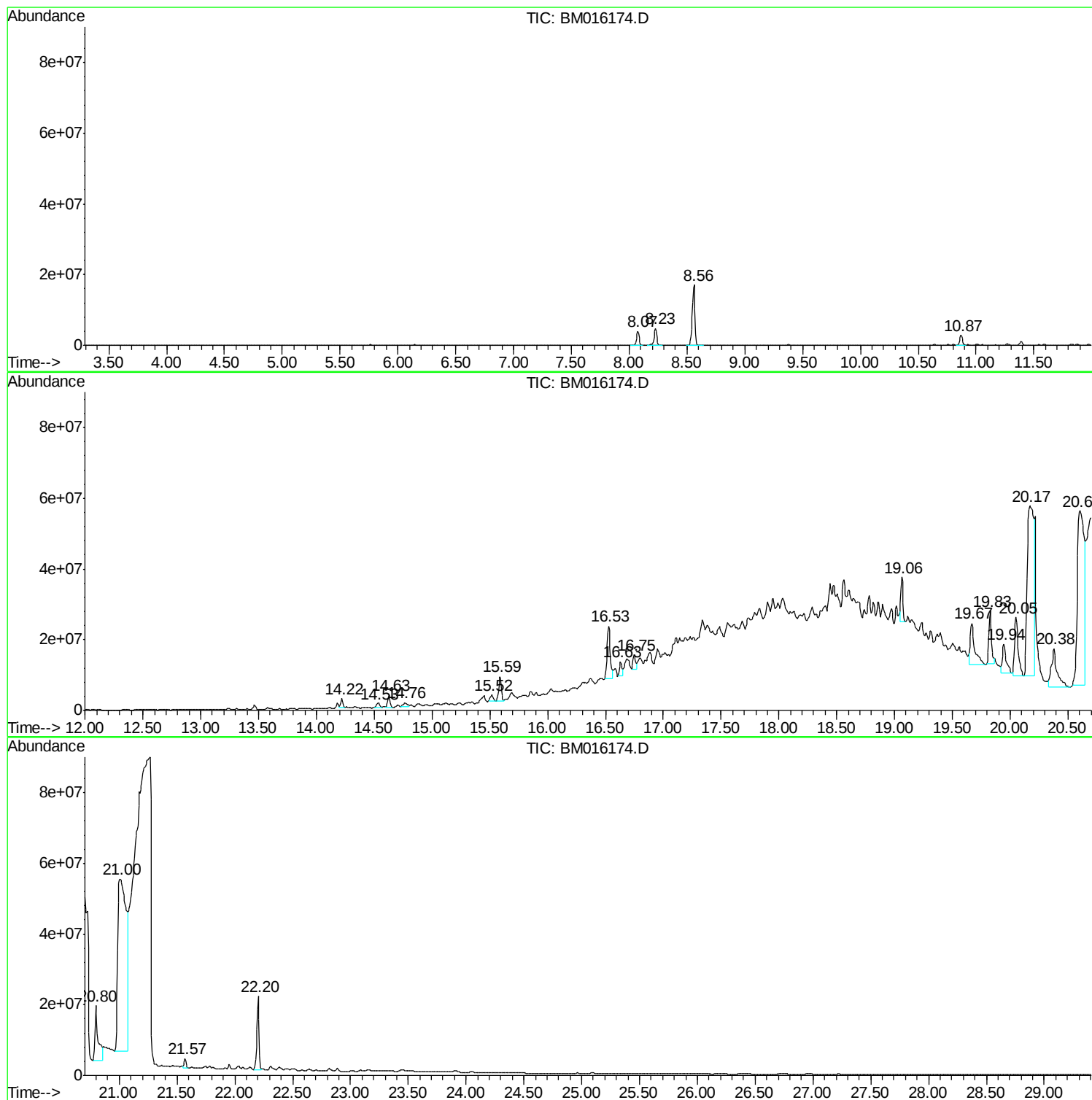
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM071318MA.M  
Quant Title : SVOA CALIBRATION

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



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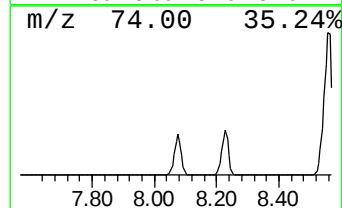
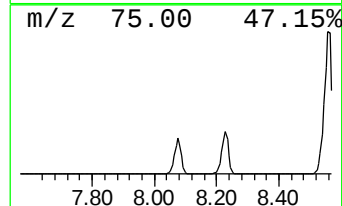
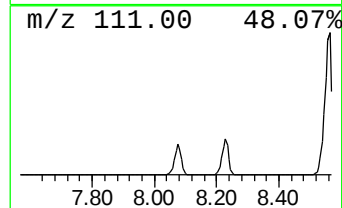
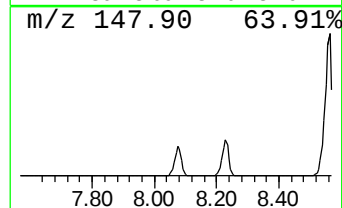
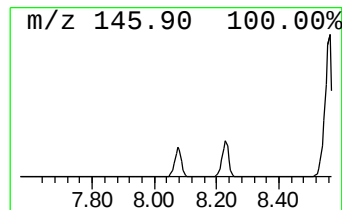
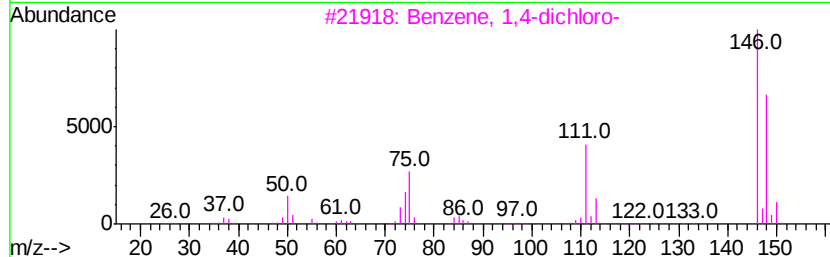
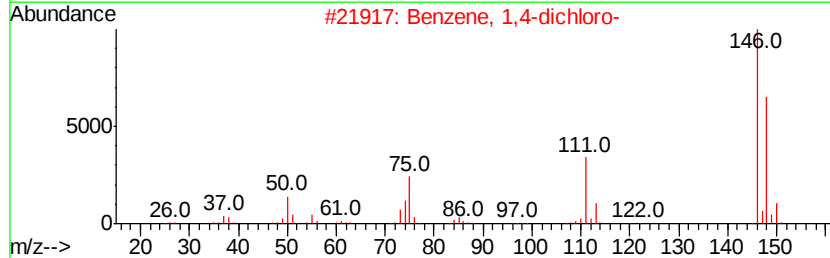
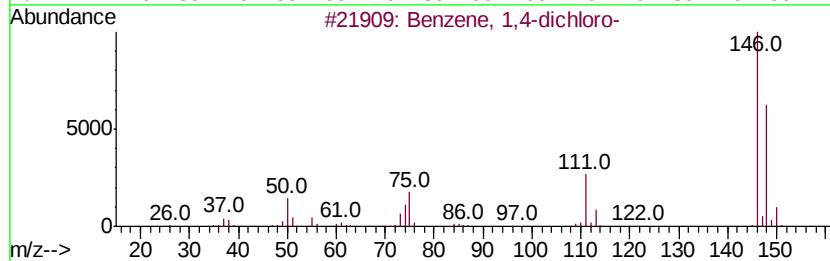
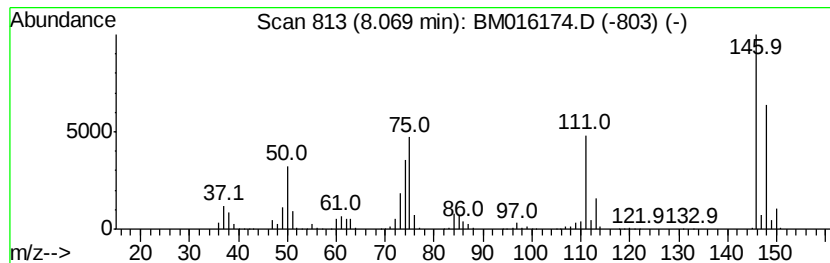
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Quant Title : SVOA CALIBRATION

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

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Peak Number 1 Benzene, 1,4-dichloro- Concentration Rank 20

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 8.07 | 16.47 ng/ul | 6158840 | 1,4-Dichlorobenzene-d4 | 8.19 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,4-dichloro- | 146 | C6H4Cl2 | 000106-46-7 | 97   |
| 2    |    | Benzene, 1,4-dichloro- | 146 | C6H4Cl2 | 000106-46-7 | 97   |
| 3    |    | Benzene, 1,4-dichloro- | 146 | C6H4Cl2 | 000106-46-7 | 96   |
| 4    |    | Benzene, 1,4-dichloro- | 146 | C6H4Cl2 | 000106-46-7 | 96   |
| 5    |    | Benzene, 1,3-dichloro- | 146 | C6H4Cl2 | 000541-73-1 | 96   |



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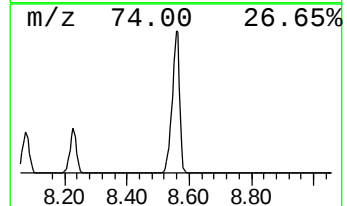
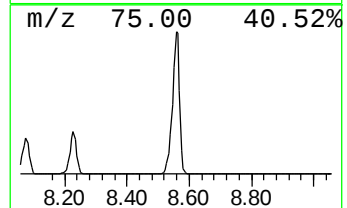
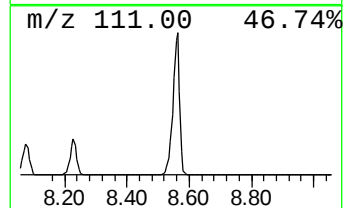
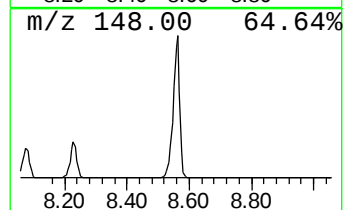
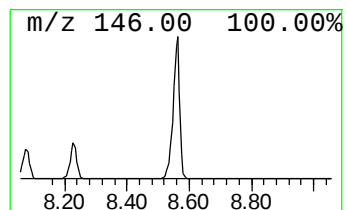
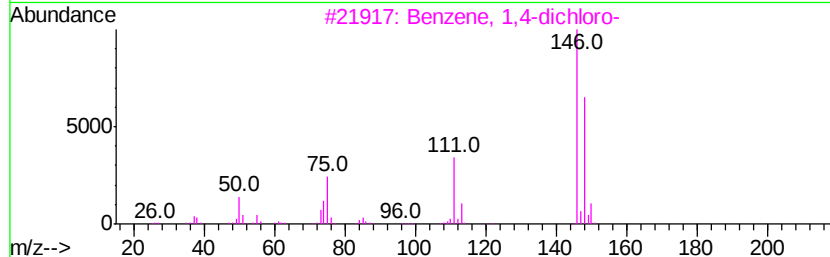
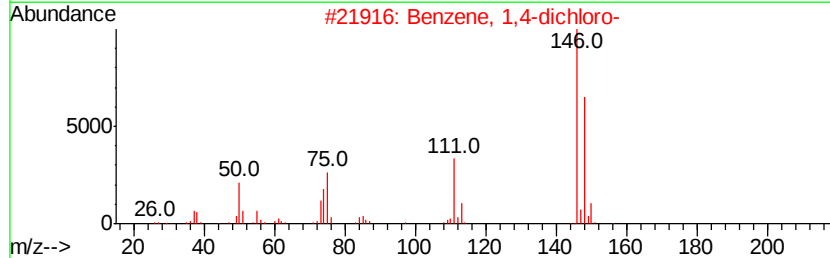
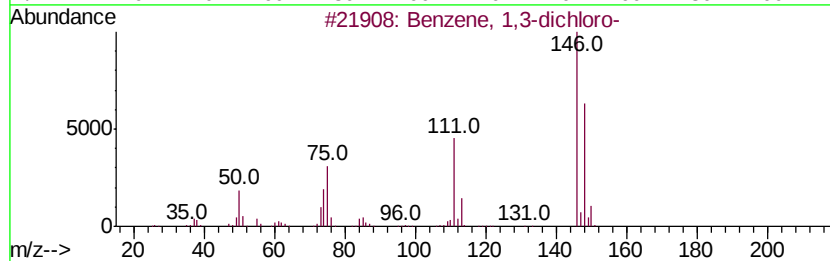
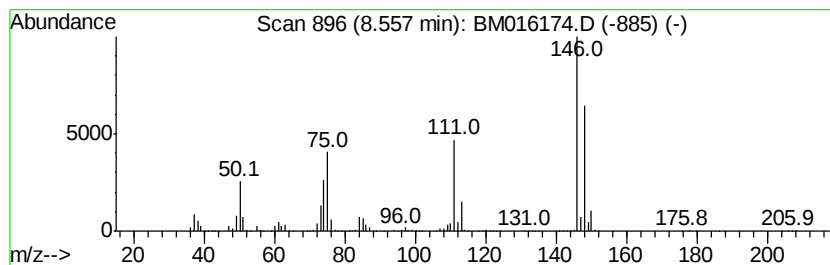
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM071318MA.M  
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TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Benzene, 1,3-dichloro- Concentration Rank 11

| R.T. | EstConc     | Area     | Relative to ISTD       | R.T. |
|------|-------------|----------|------------------------|------|
| 8.56 | 72.12 ng/ul | 26969100 | 1,4-Dichlorobenzene-d4 | 8.19 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,3-dichloro- | 146 | C6H4Cl2 | 000541-73-1 | 98   |
| 2    |    | Benzene, 1,4-dichloro- | 146 | C6H4Cl2 | 000106-46-7 | 98   |
| 3    |    | Benzene, 1,4-dichloro- | 146 | C6H4Cl2 | 000106-46-7 | 97   |
| 4    |    | Benzene, 1,2-dichloro- | 146 | C6H4Cl2 | 000095-50-1 | 97   |
| 5    |    | Benzene, 1,2-dichloro- | 146 | C6H4Cl2 | 000095-50-1 | 96   |



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ALS Vial : 5 Sample Multiplier: 1

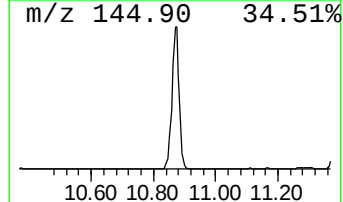
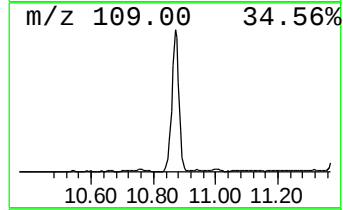
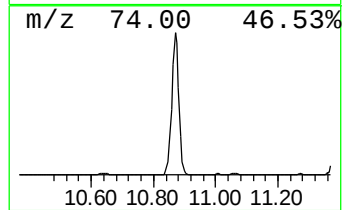
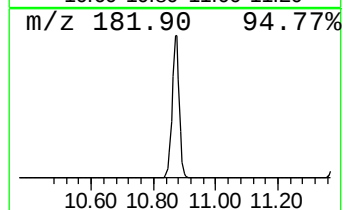
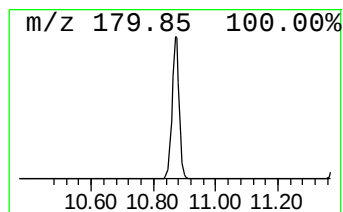
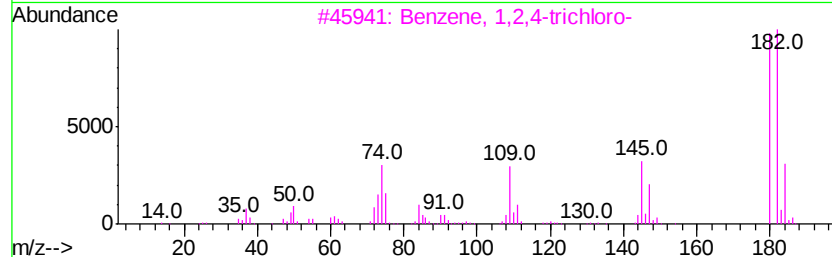
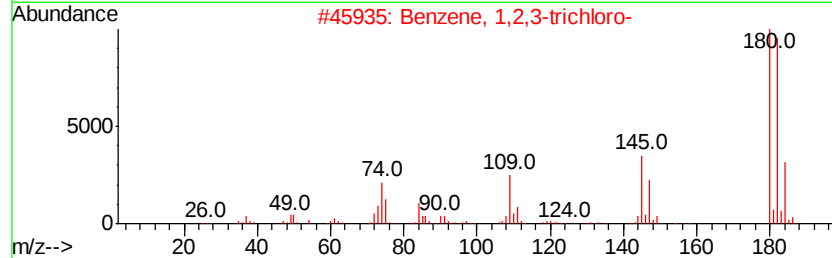
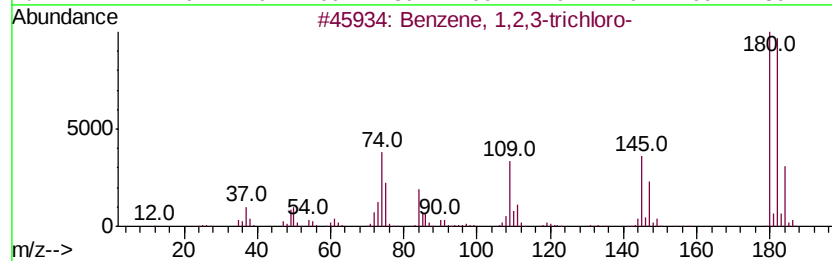
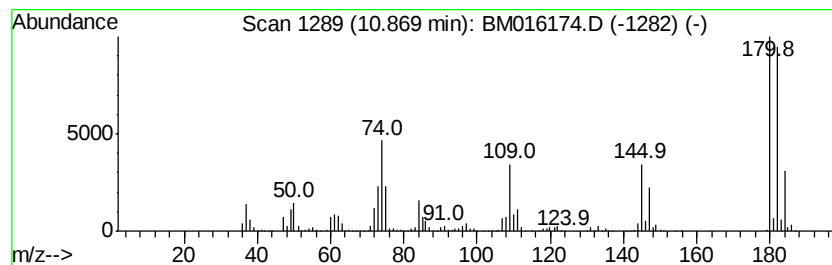
Instrument :  
BNA\_M  
ClientSampleId :  
C0AB9

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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 3 Benzene, 1,2,3-trichloro- Concentration Rank 16

| R.T.    | EstConc     | Area                      | Relative to ISTD | R.T.           |
|---------|-------------|---------------------------|------------------|----------------|
| 10.87   | 20.00 ng/ul | 4601160                   | Naphthalene-d8   | 11.01          |
| Hit# of | 5           | Tentative ID              | MW MolForm       | CAS# Qual      |
| 1       |             | Benzene, 1,2,3-trichloro- | 180 C6H3Cl3      | 000087-61-6 98 |
| 2       |             | Benzene, 1,2,3-trichloro- | 180 C6H3Cl3      | 000087-61-6 97 |
| 3       |             | Benzene, 1,2,4-trichloro- | 180 C6H3Cl3      | 000120-82-1 97 |
| 4       |             | Benzene, 1,3,5-trichloro- | 180 C6H3Cl3      | 000108-70-3 97 |
| 5       |             | Benzene, 1,2,3-trichloro- | 180 C6H3Cl3      | 000087-61-6 97 |



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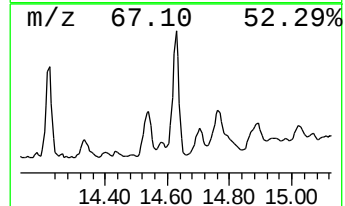
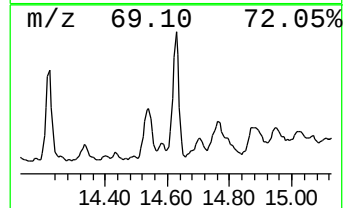
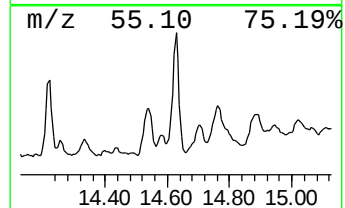
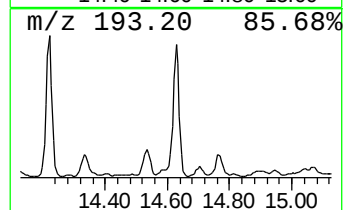
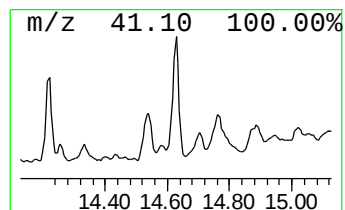
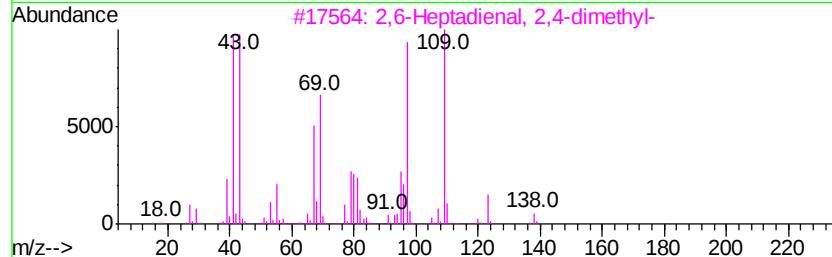
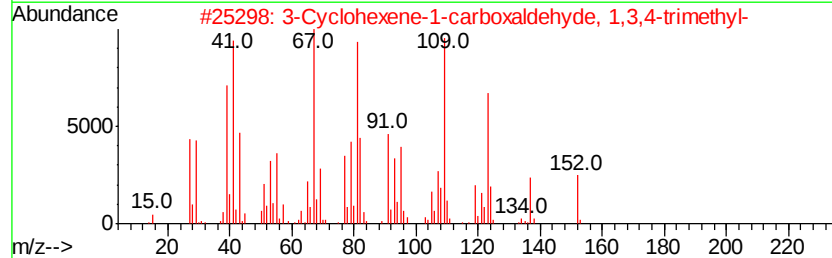
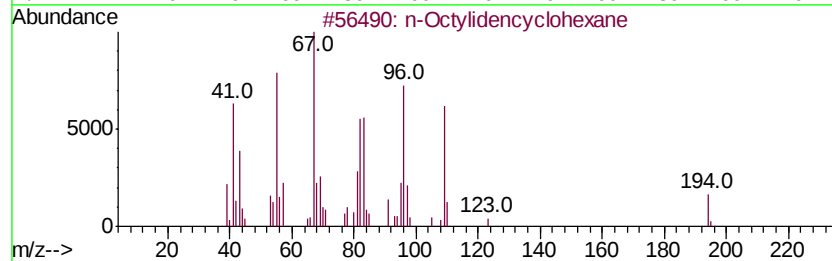
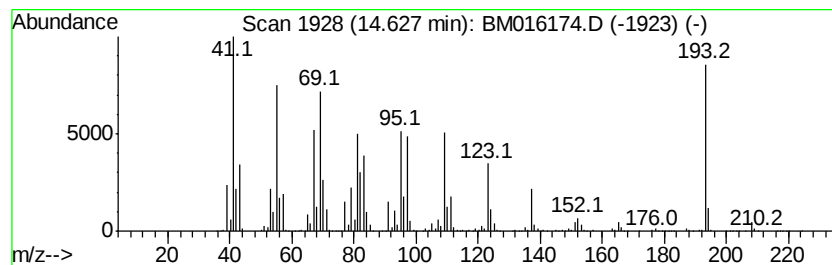
Instrument :  
BNA\_M  
ClientSampleId :  
C0AB9

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TIC Integration Parameters: LSCINT.P

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Peak Number 4 n-Octylidencyclohexane Concentration Rank 14

| R.T.    | EstConc                             | Area          | Relative to ISTD | R.T.      |
|---------|-------------------------------------|---------------|------------------|-----------|
| 14.63   | 34.87 ng/ul                         | 4977020       | Acenaphthene-d10 | 14.82     |
| Hit# of | 5                                   | Tentative ID  | MW MolForm       | CAS# Qual |
| 1       | n-Octylidencyclohexane              | 194 C14H26    | 137595-12-1      | 55        |
| 2       | 3-Cyclohexene-1-carboxaldehyde, ... | 152 C10H16O   | 040702-26-9      | 25        |
| 3       | 2,6-Heptadienal, 2,4-dimethyl-      | 138 C9H14O    | 085136-08-9      | 22        |
| 4       | 1-(p-Fluorophenyl)-4-piperidone     | 193 C11H12FNO | 1000238-56-7     | 22        |
| 5       | Neopentylidenecyclohexane           | 152 C11H20    | 039546-80-0      | 22        |



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Quant Title : SVOA CALIBRATION

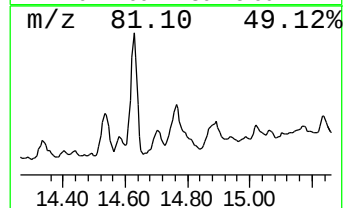
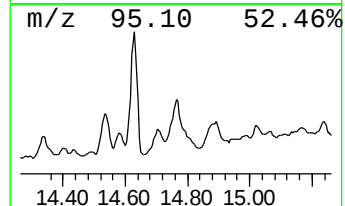
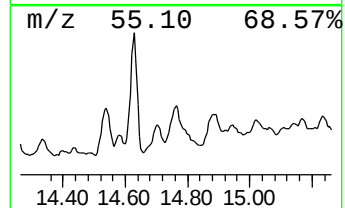
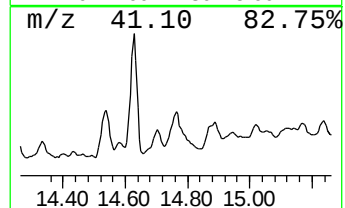
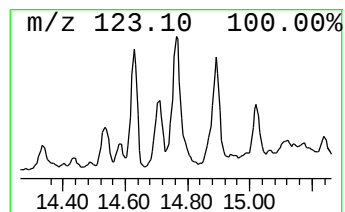
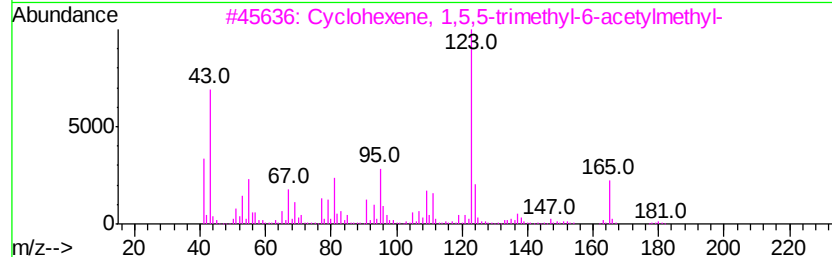
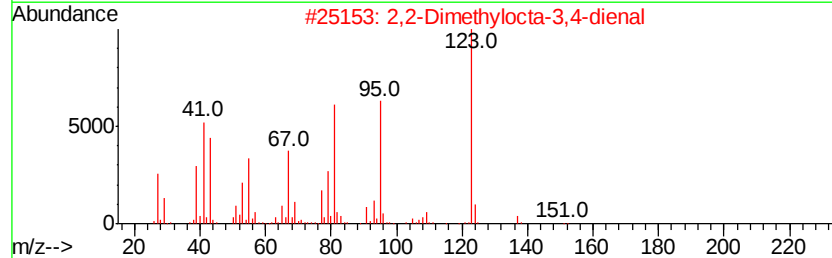
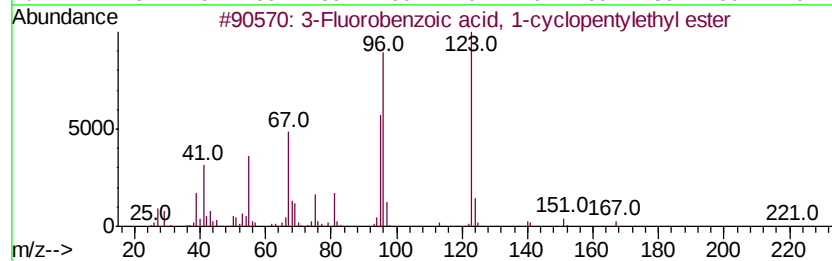
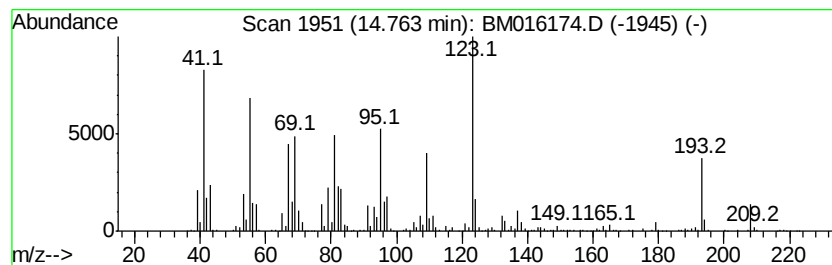
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 3-Fluorobenzoic acid, 1-cyc... Concentration Rank 17

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 14.76 | 20.00 ng/ul | 2854840 | Acenaphthene-d10 | 14.82 |

| Hit# | of | Tentative ID                           | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Fluorobenzoic acid, 1-cyclopenten... | 236 | C14H17F02 | 1000279-04-7 | 50   |
| 2    |    | 2,2-Dimethylocta-3,4-dienal            | 152 | C10H16O   | 000590-71-6  | 47   |
| 3    |    | Cyclohexene, 1,5,5-trimethyl-6-ac...   | 180 | C12H20O   | 211563-96-1  | 43   |
| 4    |    | 4-Fluorobenzoic acid, 2-pentyl e...    | 210 | C12H15F02 | 1000279-07-2 | 43   |
| 5    |    | Cyclohexene, 1,4,6,6-tetramethyl-      | 138 | C10H18    | 070092-37-4  | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080318\  
Data File : BM016174.D  
Acq On : 03 Aug 2018 11:39  
Operator : SJ/JU  
Sample : J4178-03  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0AB9

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Quant Title : SVOA CALIBRATION

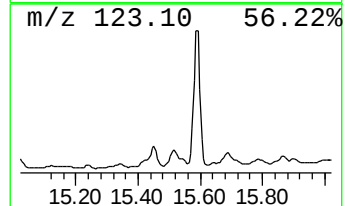
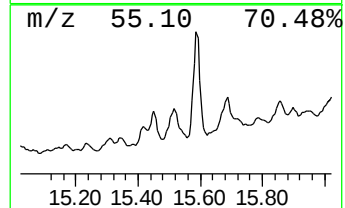
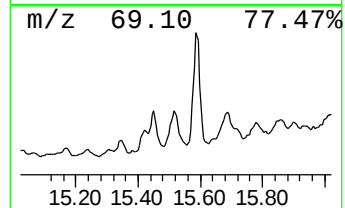
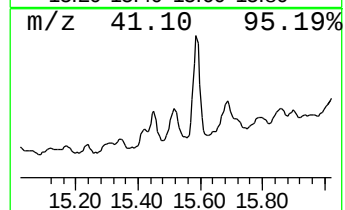
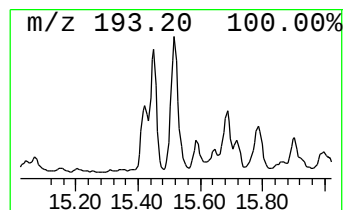
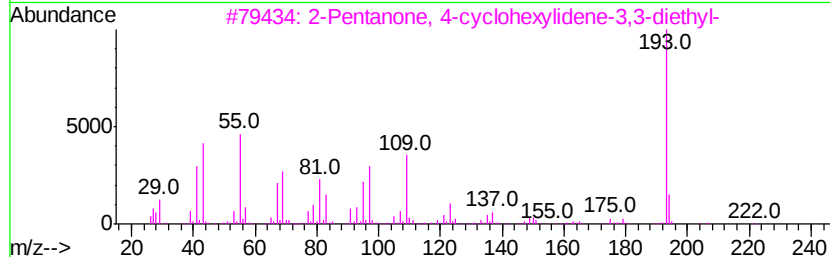
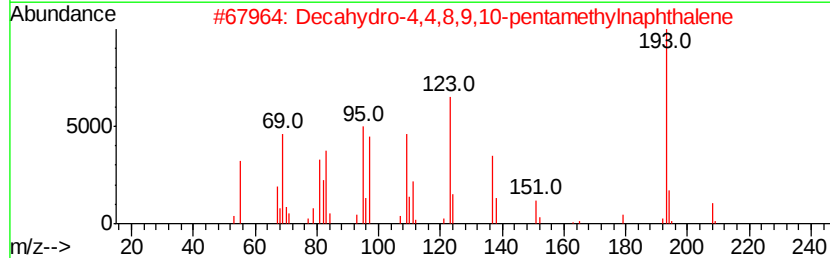
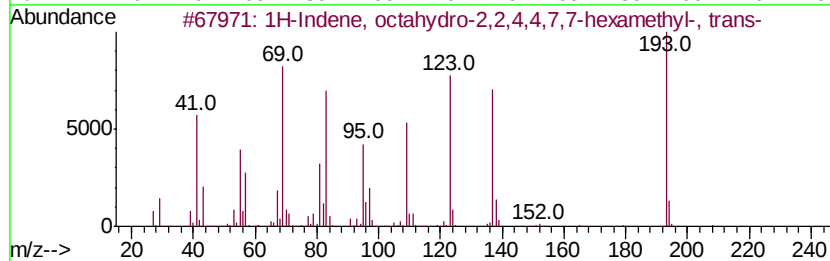
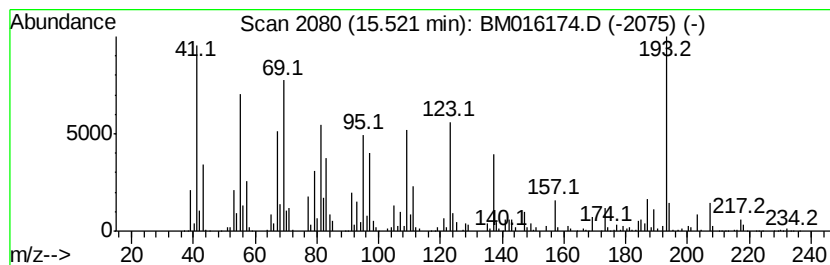
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 1H-Indene, octahydro-2,2,4,... Concentration Rank 15

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.52 | 20.13 ng/ul | 2872920 | Acenaphthene-d10 | 14.82 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | 1H-Indene, octahydro-2,2,4,4,7,7... | 208 | C15H28    | 054832-83-6 | 62   |
| 2    |    | Decahydro-4,4,8,9,10-pentamethyl... | 208 | C15H28    | 080655-44-3 | 58   |
| 3    |    | 2-Pentanone, 4-cyclohexylidene-3... | 222 | C15H26O   | 313253-65-5 | 38   |
| 4    |    | 4H-1,3,2-Dioxaborin, 4-ethenyl-2... | 208 | C12H21BO2 | 074630-04-9 | 30   |
| 5    |    | Bicyclo[4.1.0]heptane, 3,7,7-tri... | 138 | C10H18    | 006069-97-2 | 25   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080318\  
Data File : BM016174.D  
Acq On : 03 Aug 2018 11:39  
Operator : SJ/JU  
Sample : J4178-03  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

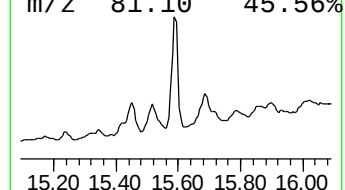
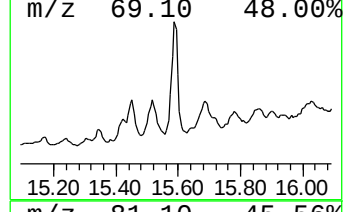
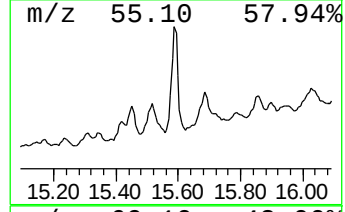
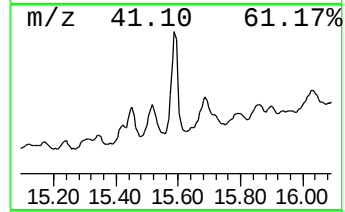
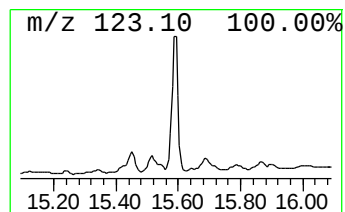
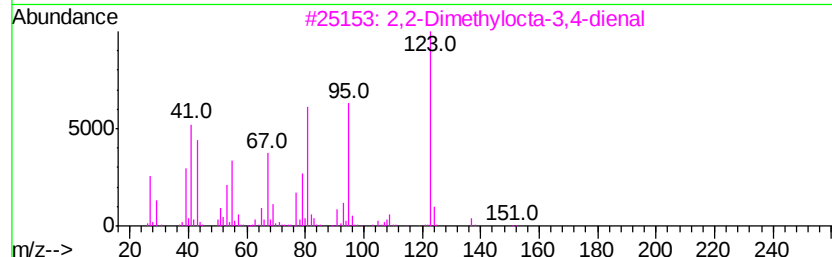
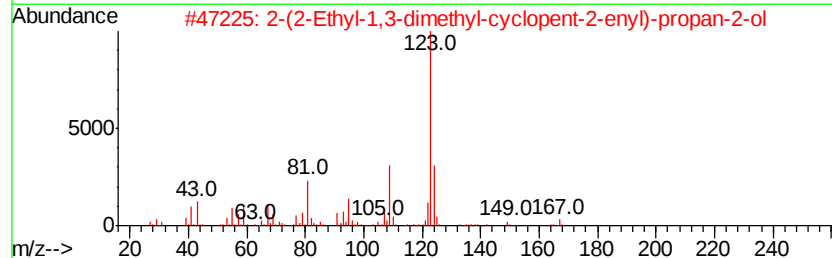
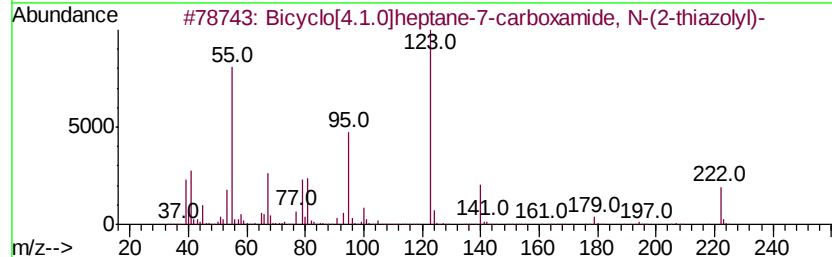
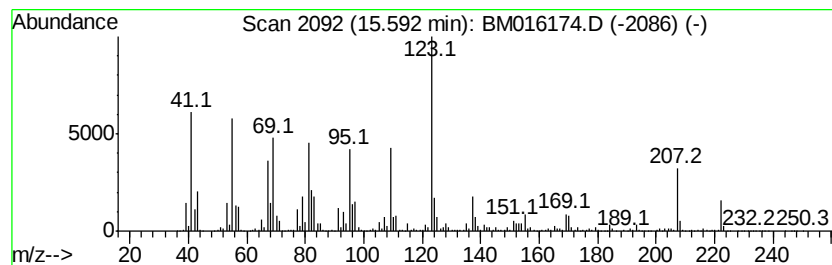
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ClientSampleId :  
C0AB9

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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 7 unknown-01 Concentration Rank 12

| R.T.    | EstConc     | Area                                 | Relative to ISTD | R.T.            |
|---------|-------------|--------------------------------------|------------------|-----------------|
| 15.59   | 70.40 ng/ul | 10049100                             | Acenaphthene-d10 | 14.82           |
| Hit# of | 5           | Tentative ID                         | MW MolForm       | CAS# Qual       |
| 1       |             | Bicyclo[4.1.0]heptane-7-carboxam...  | 222 C11H14N2O5   | 329912-72-3 46  |
| 2       |             | 2-(2-Ethyl-1,3-dimethyl-cyclopent... | 182 C12H22O      | 1000186-82-4 45 |
| 3       |             | 2,2-Dimethylocta-3,4-dienal          | 152 C10H16O      | 000590-71-6 43  |
| 4       |             | 4-Fluorobenzoic acid, 2,2,2-trif...  | 222 C9H6F4O2     | 1000325-53-2 38 |
| 5       |             | 4-Fluorobenzoic acid, pentafluor...  | 306 C13H4F6O2    | 1000355-67-4 35 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080318\  
Data File : BM016174.D  
Acq On : 03 Aug 2018 11:39  
Operator : SJ/JU  
Sample : J4178-03  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0AB9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM071318MA.M  
Quant Title : SVOA CALIBRATION

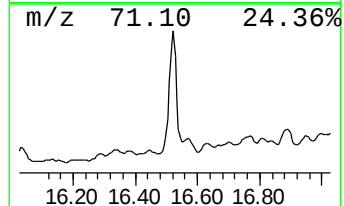
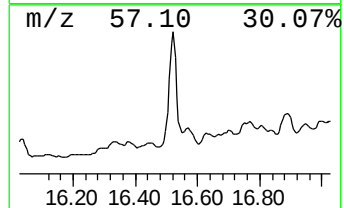
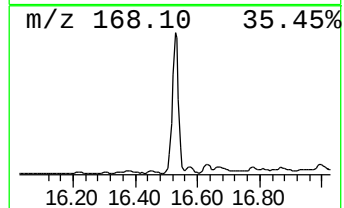
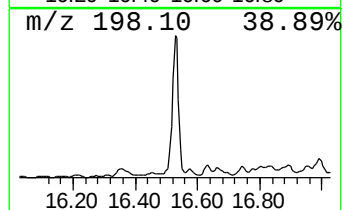
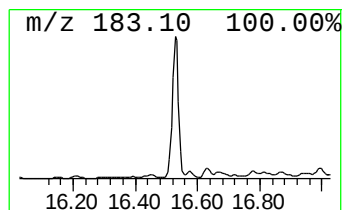
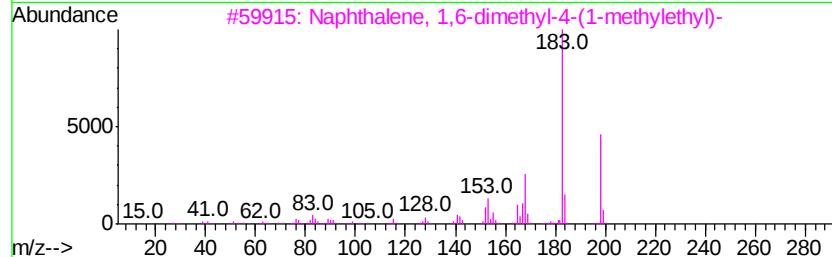
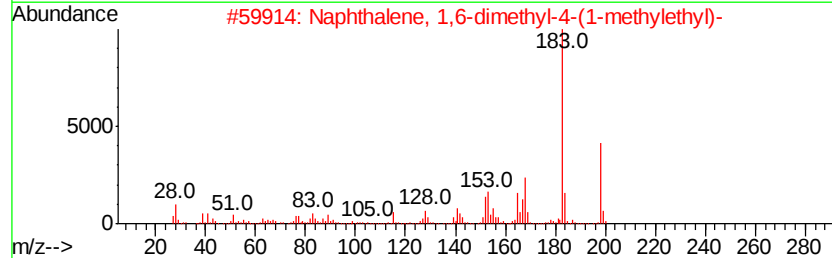
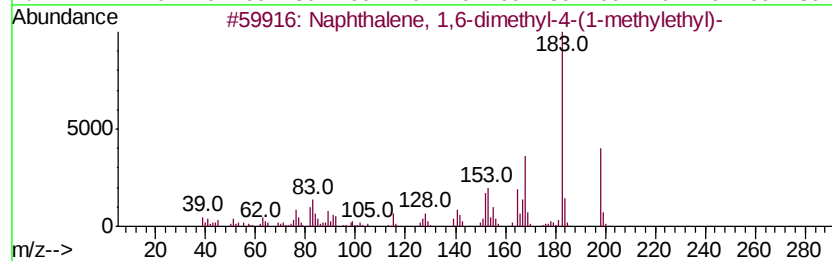
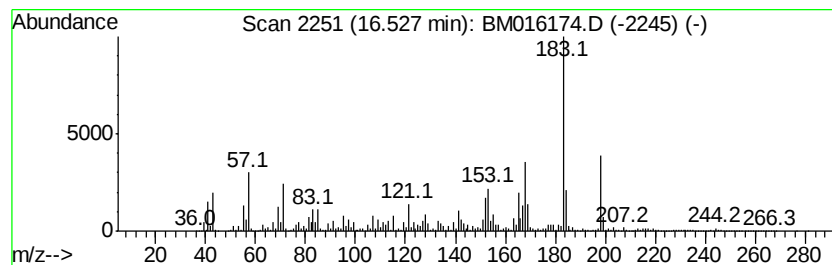
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 Naphthalene, 1,6-dimethyl-4... Concentration Rank 10

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 16.53 | 88.90 ng/ul | 28598400 | Phenanthrene-d10 | 17.57 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,6-dimethyl-4-(1-m... | 198 | C15H18  | 000483-78-3 | 98   |
| 2    |    | Naphthalene, 1,6-dimethyl-4-(1-m... | 198 | C15H18  | 000483-78-3 | 98   |
| 3    |    | Naphthalene, 1,6-dimethyl-4-(1-m... | 198 | C15H18  | 000483-78-3 | 94   |
| 4    |    | Azulene, 1,4-dimethyl-7-(1-methy... | 198 | C15H18  | 000489-84-9 | 93   |
| 5    |    | Azulene, 1,4-dimethyl-7-(1-methy... | 198 | C15H18  | 000489-84-9 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080318\  
Data File : BM016174.D  
Acq On : 03 Aug 2018 11:39  
Operator : SJ/JU  
Sample : J4178-03  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

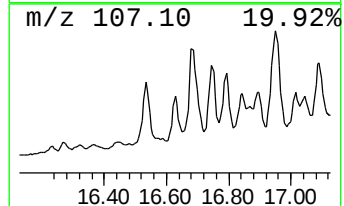
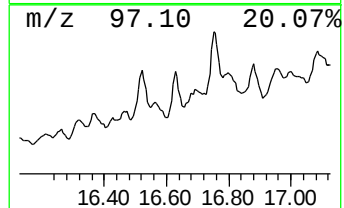
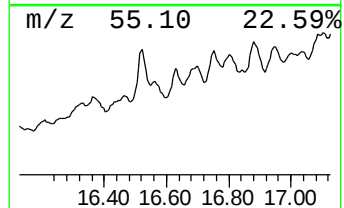
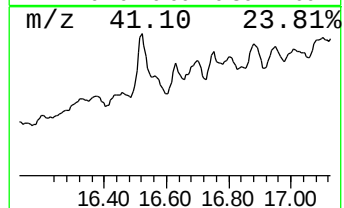
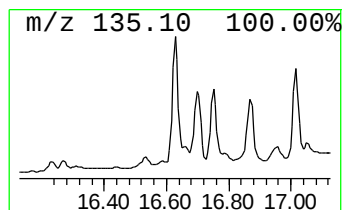
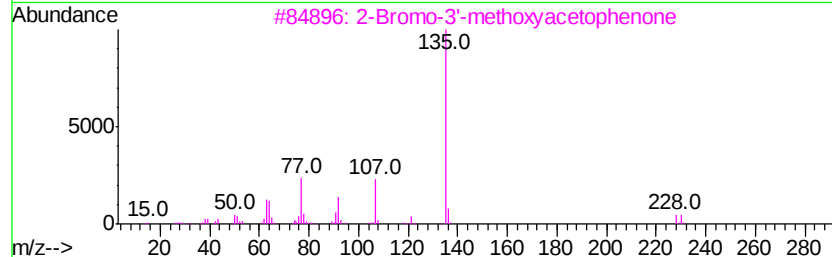
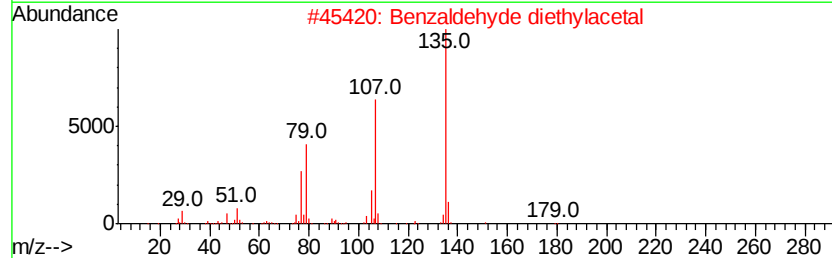
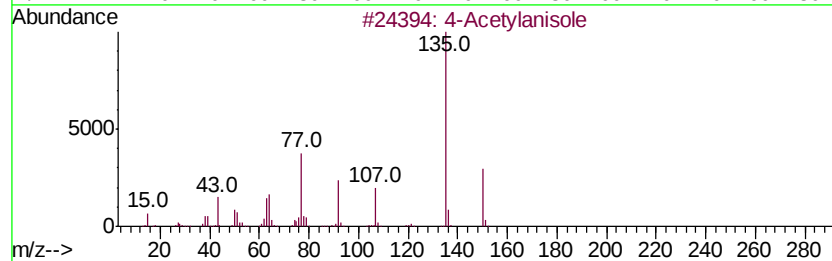
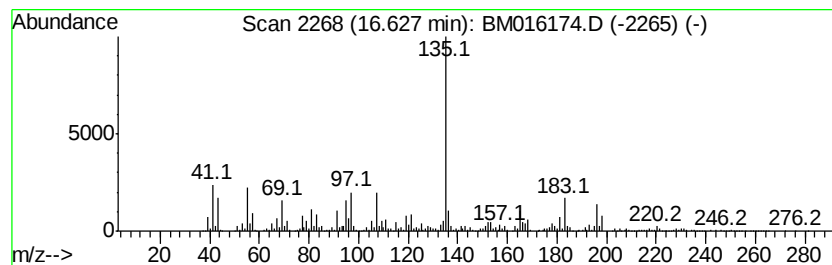
Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM071318MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 9 unknown-02 Concentration Rank 19

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.      |
|---------|-------------------------------------|--------------|------------------|-----------|
| 16.63   | 19.42 ng/ul                         | 6247630      | Phenanthrene-d10 | 17.57     |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | 4-Acetylanisole                     | 150 C9H10O2  | 000100-06-1      | 47        |
| 2       | Benzaldehyde diethylacetal          | 180 C11H16O2 | 000774-48-1      | 47        |
| 3       | 2-Bromo-3'-methoxycetophenone       | 228 C9H9BrO2 | 005000-65-7      | 47        |
| 4       | 1-Adamantaneethanol                 | 180 C12H20O  | 006240-11-5      | 43        |
| 5       | 3-Adamantan-1-yl-3-oxo-propionit... | 203 C13H17NO | 1000296-74-6     | 43        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080318\  
Data File : BM016174.D  
Acq On : 03 Aug 2018 11:39  
Operator : SJ/JU  
Sample : J4178-03  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0AB9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM071318MA.M  
Quant Title : SVOA CALIBRATION

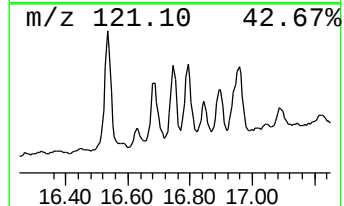
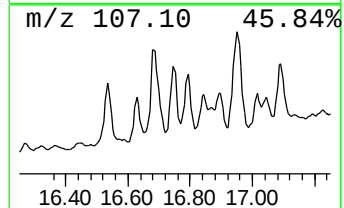
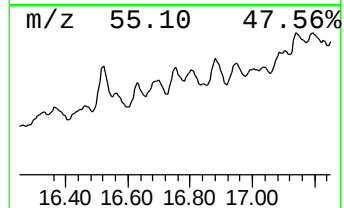
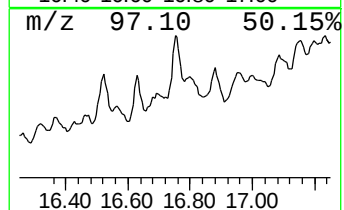
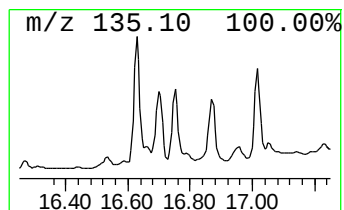
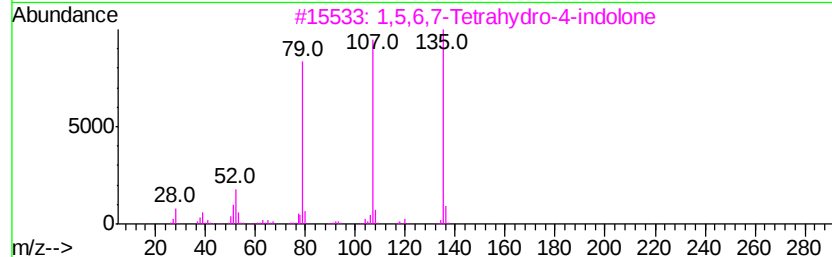
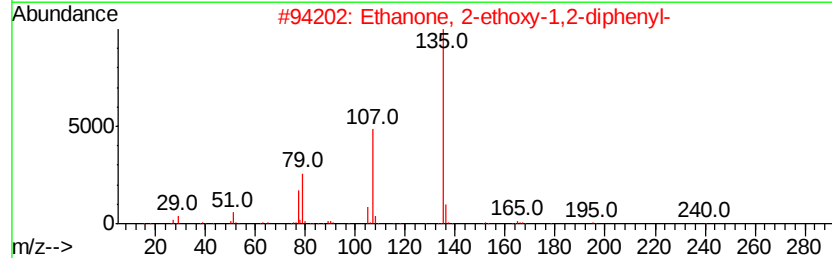
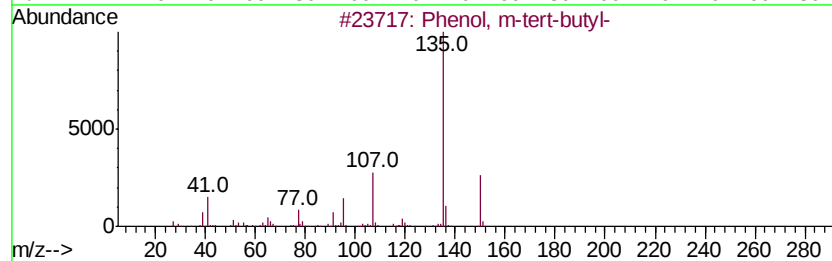
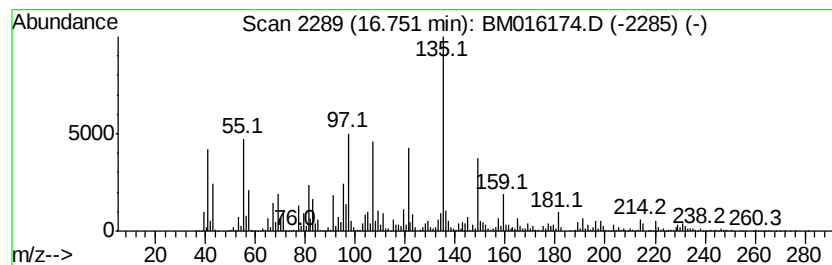
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 unknown-03 Concentration Rank 18

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 16.75 | 20.00 ng/ul | 6433870 | Phenanthrene-d10 | 17.57 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Phenol, m-tert-butyl-               | 150 | C10H14O  | 000585-34-2  | 47   |
| 2    |    | Ethanone, 2-ethoxy-1,2-diphenyl-    | 240 | C16H16O2 | 000574-09-4  | 43   |
| 3    |    | 1,5,6,7-Tetrahydro-4-indolone       | 135 | C8H9NO   | 013754-86-4  | 38   |
| 4    |    | Methanol, (cyclohexyl)(2,3-dimet... | 218 | C15H22O  | 349498-47-1  | 38   |
| 5    |    | Ethylene glycol, 2-methoxybenzoate  | 196 | C10H12O4 | 1000374-70-1 | 30   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080318\  
Data File : BM016174.D  
Acq On : 03 Aug 2018 11:39  
Operator : SJ/JU  
Sample : J4178-03  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Title : SVOA CALIBRATION

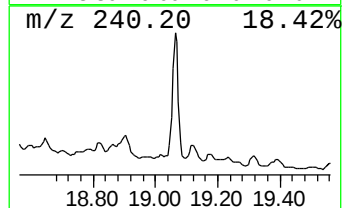
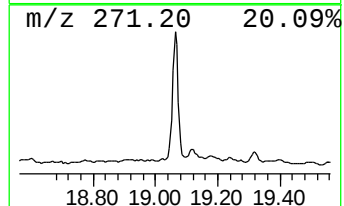
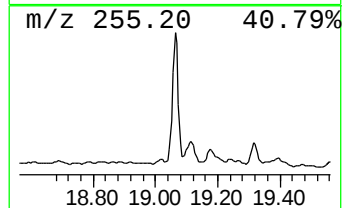
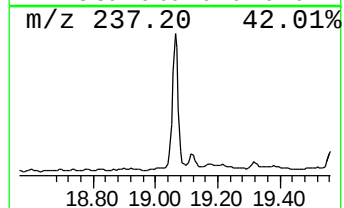
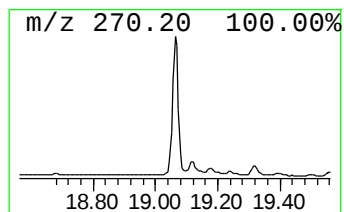
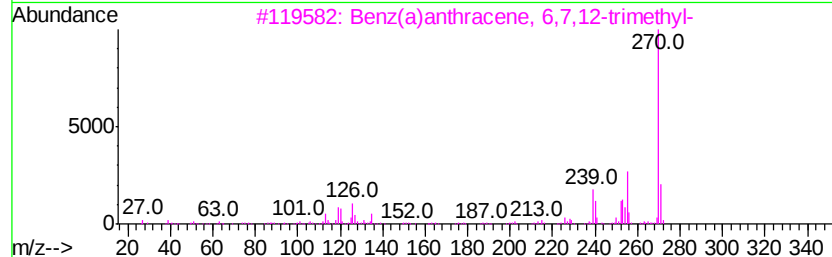
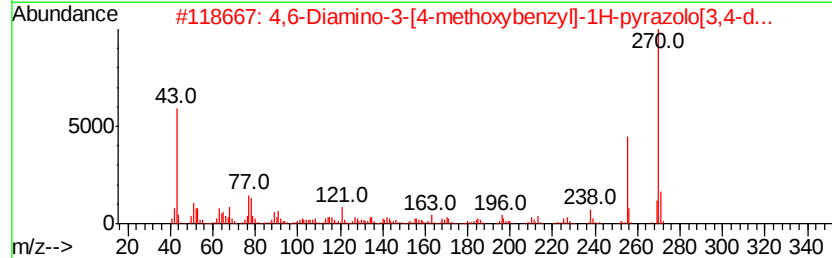
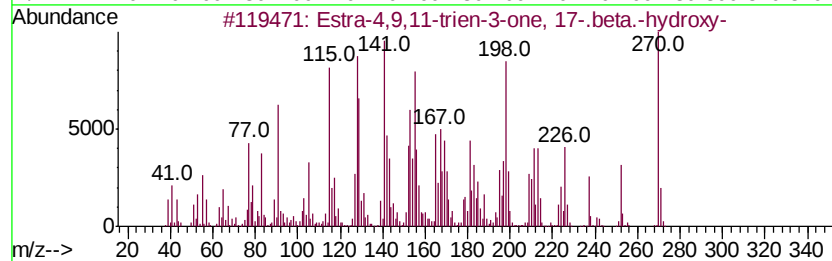
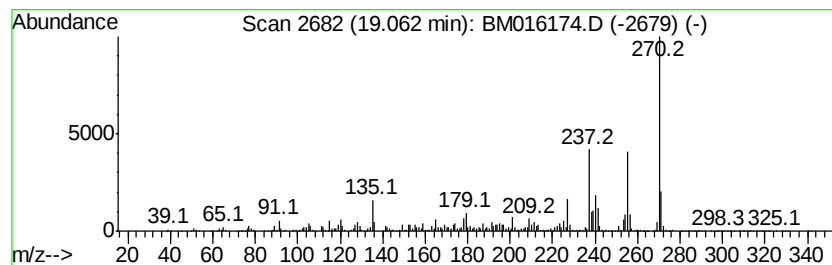
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TIC Integration Parameters: LSCINT.P

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Peak Number 11 Estra-4,9,11-trien-3-one, 1... Concentration Rank 13

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 19.06 | 49.80 ng/ul | 16019700 | Phenanthrene-d10 | 17.57 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Estra-4,9,11-trien-3-one, 17-.be... | 270 | C18H22O2  | 010161-33-8 | 92   |
| 2    |    | 4,6-Diamino-3-[4-methoxybenzyl]-... | 270 | C13H14N6O | 058791-73-4 | 78   |
| 3    |    | Benz(a)anthracene, 6,7,12-trimet... | 270 | C21H18    | 020627-33-2 | 55   |
| 4    |    | 1H-Naphtho[2,1-b]pyran-1-one, 7,... | 270 | C16H14O4  | 054889-69-9 | 50   |
| 5    |    | Benz(a)anthracene, 6,7,8-trimethyl- | 270 | C21H18    | 020627-32-1 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080318\  
Data File : BM016174.D  
Acq On : 03 Aug 2018 11:39  
Operator : SJ/JU  
Sample : J4178-03  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

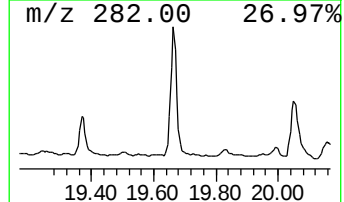
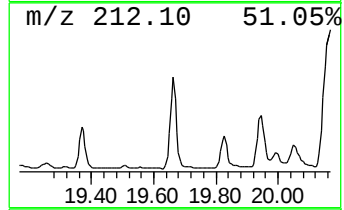
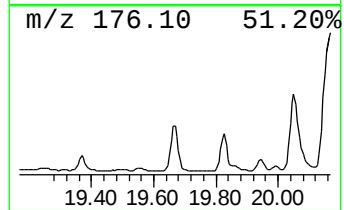
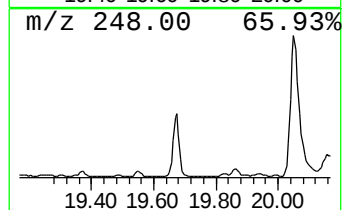
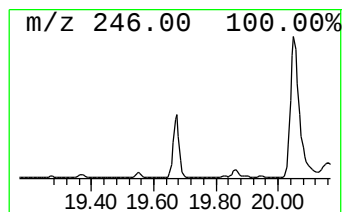
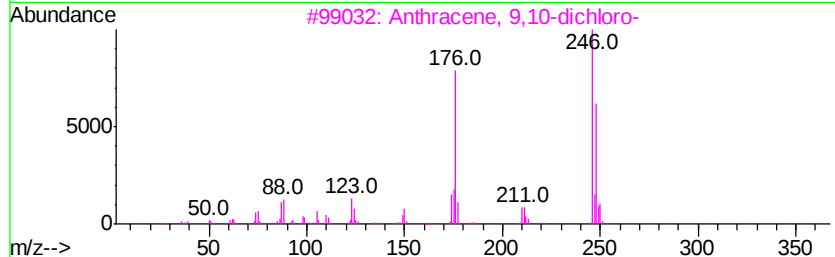
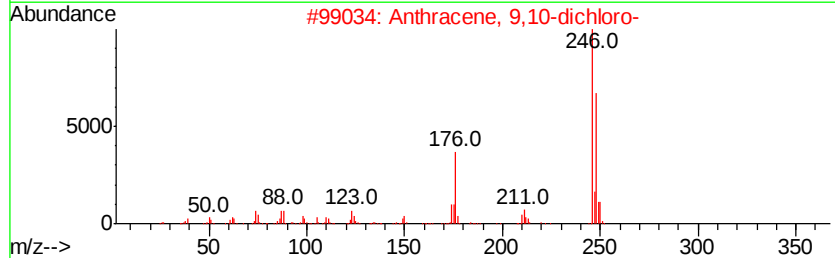
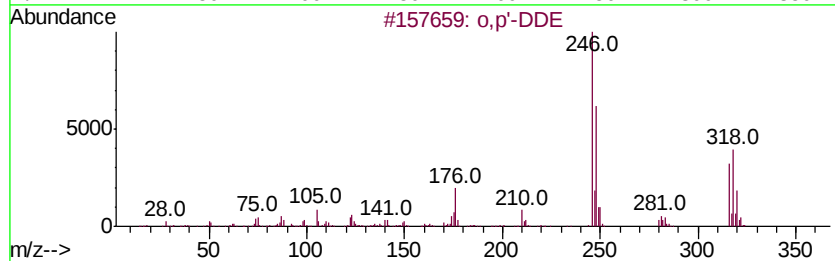
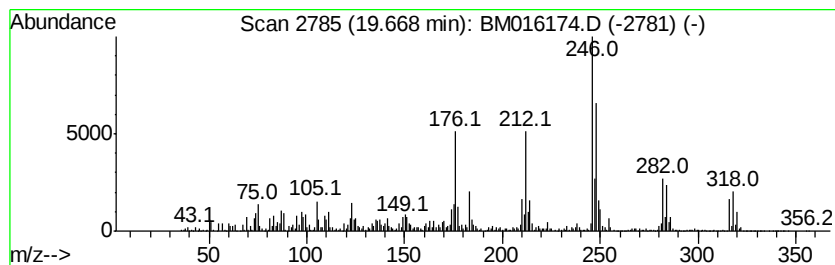
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ClientSampleId :  
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 12 o,p'-DDE Concentration Rank 6

| R.T.    | EstConc      | Area                       | Relative to ISTD | R.T.           |
|---------|--------------|----------------------------|------------------|----------------|
| 19.67   | 215.23 ng/ul | 31067000                   | Chrysene-d12     | 21.74          |
| Hit# of | 5            | Tentative ID               | MW MolForm       | CAS# Qual      |
| 1       |              | o,p'-DDE                   | 316 C14H8Cl4     | 003424-82-6 96 |
| 2       |              | Anthracene, 9,10-dichloro- | 246 C14H8Cl2     | 000605-48-1 96 |
| 3       |              | Anthracene, 9,10-dichloro- | 246 C14H8Cl2     | 000605-48-1 95 |
| 4       |              | p,p'-DDE                   | 316 C14H8Cl4     | 000072-55-9 93 |
| 5       |              | o,p'-DDE                   | 316 C14H8Cl4     | 003424-82-6 92 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080318\  
Data File : BM016174.D  
Acq On : 03 Aug 2018 11:39  
Operator : SJ/JU  
Sample : J4178-03  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM071318MA.M  
Quant Title : SVOA CALIBRATION

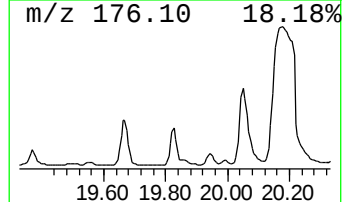
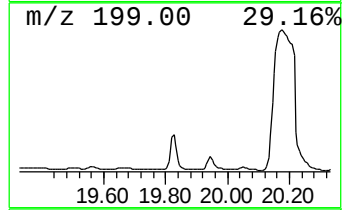
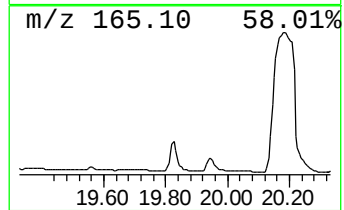
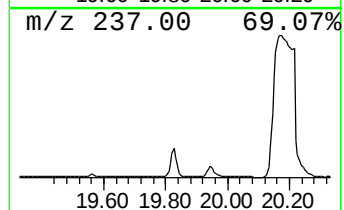
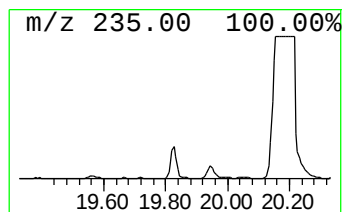
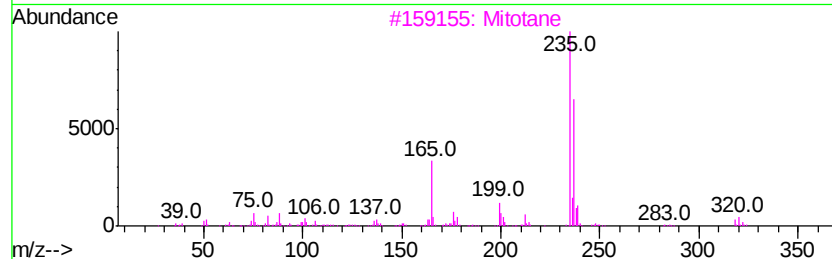
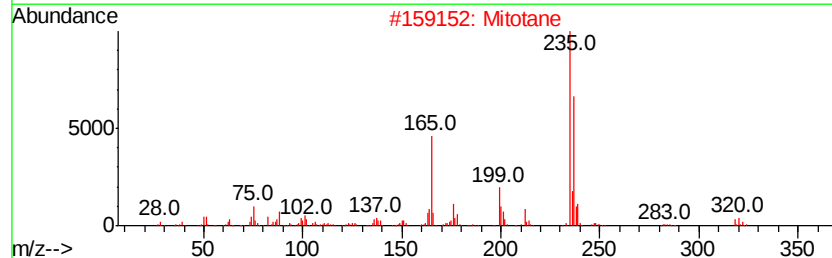
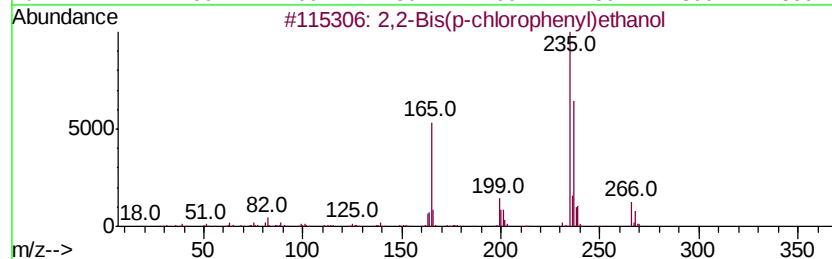
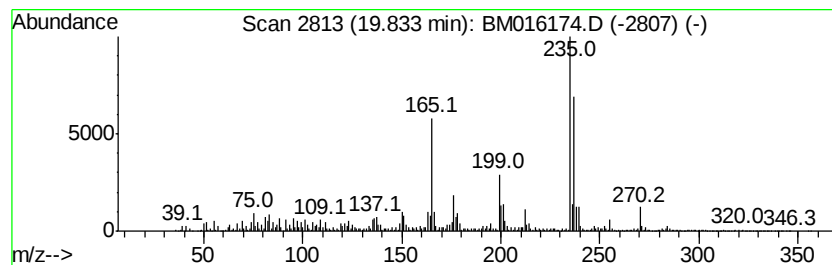
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 2,2-Bis(p-chlorophenyl)ethanol Concentration Rank 9

| R.T.  | EstConc      | Area     | Relative to ISTD | R.T.  |
|-------|--------------|----------|------------------|-------|
| 19.83 | 158.89 ng/ul | 22934400 | Chrysene-d12     | 21.74 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | 2,2-Bis(p-chlorophenyl)ethanol      | 266 | C14H12Cl2O | 002642-82-2 | 95   |
| 2    |    | Mitotane                            | 318 | C14H10Cl4  | 000053-19-0 | 91   |
| 3    |    | Mitotane                            | 318 | C14H10Cl4  | 000053-19-0 | 91   |
| 4    |    | m,p'-DDD                            | 318 | C14H10Cl4  | 004329-12-8 | 90   |
| 5    |    | 1,1-Dichloro-2,2-bis(p-chlorophe... | 318 | C14H10Cl4  | 000072-54-8 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080318\  
Data File : BM016174.D  
Acq On : 03 Aug 2018 11:39  
Operator : SJ/JU  
Sample : J4178-03  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Title : SVOA CALIBRATION

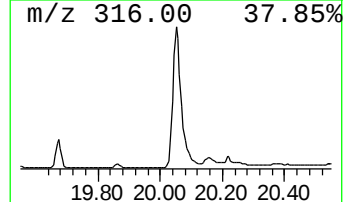
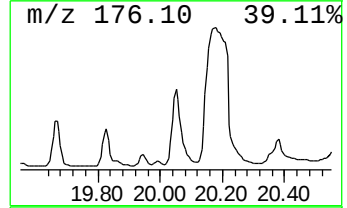
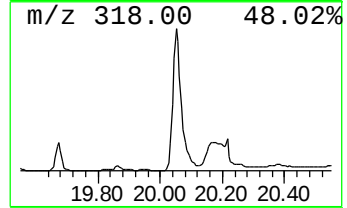
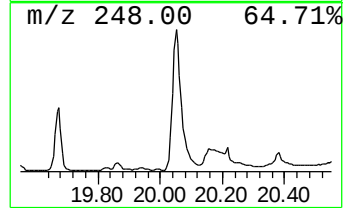
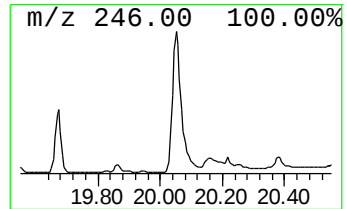
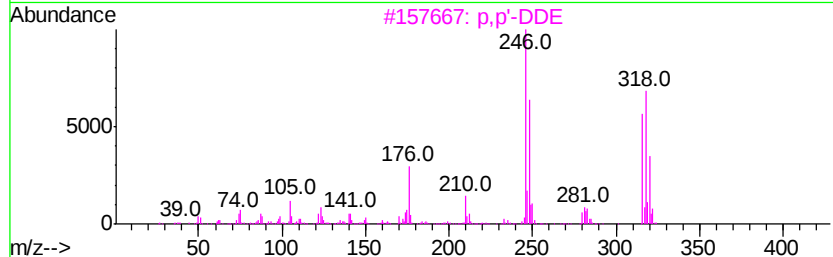
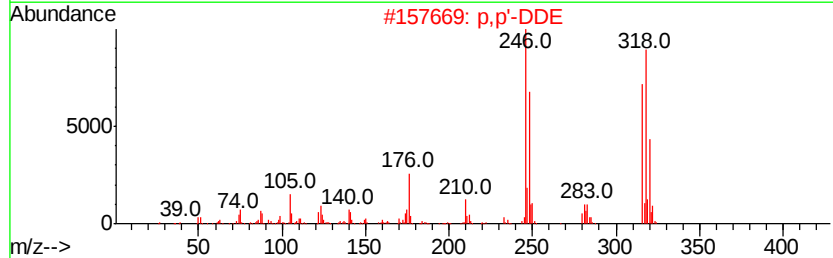
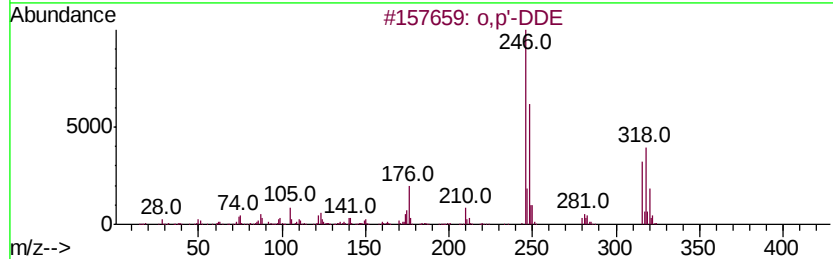
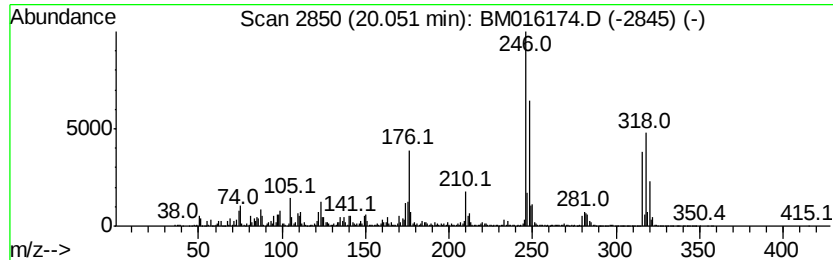
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TIC Integration Parameters: LSCINT.P

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Peak Number 14 p,p'-DDE Concentration Rank 5

| R.T.  | EstConc      | Area     | Relative to ISTD | R.T.  |
|-------|--------------|----------|------------------|-------|
| 20.05 | 253.07 ng/ul | 36529500 | Chrysene-d12     | 21.74 |

| Hit# | of | Tentative ID | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------|-----|----------|-------------|------|
| 1    | 5  | o,p'-DDE     | 316 | C14H8Cl4 | 003424-82-6 | 99   |
| 2    |    | p,p'-DDE     | 316 | C14H8Cl4 | 000072-55-9 | 99   |
| 3    |    | p,p'-DDE     | 316 | C14H8Cl4 | 000072-55-9 | 99   |
| 4    |    | p,p'-DDE     | 316 | C14H8Cl4 | 000072-55-9 | 99   |
| 5    |    | p,p'-DDE     | 316 | C14H8Cl4 | 000072-55-9 | 99   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080318\  
Data File : BM016174.D  
Acq On : 03 Aug 2018 11:39  
Operator : SJ/JU  
Sample : J4178-03  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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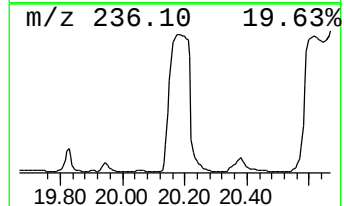
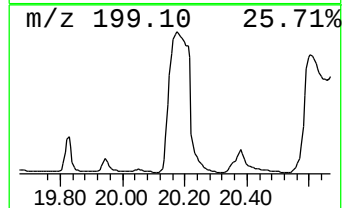
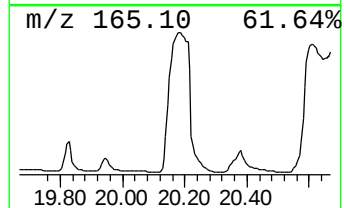
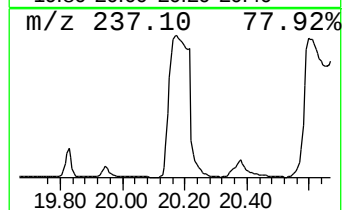
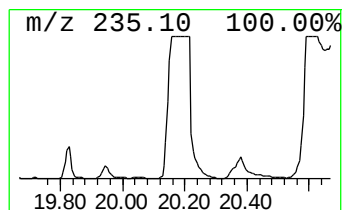
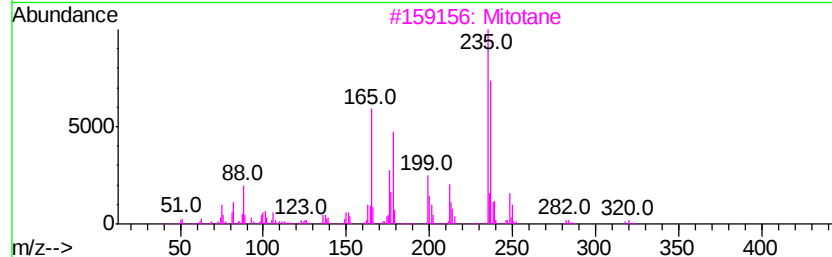
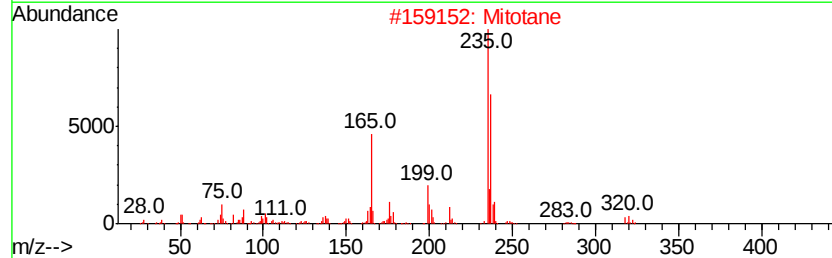
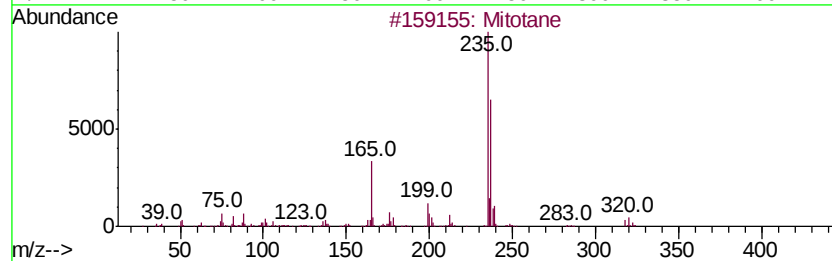
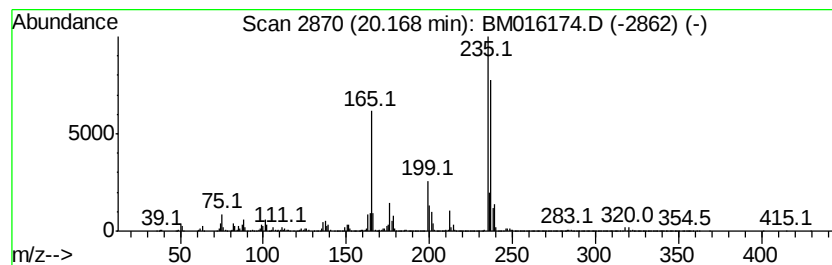
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Mitotane Concentration Rank 3

| R.T.  | EstConc       | Area      | Relative to ISTD | R.T.  |
|-------|---------------|-----------|------------------|-------|
| 20.17 | 1345.30 ng/ul | 194187000 | Chrysene-d12     | 21.74 |

| Hit# of | 5                                   | Tentative ID | MW  | MolForm    | CAS#         | Qual |
|---------|-------------------------------------|--------------|-----|------------|--------------|------|
| 1       | Mitotane                            |              | 318 | C14H10Cl4  | 000053-19-0  | 98   |
| 2       | Mitotane                            |              | 318 | C14H10Cl4  | 000053-19-0  | 95   |
| 3       | Mitotane                            |              | 318 | C14H10Cl4  | 000053-19-0  | 94   |
| 4       | Mitotane                            |              | 318 | C14H10Cl4  | 000053-19-0  | 93   |
| 5       | 3-Butanone, 1,1-bis(4-chlorophen... |              | 320 | C18H18Cl2O | 1000158-20-4 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080318\  
Data File : BM016174.D  
Acq On : 03 Aug 2018 11:39  
Operator : SJ/JU  
Sample : J4178-03  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampled :  
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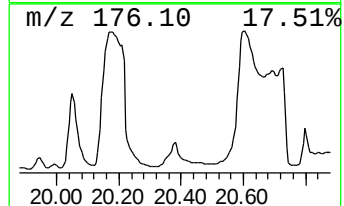
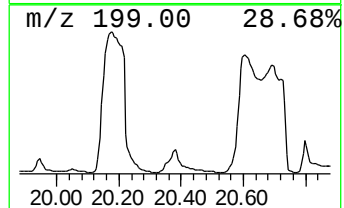
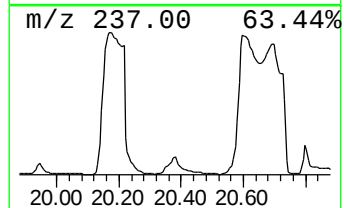
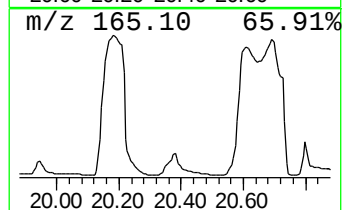
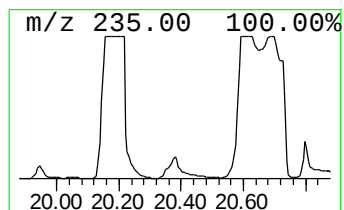
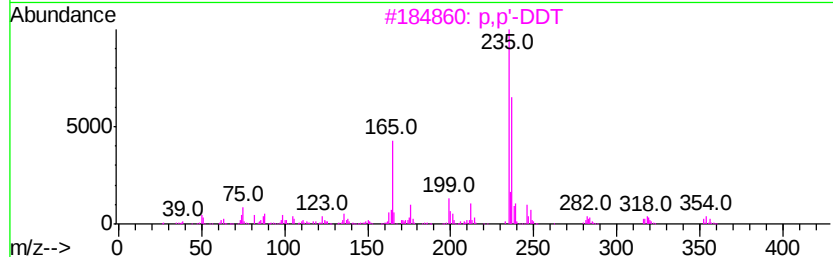
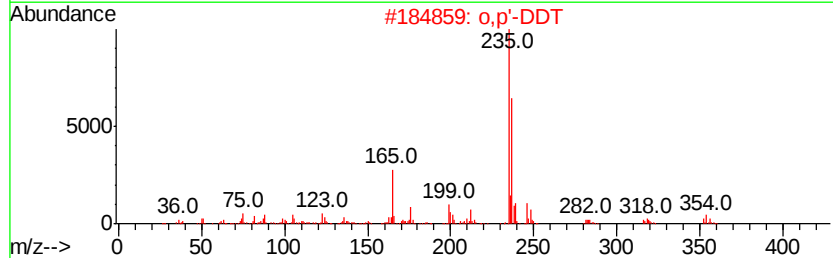
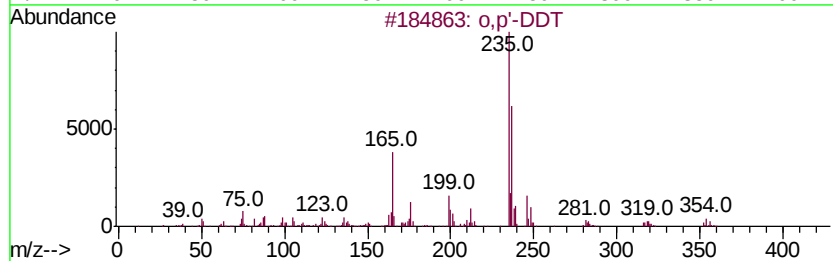
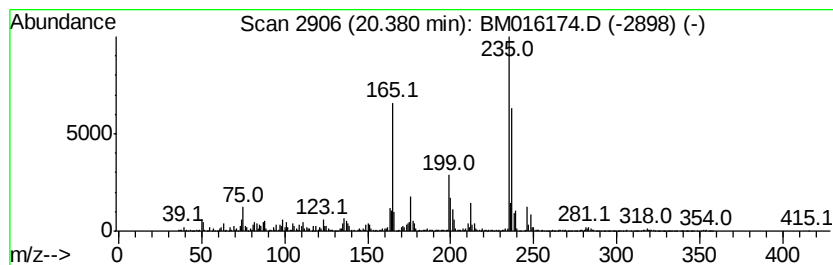
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TIC Integration Parameters: LSCINT.P

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Peak Number 16 o,p'-DDT Concentration Rank 4

| R.T.  | EstConc      | Area     | Relative to ISTD | R.T.  |
|-------|--------------|----------|------------------|-------|
| 20.38 | 260.48 ng/ul | 37598700 | Chrysene-d12     | 21.74 |

| Hit# | of | Tentative ID | MW  | MolForm   | CAS#        | Qual |
|------|----|--------------|-----|-----------|-------------|------|
| 1    | 5  | o,p'-DDT     | 352 | C14H9Cl5  | 000789-02-6 | 99   |
| 2    |    | o,p'-DDT     | 352 | C14H9Cl5  | 000789-02-6 | 97   |
| 3    |    | p,p'-DDT     | 352 | C14H9Cl5  | 000050-29-3 | 95   |
| 4    |    | Mitotane     | 318 | C14H10Cl4 | 000053-19-0 | 93   |
| 5    |    | o,p'-DDT     | 352 | C14H9Cl5  | 000789-02-6 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080318\  
Data File : BM016174.D  
Acq On : 03 Aug 2018 11:39  
Operator : SJ/JU  
Sample : J4178-03  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM071318MA.M  
Quant Title : SVOA CALIBRATION

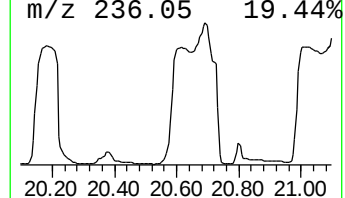
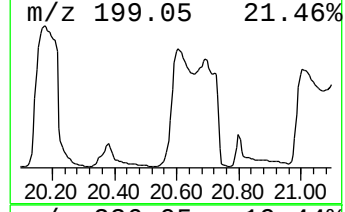
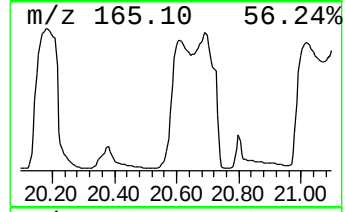
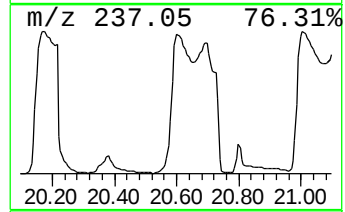
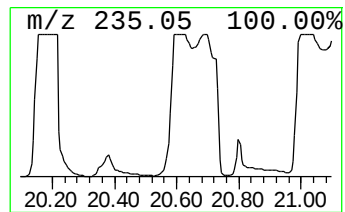
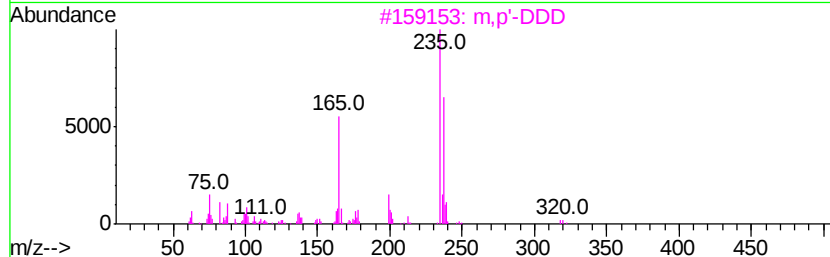
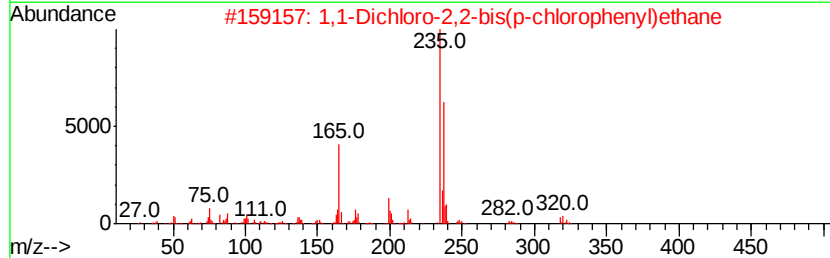
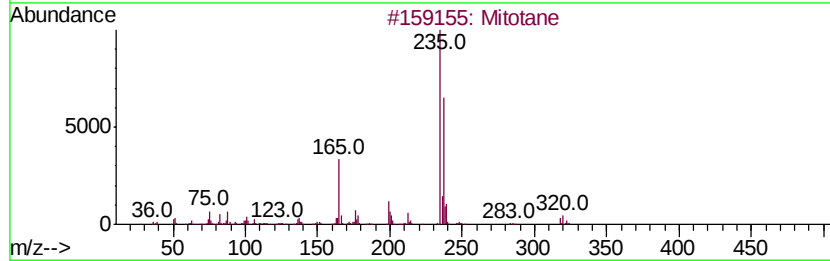
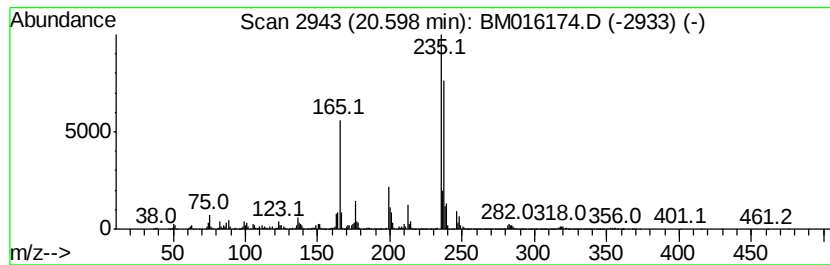
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TIC Integration Parameters: LSCINT.P

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Peak Number 17 1,1-Dichloro-2,2-bis(p-chlo... Concentration Rank 2

| R.T.  | EstConc       | Area      | Relative to ISTD | R.T.  |
|-------|---------------|-----------|------------------|-------|
| 20.60 | 1452.47 ng/ul | 209656000 | Chrysene-d12     | 21.74 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Mitotane                            | 318 | C14H10Cl4 | 000053-19-0 | 95   |
| 2    |    | 1,1-Dichloro-2,2-bis(p-chlorophe... | 318 | C14H10Cl4 | 000072-54-8 | 95   |
| 3    |    | m,p'-DDD                            | 318 | C14H10Cl4 | 004329-12-8 | 94   |
| 4    |    | o,p'-DDT                            | 352 | C14H9Cl5  | 000789-02-6 | 93   |
| 5    |    | p,p'-DDT                            | 352 | C14H9Cl5  | 000050-29-3 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080318\  
Data File : BM016174.D  
Acq On : 03 Aug 2018 11:39  
Operator : SJ/JU  
Sample : J4178-03  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM071318MA.M  
Quant Title : SVOA CALIBRATION

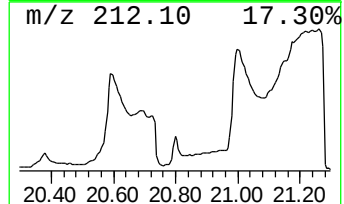
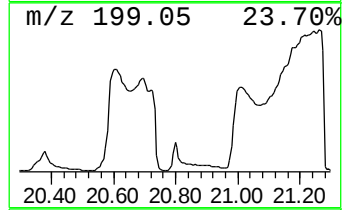
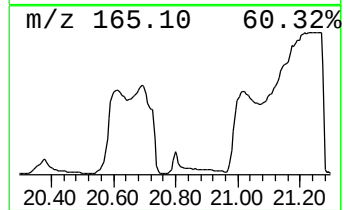
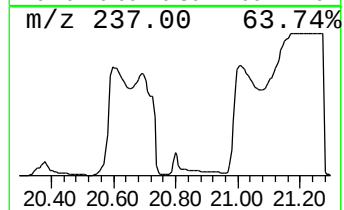
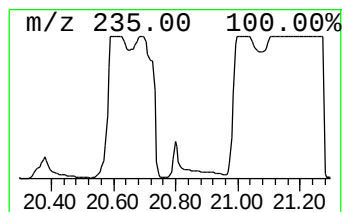
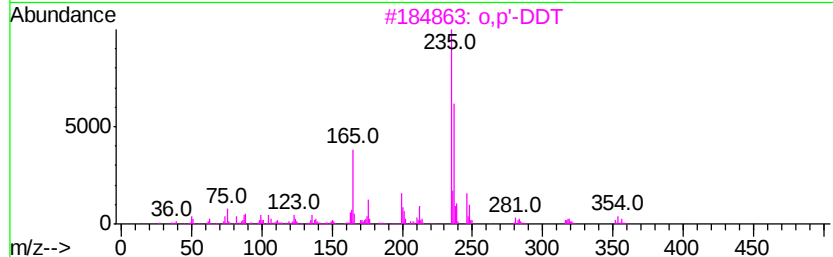
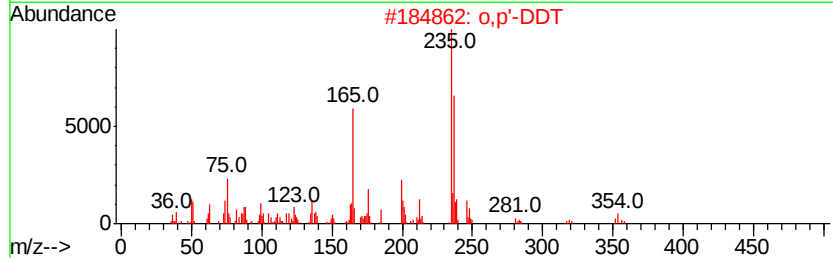
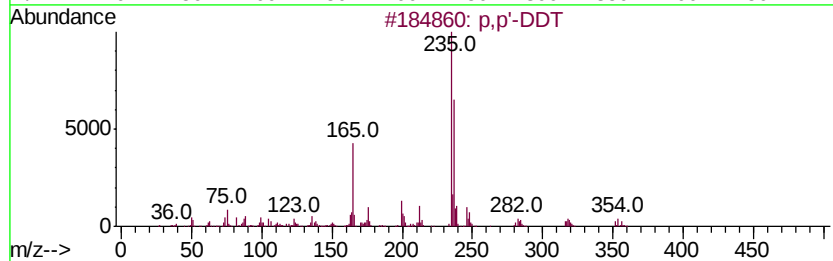
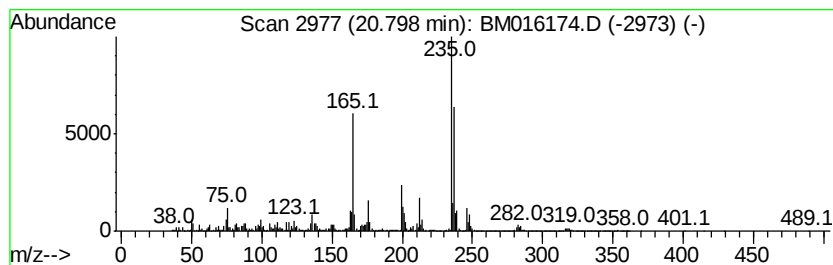
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 p,p'-DDT Concentration Rank 7

| R.T.  | EstConc      | Area     | Relative to ISTD | R.T.  |
|-------|--------------|----------|------------------|-------|
| 20.80 | 206.65 ng/ul | 29829300 | Chrysene-d12     | 21.74 |

| Hit# | of | Tentative ID | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------|-----|----------|-------------|------|
| 1    | 5  | p,p'-DDT     | 352 | C14H9Cl5 | 000050-29-3 | 99   |
| 2    |    | o,p'-DDT     | 352 | C14H9Cl5 | 000789-02-6 | 99   |
| 3    |    | o,p'-DDT     | 352 | C14H9Cl5 | 000789-02-6 | 98   |
| 4    |    | p,p'-DDT     | 352 | C14H9Cl5 | 000050-29-3 | 97   |
| 5    |    | p,p'-DDT     | 352 | C14H9Cl5 | 000050-29-3 | 97   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080318\  
Data File : BM016174.D  
Acq On : 03 Aug 2018 11:39  
Operator : SJ/JU  
Sample : J4178-03  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0AB9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM071318MA.M  
Quant Title : SVOA CALIBRATION

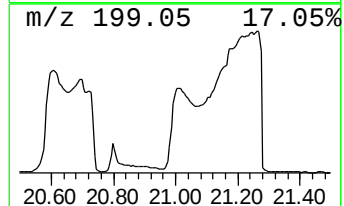
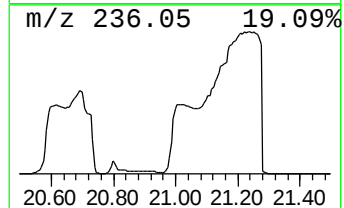
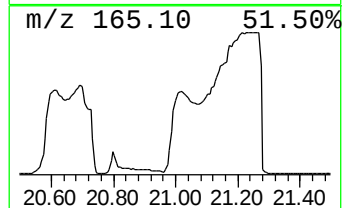
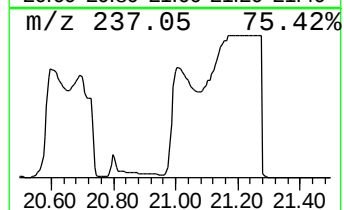
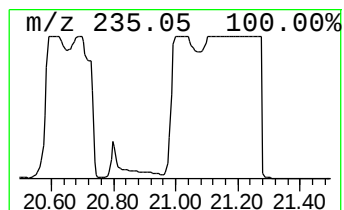
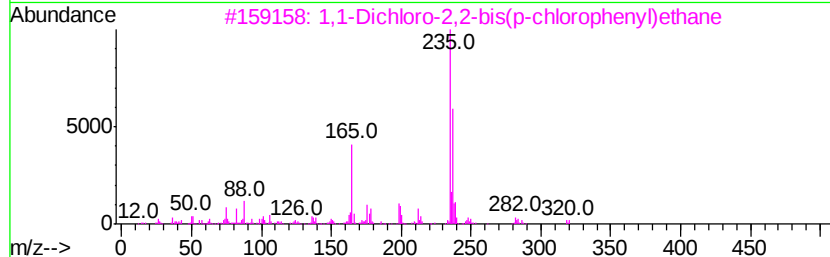
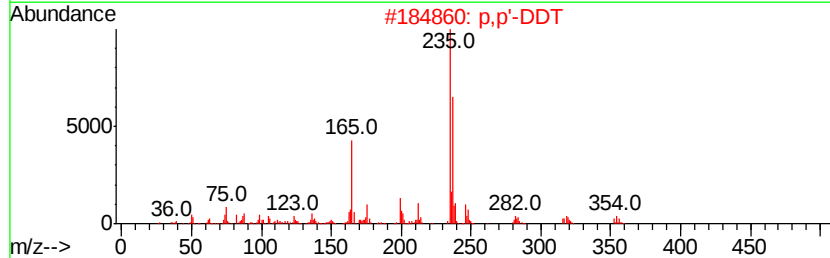
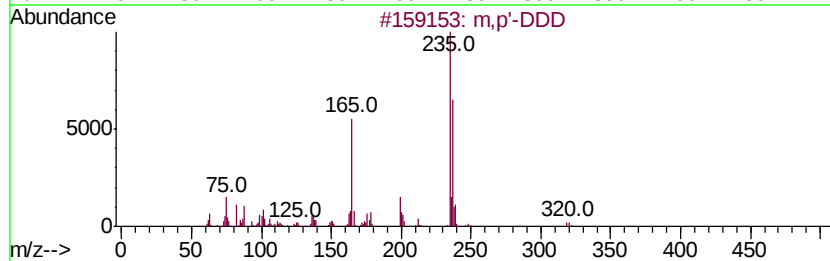
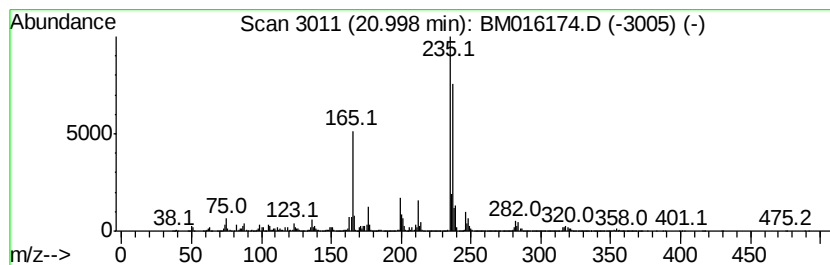
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 m,p'-DDD Concentration Rank 1

| R.T.  | EstConc       | Area      | Relative to ISTD | R.T.  |
|-------|---------------|-----------|------------------|-------|
| 21.00 | 1743.36 ng/ul | 251644000 | Chrysene-d12     | 21.74 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | m,p'-DDD                            | 318 | C14H10Cl4 | 004329-12-8 | 98   |
| 2    |    | p,p'-DDT                            | 352 | C14H9Cl5  | 000050-29-3 | 97   |
| 3    |    | 1,1-Dichloro-2,2-bis(p-chlorophe... | 318 | C14H10Cl4 | 000072-54-8 | 91   |
| 4    |    | 1,1-Dichloro-2,2-bis(p-chlorophe... | 318 | C14H10Cl4 | 000072-54-8 | 90   |
| 5    |    | p,p'-DDT                            | 352 | C14H9Cl5  | 000050-29-3 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080318\  
Data File : BM016174.D  
Acq On : 03 Aug 2018 11:39  
Operator : SJ/JU  
Sample : J4178-03  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0AB9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM071318MA.M  
Quant Title : SVOA CALIBRATION

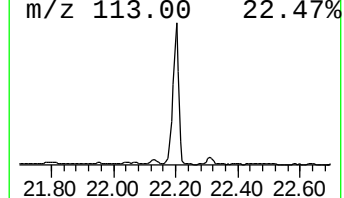
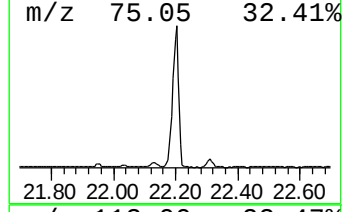
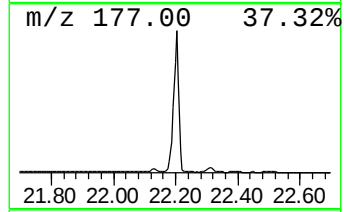
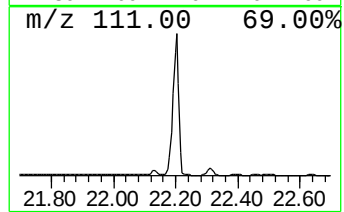
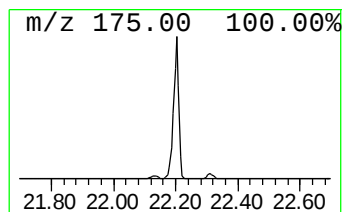
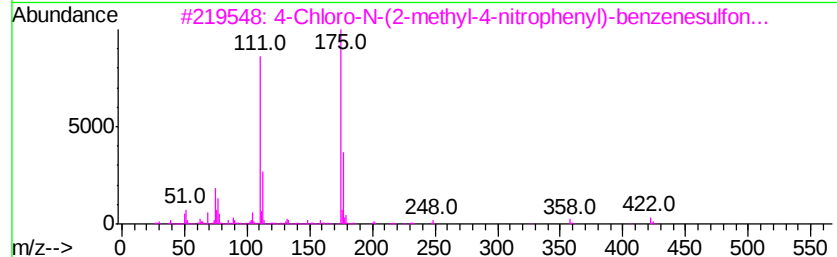
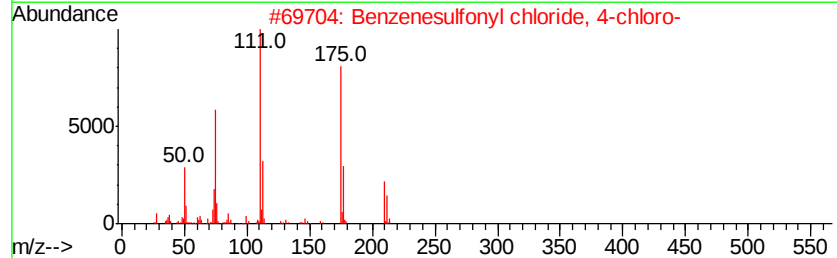
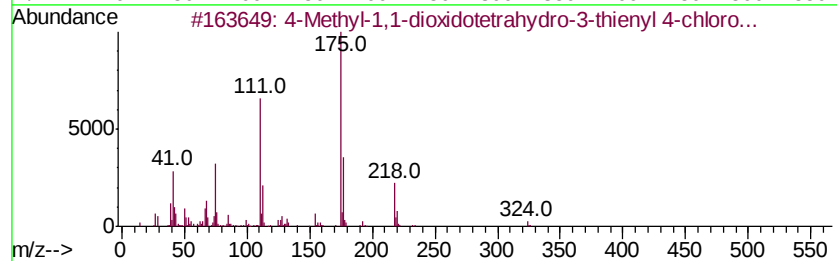
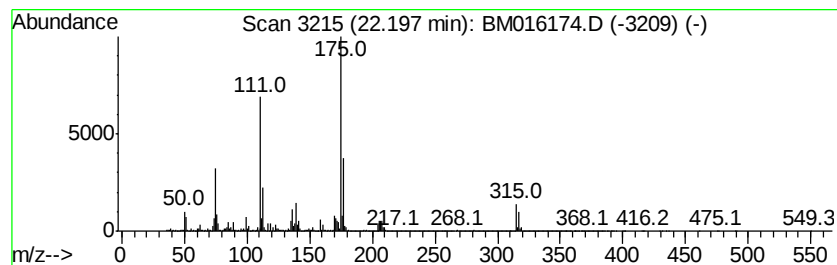
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 4-Methyl-1,1-dioxidotetrahy... Concentration Rank 8

| R.T.  | EstConc      | Area     | Relative to ISTD | R.T.  |
|-------|--------------|----------|------------------|-------|
| 22.20 | 193.15 ng/ul | 27880700 | Chrysene-d12     | 21.74 |

| Hit# | of | Tentative ID                        | MW  | MolForm         | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------------|--------------|------|
| 1    | 5  | 4-Methyl-1,1-dioxidotetrahydro-3... | 324 | C11H13ClO5S2    | 1000332-40-7 | 72   |
| 2    |    | Benzenesulfonyl chloride, 4-chloro- | 210 | C6H4Cl2O2S      | 000098-60-2  | 64   |
| 3    |    | 4-Chloro-N-(2-methyl-4-nitrophen... | 422 | C15H10ClF3N2O5S | 1000374-78-8 | 64   |
| 4    |    | Benzenemethanol, 2,4-dichloro-.a... | 224 | C8H7Cl3O        | 013692-14-3  | 53   |
| 5    |    | 4-Chlorobenzenesulfonyl isocyanate  | 217 | C7H4ClN3O3S     | 005769-15-3  | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM080318\  
 Data File : BM016174.D  
 Acq On : 03 Aug 2018 11:39  
 Operator : SJ/JU  
 Sample : J4178-03  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C0AB9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM071318MA.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 |
| Benzene, 1,4-dich... | 8.07  | 16.5    | ng/ul | 6158840   | 1                     | 8.19  | 7479360 | 20.0 |
| Benzene, 1,3-dich... | 8.56  | 72.1    | ng/ul | 26969100  | 1                     | 8.19  | 7479360 | 20.0 |
| Benzene, 1,2,3-tr... | 10.87 | 20.0    | ng/ul | 4601160   | 2                     | 11.01 | 4601160 | 20.0 |
| n-Octylidencycloh... | 14.63 | 34.9    | ng/ul | 4977020   | 3                     | 14.82 | 2854840 | 20.0 |
| 3-Fluorobenzoic a... | 14.76 | 20.0    | ng/ul | 2854840   | 3                     | 14.82 | 2854840 | 20.0 |
| 1H-Indene, octahy... | 15.52 | 20.1    | ng/ul | 2872920   | 3                     | 14.82 | 2854840 | 20.0 |
| unknown-01           | 15.59 | 70.4    | ng/ul | 10049100  | 3                     | 14.82 | 2854840 | 20.0 |
| Naphthalene, 1,6-... | 16.53 | 88.9    | ng/ul | 28598400  | 4                     | 17.57 | 6433870 | 20.0 |
| unknown-02           | 16.63 | 19.4    | ng/ul | 6247630   | 4                     | 17.57 | 6433870 | 20.0 |
| unknown-03           | 16.75 | 20.0    | ng/ul | 6433870   | 4                     | 17.57 | 6433870 | 20.0 |
| Estra-4,9,11-trie... | 19.06 | 49.8    | ng/ul | 16019700  | 4                     | 17.57 | 6433870 | 20.0 |
| o,p'-DDE             | 19.67 | 215.2   | ng/ul | 31067000  | 5                     | 21.74 | 2886890 | 20.0 |
| 2,2-Bis(p-chlorop... | 19.83 | 158.9   | ng/ul | 22934400  | 5                     | 21.74 | 2886890 | 20.0 |
| p,p'-DDE             | 20.05 | 253.1   | ng/ul | 36529500  | 5                     | 21.74 | 2886890 | 20.0 |
| Mitotane             | 20.17 | 1345.3  | ng/ul | 194187000 | 5                     | 21.74 | 2886890 | 20.0 |
| o,p'-DDT             | 20.38 | 260.5   | ng/ul | 37598700  | 5                     | 21.74 | 2886890 | 20.0 |
| 1,1-Dichloro-2,2-... | 20.60 | 1452.5  | ng/ul | 209656000 | 5                     | 21.74 | 2886890 | 20.0 |
| p,p'-DDT             | 20.80 | 206.7   | ng/ul | 29829300  | 5                     | 21.74 | 2886890 | 20.0 |
| m,p'-DDD             | 21.00 | 1743.4  | ng/ul | 251644000 | 5                     | 21.74 | 2886890 | 20.0 |
| 4-Methyl-1,1-diox... | 22.20 | 193.2   | ng/ul | 27880700  | 5                     | 21.74 | 2886890 | 20.0 |