

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\  
 Data File : BM019372.D  
 Acq On : 24 Mar 2019 03:11  
 Operator : JU/SJ  
 Sample : K2025-10  
 Misc :  
 ALS Vial : 30 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C0967

## Integration Parameters: LSCINT.P

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

Filtering: 5  
 Min Area: 0.1 % 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-BM032219MA.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         | 4.957       | 327           | 335         | 353          | rBV      | 21736708       | 31124017      | 37.87%          | 11.142%       |
| 2         | 6.793       | 640           | 647         | 660          | rBV2     | 131673         | 289820        | 0.35%           | 0.104%        |
| 3         | 6.963       | 670           | 676         | 691          | rBV      | 479299         | 801344        | 0.97%           | 0.287%        |
| 4         | 7.145       | 700           | 707         | 723          | rBV2     | 537570         | 1014973       | 1.23%           | 0.363%        |
| 5         | 7.498       | 759           | 767         | 780          | rBV      | 10793709       | 16588891      | 20.18%          | 5.939%        |
| 6         | 7.663       | 780           | 795         | 801          | rBV      | 38429350       | 72500969      | 88.20%          | 25.954%       |
| 7         | 7.981       | 837           | 849         | 854          | rBV      | 39079440       | 82196970      | 100.00%         | 29.425%       |
| 8         | 8.322       | 901           | 907         | 921          | rBV      | 341916         | 619264        | 0.75%           | 0.222%        |
| 9         | 8.781       | 976           | 985         | 987          | rBV      | 586799         | 927031        | 1.13%           | 0.332%        |
| 10        | 8.828       | 987           | 993         | 1010         | rVB      | 3717320        | 6231720       | 7.58%           | 2.231%        |
| 11        | 9.104       | 1031          | 1040        | 1050         | rVB2     | 95746          | 181875        | 0.22%           | 0.065%        |
| 12        | 9.492       | 1100          | 1106        | 1123         | rVB      | 464712         | 831617        | 1.01%           | 0.298%        |
| 13        | 10.022      | 1189          | 1196        | 1204         | rBV      | 706227         | 1345749       | 1.64%           | 0.482%        |
| 14        | 10.086      | 1204          | 1207        | 1217         | rVV      | 172373         | 327695        | 0.40%           | 0.117%        |
| 15        | 10.181      | 1217          | 1223        | 1228         | rVV      | 155835         | 281852        | 0.34%           | 0.101%        |
| 16        | 10.257      | 1228          | 1236        | 1244         | rVV      | 10438478       | 16823444      | 20.47%          | 6.023%        |
| 17        | 10.333      | 1244          | 1249        | 1253         | rVV3     | 55205          | 140093        | 0.17%           | 0.050%        |
| 18        | 10.392      | 1253          | 1259        | 1268         | rVB      | 945197         | 1520726       | 1.85%           | 0.544%        |
| 19        | 10.775      | 1316          | 1324        | 1335         | rBV      | 4598916        | 7476601       | 9.10%           | 2.676%        |
| 20        | 11.010      | 1357          | 1364        | 1387         | rVB      | 1530847        | 2670569       | 3.25%           | 0.956%        |
| 21        | 11.228      | 1393          | 1401        | 1420         | rBV2     | 294009         | 653985        | 0.80%           | 0.234%        |
| 22        | 12.439      | 1596          | 1607        | 1612         | rBV      | 82247          | 151834        | 0.18%           | 0.054%        |
| 23        | 13.057      | 1705          | 1712        | 1729         | rBV3     | 334399         | 990883        | 1.21%           | 0.355%        |
| 24        | 13.298      | 1747          | 1753        | 1769         | rVB2     | 363915         | 745664        | 0.91%           | 0.267%        |
| 25        | 13.686      | 1811          | 1819        | 1825         | rBV      | 1443157        | 2035343       | 2.48%           | 0.729%        |
| 26        | 13.745      | 1825          | 1829        | 1841         | rVB4     | 55492          | 99548         | 0.12%           | 0.036%        |
| 27        | 13.957      | 1858          | 1865        | 1875         | rBV      | 1643701        | 2356172       | 2.87%           | 0.843%        |
| 28        | 14.263      | 1910          | 1917        | 1925         | rBV2     | 1274613        | 1972527       | 2.40%           | 0.706%        |
| 29        | 14.533      | 1957          | 1963        | 1969         | rBV3     | 38378          | 92246         | 0.11%           | 0.033%        |
| 30        | 15.268      | 2081          | 2088        | 2100         | rBV      | 2044538        | 2955683       | 3.60%           | 1.058%        |
| 31        | 15.404      | 2106          | 2111        | 2125         | rBV      | 591206         | 924982        | 1.13%           | 0.331%        |
| 32        | 17.021      | 2380          | 2386        | 2397         | rBV2     | 1621562        | 2280197       | 2.77%           | 0.816%        |
| 33        | 17.121      | 2397          | 2403        | 2422         | rBV2     | 2197600        | 3253858       | 3.96%           | 1.165%        |
| 34        | 17.915      | 2531          | 2538        | 2550         | rBV      | 180829         | 299609        | 0.36%           | 0.107%        |

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Instrument :  
BNA\_M  
ClientSampleId :  
C0967

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

Filtering: 5  
Min Area: 0.1 % 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-BM032219MA.M  
Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |       |        |
|----|--------|------|------|------|------|---------|---------|-------|--------|
| 35 | 19.209 | 2753 | 2758 | 2770 | rBV  | 368795  | 590456  | 0.72% | 0.211% |
| 36 | 19.427 | 2789 | 2795 | 2807 | rBV2 | 2796674 | 3889260 | 4.73% | 1.392% |
| 37 | 20.927 | 3047 | 3050 | 3055 | rVB  | 94625   | 97808   | 0.12% | 0.035% |
| 38 | 21.227 | 3096 | 3101 | 3114 | rBV  | 2265454 | 2915870 | 3.55% | 1.044% |
| 39 | 21.362 | 3121 | 3124 | 3129 | rBV  | 120107  | 132062  | 0.16% | 0.047% |
| 40 | 21.792 | 3194 | 3197 | 3203 | rBV  | 144860  | 161711  | 0.20% | 0.058% |
| 41 | 22.244 | 3271 | 3274 | 3279 | rBV  | 190738  | 228075  | 0.28% | 0.082% |
| 42 | 22.739 | 3355 | 3358 | 3363 | rVB  | 200473  | 280705  | 0.34% | 0.100% |
| 43 | 23.303 | 3447 | 3454 | 3469 | rVB2 | 2577609 | 4669363 | 5.68% | 1.672% |
| 44 | 23.444 | 3471 | 3478 | 3488 | rVV  | 1591923 | 3018648 | 3.67% | 1.081% |
| 45 | 23.927 | 3555 | 3560 | 3567 | rVB2 | 181811  | 325607  | 0.40% | 0.117% |
| 46 | 24.656 | 3680 | 3684 | 3693 | rVB  | 162894  | 325708  | 0.40% | 0.117% |

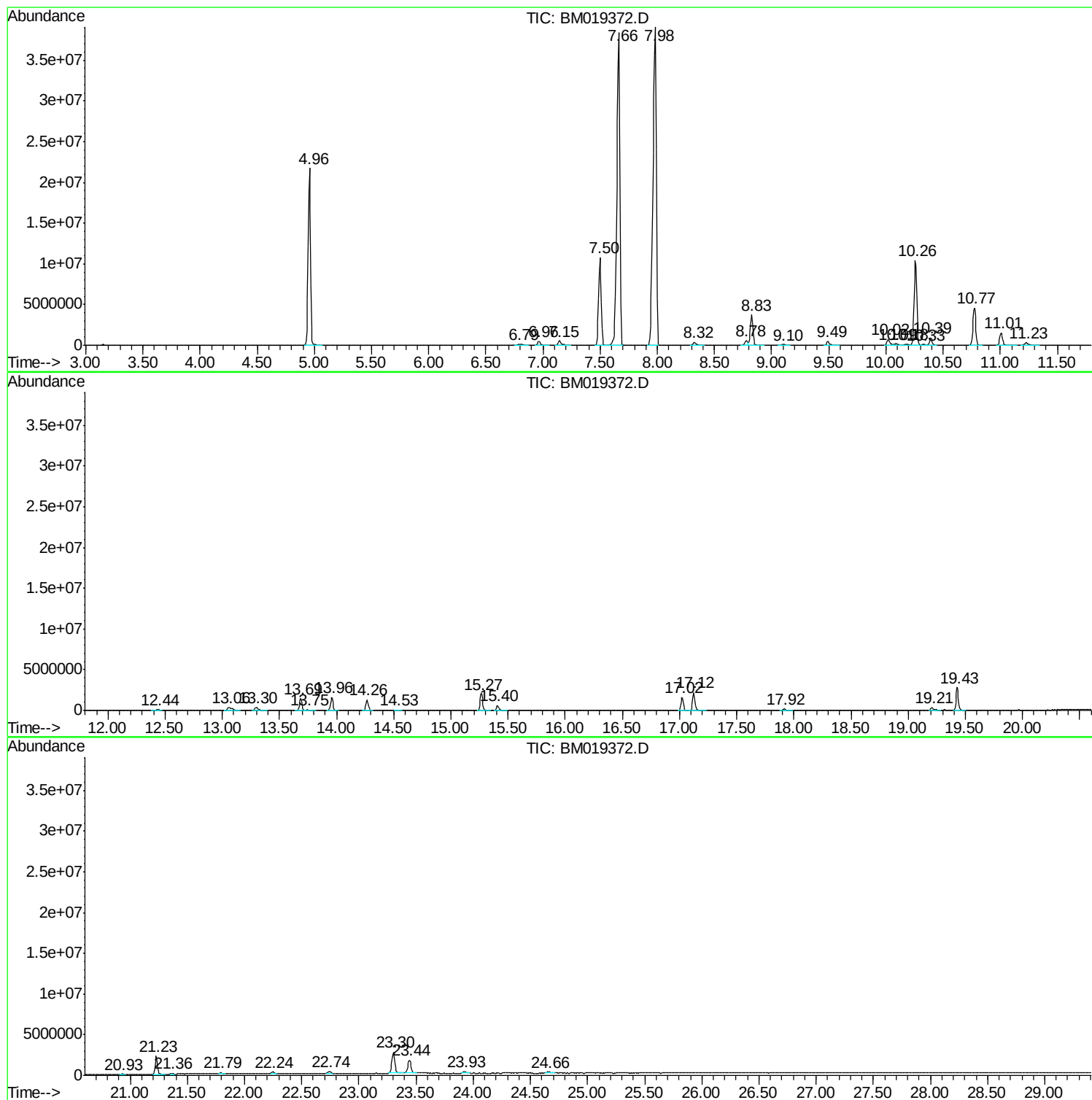
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

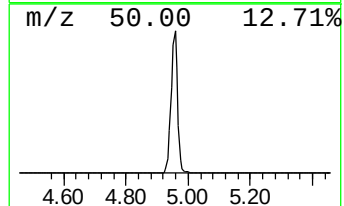
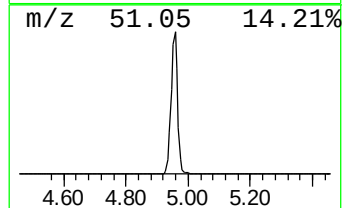
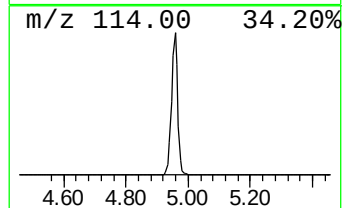
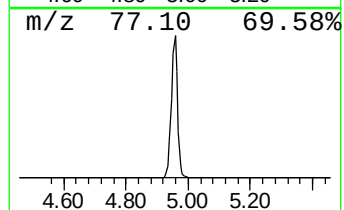
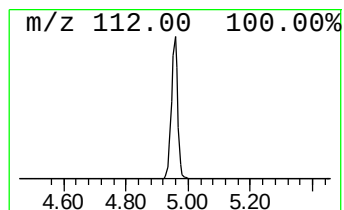
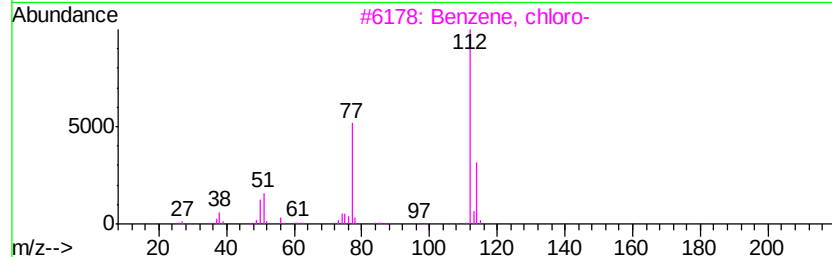
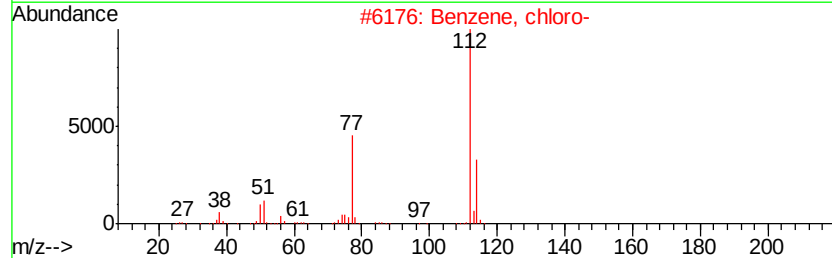
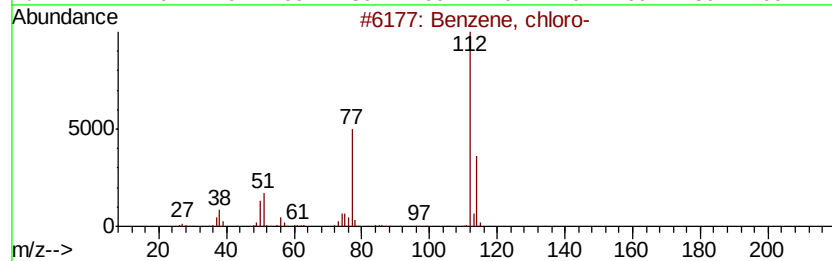
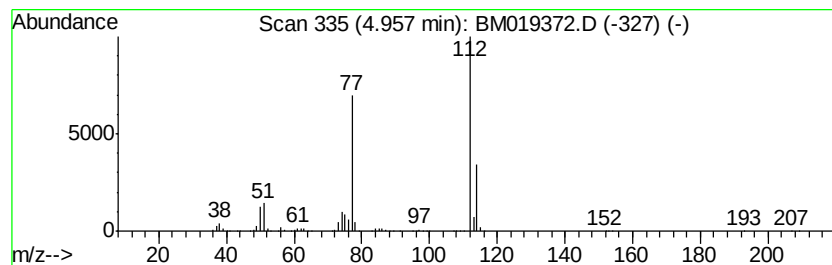
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Benzene, chloro- Concentration Rank 6

| R.T. | EstConc    | Area     | Relative to ISTD       | R.T. |
|------|------------|----------|------------------------|------|
| 4.96 | 8.59 ng/ul | 31124000 | 1,4-Dichlorobenzene-d4 | 7.62 |

| Hit# | of | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, chloro-  | 112 | C6H5Cl  | 000108-90-7 | 95   |
| 2    |    | Benzene, chloro-  | 112 | C6H5Cl  | 000108-90-7 | 95   |
| 3    |    | Benzene, chloro-  | 112 | C6H5Cl  | 000108-90-7 | 95   |
| 4    |    | Benzene, chloro-  | 112 | C6H5Cl  | 000108-90-7 | 94   |
| 5    |    | Phenol, 4-fluoro- | 112 | C6H5FO  | 000371-41-5 | 38   |



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

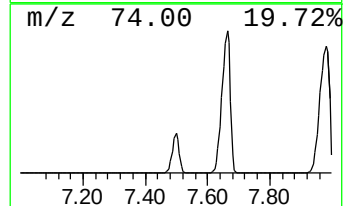
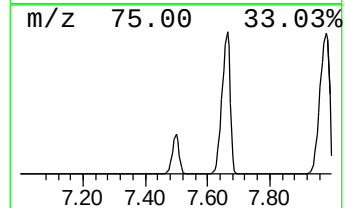
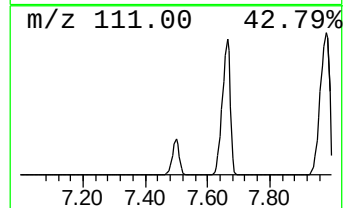
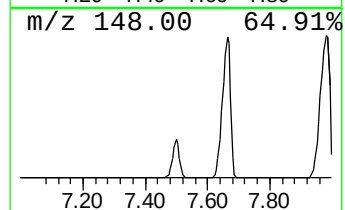
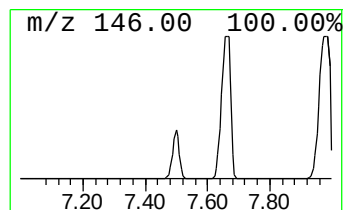
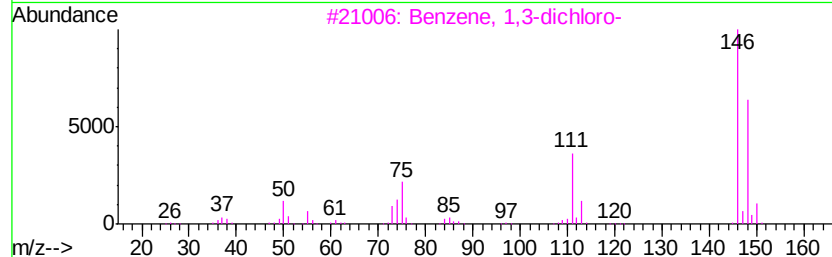
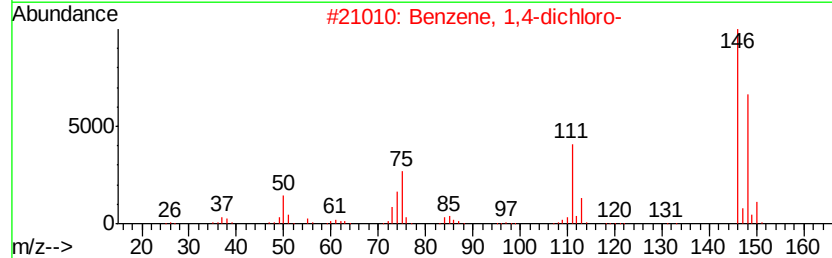
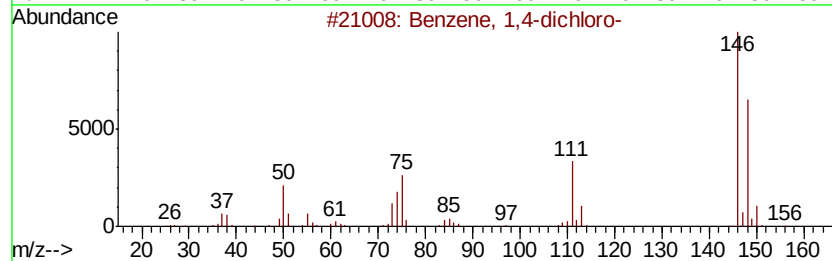
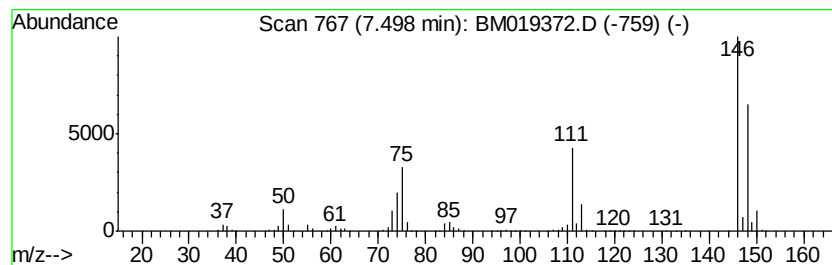
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Benzene, 1,4-dichloro- Concentration Rank 8

| R.T. | EstConc    | Area     | Relative to ISTD       | R.T. |
|------|------------|----------|------------------------|------|
| 7.50 | 4.58 ng/ul | 16588900 | 1,4-Dichlorobenzene-d4 | 7.62 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,4-dichloro- | 146 | C6H4Cl2 | 000106-46-7 | 96   |
| 2    |    | Benzene, 1,4-dichloro- | 146 | C6H4Cl2 | 000106-46-7 | 96   |
| 3    |    | Benzene, 1,3-dichloro- | 146 | C6H4Cl2 | 000541-73-1 | 95   |
| 4    |    | Benzene, 1,3-dichloro- | 146 | C6H4Cl2 | 000541-73-1 | 95   |
| 5    |    | Benzene, 1,2-dichloro- | 146 | C6H4Cl2 | 000095-50-1 | 95   |



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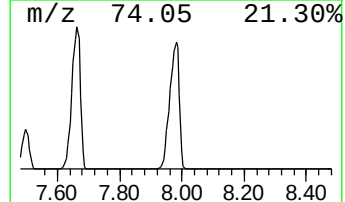
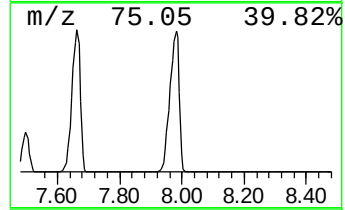
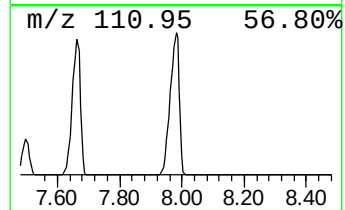
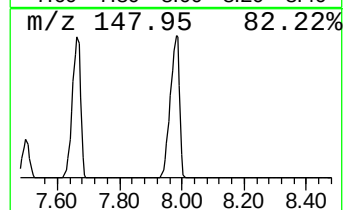
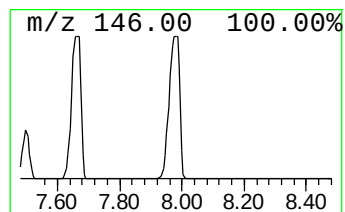
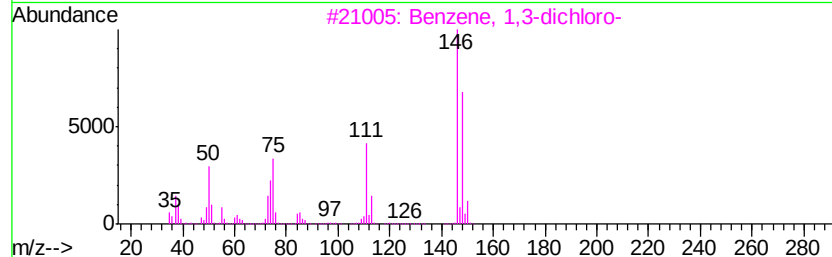
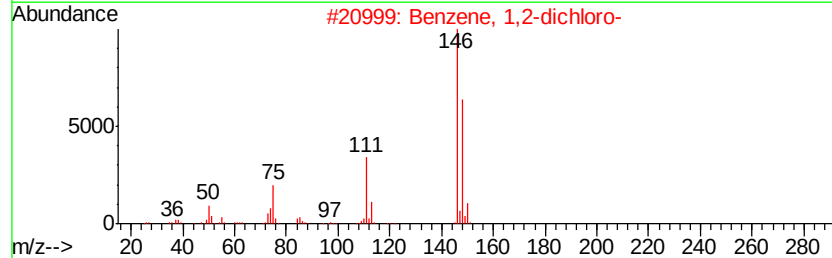
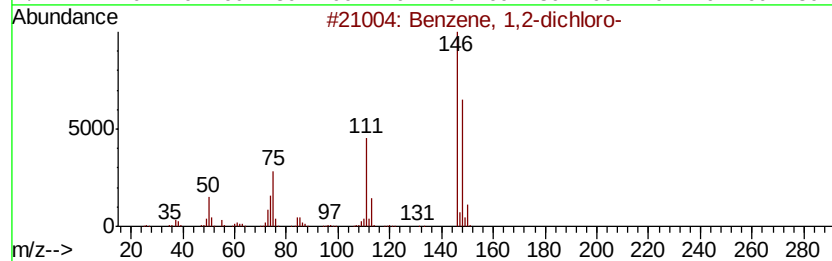
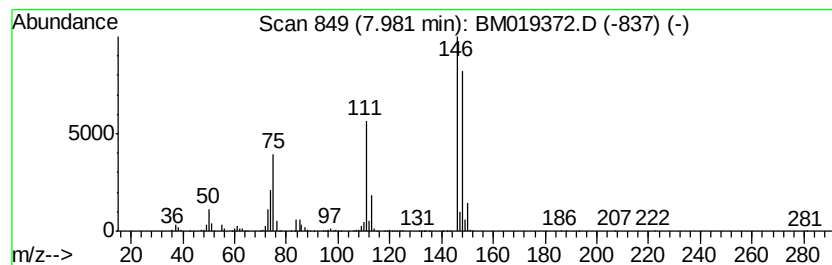
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 Benzene, 1,2-dichloro- Concentration Rank 4

| R.T. | EstConc     | Area     | Relative to ISTD       | R.T. |
|------|-------------|----------|------------------------|------|
| 7.98 | 22.67 ng/ul | 82197000 | 1,4-Dichlorobenzene-d4 | 7.62 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2-dichloro- | 146 | C6H4Cl2 | 000095-50-1 | 95   |
| 2    |    | Benzene, 1,2-dichloro- | 146 | C6H4Cl2 | 000095-50-1 | 94   |
| 3    |    | Benzene, 1,3-dichloro- | 146 | C6H4Cl2 | 000541-73-1 | 94   |
| 4    |    | Benzene, 1,3-dichloro- | 146 | C6H4Cl2 | 000541-73-1 | 94   |
| 5    |    | Benzene, 1,4-dichloro- | 146 | C6H4Cl2 | 000106-46-7 | 94   |



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Quant Title : SVOA CALIBRATION

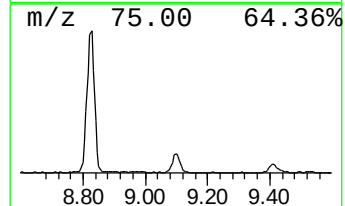
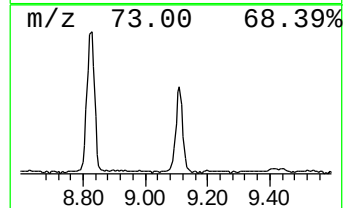
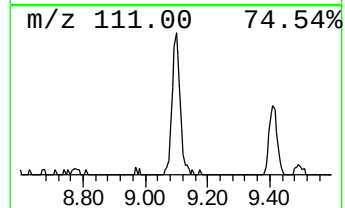
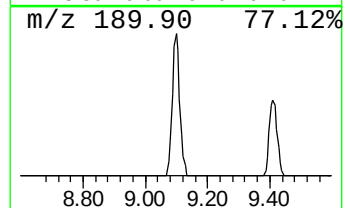
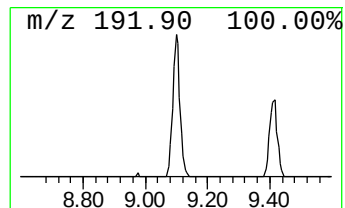
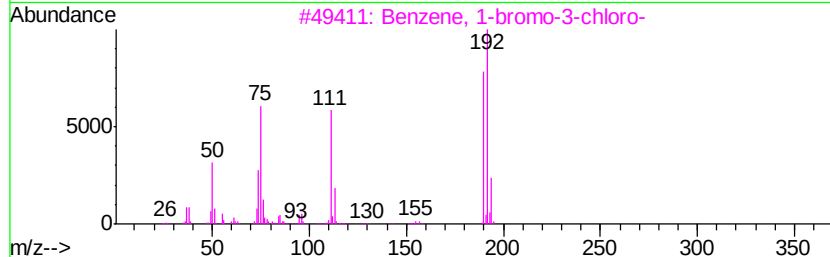
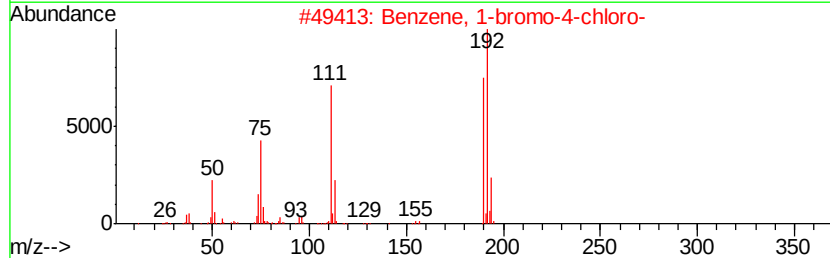
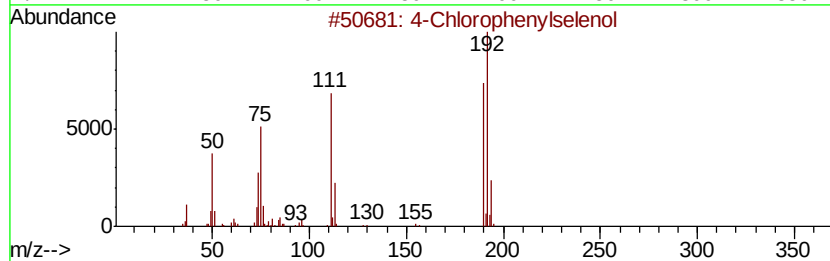
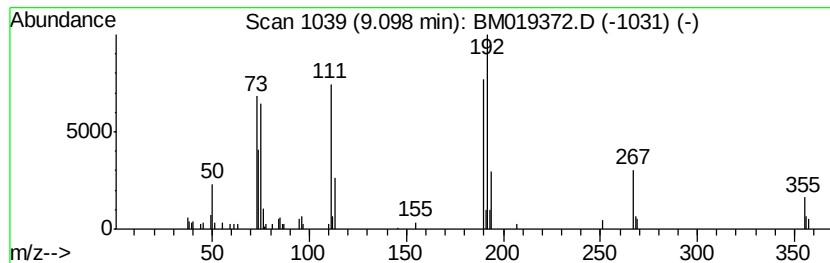
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 4-Chlorophenylselenol Concentration Rank 12

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.10 | 2.39 ng/ul | 181875 | Napthalene-d8    | 10.39 |

| Hit# | of | Tentative ID               | MW  | MolForm  | CAS#        | Qual |
|------|----|----------------------------|-----|----------|-------------|------|
| 1    | 5  | 4-Chlorophenylselenol      | 192 | C6H5ClSe | 016645-10-6 | 86   |
| 2    |    | Benzene, 1-bromo-4-chloro- | 190 | C6H4BrCl | 000106-39-8 | 86   |
| 3    |    | Benzene, 1-bromo-3-chloro- | 190 | C6H4BrCl | 000108-37-2 | 83   |
| 4    |    | Benzene, 1-bromo-2-chloro- | 190 | C6H4BrCl | 000694-80-4 | 83   |
| 5    |    | Benzene, 1-bromo-2-chloro- | 190 | C6H4BrCl | 000694-80-4 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\  
Data File : BM019372.D  
Acq On : 24 Mar 2019 03:11  
Operator : JU/SJ  
Sample : K2025-10  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0967

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

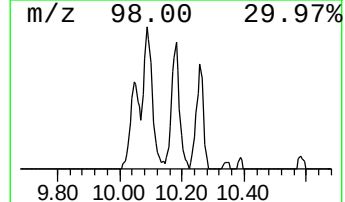
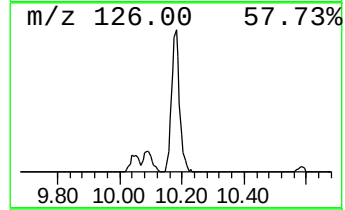
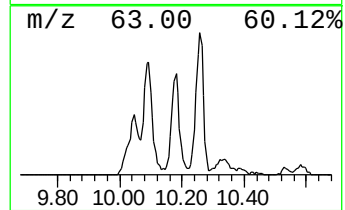
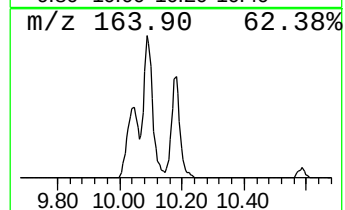
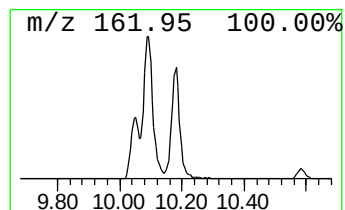
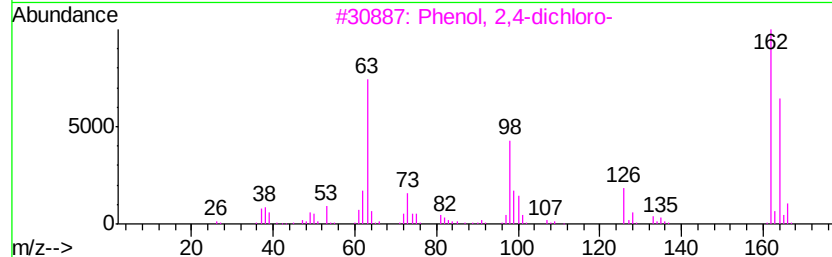
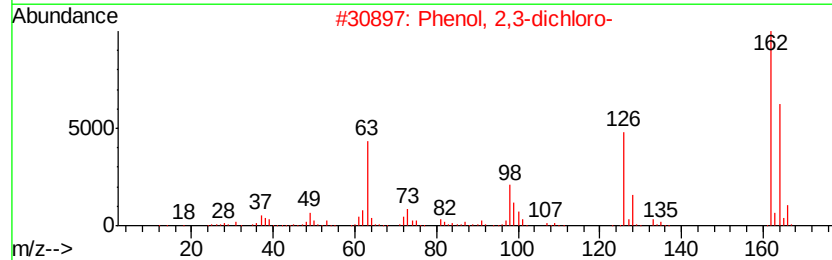
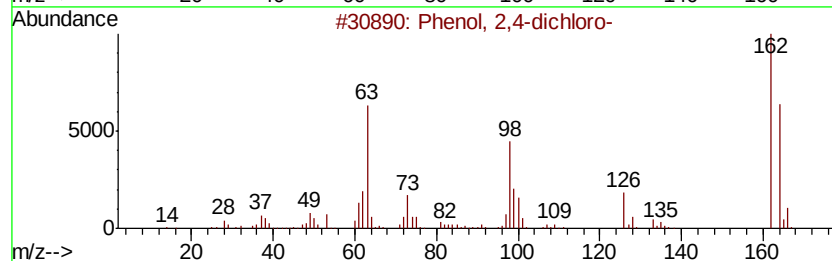
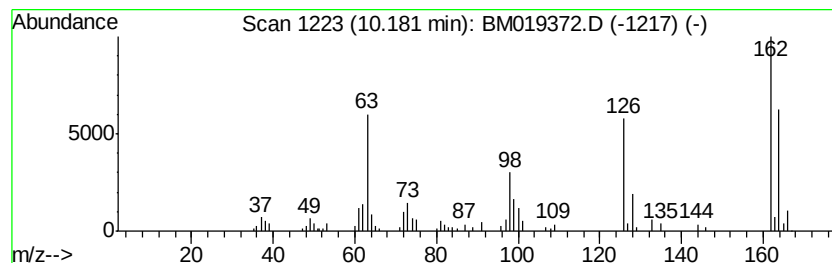
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Phenol, 2,4-dichloro- Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 10.18 | 3.71 ng/ul | 281852 | Naphthalene-d8   | 10.39 |

| Hit# of | 5                     | Tentative ID | MW       | MolForm     | CAS# | Qual |
|---------|-----------------------|--------------|----------|-------------|------|------|
| 1       | Phenol, 2,4-dichloro- | 162          | C6H4Cl2O | 000120-83-2 | 98   |      |
| 2       | Phenol, 2,3-dichloro- | 162          | C6H4Cl2O | 000576-24-9 | 98   |      |
| 3       | Phenol, 2,4-dichloro- | 162          | C6H4Cl2O | 000120-83-2 | 97   |      |
| 4       | Phenol, 2,6-dichloro- | 162          | C6H4Cl2O | 000087-65-0 | 96   |      |
| 5       | Phenol, 2,4-dichloro- | 162          | C6H4Cl2O | 000120-83-2 | 96   |      |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\  
Data File : BM019372.D  
Acq On : 24 Mar 2019 03:11  
Operator : JU/SJ  
Sample : K2025-10  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

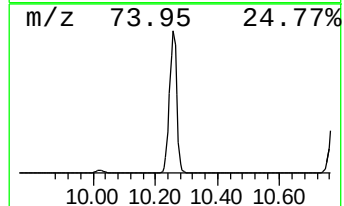
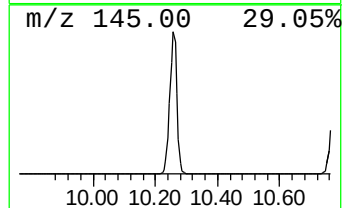
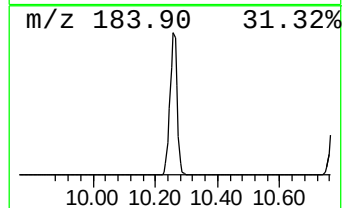
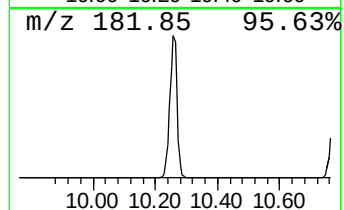
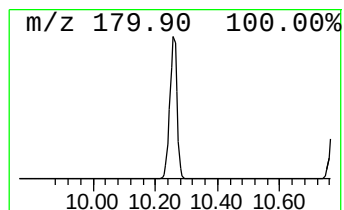
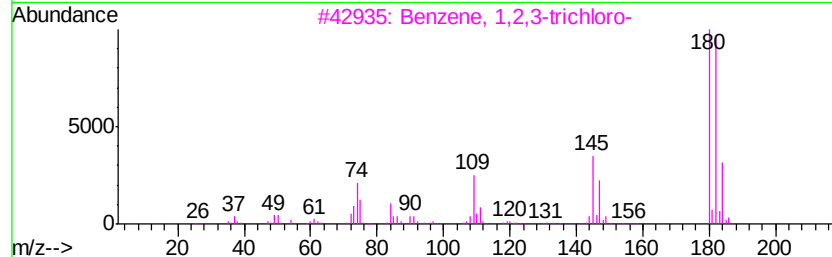
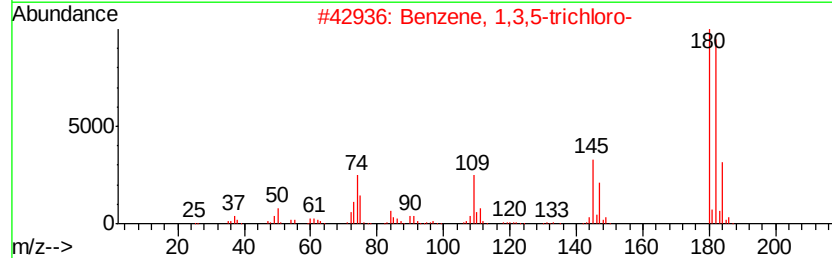
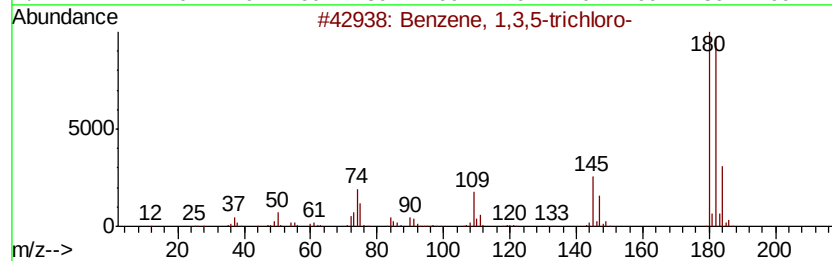
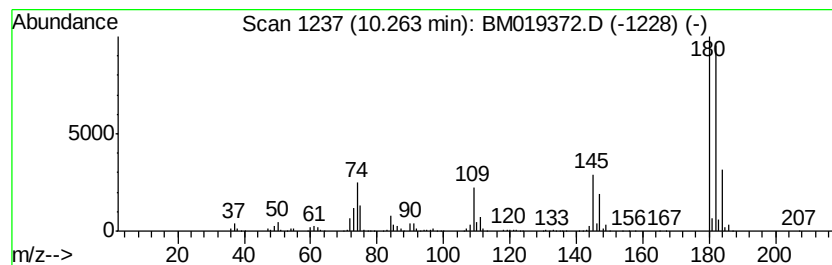
Instrument :  
BNA\_M  
ClientSampleId :  
C0967

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Benzene, 1,3,5-trichloro- Concentration Rank 1

| R.T.    | EstConc      | Area                      | Relative to ISTD | R.T.           |
|---------|--------------|---------------------------|------------------|----------------|
| 10.26   | 221.26 ng/ul | 16823400                  | Naphthalene-d8   | 10.39          |
| Hit# of | 5            | Tentative ID              | MW MolForm       | CAS# Qual      |
| 1       |              | Benzene, 1,3,5-trichloro- | 180 C6H3Cl3      | 000108-70-3 98 |
| 2       |              | Benzene, 1,3,5-trichloro- | 180 C6H3Cl3      | 000108-70-3 98 |
| 3       |              | Benzene, 1,2,3-trichloro- | 180 C6H3Cl3      | 000087-61-6 98 |
| 4       |              | Benzene, 1,2,3-trichloro- | 180 C6H3Cl3      | 000087-61-6 98 |
| 5       |              | Benzene, 1,3,5-trichloro- | 180 C6H3Cl3      | 000108-70-3 97 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\  
Data File : BM019372.D  
Acq On : 24 Mar 2019 03:11  
Operator : JU/SJ  
Sample : K2025-10  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0967

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

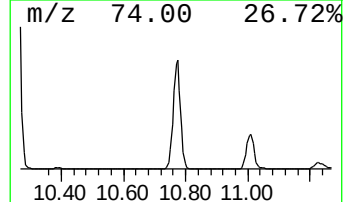
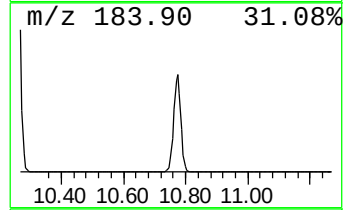
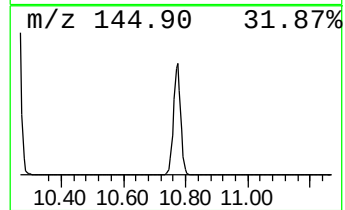
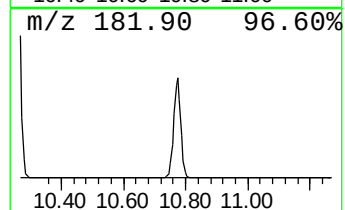
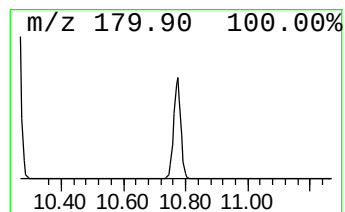
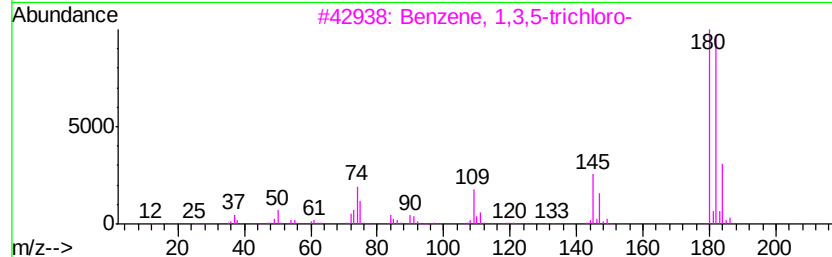
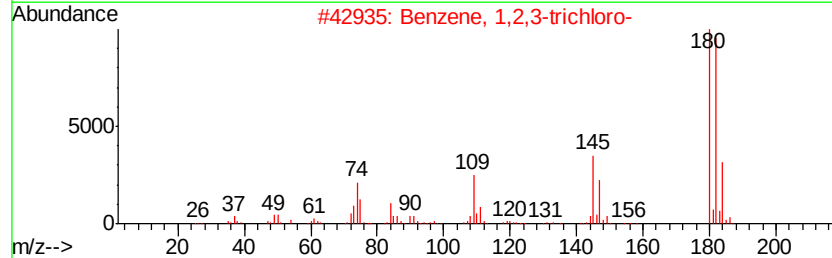
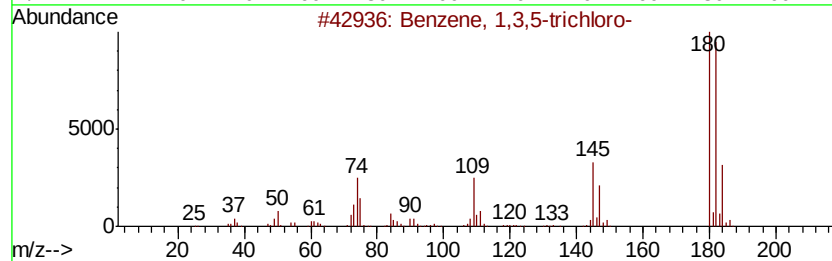
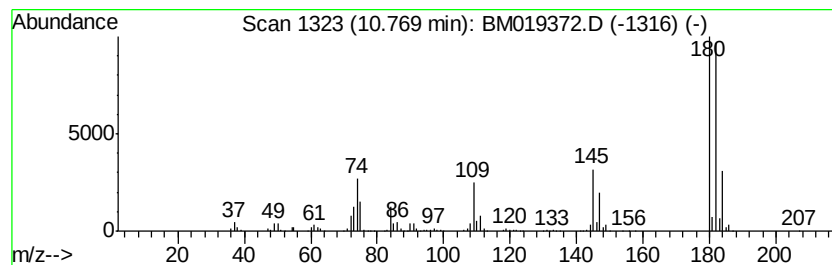
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Benzene, 1,2,3-trichloro- Concentration Rank 2

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 10.77 | 98.33 ng/ul | 7476600 | Naphthalene-d8   | 10.39 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,3,5-trichloro- | 180 | C6H3Cl3 | 000108-70-3 | 98   |
| 2    |    | Benzene, 1,2,3-trichloro- | 180 | C6H3Cl3 | 000087-61-6 | 98   |
| 3    |    | Benzene, 1,3,5-trichloro- | 180 | C6H3Cl3 | 000108-70-3 | 98   |
| 4    |    | Benzene, 1,2,3-trichloro- | 180 | C6H3Cl3 | 000087-61-6 | 98   |
| 5    |    | Benzene, 1,3,5-trichloro- | 180 | C6H3Cl3 | 000108-70-3 | 97   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\  
Data File : BM019372.D  
Acq On : 24 Mar 2019 03:11  
Operator : JU/SJ  
Sample : K2025-10  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0967

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

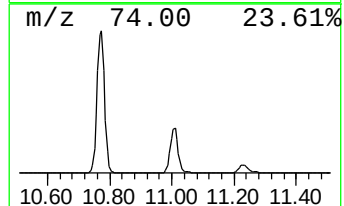
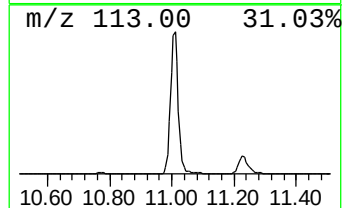
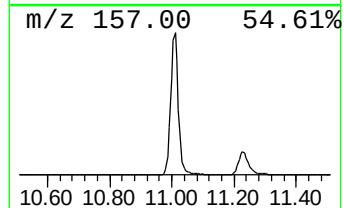
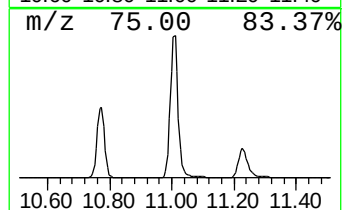
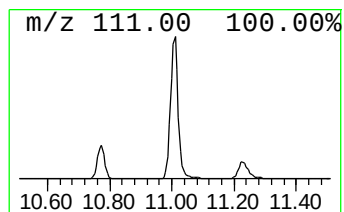
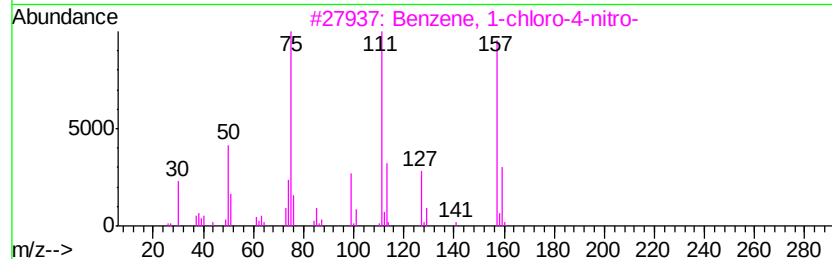
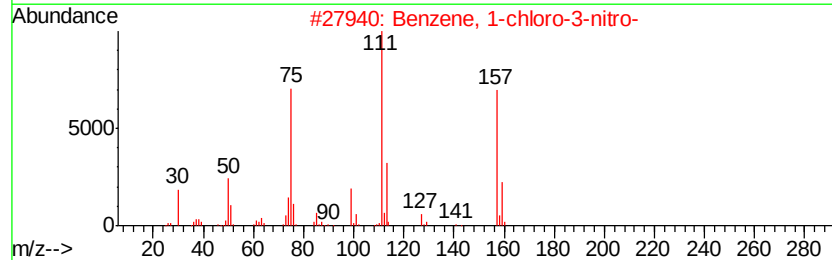
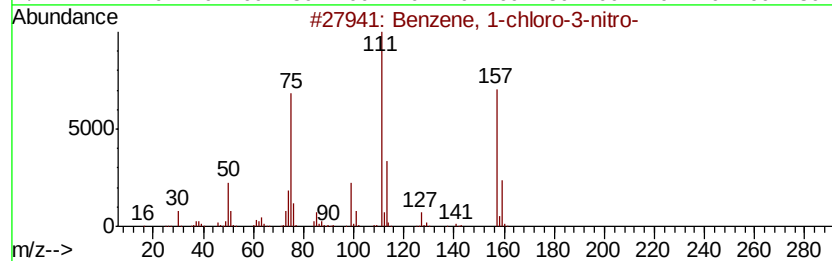
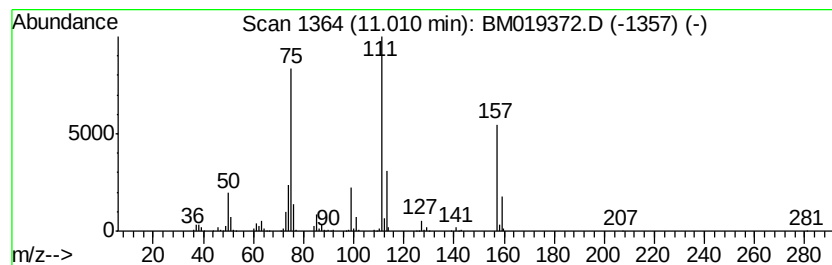
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Benzene, 1-chloro-3-nitro- Concentration Rank 3

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 11.01 | 35.12 ng/ul | 2670570 | Naphthalene-d8   | 10.39 |

| Hit# | of | Tentative ID               | MW  | MolForm   | CAS#        | Qual |
|------|----|----------------------------|-----|-----------|-------------|------|
| 1    | 5  | Benzene, 1-chloro-3-nitro- | 157 | C6H4ClNO2 | 000121-73-3 | 96   |
| 2    |    | Benzene, 1-chloro-3-nitro- | 157 | C6H4ClNO2 | 000121-73-3 | 95   |
| 3    |    | Benzene, 1-chloro-4-nitro- | 157 | C6H4ClNO2 | 000100-00-5 | 94   |
| 4    |    | Benzene, 1-chloro-3-nitro- | 157 | C6H4ClNO2 | 000121-73-3 | 93   |
| 5    |    | Benzene, 1-chloro-2-nitro- | 157 | C6H4ClNO2 | 000088-73-3 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\  
Data File : BM019372.D  
Acq On : 24 Mar 2019 03:11  
Operator : JU/SJ  
Sample : K2025-10  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

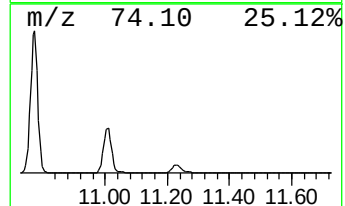
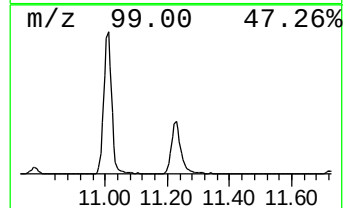
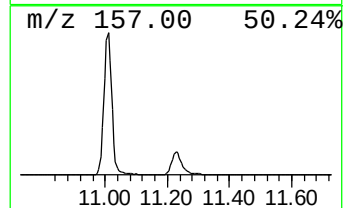
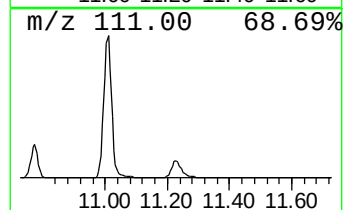
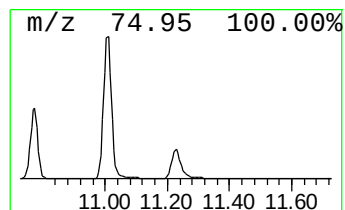
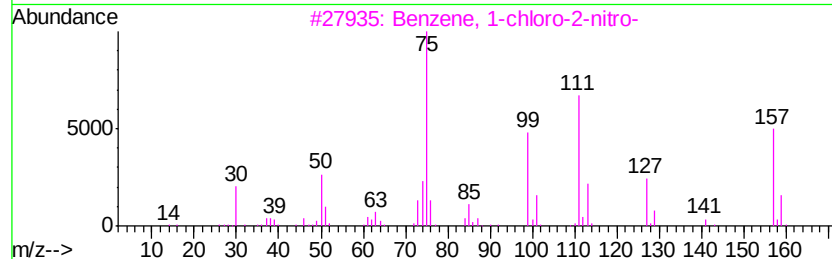
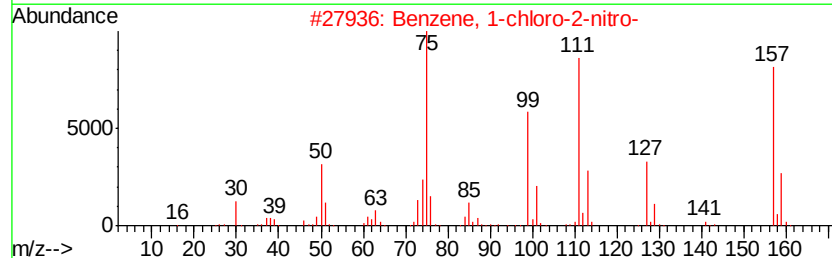
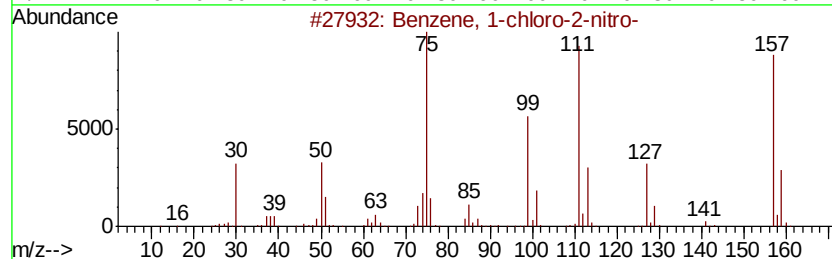
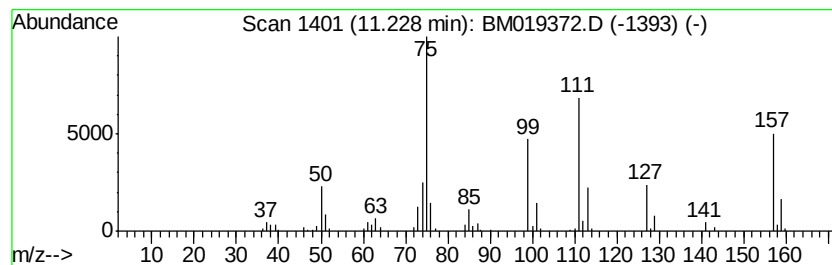
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Benzene, 1-chloro-2-nitro- Concentration Rank 5

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.23 | 8.60 ng/ul | 653985 | Naphthalene-d8   | 10.39 |

| Hit# | of | Tentative ID               | MW  | MolForm   | CAS#        | Qual |
|------|----|----------------------------|-----|-----------|-------------|------|
| 1    | 5  | Benzene, 1-chloro-2-nitro- | 157 | C6H4ClNO2 | 000088-73-3 | 97   |
| 2    |    | Benzene, 1-chloro-2-nitro- | 157 | C6H4ClNO2 | 000088-73-3 | 96   |
| 3    |    | Benzene, 1-chloro-2-nitro- | 157 | C6H4ClNO2 | 000088-73-3 | 96   |
| 4    |    | Benzene, 1-chloro-2-nitro- | 157 | C6H4ClNO2 | 000088-73-3 | 96   |
| 5    |    | Benzene, 1-chloro-4-nitro- | 157 | C6H4ClNO2 | 000100-00-5 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\  
Data File : BM019372.D  
Acq On : 24 Mar 2019 03:11  
Operator : JU/SJ  
Sample : K2025-10  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0967

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

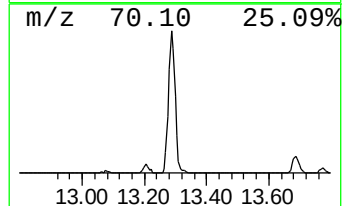
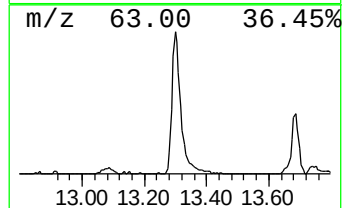
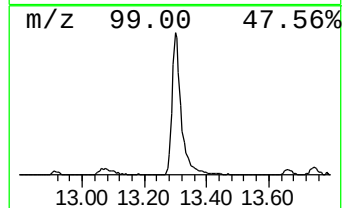
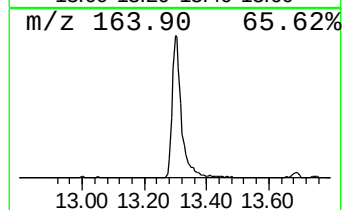
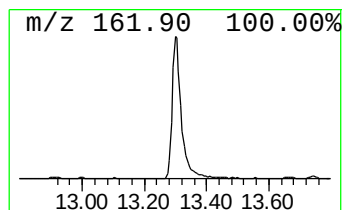
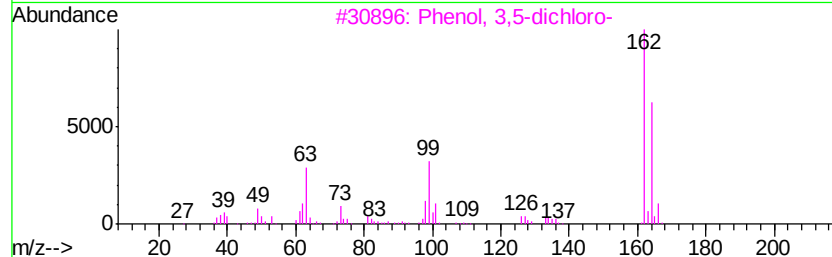
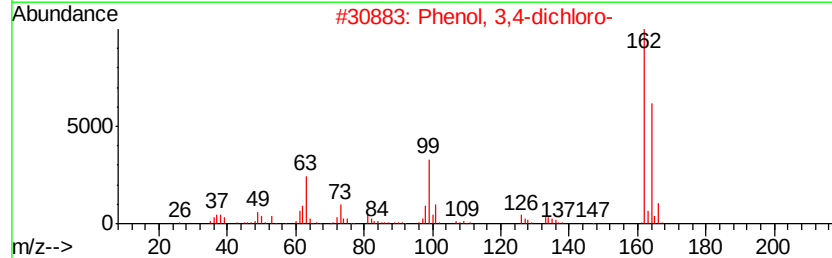
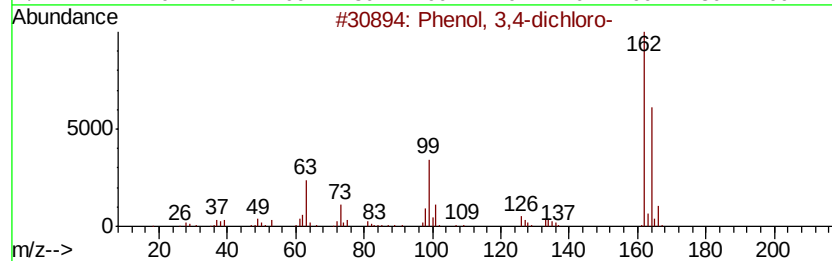
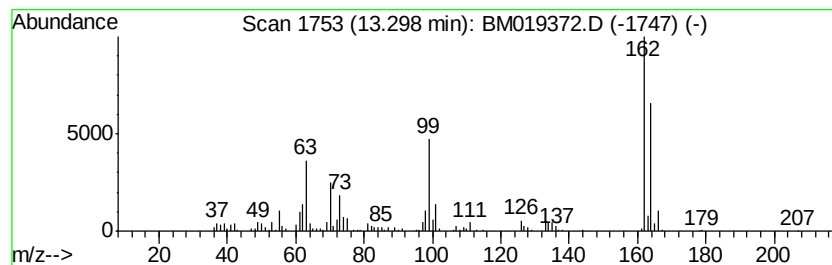
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Phenol, 3,4-dichloro- Concentration Rank 7

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.30 | 7.56 ng/ul | 745664 | Acenaphthene-d10 | 14.26 |

| Hit# | of                    | 5 | Tentative ID | MW  | MolForm  | CAS#        | Qual |
|------|-----------------------|---|--------------|-----|----------|-------------|------|
| 1    | Phenol, 3,4-dichloro- |   |              | 162 | C6H4Cl2O | 000095-77-2 | 95   |
| 2    | Phenol, 3,4-dichloro- |   |              | 162 | C6H4Cl2O | 000095-77-2 | 93   |
| 3    | Phenol, 3,5-dichloro- |   |              | 162 | C6H4Cl2O | 000591-35-5 | 90   |
| 4    | Phenol, 3,5-dichloro- |   |              | 162 | C6H4Cl2O | 000591-35-5 | 90   |
| 5    | Phenol, 3,5-dichloro- |   |              | 162 | C6H4Cl2O | 000591-35-5 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\  
Data File : BM019372.D  
Acq On : 24 Mar 2019 03:11  
Operator : JU/SJ  
Sample : K2025-10  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

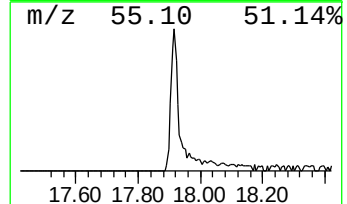
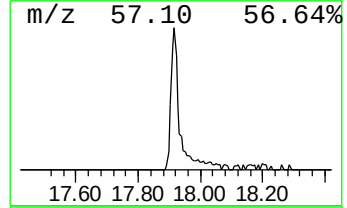
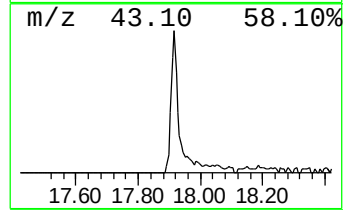
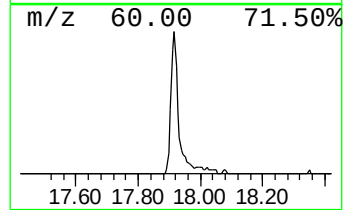
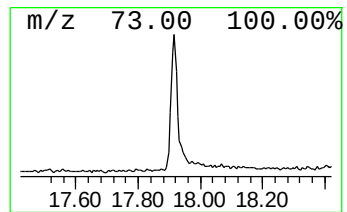
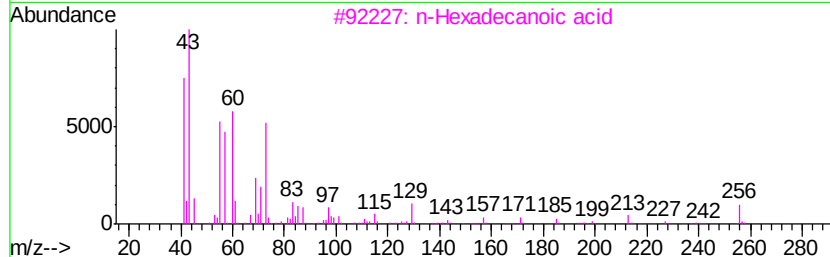
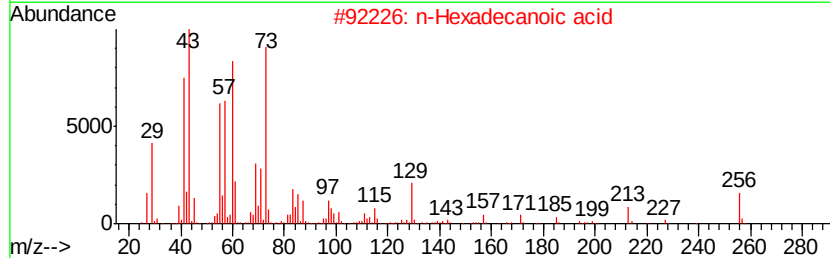
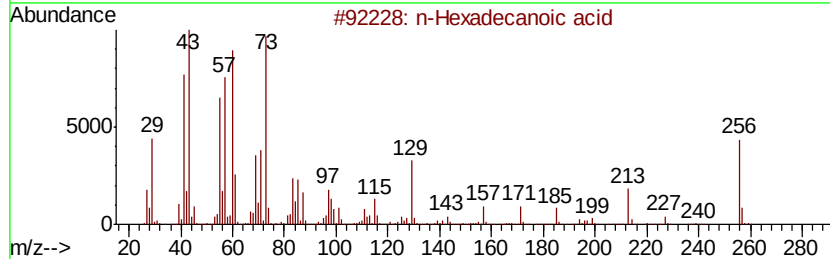
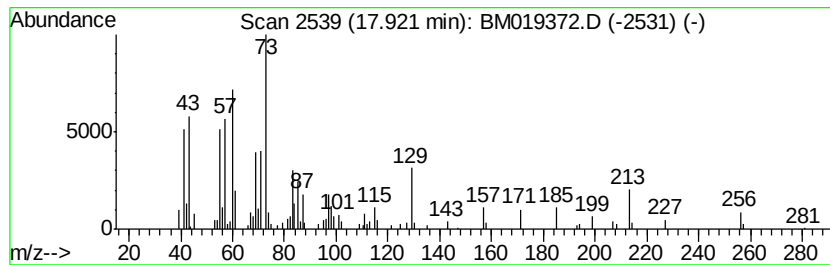
Instrument :  
BNA\_M  
ClientSampleId :  
C0967

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 n-Hexadecanoic acid Concentration Rank 11

| R.T.      | EstConc             | Area   | Relative to ISTD | R.T.        |      |
|-----------|---------------------|--------|------------------|-------------|------|
| 17.92     | 2.63 ng/ul          | 299609 | Phenanthrene-d10 | 17.02       |      |
| Hit# of 5 | Tentative ID        | MW     | MolForm          | CAS#        | Qual |
| 1         | n-Hexadecanoic acid | 256    | C16H32O2         | 000057-10-3 | 98   |
| 2         | n-Hexadecanoic acid | 256    | C16H32O2         | 000057-10-3 | 98   |
| 3         | n-Hexadecanoic acid | 256    | C16H32O2         | 000057-10-3 | 90   |
| 4         | Pentadecanoic acid  | 242    | C15H30O2         | 001002-84-2 | 72   |
| 5         | Undecanoic acid     | 186    | C11H22O2         | 000112-37-8 | 46   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\  
Data File : BM019372.D  
Acq On : 24 Mar 2019 03:11  
Operator : JU/SJ  
Sample : K2025-10  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0967

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

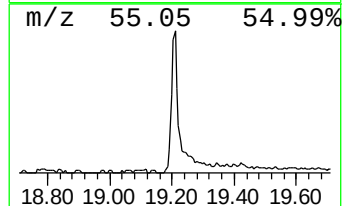
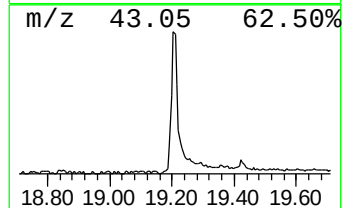
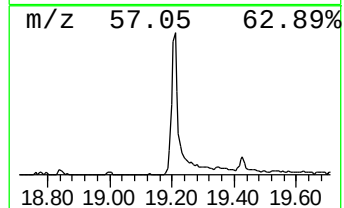
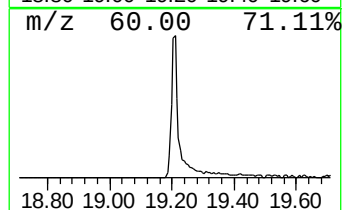
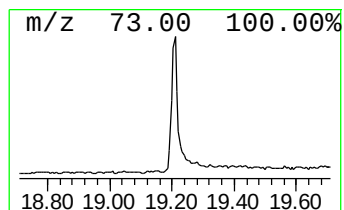
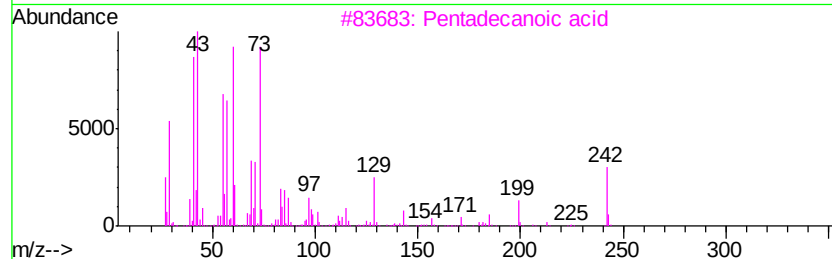
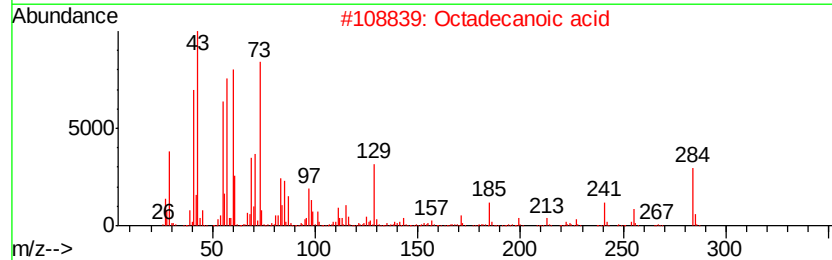
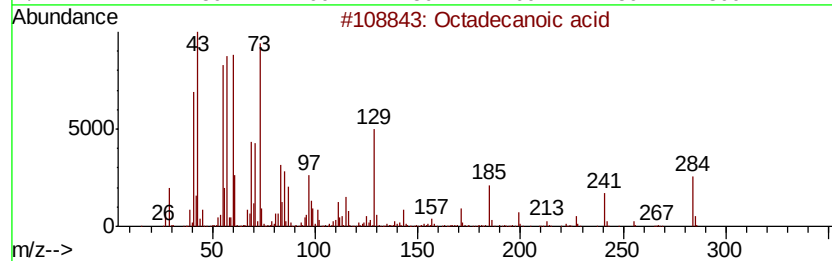
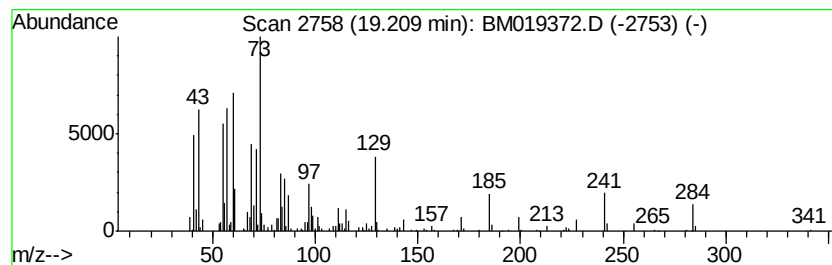
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Octadecanoic acid Concentration Rank 9

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.21 | 4.05 ng/ul | 590456 | Chrysene-d12     | 21.23 |

| Hit# | of | Tentative ID       | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------|-----|----------|-------------|------|
| 1    | 5  | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 99   |
| 2    |    | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 95   |
| 3    |    | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 90   |
| 4    |    | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 89   |
| 5    |    | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\  
Data File : BM019372.D  
Acq On : 24 Mar 2019 03:11  
Operator : JU/SJ  
Sample : K2025-10  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0967

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

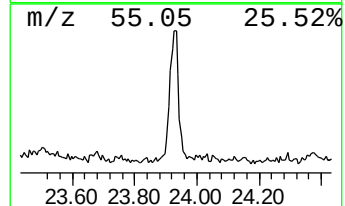
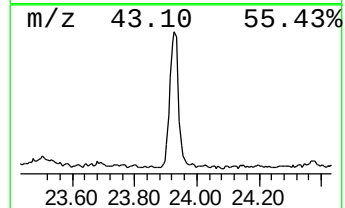
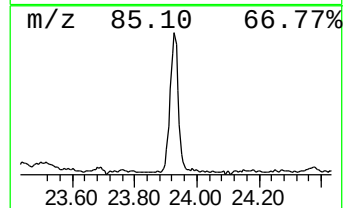
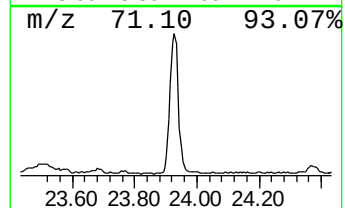
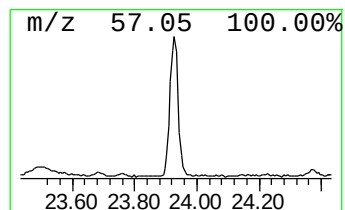
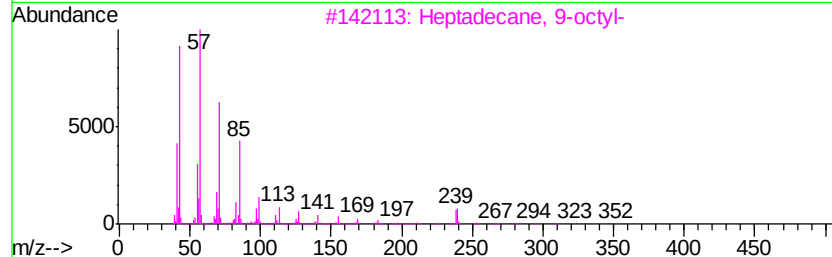
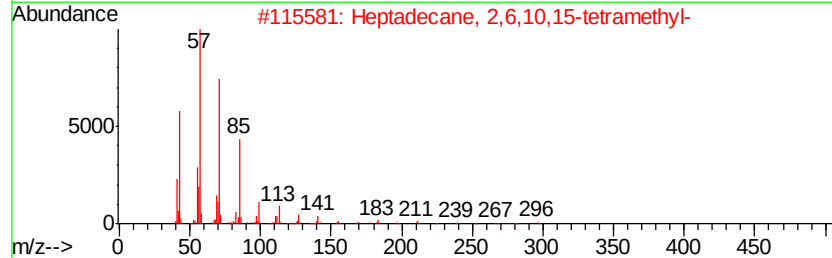
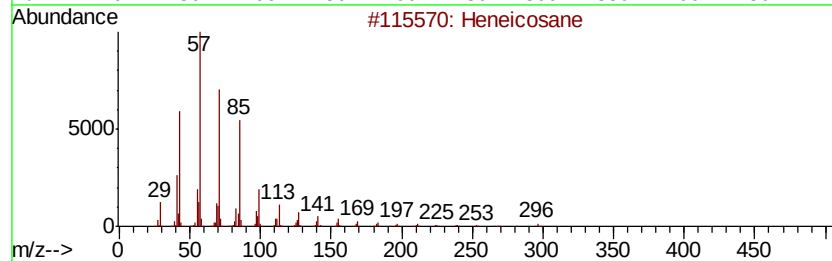
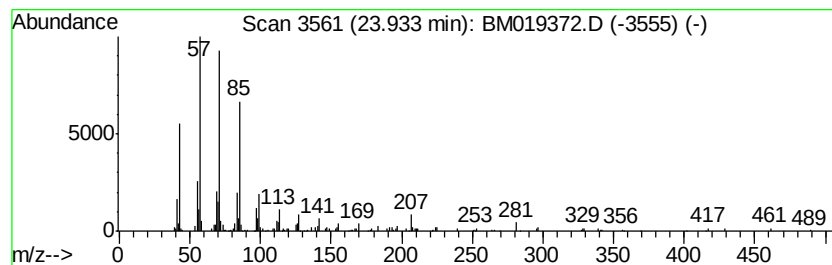
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 23.93 | 2.16 ng/ul | 325607 | Perylene-d12     | 23.44 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Heneicosane                         | 296 | C21H44  | 000629-94-7 | 98   |
| 2    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 93   |
| 3    |    |   | Heptadecane, 9-octyl-               | 352 | C25H52  | 007225-64-1 | 87   |
| 4    |    |   | Heptacosane                         | 380 | C27H56  | 000593-49-7 | 87   |
| 5    |    |   | Docosane                            | 310 | C22H46  | 000629-97-0 | 86   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\  
Data File : BM019372.D  
Acq On : 24 Mar 2019 03:11  
Operator : JU/SJ  
Sample : K2025-10  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0967

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

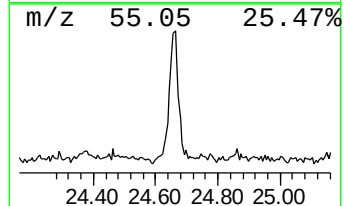
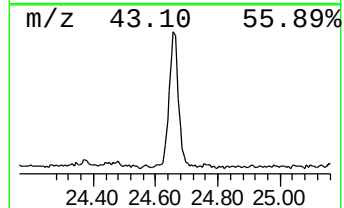
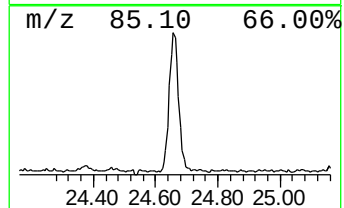
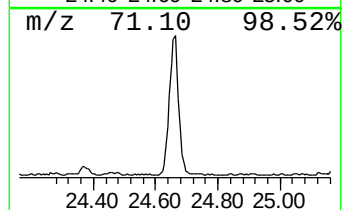
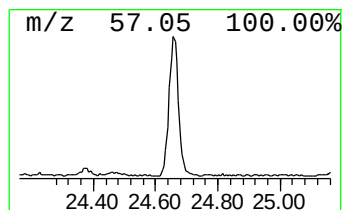
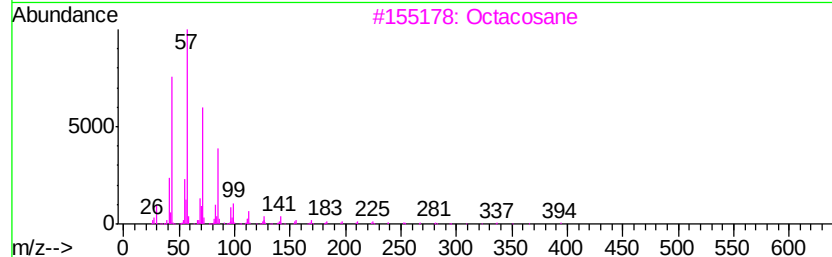
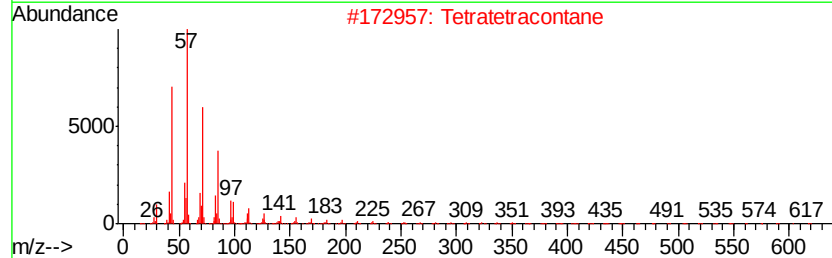
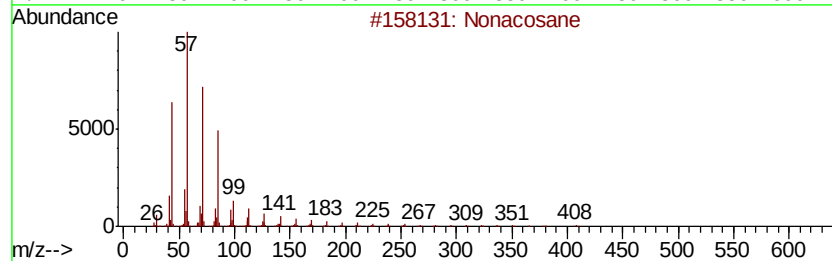
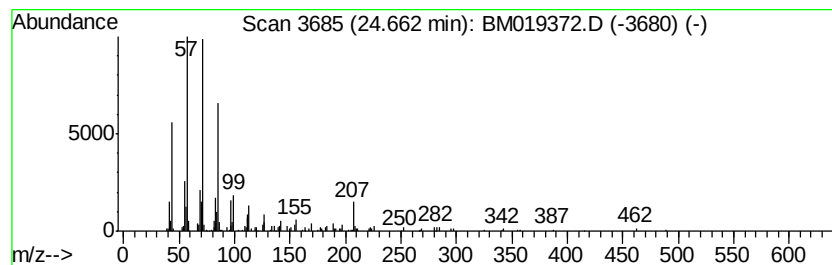
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 13

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 24.66 | 2.16 ng/ul | 325708 | Perylene-d12     | 23.44 |

| Hit# | of | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------|-----|---------|-------------|------|
| 1    | 5  | Nonacosane        | 408 | C29H60  | 000630-03-5 | 83   |
| 2    |    | Tetratetracontane | 619 | C44H90  | 007098-22-8 | 83   |
| 3    |    | Octacosane        | 394 | C28H58  | 000630-02-4 | 83   |
| 4    |    | triacontane       | 422 | C30H62  | 000638-68-6 | 83   |
| 5    |    | Eicosane          | 282 | C20H42  | 000112-95-8 | 81   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM032319\  
 Data File : BM019372.D  
 Acq On : 24 Mar 2019 03:11  
 Operator : JU/SJ  
 Sample : K2025-10  
 Misc :  
 ALS Vial : 30 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C0967

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |          |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|----------|------|
|                      |       |         |       |          | #                     | RT    | Resp     | Conc |
| Benzene, chloro-     | 4.96  | 8.6     | ng/ul | 31124000 | 1                     | 7.62  | 72501000 | 20.0 |
| Benzene, 1,4-dich... | 7.50  | 4.6     | ng/ul | 16588900 | 1                     | 7.62  | 72501000 | 20.0 |
| Benzene, 1,2-dich... | 7.98  | 22.7    | ng/ul | 82197000 | 1                     | 7.62  | 72501000 | 20.0 |
| 4-Chlorophenylsel... | 9.10  | 2.4     | ng/ul | 181875   | 2                     | 10.39 | 1520730  | 20.0 |
| Phenol, 2,4-dichl... | 10.18 | 3.7     | ng/ul | 281852   | 2                     | 10.39 | 1520730  | 20.0 |
| Benzene, 1,3,5-tr... | 10.26 | 221.3   | ng/ul | 16823400 | 2                     | 10.39 | 1520730  | 20.0 |
| Benzene, 1,2,3-tr... | 10.77 | 98.3    | ng/ul | 7476600  | 2                     | 10.39 | 1520730  | 20.0 |
| Benzene, 1-chloro... | 11.01 | 35.1    | ng/ul | 2670570  | 2                     | 10.39 | 1520730  | 20.0 |
| Benzene, 1-chloro... | 11.23 | 8.6     | ng/ul | 653985   | 2                     | 10.39 | 1520730  | 20.0 |
| Phenol, 3,4-dichl... | 13.30 | 7.6     | ng/ul | 745664   | 3                     | 14.26 | 1972530  | 20.0 |
| n-Hexadecanoic acid  | 17.92 | 2.6     | ng/ul | 299609   | 4                     | 17.02 | 2280200  | 20.0 |
| Octadecanoic acid    | 19.21 | 4.0     | ng/ul | 590456   | 5                     | 21.23 | 2915870  | 20.0 |
| (DEL) Alkane: Str... | 23.93 | 2.2     | ng/ul | 325607   | 6                     | 23.44 | 3018650  | 20.0 |
| (DEL) Alkane: Str... | 24.66 | 2.2     | ng/ul | 325708   | 6                     | 23.44 | 3018650  | 20.0 |