

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081720\  
 Data File : BG046320.D  
 Acq On : 18 Aug 2020 9:46  
 Operator : CG/JU  
 Sample : PB130999BL  
 Misc :  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SBLK99

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\_G\METHODS\SOM-EPA-BG081320PAH.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         | 3.140       | 5             | 9           | 18           | rVB2     | 30339          | 46756         | 1.93%           | 0.205%        |
| 2         | 3.399       | 49            | 53          | 61           | rVB      | 19955          | 29754         | 1.23%           | 0.131%        |
| 3         | 4.145       | 177           | 180         | 184          | rVB2     | 3246           | 4648          | 0.19%           | 0.020%        |
| 4         | 4.269       | 198           | 201         | 206          | rBV4     | 1573           | 2441          | 0.10%           | 0.011%        |
| 5         | 7.106       | 677           | 684         | 696          | rBV      | 295872         | 522796        | 21.54%          | 2.295%        |
| 6         | 7.265       | 703           | 711         | 721          | rBV      | 245679         | 408320        | 16.83%          | 1.792%        |
| 7         | 7.471       | 738           | 746         | 755          | rBV      | 315554         | 540511        | 22.27%          | 2.373%        |
| 8         | 7.941       | 818           | 826         | 834          | rBV      | 258628         | 442921        | 18.25%          | 1.944%        |
| 9         | 8.646       | 938           | 946         | 958          | rBV      | 377007         | 666710        | 27.47%          | 2.927%        |
| 10        | 9.092       | 1014          | 1022        | 1039         | rBV      | 283446         | 511540        | 21.08%          | 2.246%        |
| 11        | 9.815       | 1137          | 1145        | 1155         | rBV      | 241005         | 431675        | 17.79%          | 1.895%        |
| 12        | 10.356      | 1229          | 1237        | 1249         | rBV      | 381108         | 701810        | 28.92%          | 3.081%        |
| 13        | 10.737      | 1292          | 1302        | 1310         | rBV      | 371174         | 666780        | 27.48%          | 2.927%        |
| 14        | 10.861      | 1316          | 1323        | 1334         | rBV      | 356790         | 675376        | 27.83%          | 2.965%        |
| 15        | 12.324      | 1567          | 1572        | 1579         | rBV3     | 5526           | 9102          | 0.38%           | 0.040%        |
| 16        | 13.969      | 1846          | 1852        | 1870         | rBV      | 738353         | 1092498       | 45.02%          | 4.796%        |
| 17        | 14.257      | 1894          | 1901        | 1914         | rBV      | 865504         | 1362339       | 56.14%          | 5.981%        |
| 18        | 14.568      | 1945          | 1954        | 1962         | rBV      | 622112         | 969550        | 39.95%          | 4.256%        |
| 19        | 14.762      | 1980          | 1987        | 2003         | rBV      | 242270         | 456188        | 18.80%          | 2.003%        |
| 20        | 15.555      | 2116          | 2122        | 2131         | rBV      | 1149030        | 1742642       | 71.81%          | 7.650%        |
| 21        | 15.673      | 2136          | 2142        | 2159         | rVV      | 281565         | 442623        | 18.24%          | 1.943%        |
| 22        | 15.796      | 2159          | 2163        | 2169         | rVB      | 13143          | 19066         | 0.79%           | 0.084%        |
| 23        | 16.343      | 2250          | 2256        | 2261         | rBV3     | 4798           | 6889          | 0.28%           | 0.030%        |
| 24        | 16.995      | 2362          | 2367        | 2370         | rBV4     | 2197           | 2532          | 0.10%           | 0.011%        |
| 25        | 17.095      | 2379          | 2384        | 2388         | rBV2     | 2002           | 3583          | 0.15%           | 0.016%        |
| 26        | 17.306      | 2414          | 2420        | 2430         | rBV      | 796936         | 1204861       | 49.65%          | 5.289%        |
| 27        | 17.406      | 2430          | 2437        | 2454         | rVB      | 1307018        | 1909510       | 78.69%          | 8.383%        |
| 28        | 17.594      | 2468          | 2469        | 2477         | rVB4     | 1343           | 2779          | 0.11%           | 0.012%        |
| 29        | 19.333      | 2759          | 2765        | 2769         | rBV      | 14599          | 23586         | 0.97%           | 0.104%        |
| 30        | 19.580      | 2802          | 2807        | 2811         | rBV      | 13923          | 18580         | 0.77%           | 0.082%        |
| 31        | 19.692      | 2820          | 2826        | 2838         | rBV      | 1594748        | 2320023       | 95.60%          | 10.185%       |
| 32        | 20.238      | 2917          | 2919        | 2924         | rBV6     | 2920           | 4355          | 0.18%           | 0.019%        |
| 33        | 20.373      | 2940          | 2942        | 2943         | rBV2     | 3677           | 2623          | 0.11%           | 0.012%        |
| 34        | 20.608      | 2980          | 2982        | 2988         | rBV5     | 5564           | 9459          | 0.39%           | 0.042%        |

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Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SBLK99

## 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\_G\METHODS\SOM-EPA-BG081320PAH.M  
Title : SVOA CALIBRATION

|    |        |      |      |      |      |        |         |         |         |
|----|--------|------|------|------|------|--------|---------|---------|---------|
| 35 | 20.890 | 3028 | 3030 | 3032 | rBV3 | 5565   | 3201    | 0.13%   | 0.014%  |
| 36 | 21.554 | 3136 | 3143 | 3154 | rBV  | 880130 | 1396632 | 57.55%  | 6.131%  |
| 37 | 23.023 | 3391 | 3393 | 3399 | rVB6 | 19256  | 28980   | 1.19%   | 0.127%  |
| 38 | 23.799 | 3522 | 3525 | 3532 | rVB4 | 23504  | 44784   | 1.85%   | 0.197%  |
| 39 | 24.445 | 3624 | 3635 | 3647 | rBV  | 895250 | 2426692 | 100.00% | 10.653% |
| 40 | 24.656 | 3661 | 3671 | 3690 | rVB2 | 510237 | 1532516 | 63.15%  | 6.728%  |
| 41 | 25.808 | 3862 | 3867 | 3878 | rVB4 | 34653  | 91763   | 3.78%   | 0.403%  |

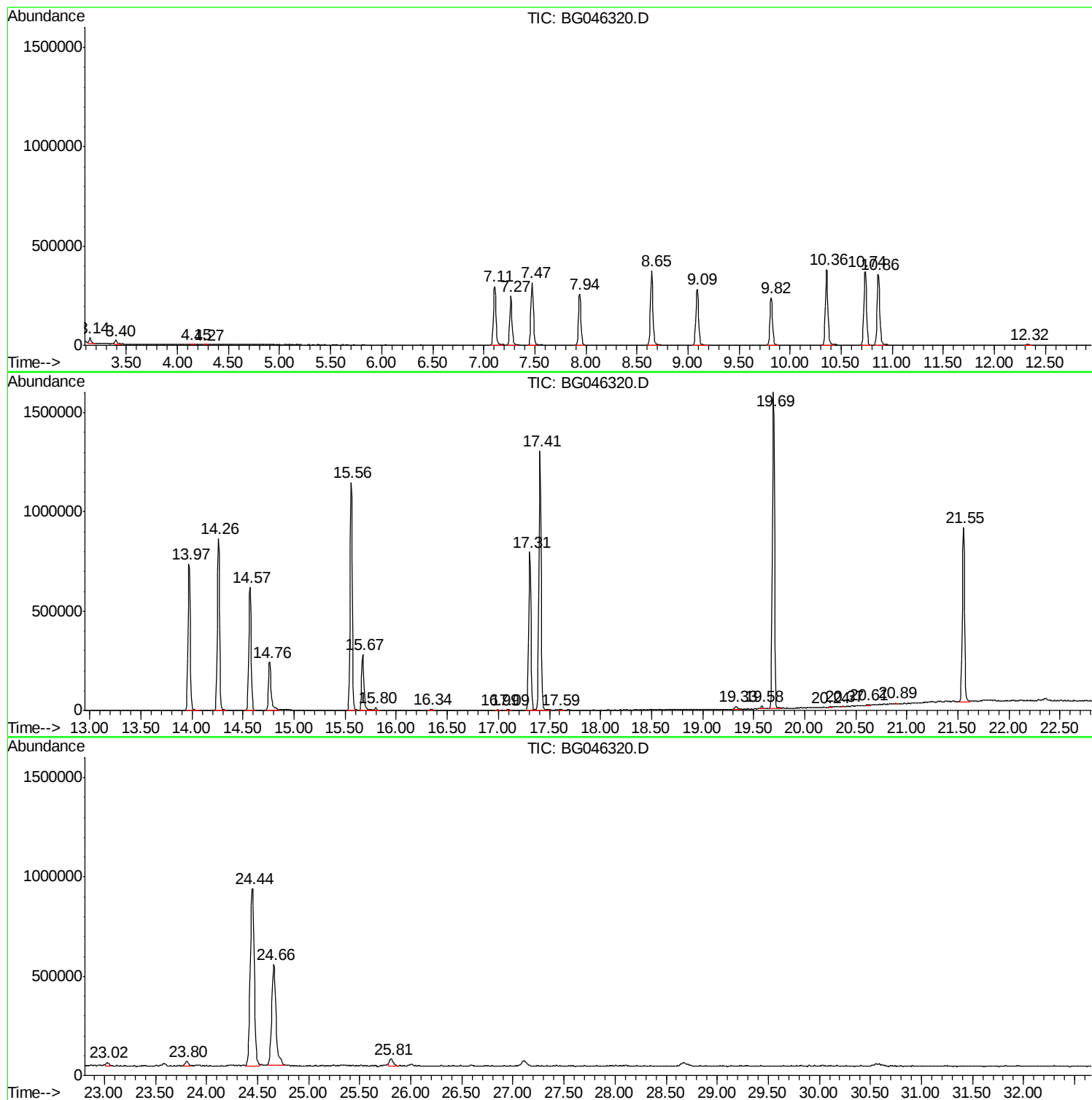
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

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



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Sample : PB130999BL  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
SBLK99

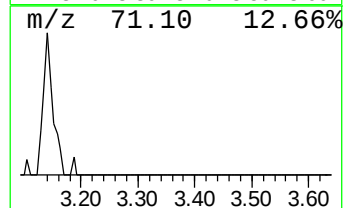
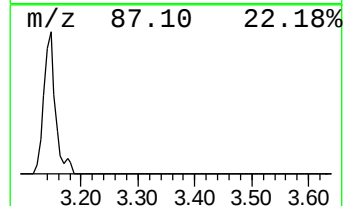
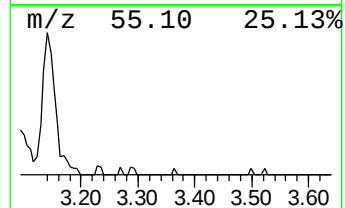
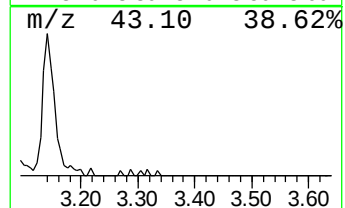
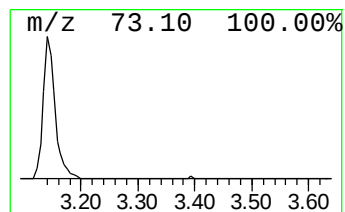
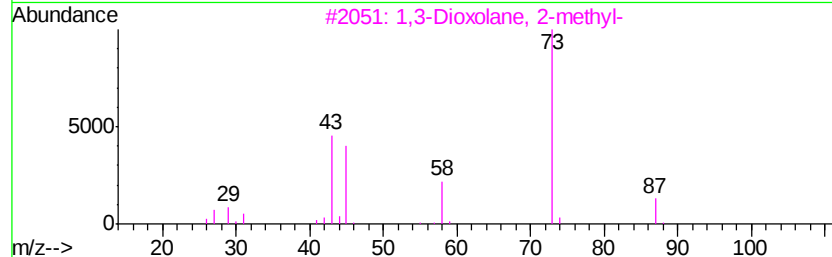
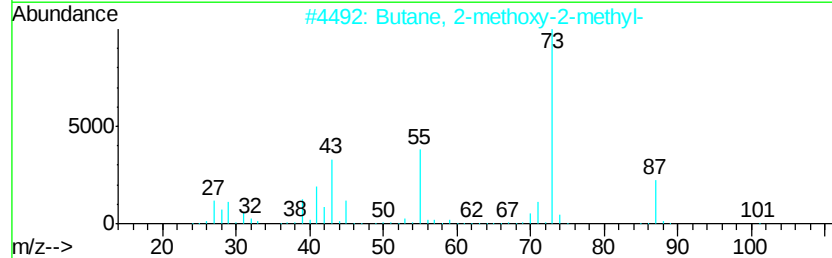
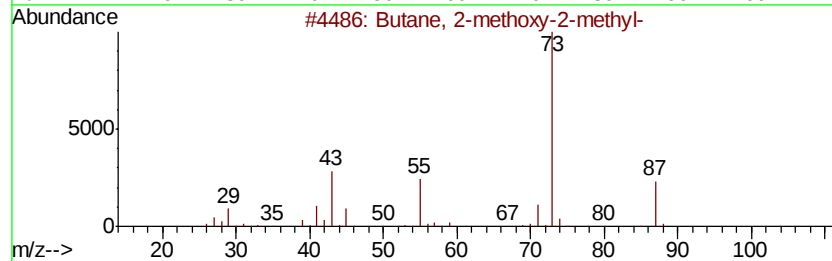
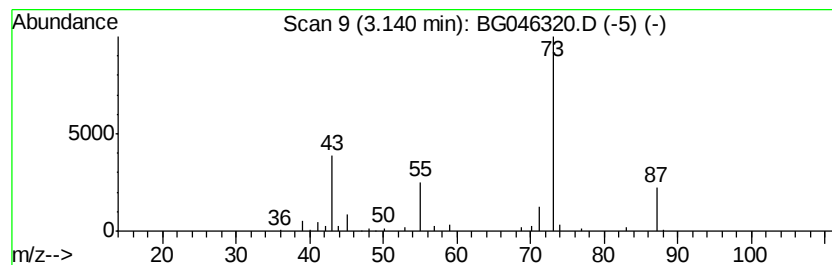
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 11

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 3.14 | 2.11 ng/ul | 46756 | 1,4-Dichlorobenzene-d4 | 7.94 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 56   |
| 2    |    |   | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 50   |
| 3    |    |   | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 23   |
| 4    |    |   | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 12   |
| 5    |    |   | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 9    |



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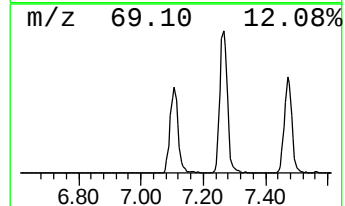
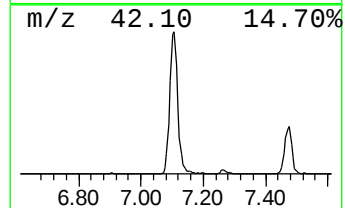
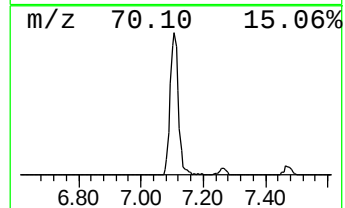
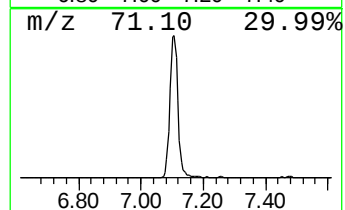
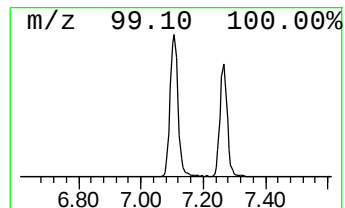
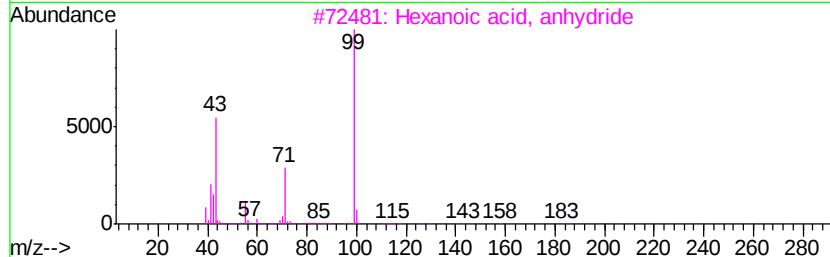
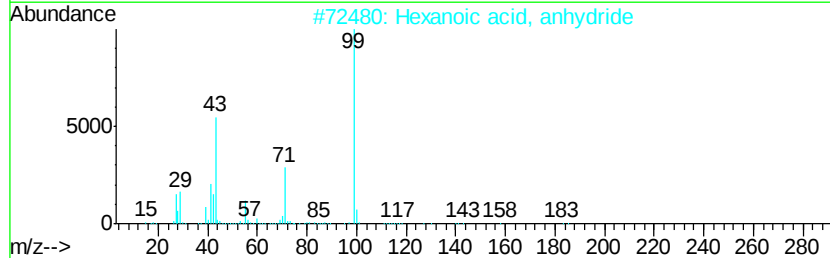
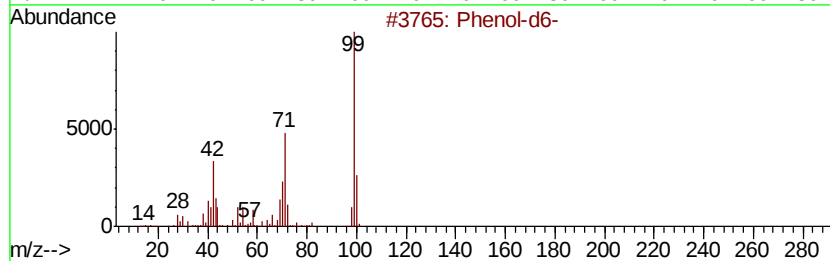
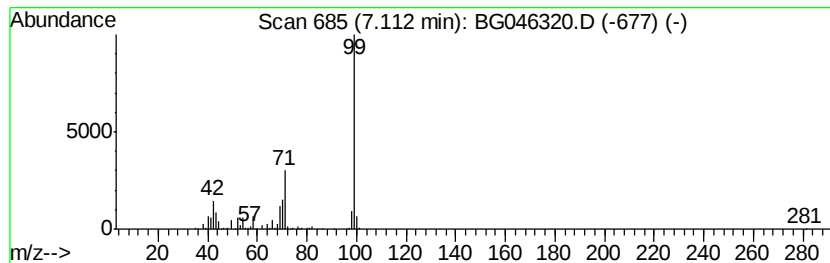
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 Hexanoic acid, anhydride Concentration Rank 3

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 7.11 | 23.61 ng/ul | 522796 | 1,4-Dichlorobenzene-d4 | 7.94 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Phenol-d6-                          | 100 | C6D6O    | 013127-88-3 | 64   |
| 2    |    | Hexanoic acid, anhydride            | 214 | C12H22O3 | 002051-49-2 | 50   |
| 3    |    | Hexanoic acid, anhydride            | 214 | C12H22O3 | 002051-49-2 | 50   |
| 4    |    | 4H-Imidazol-4-one, 2-amino-1,5-d... | 99  | C3H5N3O  | 000503-86-6 | 45   |
| 5    |    | 3,4-Diamino-1,2,4(4H)-triazole      | 99  | C2H5N5   | 038104-45-9 | 38   |



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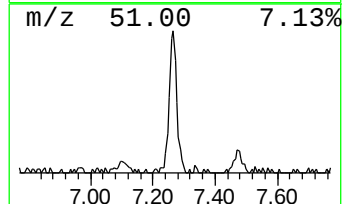
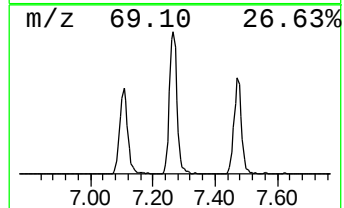
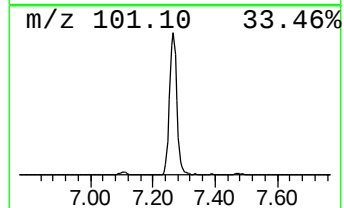
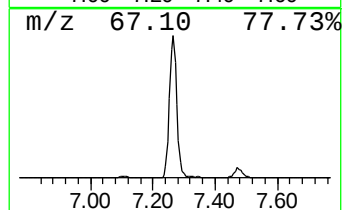
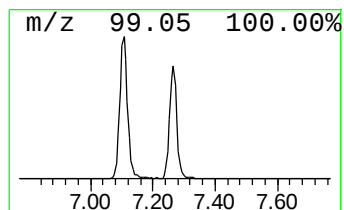
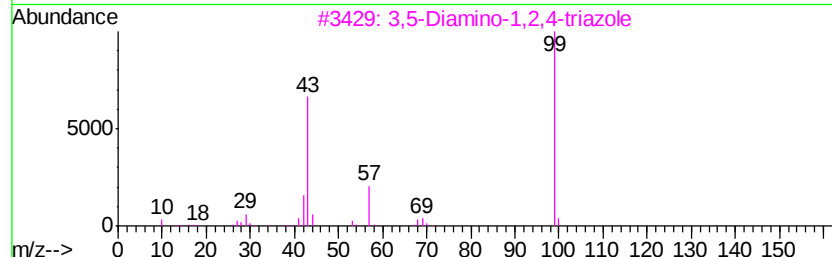
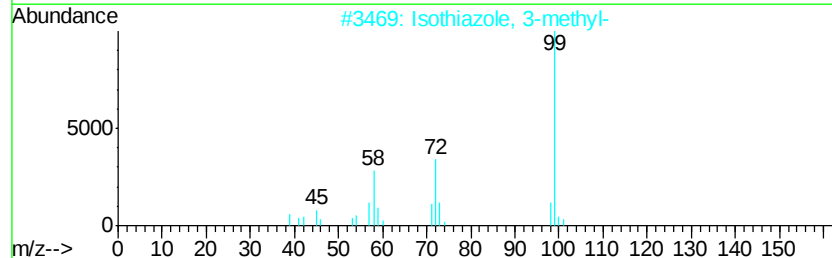
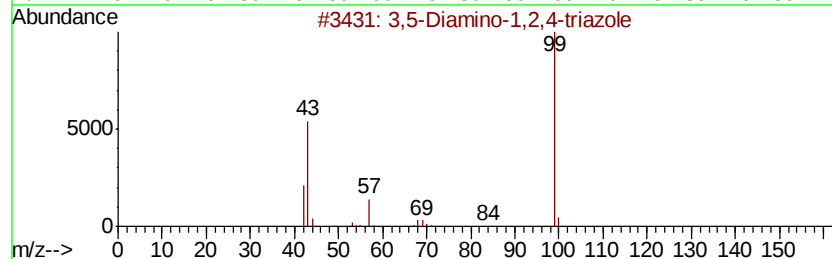
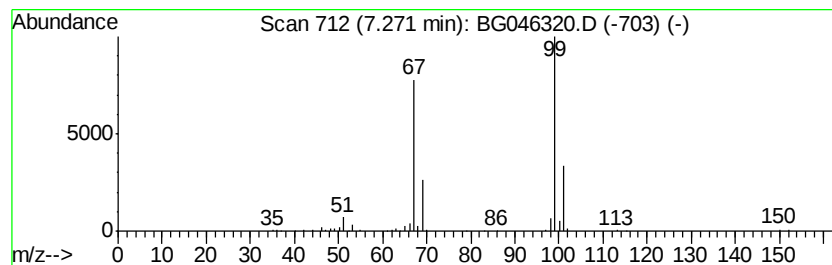
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 unknown-01 Concentration Rank 7

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 7.27 | 18.44 ng/ul | 408320 | 1,4-Dichlorobenzene-d4 | 7.94 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | 3,5-Diamino-1,2,4-triazole     | 99  | C2H5N5  | 001455-77-2 | 9    |
| 2    |    | Isothiazole, 3-methyl-         | 99  | C4H5NS  | 000693-92-5 | 9    |
| 3    |    | 3,5-Diamino-1,2,4-triazole     | 99  | C2H5N5  | 001455-77-2 | 9    |
| 4    |    | 1-Pentanamine, 5-(methylthio)- | 133 | C6H15NS | 055021-78-8 | 9    |
| 5    |    | Isothiazole, 3-methyl-         | 99  | C4H5NS  | 000693-92-5 | 7    |



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Quant Title : SVOA CALIBRATION

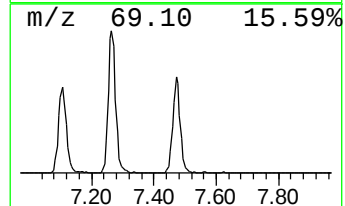
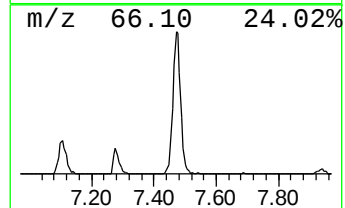
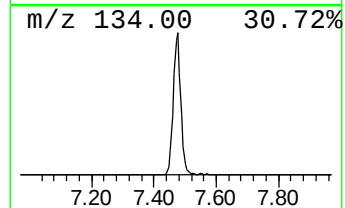
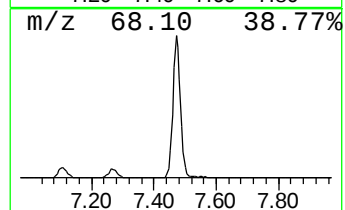
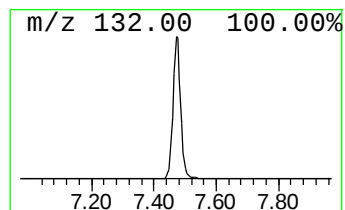
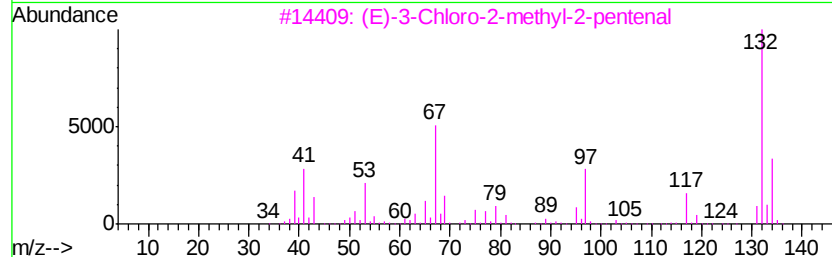
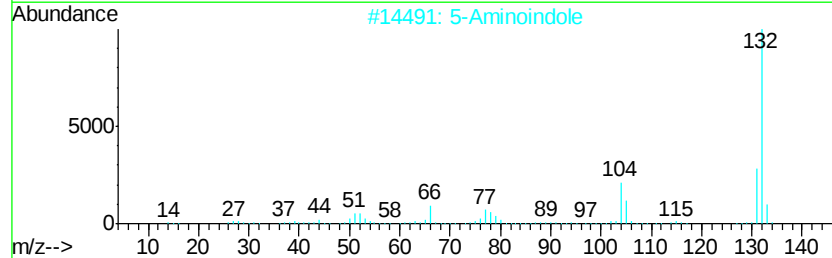
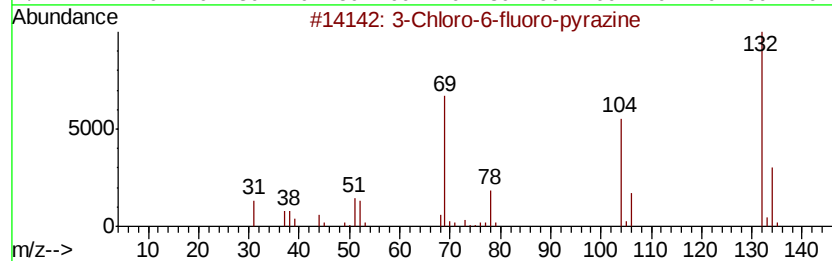
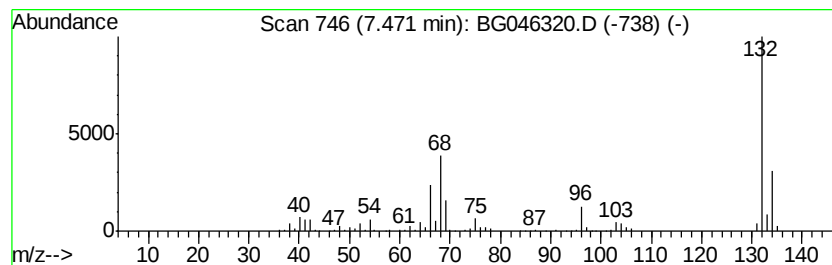
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 unknown-02 Concentration Rank 2

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 7.47 | 24.41 ng/ul | 540511 | 1,4-Dichlorobenzene-d4 | 7.94 |

| Hit# | of                               | 5 | Tentative ID | MW  | MolForm   | CAS#         | Qual |
|------|----------------------------------|---|--------------|-----|-----------|--------------|------|
| 1    | 3-Chloro-6-fluoro-pyrazine       |   |              | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    | 5-Aminoindole                    |   |              | 132 | C8H8N2    | 005192-03-0  | 22   |
| 3    | (E)-3-Chloro-2-methyl-2-pentenal |   |              | 132 | C6H9ClO   | 031357-76-3  | 16   |
| 4    | Tranylcypromine-propionyl        |   |              | 189 | C12H15NO  | 1000123-86-3 | 12   |
| 5    | Amphetaminil                     |   |              | 250 | C17H18N2  | 017590-01-1  | 12   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081720\  
Data File : BG046320.D  
Acq On : 18 Aug 2020 9:46  
Operator : CG/JU  
Sample : PB130999BL  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SBLK99

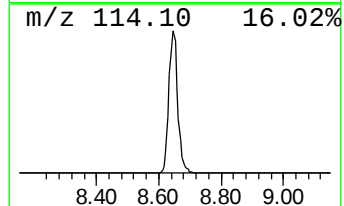
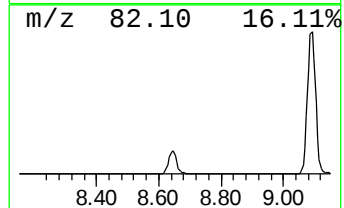
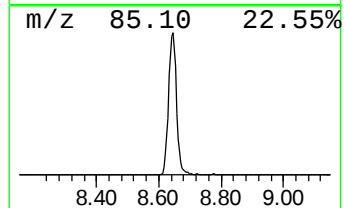
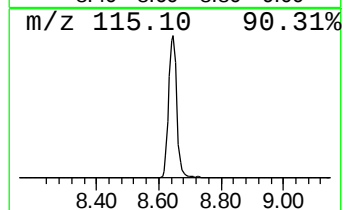
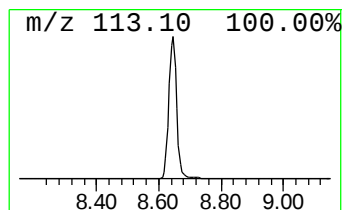
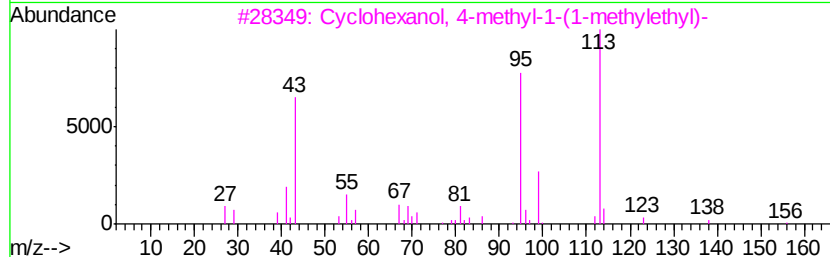
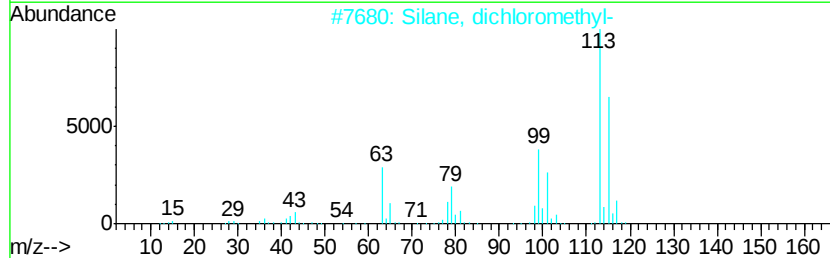
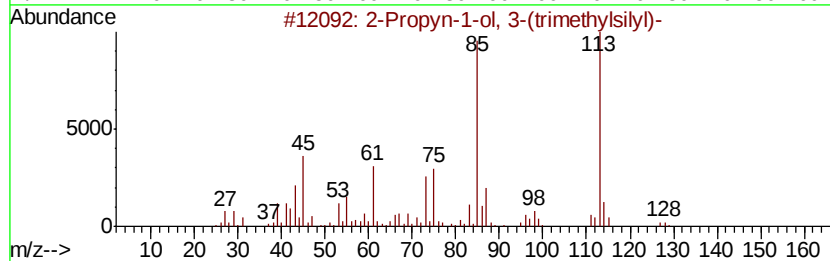
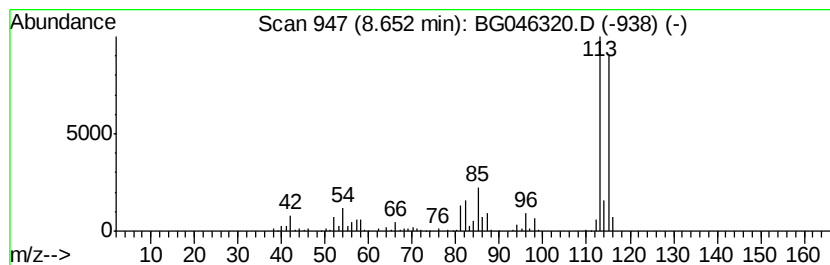
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 unknown-03 Concentration Rank 1

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 8.65 | 30.11 ng/ul | 666710 | 1,4-Dichlorobenzene-d4 | 7.94 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm   | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | 2-Propyn-1-ol, 3-(trimethylsilyl)-   | 128 | C6H12OSi  | 005272-36-6 | 43   |
| 2    |    |   | Silane, dichloromethyl-              | 114 | CH4Cl2Si  | 000075-54-7 | 38   |
| 3    |    |   | Cyclohexanol, 4-methyl-1-(1-meth...  | 156 | C10H20O   | 000470-65-5 | 35   |
| 4    |    |   | Butanedioyl dihydrazide              | 146 | C4H10N4O2 | 004146-43-4 | 35   |
| 5    |    |   | 2H-Pyrrrol-2-one, 1,5-dihydro-4-m... | 113 | C5H7NO2   | 069778-83-2 | 32   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081720\  
Data File : BG046320.D  
Acq On : 18 Aug 2020 9:46  
Operator : CG/JU  
Sample : PB130999BL  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SBLK99

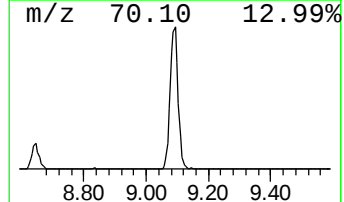
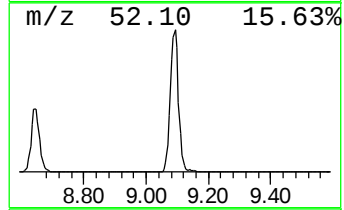
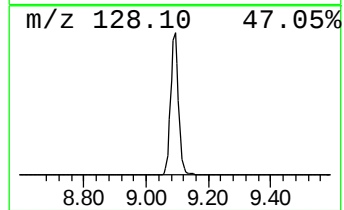
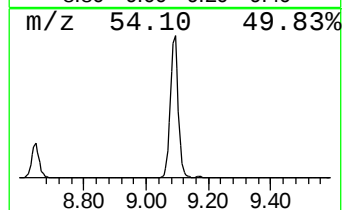
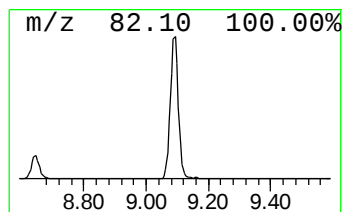
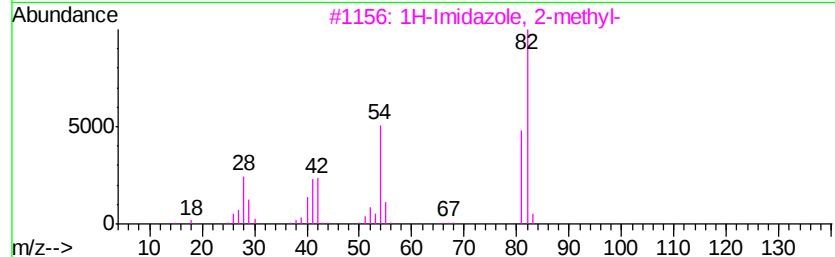
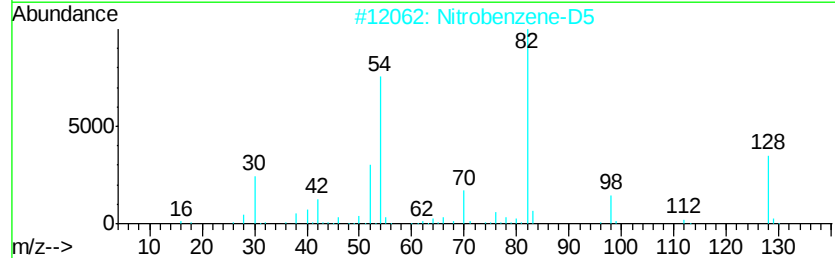
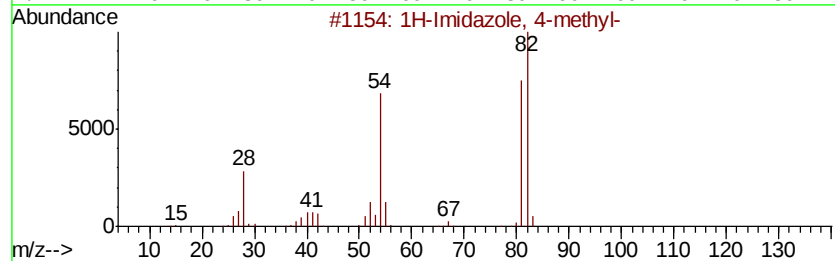
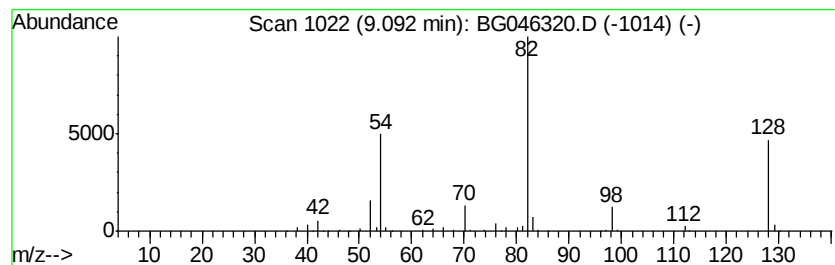
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Quant Title : SVOA CALIBRATION

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

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Peak Number 6 1H-Imidazole, 4-methyl- Concentration Rank 4

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 9.09 | 23.10 ng/ul | 511540 | 1,4-Dichlorobenzene-d4 | 7.94 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | 1H-Imidazole, 4-methyl- | 82  | C4H6N2  | 000822-36-6 | 50   |
| 2    |    | Nitrobenzene-D5         | 128 | C6D5NO2 | 004165-60-0 | 47   |
| 3    |    | 1H-Imidazole, 2-methyl- | 82  | C4H6N2  | 000693-98-1 | 40   |
| 4    |    | 1H-Imidazole, 2-methyl- | 82  | C4H6N2  | 000693-98-1 | 38   |
| 5    |    | Nitrobenzene-D5         | 128 | C6D5NO2 | 004165-60-0 | 37   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081720\  
Data File : BG046320.D  
Acq On : 18 Aug 2020 9:46  
Operator : CG/JU  
Sample : PB130999BL  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SBLK99

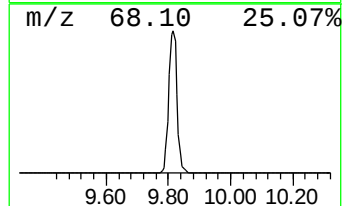
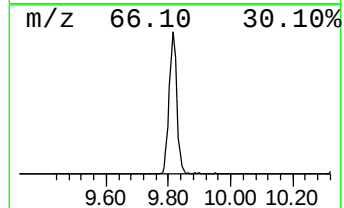
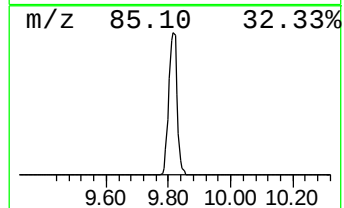
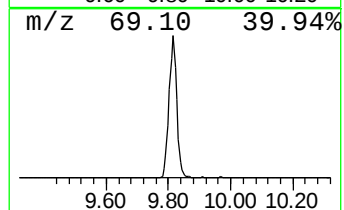
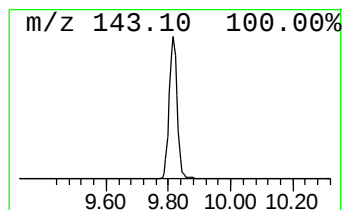
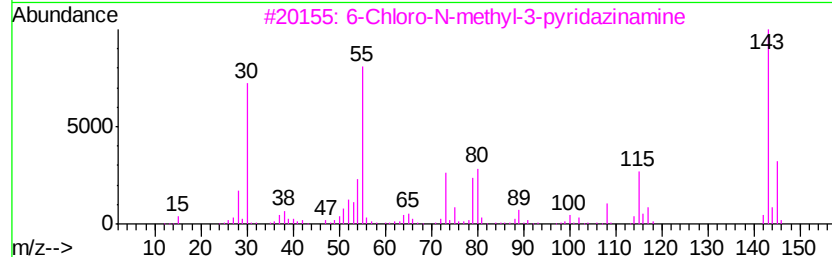
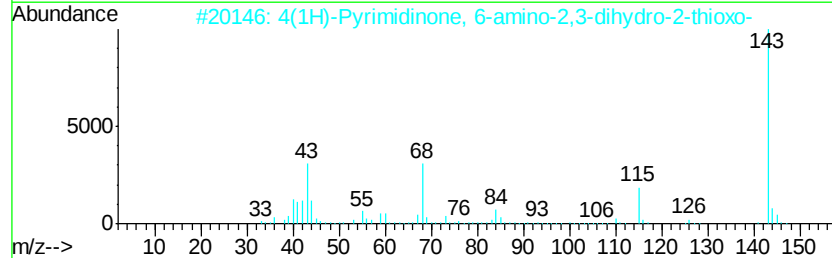
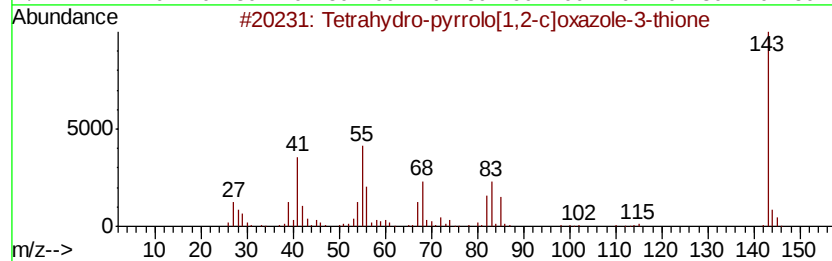
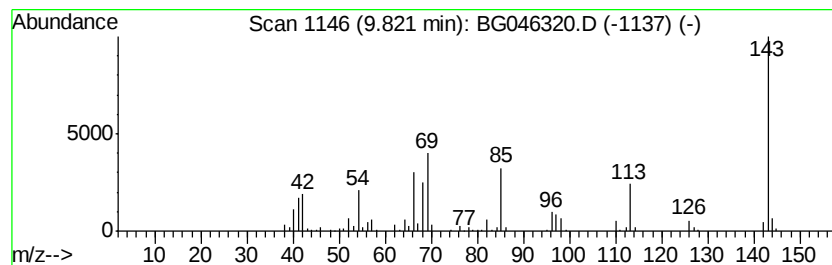
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 7 unknown-04 Concentration Rank 8

| R.T. | EstConc     | Area   | Relative to ISTD | R.T.  |
|------|-------------|--------|------------------|-------|
| 9.82 | 12.95 ng/ul | 431675 | Naphthalene-d8   | 10.74 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Tetrahydro-pyrrolo[1,2-c]oxazole... | 143 | C6H9NOS  | 1000190-53-6 | 35   |
| 2    |    | 4(1H)-Pyrimidinone, 6-amino-2,3-... | 143 | C4H5N3OS | 001004-40-6  | 16   |
| 3    |    | 6-Chloro-N-methyl-3-pyridazinamine  | 143 | C5H6ClN3 | 014959-32-1  | 12   |
| 4    |    | 1,2-Benzenedicarbonitrile, 4-amino- | 143 | C8H5N3   | 056765-79-8  | 10   |
| 5    |    | Glutaric acid, ethyl 3-heptyl ester | 258 | C14H26O4 | 1000359-73-0 | 10   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081720\  
Data File : BG046320.D  
Acq On : 18 Aug 2020 9:46  
Operator : CG/JU  
Sample : PB130999BL  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

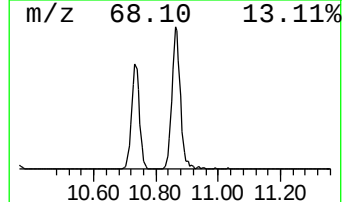
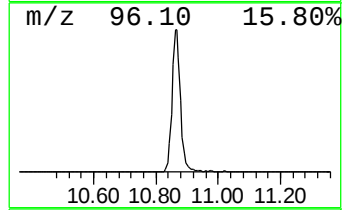
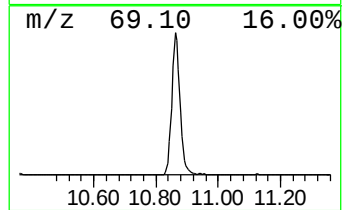
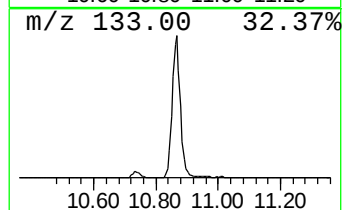
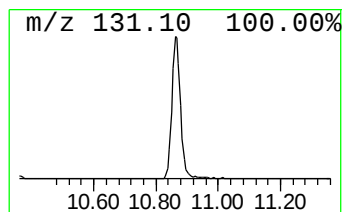
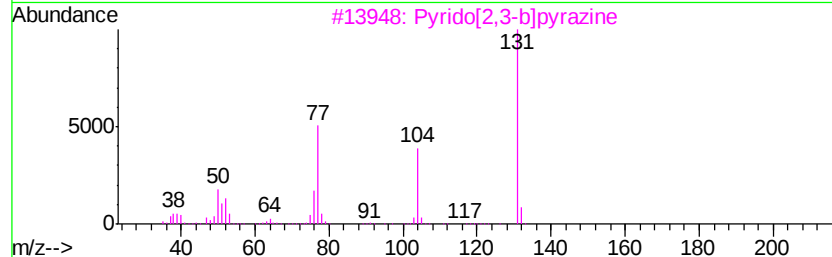
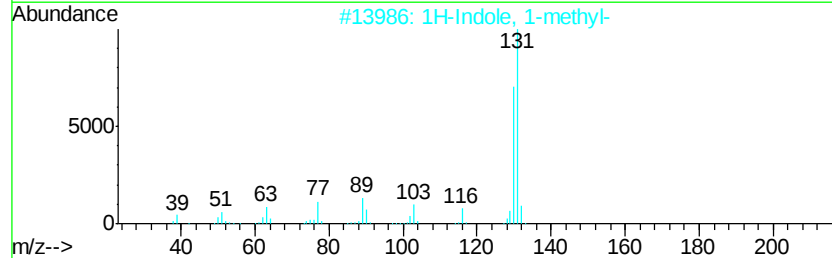
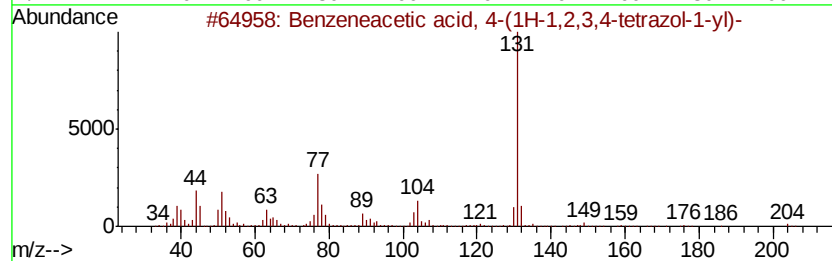
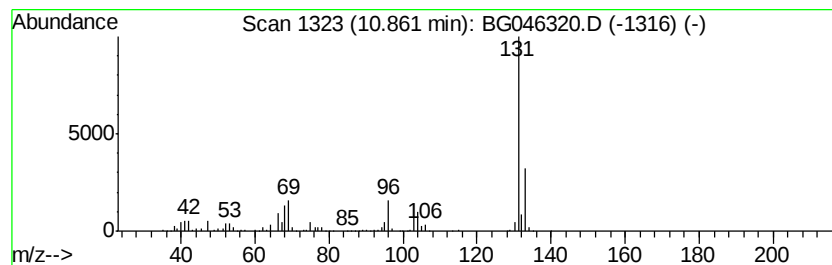
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 unknown-05 Concentration Rank 6

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 10.86 | 20.26 ng/ul | 675376 | Naphthalene-d8   | 10.74 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Benzeneacetic acid, 4-(1H-1,2,3,... | 204 | C9H8N4O2 | 1000351-56-5 | 33   |
| 2    |    | 1H-Indole, 1-methyl-                | 131 | C9H9N    | 000603-76-9  | 28   |
| 3    |    | Pyrido[2,3-b]pyrazine               | 131 | C7H5N3   | 000322-46-3  | 9    |
| 4    |    | 4-Pyrimidinamine, 2,6-difluoro-     | 131 | C4H3F2N3 | 000675-12-7  | 9    |
| 5    |    | 1H-Indole, 5-methyl-                | 131 | C9H9N    | 000614-96-0  | 9    |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081720\  
Data File : BG046320.D  
Acq On : 18 Aug 2020 9:46  
Operator : CG/JU  
Sample : PB130999BL  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

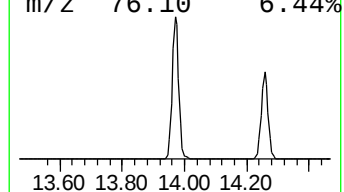
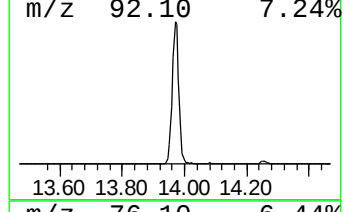
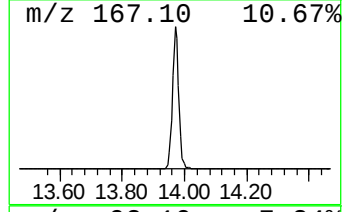
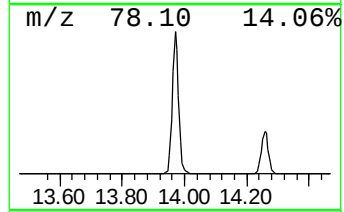
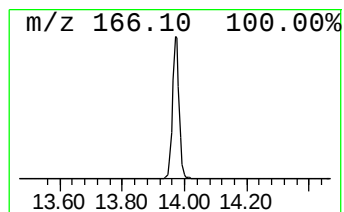
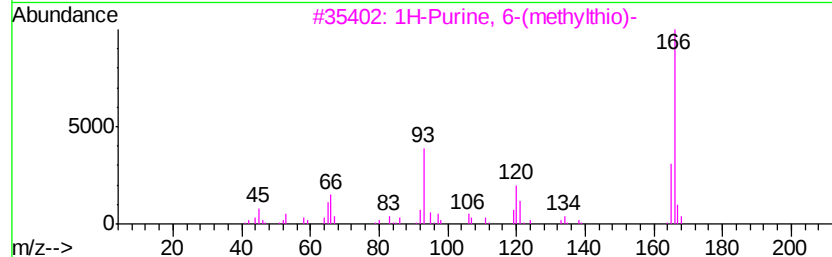
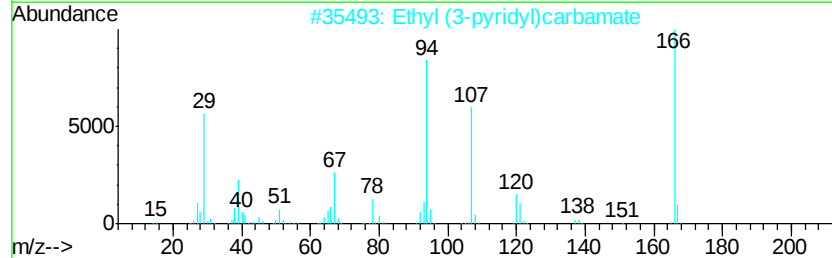
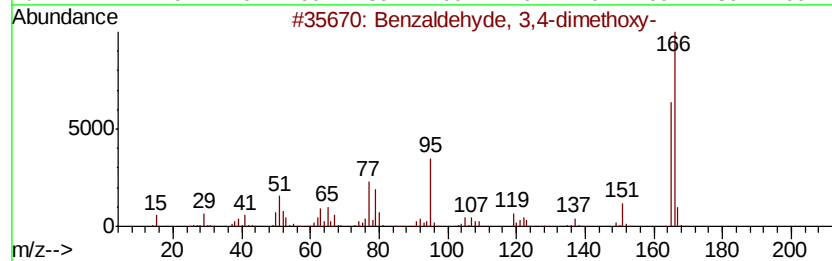
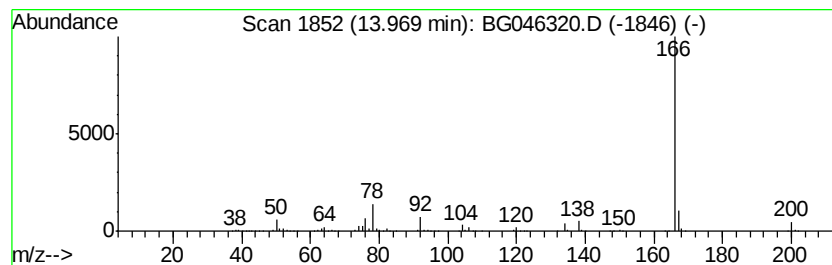
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ClientSampleId :  
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Quant Title : SVOA CALIBRATION

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

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Peak Number 9 unknown-06 Concentration Rank 5

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 13.97     | 22.54 ng/ul                         | 1092500 | Acenaphthene-d10 | 14.57       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | Benzaldehyde, 3,4-dimethoxy-        | 166     | C9H10O3          | 000120-14-9 | 45   |
| 2         | Ethyl (3-pyridyl)carbamate          | 166     | C8H10N2O2        | 006276-11-5 | 42   |
| 3         | 1H-Purine, 6-(methylthio)-          | 166     | C6H6N4S          | 000050-66-8 | 38   |
| 4         | 1H-Benzimidazole-2-ethanamine, 5... | 195     | C9H10ClN3        | 135875-16-0 | 38   |
| 5         | Benzenamine, N,N-dimethyl-4-nitro-  | 166     | C8H10N2O2        | 000100-23-2 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081720\  
Data File : BG046320.D  
Acq On : 18 Aug 2020 9:46  
Operator : CG/JU  
Sample : PB130999BL  
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ALS Vial : 27 Sample Multiplier: 1

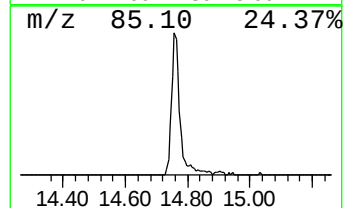
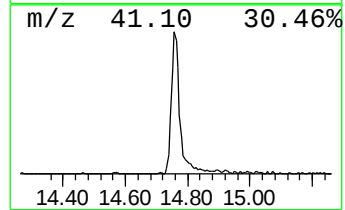
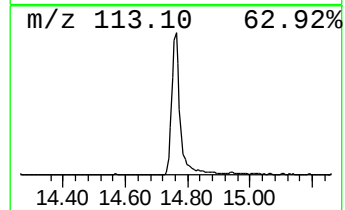
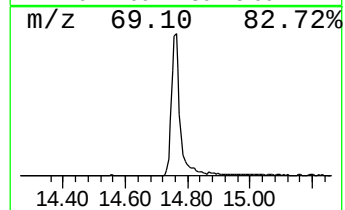
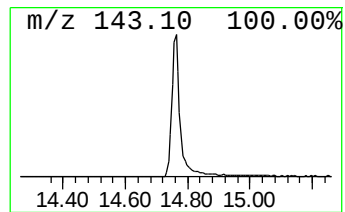
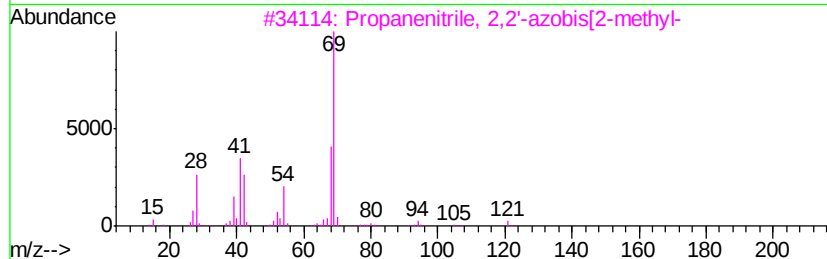
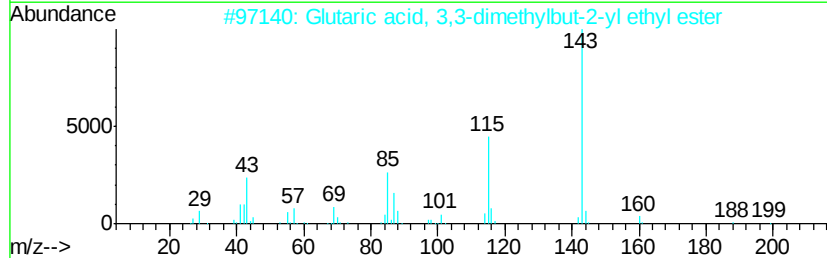
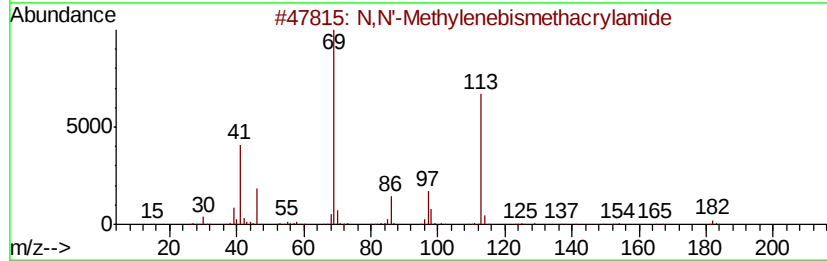
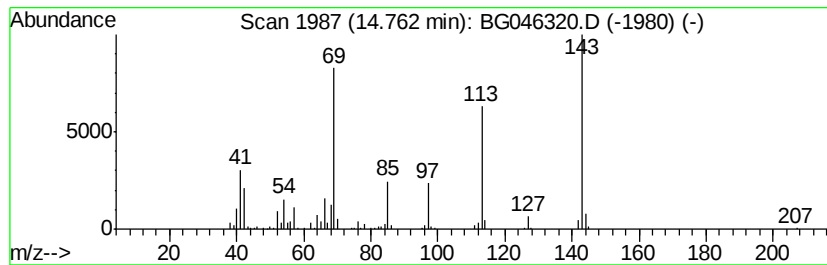
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ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 10 unknown-07 Concentration Rank 9

| R.T.      | EstConc                                                        | Area   | Relative to ISTD |              | R.T.  |
|-----------|----------------------------------------------------------------|--------|------------------|--------------|-------|
| 14.76     | 9.41 ng/ul                                                     | 456188 | Acenaphthene-d10 |              | 14.57 |
| Hit# of 5 | Tentative ID                                                   | MW     | MolForm          | CAS#         | Qual  |
| 1         | N,N'-Methylenebismethacrylamide                                | 182    | C9H14N2O2        | 002359-15-1  | 35    |
| 2         | Glutaric acid, 3,3-dimethylbut-2-ylidenehydrazide              | 244    | C13H24O4         | 1000359-74-7 | 25    |
| 3         | Propanenitrile, 2,2'-azobis[2-methylpropan-2-ylidenehydrazide] | 164    | C8H12N4          | 000078-67-1  | 22    |
| 4         | 2-Propenoic acid, 2-methyl-, 1,2-dichloroethyl ester           | 286    | C14H22O6         | 000109-16-0  | 17    |
| 5         | 6-Chloro-N-methyl-3-pyridazinamine                             | 143    | C5H6ClN3         | 014959-32-1  | 14    |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081720\  
Data File : BG046320.D  
Acq On : 18 Aug 2020 9:46  
Operator : CG/JU  
Sample : PB130999BL  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SBLK99

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

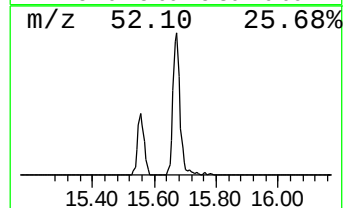
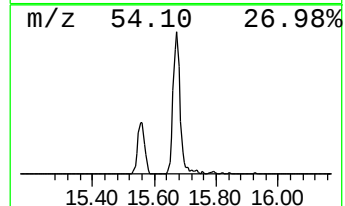
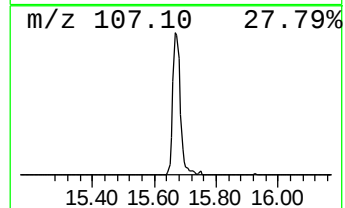
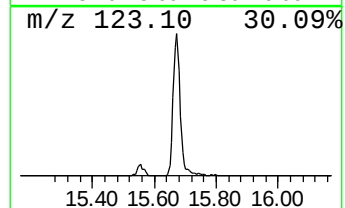
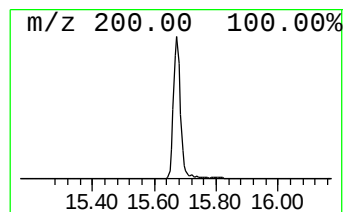
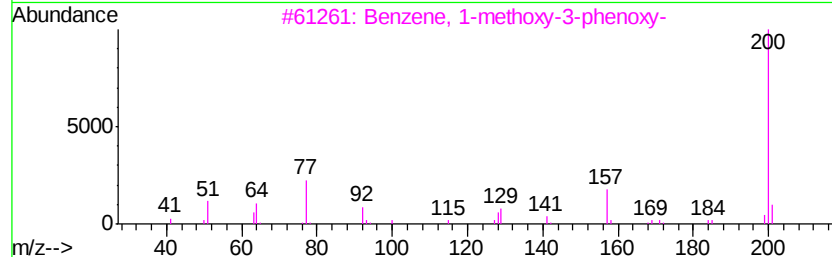
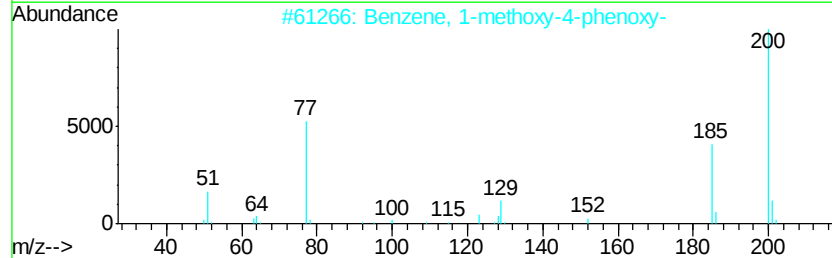
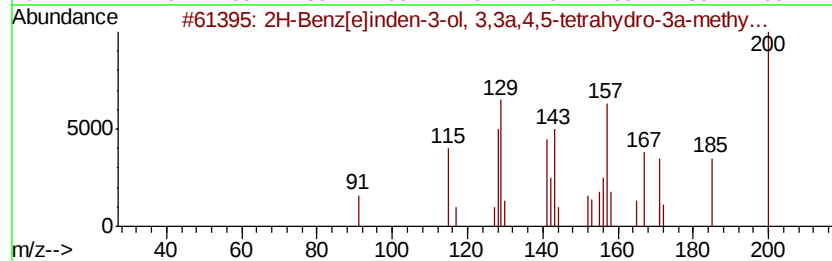
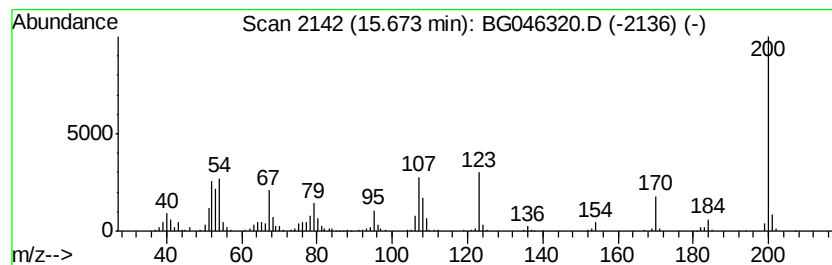
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 unknown-08 Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.67 | 9.13 ng/ul | 442623 | Acenaphthene-d10 | 14.57 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | 2H-Benz[e]inden-3-ol, 3,3a,4,5-t... | 200 | C14H16O   | 071805-91-9 | 38   |
| 2    |    | Benzene, 1-methoxy-4-phenoxy-       | 200 | C13H12O2  | 001655-69-2 | 14   |
| 3    |    | Benzene, 1-methoxy-3-phenoxy-       | 200 | C13H12O2  | 001655-68-1 | 14   |
| 4    |    | Benzenamine, 4,4'-oxybis-           | 200 | C12H12N2O | 000101-80-4 | 10   |
| 5    |    | 5-Hexyn-1-ol                        | 98  | C6H10O    | 000928-90-5 | 10   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG081720\  
Data File : BG046320.D  
Acq On : 18 Aug 2020 9:46  
Operator : CG/JU  
Sample : PB130999BL  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SBLK99

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.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 |
| Butane, 2-methoxy... | 3.14  | 2.1     | ng/ul | 46756    | 1                     | 7.94  | 442921 | 20.0 |
| Hexanoic acid, an... | 7.11  | 23.6    | ng/ul | 522796   | 1                     | 7.94  | 442921 | 20.0 |
| unknown-01           | 7.27  | 18.4    | ng/ul | 408320   | 1                     | 7.94  | 442921 | 20.0 |
| unknown-02           | 7.47  | 24.4    | ng/ul | 540511   | 1                     | 7.94  | 442921 | 20.0 |
| unknown-03           | 8.65  | 30.1    | ng/ul | 666710   | 1                     | 7.94  | 442921 | 20.0 |
| 1H-Imidazole, 4-m... | 9.09  | 23.1    | ng/ul | 511540   | 1                     | 7.94  | 442921 | 20.0 |
| unknown-04           | 9.82  | 12.9    | ng/ul | 431675   | 2                     | 10.74 | 666780 | 20.0 |
| unknown-05           | 10.86 | 20.3    | ng/ul | 675376   | 2                     | 10.74 | 666780 | 20.0 |
| unknown-06           | 13.97 | 22.5    | ng/ul | 1092500  | 3                     | 14.57 | 969550 | 20.0 |
| unknown-07           | 14.76 | 9.4     | ng/ul | 456188   | 3                     | 14.57 | 969550 | 20.0 |
| unknown-08           | 15.67 | 9.1     | ng/ul | 442623   | 3                     | 14.57 | 969550 | 20.0 |