

(QT Reviewed)

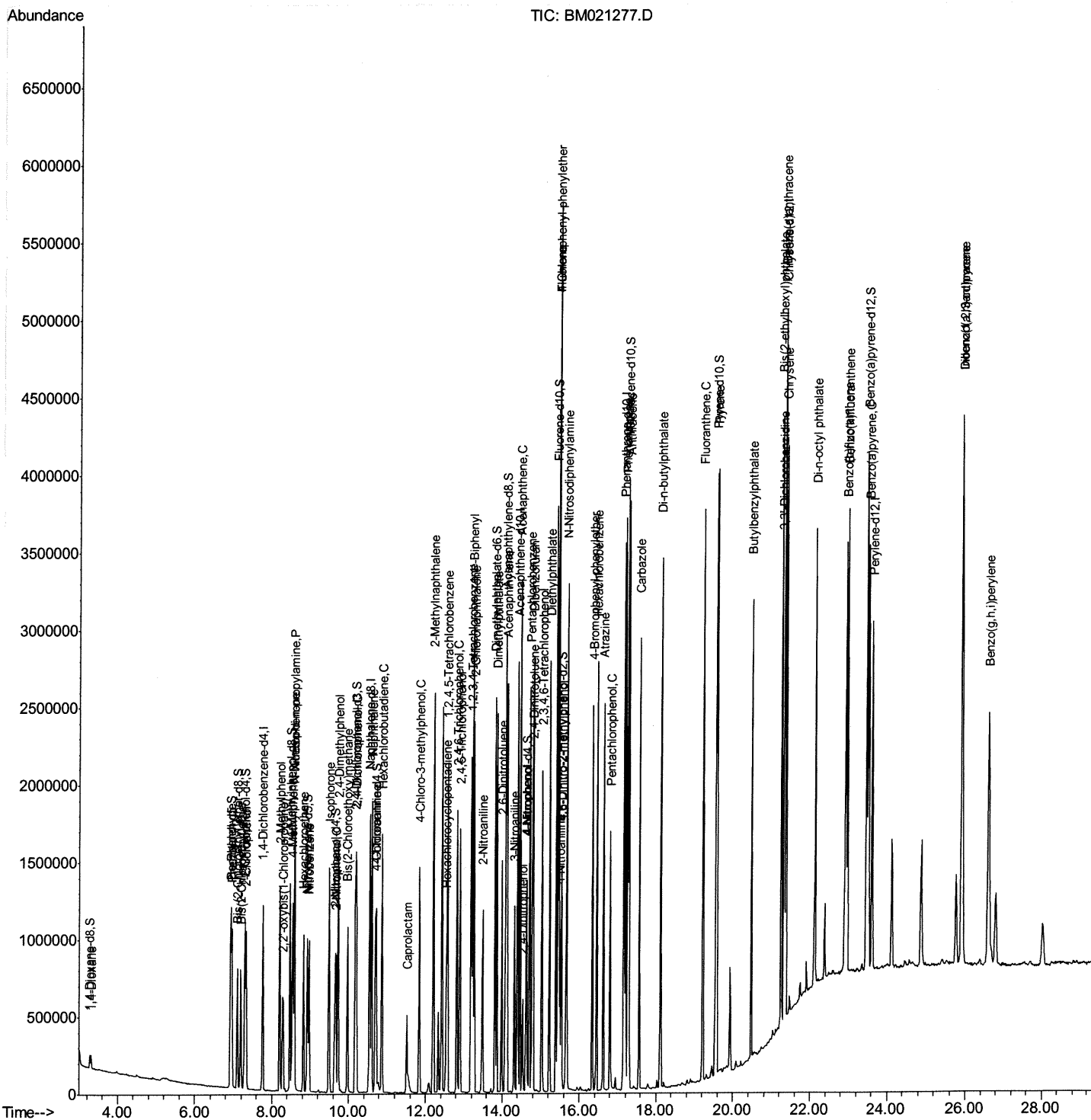
```
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062919\  
Data File : BM021277.D  
Acq On    : 30 Jun 2019   00:18  
Operator  : HP/JU  
Sample    : SSTDCCC020  
Misc      :  
ALS Vial  : 2      Sample Multiplier: 1
```

**Instrument :**  
BNA\_M  
**LabSampleId :**  
SSTD02054

## Manual Integrations

mohammad  
7/1/2019 2:43:34 PM

Quant Time: Jun 30 03:36:18 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sat Jun 29 23:19:36 2019  
Response via : Initial Calibration



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM062919\

Data File : BM021277.D

Acq On : 30 Jun 2019 00:18

Operator : HP/JU

Sample : SSTDCCC020

Misc :

ALS Vial : 2 Sample Multiplier: 1

Quant Time: Jun 30 00:50:42 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Sat Jun 29 23:19:36 2019

Response via : Initial Calibration

Instrument :

BNA\_M

LabSampleId :

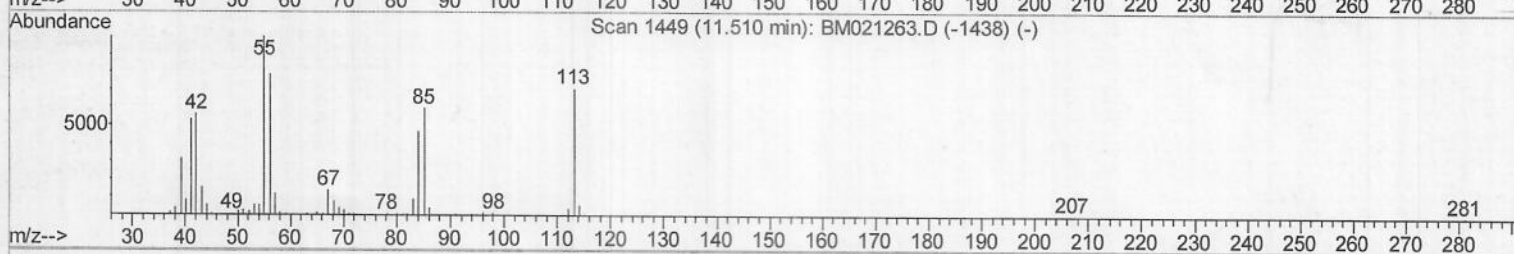
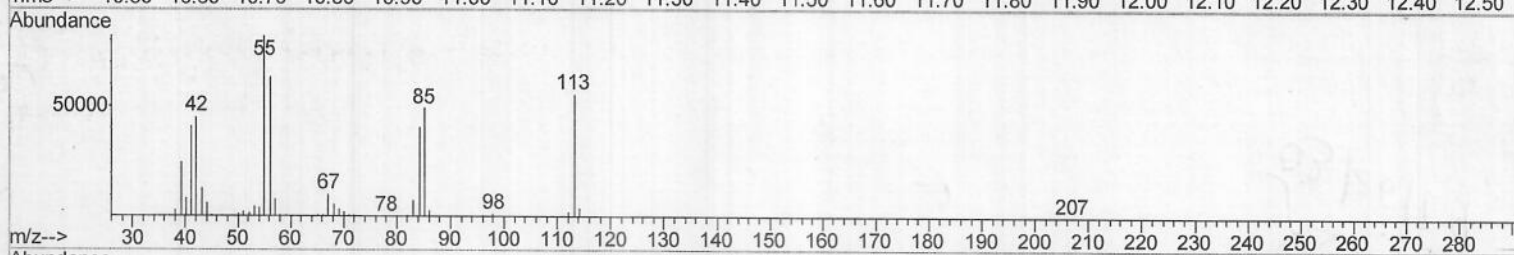
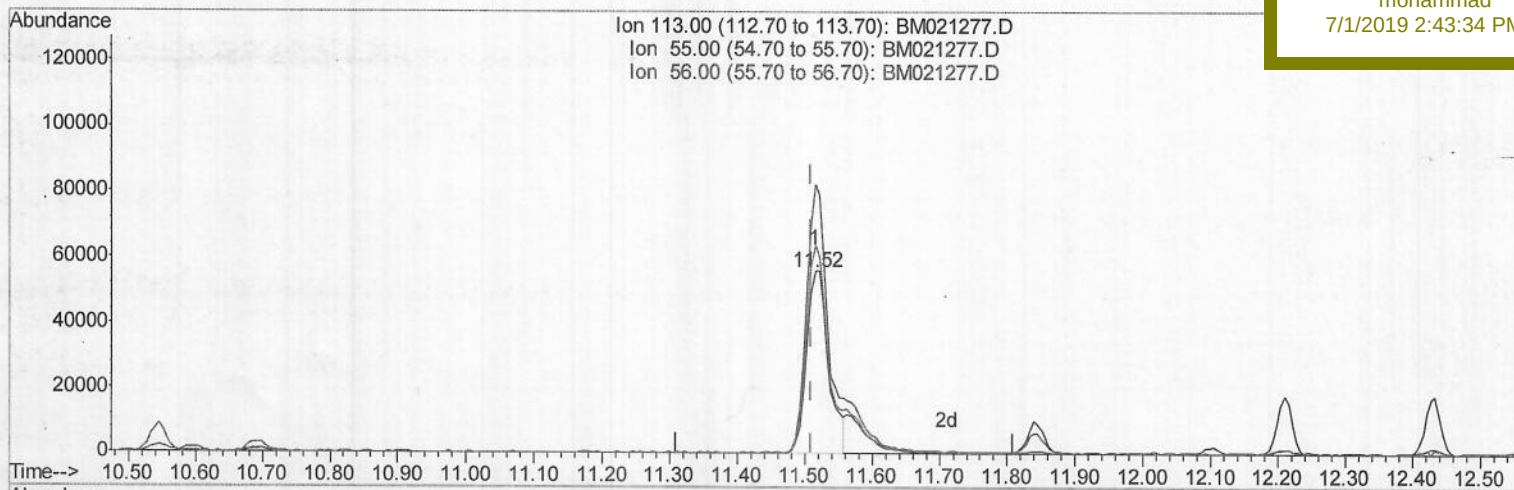
SSTD02054

Manual Integrations

APPROVED

mohammad

7/1/2019 2:43:34 PM



TIC: BM021277.D

(32) Caprolactam

11.516min (+0.006) 18.10ng/ul

response 127166

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 144.70 | 146.88 |
| 56.00  | 113.80 | 114.14 |
| 0.00   | 0.00   | 0.00   |

# Quantitation Report (Qedit)

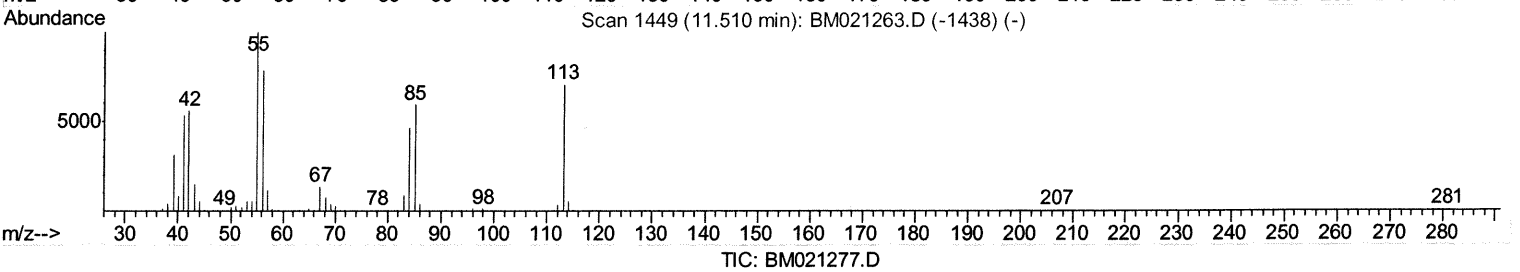
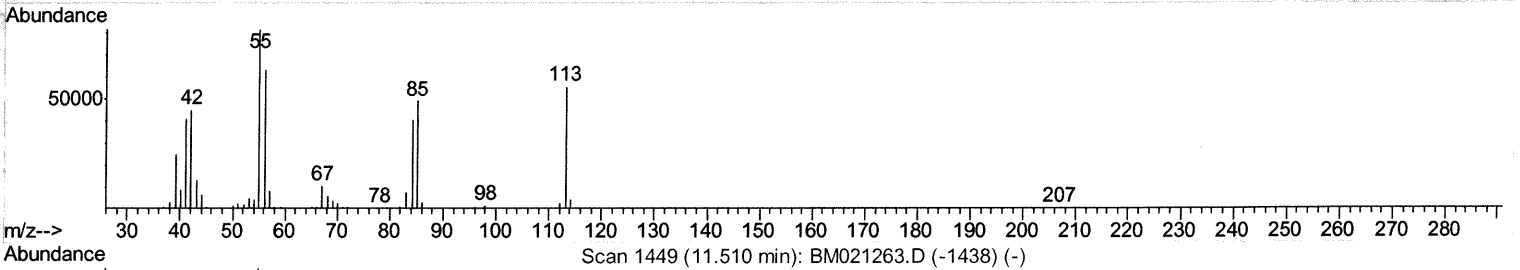
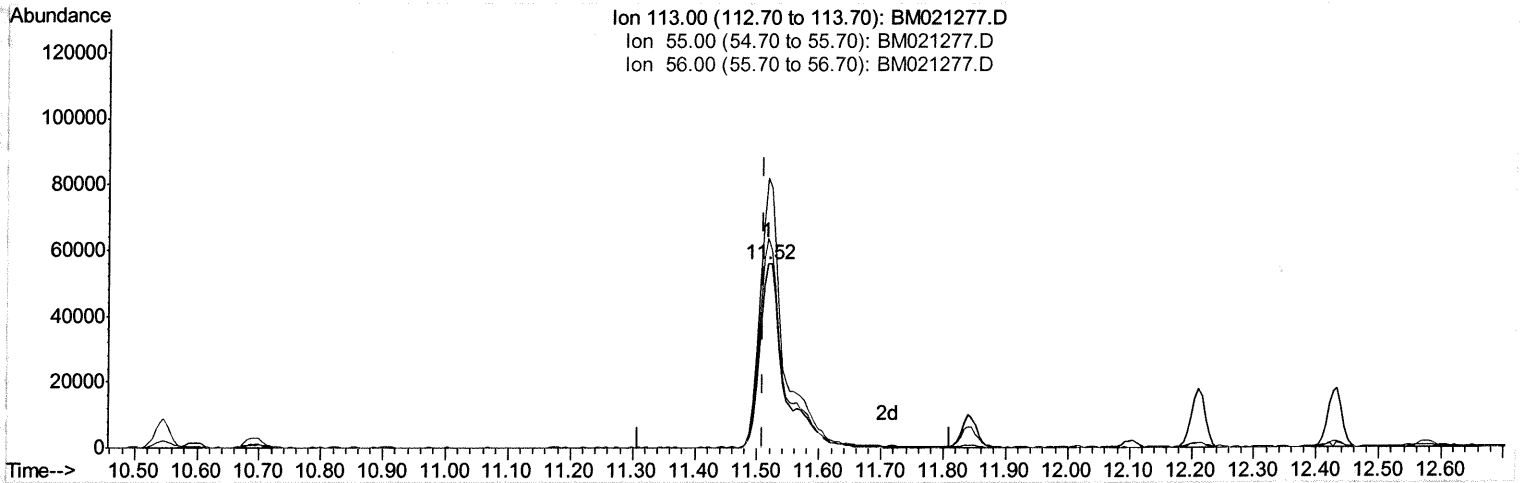
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM062919\  
 Data File : BM021277.D  
 Acq On : 30 Jun 2019 00:18  
 Operator : HP/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02054

Manual Integrations  
 APPROVED

mohammad  
 7/1/2019 2:43:34 PM

Quant Time: Jun 30 00:50:42 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Jun 29 23:19:36 2019  
 Response via : Initial Calibration



(32) Caprolactam

11.516min (+0.006) 21.99ng/ul m *Ju 07/01/19*

response 154500

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 144.70 | 146.88 |
| 56.00  | 113.80 | 114.14 |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM062919\  
 Data File : BM021277.D  
 Acq On : 30 Jun 2019 00:18  
 Operator : HP/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02054

Manual Integrations  
 APPROVED

mohammad  
 7/1/2019 2:43:34 PM

Quant Time: Jun 30 03:36:18 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Jun 29 23:19:36 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.76  | 152  | 336440   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.55 | 136  | 1497931  | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.40 | 164  | 947347   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.15 | 188  | 2042612  | 20.00 | ng/ul | 0.00     |
| 77) Chrysene-d12          | 21.33 | 240  | 1465171  | 20.00 | ng/ul | 0.00     |
| 85) Perylene-d12          | 23.60 | 264  | 1688413  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.28  | 96  | 44026   | 6.21  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.93  | 99  | 596150  | 20.36 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.10  | 67  | 338672  | 19.41 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.29  | 132 | 494296  | 20.77 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.47  | 113 | 506129  | 20.70 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.92  | 128 | 256496  | 21.09 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.64  | 143 | 309609  | 22.93 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.17 | 165 | 605296  | 22.05 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.69 | 131 | 560278  | 19.57 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.82 | 166 | 1739492 | 20.32 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.09 | 160 | 2174629 | 20.75 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.62 | 143 | 258394  | 18.56 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.40 | 176 | 1496200 | 20.09 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.53 | 200 | 253612  | 20.20 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.25 | 188 | 2165740 | 20.40 | ng/ul | 0.00 |
| 78) Pyrene-d10                 | 19.54 | 212 | 1932694 | 21.98 | ng/ul | 0.00 |
| 89) Benzo(a)pyrene-d12         | 23.46 | 264 | 2035943 | 20.44 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |         |              | Qvalue |
|--------------------------------|-------|-----|---------|--------------|--------|
| 2) 1,4-Dioxane                 | 3.31  | 88  | 46382   | 6.469 ng/uL  | 96     |
| 4) Benzaldehyde                | 6.92  | 77  | 225982  | 16.927 ng/ul | 99     |
| 6) Phenol                      | 6.96  | 94  | 553598  | 19.986 ng/ul | 98     |
| 8) Bis(2-Chloroethyl)ether     | 7.19  | 93  | 418074  | 19.776 ng/ul | 99     |
| 10) 2-Chlorophenol             | 7.32  | 128 | 461681  | 20.589 ng/ul | 98     |
| 11) 2-Methylphenol             | 8.20  | 108 | 444668  | 20.748 ng/ul | 97     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.28  | 45  | 519214  | 19.234 ng/ul | 99     |
| 14) Acetophenone               | 8.59  | 105 | 729856  | 20.746 ng/ul | 98     |
| 15) N-Nitroso-di-n-propylamine | 8.57  | 70  | 374701  | 20.464 ng/ul | 97     |
| 16) 4-Methylphenol             | 8.53  | 108 | 481138  | 20.630 ng/ul | 96     |
| 17) Hexachloroethane           | 8.82  | 117 | 184912  | 19.784 ng/ul | 97     |
| 20) Nitrobenzene               | 8.96  | 77  | 541900  | 19.832 ng/ul | 100    |
| 21) Isophorone                 | 9.49  | 82  | 1039265 | 20.196 ng/ul | 98     |
| 23) 2-Nitrophenol              | 9.67  | 139 | 293747  | 21.637 ng/ul | 99     |
| 24) 2,4-Dimethylphenol         | 9.73  | 107 | 565444  | 20.387 ng/ul | 99     |
| 25) Bis(2-Chloroethoxy)methane | 9.96  | 93  | 628190  | 19.809 ng/ul | 99     |
| 27) 2,4-Dichlorophenol         | 10.20 | 162 | 540880  | 21.641 ng/ul | 96     |
| 28) Naphthalene                | 10.60 | 128 | 1554909 | 20.218 ng/ul | 99     |
| 30) 4-Chloroaniline            | 10.72 | 127 | 528817  | 19.635 ng/ul | 99     |
| 31) Hexachlorobutadiene        | 10.86 | 225 | 354802  | 20.835 ng/ul | 98     |
| 32) Caprolactam                | 11.52 | 113 | 154500m | 21.989 ng/ul |        |
| 33) 4-Chloro-3-methylphenol    | 11.84 | 107 | 539866  | 21.187 ng/ul | 99     |
| 34) 2-Methylnaphthalene        | 12.21 | 142 | 1210477 | 20.527 ng/ul | 98     |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM062919\  
 Data File : BM021277.D  
 Acq On : 30 Jun 2019 00:18  
 Operator : HP/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02054

Manual Integrations  
 APPROVED

mohammad  
 7/1/2019 2:43:34 PM

Quant Time: Jun 30 03:36:18 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Jun 29 23:19:36 2019  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|-------|------|----------|--------|-------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 12.58 | 216  | 719914   | 21.287 | ng/ul | 99       |
| 37) Hexachlorocyclopentadiene  | 12.55 | 237  | 291898   | 14.508 | ng/ul | 97       |
| 38) 2,4,6-Trichlorophenol      | 12.83 | 196  | 471870   | 22.452 | ng/ul | 97       |
| 39) 2,4,5-Trichlorophenol      | 12.90 | 196  | 496156   | 22.101 | ng/ul | 98       |
| 40) 1,1'-Biphenyl              | 13.23 | 154  | 1595703  | 20.419 | ng/ul | 99       |
| 41) 2-Chloronaphthalene        | 13.27 | 162  | 1274278  | 20.639 | ng/ul | 100      |
| 42) 2-Nitroaniline             | 13.49 | 65   | 312181   | 20.805 | ng/ul | 98       |
| 44) Dimethylphthalate          | 13.86 | 163  | 1565412  | 20.090 | ng/ul | 99       |
| 45) 2,6-Dinitrotoluene         | 13.99 | 165  | 325181   | 21.865 | ng/ul | 92       |
| 47) Acenaphthylene             | 14.12 | 152  | 1862534  | 20.761 | ng/ul | 99       |
| 48) 3-Nitroaniline             | 14.32 | 138  | 301528   | 23.160 | ng/ul | 98       |
| 49) Acenaphthene               | 14.46 | 153  | 1334233  | 20.286 | ng/ul | 99       |
| 50) 2,4-Dinitrophenol          | 14.53 | 184  | 139982   | 18.736 | ng/ul | 95       |
| 52) 4-Nitrophenol              | 14.64 | 109  | 188095   | 18.264 | ng/ul | 97       |
| 53) Dibenzofuran               | 14.80 | 168  | 1881391  | 20.090 | ng/ul | 97       |
| 54) 2,4-Dinitrotoluene         | 14.78 | 165  | 461197   | 21.370 | ng/ul | 100      |
| 55) 2,3,4,6-Tetrachlorophenol  | 15.03 | 232  | 425616   | 21.741 | ng/ul | 99       |
| 56) Diethylphthalate           | 15.23 | 149  | 1529317  | 20.342 | ng/ul | 99       |
| 58) Fluorene                   | 15.45 | 166  | 1536420  | 20.193 | ng/ul | 99       |
| 59) 4-Chlorophenyl-phenylether | 15.45 | 204  | 845044   | 20.371 | ng/ul | 100      |
| 60) 4-Nitroaniline             | 15.49 | 138  | 280687   | 20.283 | ng/ul | 99       |
| 63) 4,6-Dinitro-2-methylphenol | 15.55 | 198  | 249820   | 19.820 | ng/ul | 95       |
| 64) N-Nitrosodiphenylamine     | 15.67 | 169  | 1318182  | 21.308 | ng/ul | 100      |
| 65) 4-Bromophenyl-phenylether  | 16.34 | 248  | 541888   | 21.634 | ng/ul | 99       |
| 66) Hexachlorobenzene          | 16.45 | 284  | 613377   | 21.366 | ng/ul | 97       |
| 67) Atrazine                   | 16.62 | 200  | 474735   | 21.153 | ng/ul | 97       |
| 68) Pentachlorophenol          | 16.80 | 266  | 321772   | 22.354 | ng/ul | 98       |
| 69) Phenanthrene               | 17.19 | 178  | 2277912  | 20.164 | ng/ul | 99       |
| 71) Anthracene                 | 17.29 | 178  | 2333137  | 20.267 | ng/ul | 99       |
| 72) 1,2,3,4-Tetrachlorobenzene | 13.19 | 216  | 716119   | 22.286 | ng/uL | 98       |
| 73) Pentachlorobenzene         | 14.72 | 250  | 723272   | 21.875 | ng/uL | 98       |
| 74) Carbazole                  | 17.56 | 167  | 1876713  | 19.781 | ng/ul | 100      |
| 75) Di-n-butylphthalate        | 18.12 | 149  | 2449215  | 21.502 | ng/ul | 100      |
| 76) Fluoranthene               | 19.21 | 202  | 2328407  | 19.076 | ng/ul | 100      |
| 79) Pyrene                     | 19.57 | 202  | 2269463  | 21.910 | ng/ul | 100      |
| 80) Butylbenzylphthalate       | 20.47 | 149  | 906041   | 23.630 | ng/ul | 95       |
| 81) 3,3'-Dichlorobenzidine     | 21.26 | 252  | 596912   | 18.177 | ng/ul | 99       |
| 82) Benzo(a)anthracene         | 21.32 | 228  | 2055707  | 20.580 | ng/ul | 99       |
| 83) Bis(2-ethylhexyl)phthalate | 21.24 | 149  | 1265201  | 23.176 | ng/ul | 100      |
| 84) Chrysene                   | 21.37 | 228  | 1901734  | 20.328 | ng/ul | 99       |
| 86) Di-n-octyl phthalate       | 22.13 | 149  | 1993007  | 20.485 | ng/ul | 100      |
| 87) Benzo(b)fluoranthene       | 22.92 | 252  | 2114068  | 19.766 | ng/ul | 100      |
| 88) Benzo(k)fluoranthene       | 22.97 | 252  | 2008297  | 19.518 | ng/ul | 99       |
| 90) Benzo(a)pyrene             | 23.51 | 252  | 2027676  | 20.143 | ng/ul | 98       |
| 91) Indeno(1,2,3-cd)pyrene     | 25.90 | 276  | 2589136  | 21.842 | ng/ul | 99       |
| 92) Dibenzo(a,h)anthracene     | 25.91 | 278  | 2163514  | 21.696 | ng/ul | 99       |
| 93) Benzo(g,h,i)perylene       | 26.60 | 276  | 2128258  | 21.799 | ng/ul | 98       |

(#) = qualifier out of range (m) = manual integration (+) = signals summed