

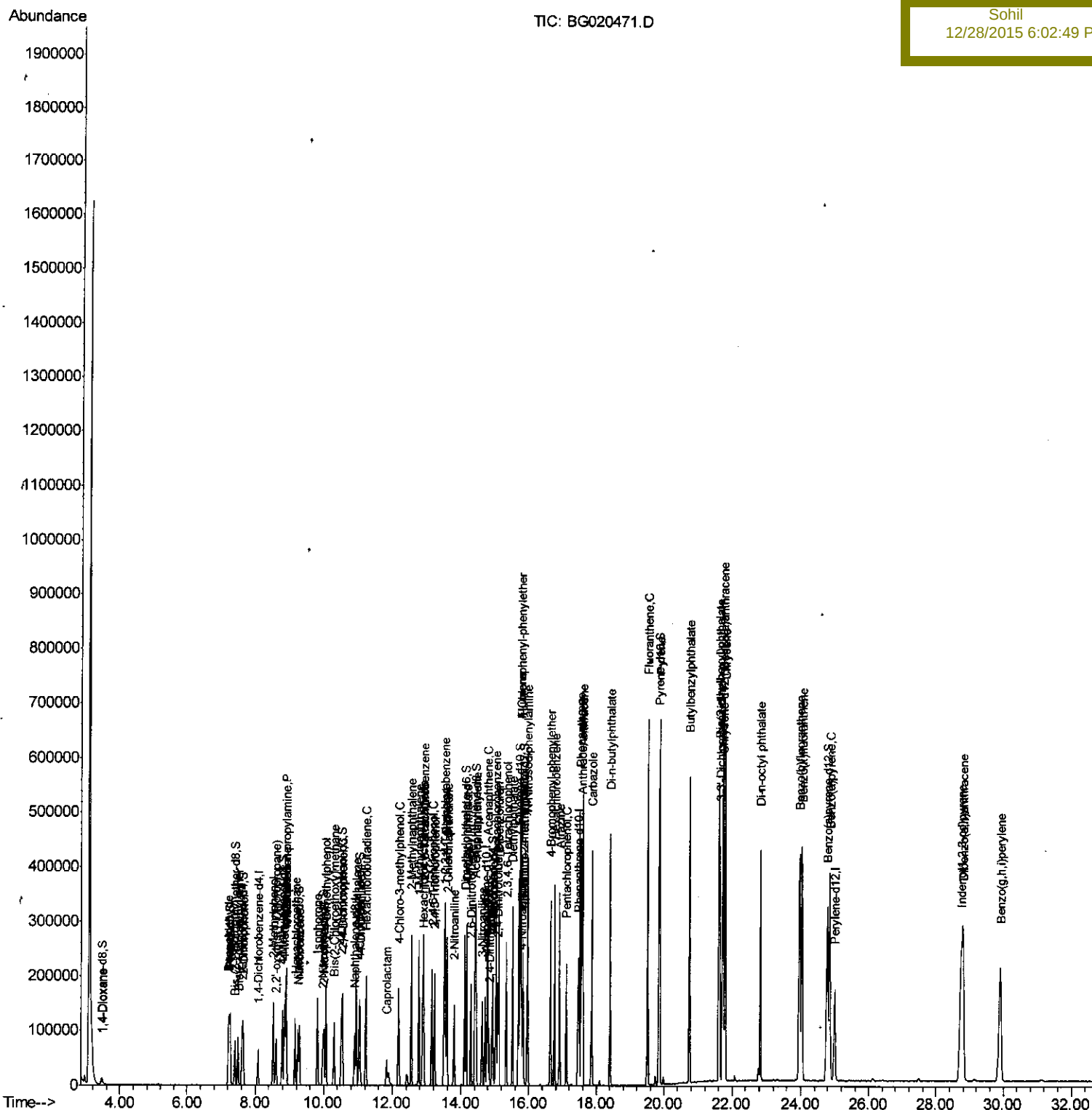
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Data Path   : Z:\HPCHEM1\BNA_G\Data\BG122415\  
Data File  : BG020471.D  
Acq On     : 24 Dec 2015   12:09  
Operator   : UM/SJ  
Sample     : SSTD04004  
Misc       :  
ALS Vial   : 5      Sample Multiplier: 1
```

**Instrument :**  
BNA\_G  
**ClientSampleId :**  
SSTD04004

## Manual Integrations

Sohil  
12/28/2015 6:02:49 PM

Quant Time: Dec 24 12:49:28 2015  
Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG122415.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Dec 24 12:16:41 2015  
Response via : Initial Calibration



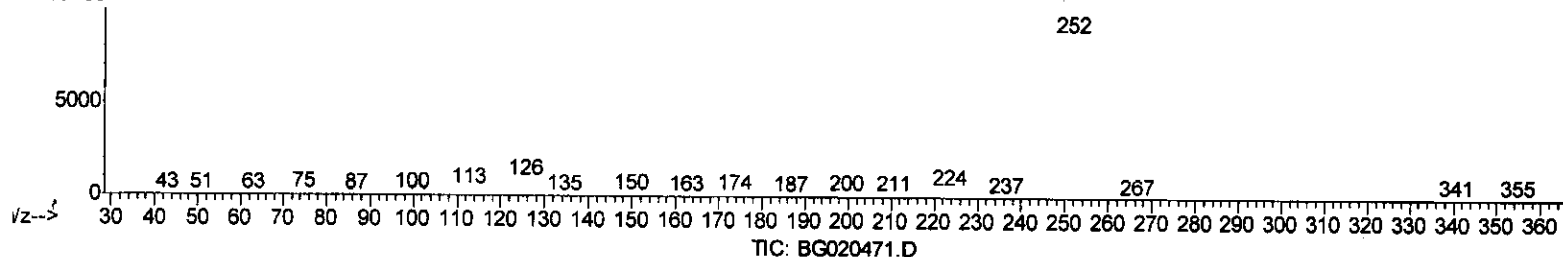
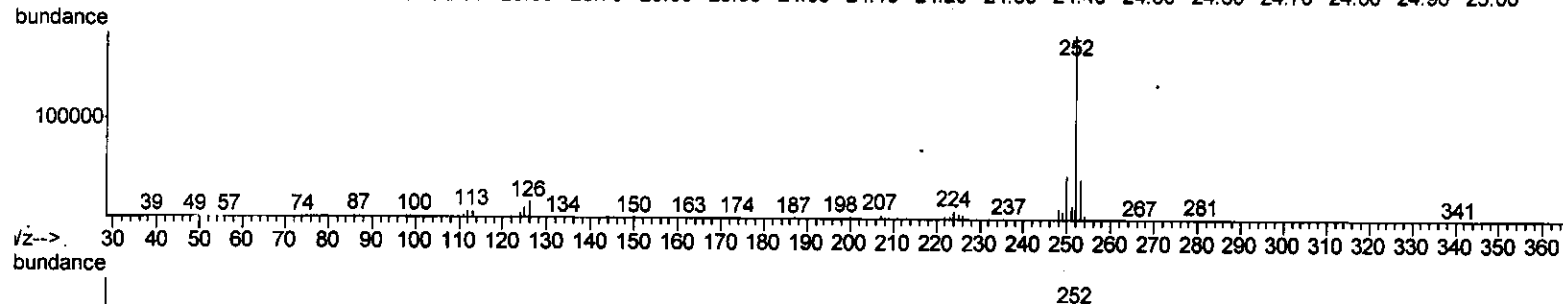
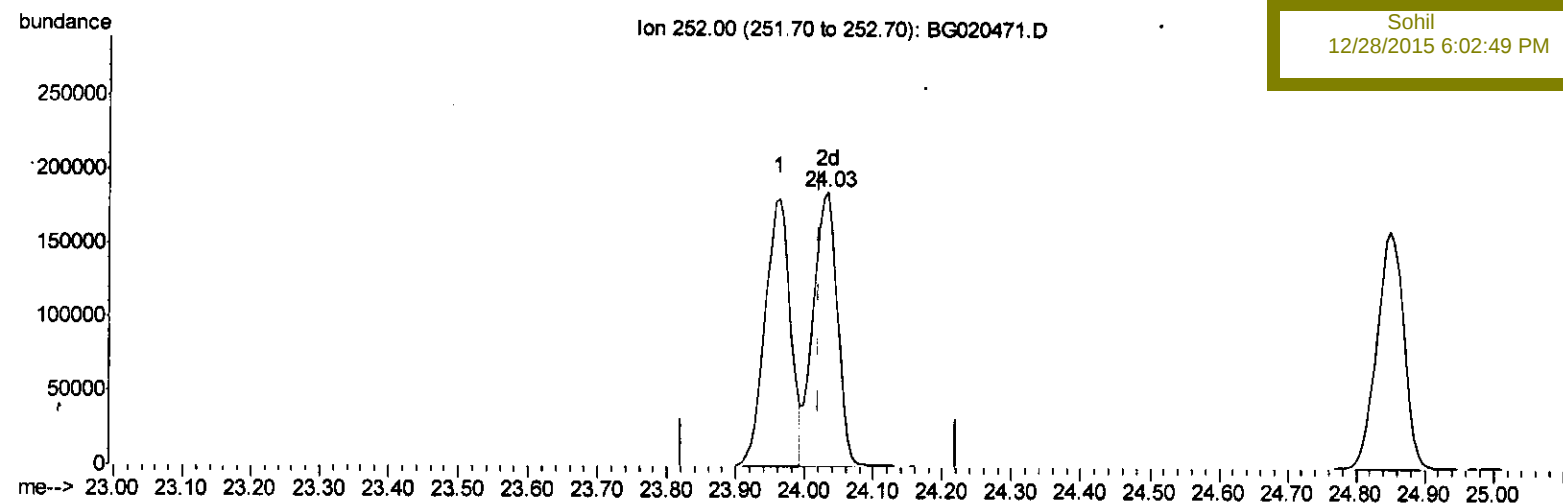
Data Path : Z:\HPCHEM1\BNA\_G\Data\BG122415\  
 Data File : BG020471.D  
 Acq On : 24 Dec 2015 12:09  
 Operator : UM/SJ  
 Sample : SST04004  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Dec 24 12:47:08 2015  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG122415.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Dec 24 12:16:41 2015  
 Response via : Initial Calibration

Instrument :  
 BNA\_G  
 ClientSampled :  
 SST04004

Manual Integrations  
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 12/28/2015 6:02:49 PM



(89) Benzo(k)fluoranthene

24.033min (+0.012) 39.19ng/ul m

response 455699

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 252.00 | 100   | 100   |
| 253.00 | 22.40 | 22.09 |
| 125.00 | 7.20  | 5.66# |
| 0.00   | 0.00  | 0.00  |

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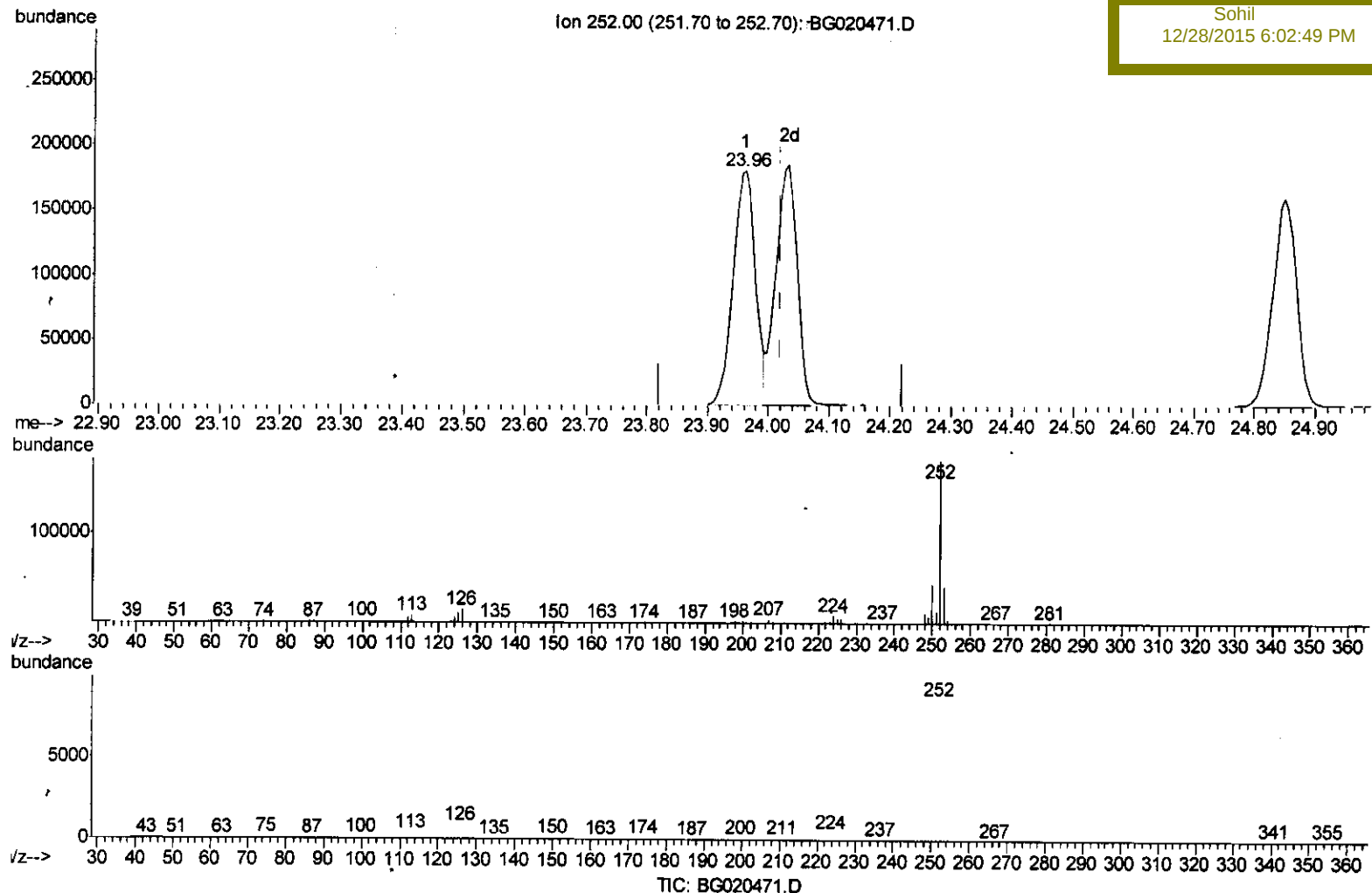
Data Path : Z:\HPCHEM1\BNA\_G\Data\BG122415\  
 Data File : BG020471.D  
 Acq On : 24 Dec 2015 12:09  
 Operator : UM/SJ  
 Sample : SST04004  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Dec 24 12:47:08 2015  
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 QLast Update : Thu Dec 24 12:16:41 2015  
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Instrument :  
 BNA\_G  
 ClientSampleId :  
 SST04004

Manual Integrations  
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 12/28/2015 6:02:49 PM



(89) Benzo(k)fluoranthene

23.962min (-0.059) 39.71ng/ul

response 461747

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 252.00 | 100   | 100   |
| 253.00 | 22.40 | 22.72 |
| 125.00 | 7.20  | 5.95  |
| 0.00   | 0.00  | 0.00  |

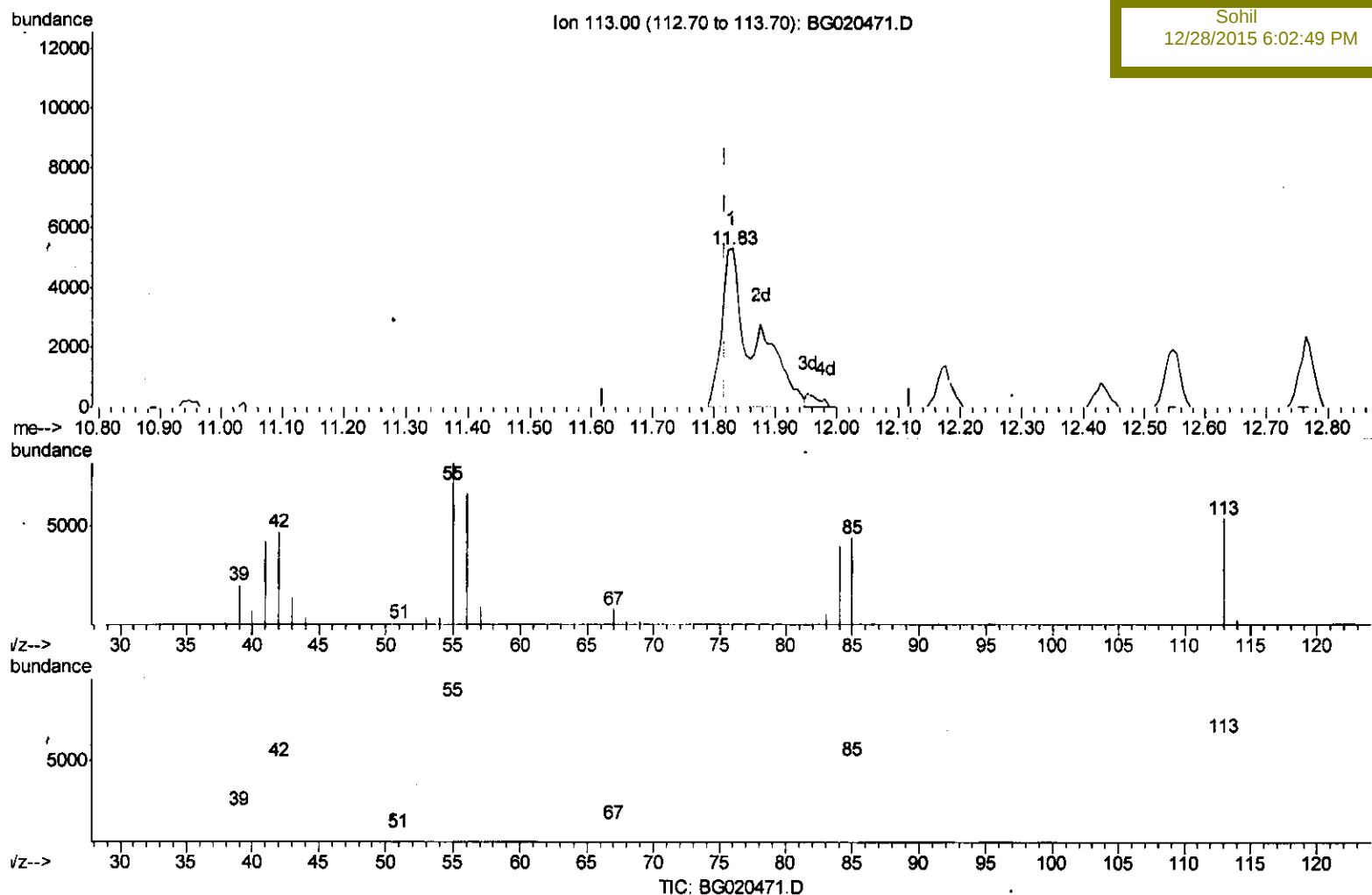
Data Path : Z:\HPCHEM1\BNA\_G\Data\BG122415\  
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 Sample : SSTD04004  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Dec 24 12:47:08 2015  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG122415.M  
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 Response via : Initial Calibration

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD04004

Manual Integrations  
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 12/28/2015 6:02:49 PM



(32) Caprolactam

11.829min (+0.012) 43.56ng/ul m

response 19105

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 165.10 | 151.58 |
| 56.00  | 129.10 | 123.02 |
| 0.00   | 0.00   | 0.00   |

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 12/25/15

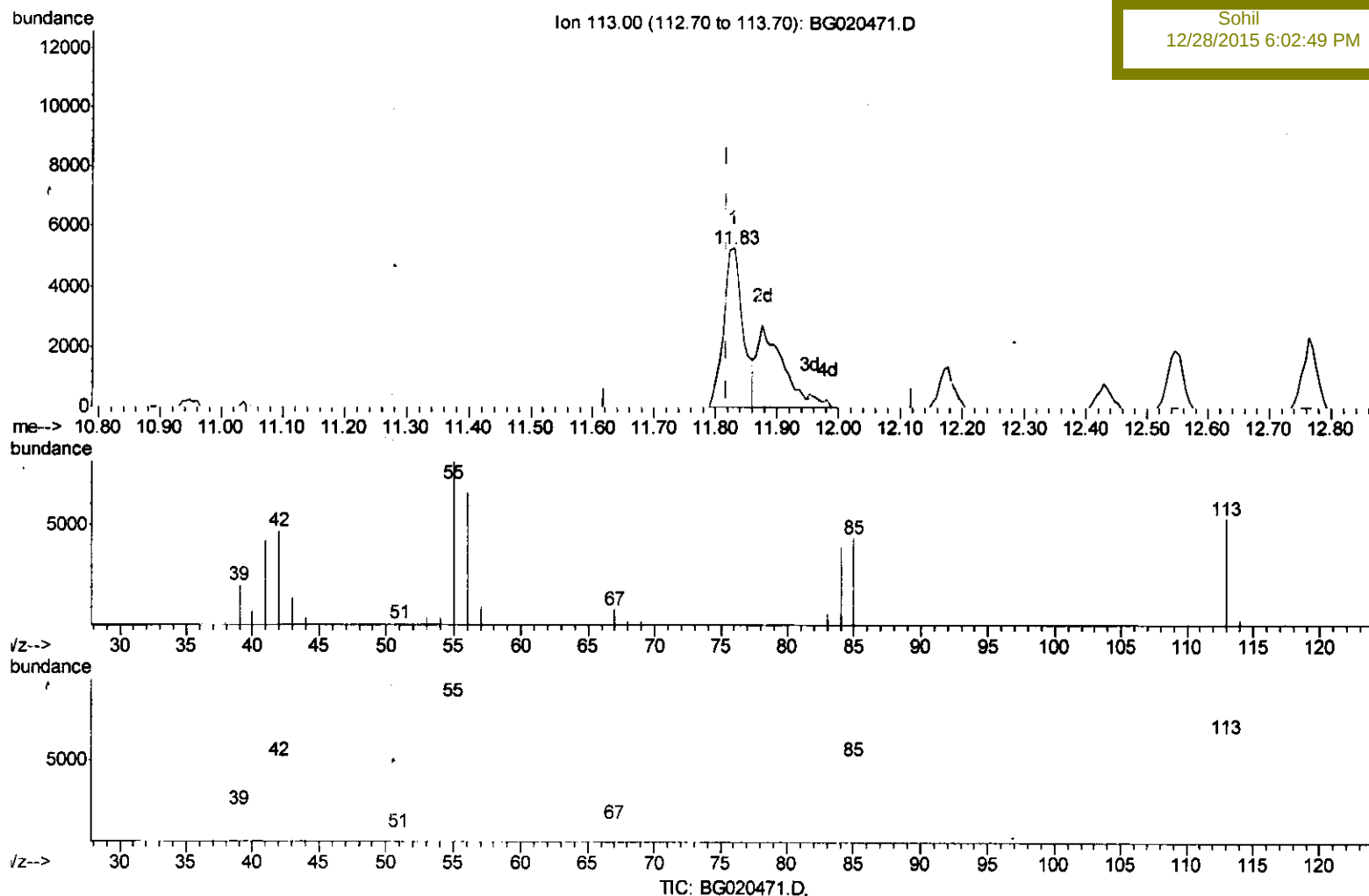
Data Path : Z:\HPCHEM1\BNA\_G\Data\BG122415\  
 Data File : BG020471.D  
 Acq On : 24 Dec 2015 12:09  
 Operator : UM/SJ  
 Sample : SSTD04004  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Dec 24 12:47:08 2015  
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Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD04004

Manual Integrations  
 APPROVED

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 12/28/2015 6:02:49 PM



(32) Caprolactam

11.829min (+0.012) 26.11ng/ul

response 11450

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 165.10 | 151.58 |
| 56.00  | 129.10 | 123.02 |
| 0.00   | 0.00   | 0.00   |

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 BNA\_G  
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 SSTD04004

Manual Integrations  
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 12/28/2015 6:02:49 PM

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 35) 1-Methylnaphthalene        | 12.76 | 142  | 126335   | 39.69 | ng/ul  | 97       |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.91 | 216  | 95979    | 39.89 | ng/ul  | 98       |
| 38) Hexachlorocyclopentadiene  | 12.89 | 237  | 53633    | 53.62 | ng/ul  | 95       |
| 39) 2,4,6-Trichlorophenol      | 13.15 | 196  | 58382    | 41.93 | ng/ul  | 98       |
| 40) 2,4,5-Trichlorophenol      | 13.23 | 196  | 63641    | 50.83 | ng/ul  | 95       |
| 41) 1,1'-Biphenyl              | 13.55 | 154  | 182951   | 39.94 | ng/ul  | 96       |
| 42) 2-Chloronaphthalene        | 13.59 | 162  | 146764   | 39.42 | ng/ul  | 98       |
| 43) 2-Nitroaniline             | 13.80 | 65   | 44360    | 45.03 | ng/ul  | 95       |
| 45) Dimethylphthalate          | 14.16 | 163  | 197922   | 39.41 | ng/ul  | 99       |
| 46) 2,6-Dinitrotoluene         | 14.29 | 165  | 42623    | 43.67 | ng/ul  | 98       |
| 48) Acenaphthylene             | 14.44 | 152  | 238005   | 39.74 | ng/ul  | 99       |
| 49) 3-Nitroaniline             | 14.62 | 138  | 40584    | 44.91 | ng/ul  | 90       |
| 50) Acenaphthene               | 14.78 | 153  | 160884   | 39.77 | ng/ul  | 99       |
| 51) 2,4-Dinitrophenol          | 14.83 | 184  | 27461    | 61.20 | ng/ul  | 93       |
| 53) 4-Nitrophenol              | 14.94 | 109  | 31257    | 49.18 | ng/ul  | 94       |
| 54) Dibenzofuran               | 15.11 | 168  | 237912   | 39.34 | ng/ul  | 97       |
| 55) 2,4-Dinitrotoluene         | 15.08 | 165  | 62342    | 41.88 | ng/ul  | 92       |
| 56) 2,3,4,6-Tetrachlorophenol  | 15.34 | 232  | 63806    | 42.81 | ng/ul# | 98       |
| 57) Diethylphthalate           | 15.53 | 149  | 203660   | 39.60 | ng/ul  | 98       |
| 59) Fluorene                   | 15.76 | 166  | 192968   | 39.29 | ng/ul  | 99       |
| 60) 4-Chlorophenyl-phenylether | 15.75 | 204  | 110745   | 39.05 | ng/ul  | 96       |
| 61) 4-Nitroaniline             | 15.79 | 138  | 44211    | 45.04 | ng/ul  | 97       |
| 64) 4,6-Dinitro-2-methylphenol | 15.84 | 198  | 45137    | 46.82 | ng/ul  | 95       |
| 65) N-Nitrosodiphenylamine     | 15.96 | 169  | 173335   | 39.82 | ng/ul  | 98       |
| 66) 4-Bromophenyl-phenylether  | 16.64 | 248  | 78558    | 41.19 | ng/ul  | 98       |
| 67) Hexachlorobenzene          | 16.76 | 284  | 85639    | 39.41 | ng/ul# | 89       |
| 68) Atrazine                   | 16.91 | 200  | 76733    | 40.34 | ng/ul  | 99       |
| 69) Pentachlorophenol          | 17.11 | 266  | 46933    | 47.20 | ng/ul  | 95       |
| 70) Phenanthrene               | 17.51 | 178  | 339993   | 38.01 | ng/ul  | 100      |
| 72) Anthracene                 | 17.59 | 178  | 346978   | 38.29 | ng/ul  | 99       |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.52 | 216  | 104859   | 40.81 | ng/uL  | 99       |
| 74) Pentachlorobenzene         | 15.03 | 250  | 95488    | 39.87 | ng/uL  | 90       |
| 75) Carbazole                  | 17.86 | 167  | 302475   | 39.99 | ng/ul  | 99       |
| 76) Di-n-butylphthalate        | 18.41 | 149  | 347653   | 38.17 | ng/ul  | 100      |
| 77) Fluoranthene               | 19.51 | 202  | 438925   | 37.98 | ng/ul  | 97       |
| 80) Pyrene                     | 19.87 | 202  | 442648   | 37.89 | ng/ul  | 97       |
| 81) Butylbenzylphthalate       | 20.74 | 149  | 162869   | 40.57 | ng/ul  | 93       |
| 82) 3,3'-Dichlorobenzidine     | 21.63 | 252  | 166416   | 41.46 | ng/ul  | 96       |
| 83) Benzo(a)anthracene         | 21.72 | 228  | 465743   | 39.00 | ng/ul  | 98       |
| 84) Bis(2-ethylhexyl)phthalate | 21.60 | 149  | 233612   | 39.47 | ng/ul  | 99       |
| 85) Chrysene                   | 21.78 | 228  | 437517   | 38.05 | ng/ul  | 98       |
| 87) Di-n-octyl phthalate       | 22.83 | 149  | 406305   | 39.33 | ng/ul  | 100      |
| 88) Benzo(b)fluoranthene       | 23.96 | 252  | 461747   | 38.83 | ng/ul# | 96       |
| 89) Benzo(k)fluoranthene       | 24.03 | 252  | 455699m  | 39.19 | ng/ul  |          |
| 91) Benzo(a)pyrene             | 24.85 | 252  | 444507   | 38.82 | ng/ul  | 97       |
| 92) Indeno(1,2,3-cd)pyrene     | 28.73 | 276  | 540553   | 39.99 | ng/ul  | 95       |
| 93) Dibenzo(a,h)anthracene     | 28.80 | 278  | 453184   | 40.57 | ng/ul  | 98       |
| 94) Benzo(g,h,i)perylene       | 29.91 | 276  | 443372   | 39.94 | ng/ul  | 98       |

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 QLast Update : Thu Dec 24 12:16:41 2015  
 Response via : Initial Calibration

Instrument :  
 BNA\_G  
 Client Sample Id :  
 SSTD04004

Manual Integrations  
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 12/28/2015 6:02:49 PM

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.08  | 152  | 20770    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.90 | 136  | 84747    | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.71 | 164  | 58975    | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.46 | 188  | 155103   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.74 | 240  | 199382   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 25.00 | 264  | 195663   | 20.00 | ng/ul | 0.00     |

#### System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.45  | 96  | 6243   | 41.03 | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.24  | 99  | 66362  | 43.80 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.40  | 67  | 40255  | 42.88 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.61  | 132 | 52643  | 42.33 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.79  | 113 | 56155  | 43.56 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.25  | 128 | 27466  | 41.93 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.97  | 143 | 32324  | 44.71 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.52 | 165 | 60541  | 42.35 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.03 | 131 | 71834  | 44.76 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 14.11 | 166 | 193618 | 39.16 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.41 | 160 | 224540 | 39.39 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.92 | 143 | 34382  | 49.09 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.71 | 176 | 175176 | 39.19 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.82 | 200 | 41925  | 47.24 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.56 | 188 | 288367 | 38.98 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.84 | 212 | 347148 | 39.67 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 24.78 | 264 | 394721 | 39.71 | ng/ul | 0.01 |

#### Target Compounds

|                                |       |     |        |       |       | Qvalue |
|--------------------------------|-------|-----|--------|-------|-------|--------|
| 2) 1,4-Dioxane                 | 3.49  | 88  | 6881   | 39.86 | ng/uL | 97     |
| 4) Benzaldehyde                | 7.21  | 77  | 44902  | 42.91 | ng/ul | 97     |
| 6) Phenol                      | 7.26  | 94  | 70219  | 44.98 | ng/ul | 99     |
| 8) Bis(2-Chloroethyl)ether     | 7.49  | 93  | 52890  | 43.33 | ng/ul | 95     |
| 10) 2-Chlorophenol             | 7.64  | 128 | 54572  | 41.69 | ng/ul | 97     |
| 11) 2-Methylphenol             | 8.52  | 108 | 54194  | 44.52 | ng/ul | 97     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.61  | 40  | 66027  | 40.70 | ng/ul | 96     |
| 14) Acetophenone               | 8.91  | 105 | 89976  | 42.20 | ng/ul | 95     |
| 15) N-Nitroso-di-n-propylamine | 8.89  | 70  | 46433  | 41.86 | ng/ul | 97     |
| 16) 4-Methylphenol             | 8.86  | 108 | 58325  | 43.09 | ng/ul | 94     |
| 17) Hexachloroethane           | 9.17  | 117 | 22941  | 41.06 | ng/ul | 99     |
| 20) Nitrobenzene               | 9.29  | 77  | 67998  | 41.29 | ng/ul | 98     |
| 21) Isophorone                 | 9.81  | 82  | 126451 | 40.52 | ng/ul | 99     |
| 23) 2-Nitrophenol              | 10.01 | 139 | 33498  | 43.37 | ng/ul | 93     |
| 24) 2,4-Dimethylphenol         | 10.06 | 107 | 70109  | 42.18 | ng/ul | 98     |
| 25) Bis(2-Chloroethoxy)methane | 10.30 | 93  | 69121  | 39.64 | ng/ul | 95     |
| 27) 2,4-Dichlorophenol         | 10.54 | 162 | 61958  | 42.28 | ng/ul | 96     |
| 28) Naphthalene                | 10.95 | 128 | 179624 | 39.39 | ng/ul | 99     |
| 30) 4-Chloroaniline            | 11.05 | 127 | 71618  | 43.78 | ng/ul | 97     |
| 31) Hexachlorobutadiene        | 11.22 | 225 | 51231  | 40.83 | ng/ul | 96     |
| 32) Caprolactam                | 11.83 | 113 | 19105m | 43.56 | ng/ul | 93     |
| 33) 4-Chloro-3-methylphenol    | 12.18 | 107 | 65509  | 42.77 | ng/ul | 93     |
| 34) 2-Methylnaphthalene        | 12.55 | 142 | 133627 | 40.01 | ng/ul | 96     |

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Instrument :  
BNA\_G  
ClientSampleId :  
SSTD04004

Manual Integrations  
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Sohil  
12/28/2015 6:02:49 PM

| Internal Standards                                                          | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------------------------------------------------------|------|------|----------|------|-------|----------|
| -----                                                                       |      |      |          |      |       |          |
| (# ) = qualifier out of range (m) = manual integration (+) = signals summed |      |      |          |      |       |          |