

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050716\  
 Data File : BF086815.D  
 Acq On : 8 May 2016 11:03  
 Operator : UM/SJ  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SSTDCCC040EC

Quant Time: May 09 00:58:50 2016  
 Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF050416.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed May 04 15:12:53 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 6.70  | 152  | 248366   | 40.00 | ng    | -0.03    |
| 21) Naphthalene-d8        | 8.00  | 136  | 973235   | 40.00 | ng    | -0.03    |
| 38) Acenaphthene-d10      | 9.74  | 164  | 416361   | 40.00 | ng    | -0.03    |
| 63) Phenanthrene-d10      | 11.22 | 188  | 638598   | 40.00 | ng    | -0.03    |
| 75) Chrysene-d12          | 13.86 | 240  | 434800   | 40.00 | ng    | -0.02    |
| 86) Perylene-d12          | 15.23 | 264  | 429267   | 40.00 | ng    | -0.03    |

## System Monitoring Compounds

|                          |       |     |         |       |    |       |
|--------------------------|-------|-----|---------|-------|----|-------|
| 5) 2-Fluorophenol        | 5.31  | 112 | 1290142 | 85.99 | ng | -0.01 |
| 7) Phenol-d6             | 6.37  | 99  | 1510072 | 74.72 | ng | 0.00  |
| 23) Nitrobenzene-d5      | 7.29  | 82  | 1463159 | 77.26 | ng | -0.02 |
| 41) 2,4,6-Tribromophenol | 10.53 | 330 | 283877  | 69.70 | ng | -0.03 |
| 44) 2-Fluorobiphenyl     | 9.07  | 172 | 1945821 | 76.61 | ng | -0.03 |
| 78) Terphenyl-d14        | 12.81 | 244 | 1610907 | 74.88 | ng | -0.03 |

## Target Compounds

|                                |      |     |        |       |    | Qvalue |
|--------------------------------|------|-----|--------|-------|----|--------|
| 2) 1,4-Dioxane                 | 2.21 | 88  | 300354 | 38.73 | ng | # 71   |
| 3) Pyridine                    | 2.90 | 79  | 821469 | 43.20 | ng | # 66   |
| 4) n-Nitrosodimethylamine      | 2.86 | 42  | 549556 | 47.65 | ng | # 50   |
| 6) Aniline                     | 6.38 | 93  | 938657 | 37.63 | ng | # 59   |
| 8) 2-Chlorophenol              | 6.50 | 128 | 675239 | 37.51 | ng | 85     |
| 9) Benzaldehyde                | 6.26 | 77  | 452684 | 44.90 | ng | 95     |
| 10) Phenol                     | 6.40 | 94  | 903224 | 40.85 | ng | 92     |
| 11) bis(2-Chloroethyl)ether    | 6.45 | 93  | 655117 | 34.43 | ng | # 83   |
| 12) 1,3-Dichlorobenzene        | 6.65 | 146 | 782536 | 40.93 | ng | 96     |
| 13) 1,4-Dichlorobenzene        | 6.73 | 146 | 783240 | 39.63 | ng | 97     |
| 14) 1,2-Dichlorobenzene        | 6.73 | 146 | 785393 | 45.63 | ng | 98     |
| 15) Benzyl Alcohol             | 6.88 | 79  | 561132 | 42.92 | ng | 94     |
| 16) 2,2'-oxybis(1-Chloropropan | 6.99 | 45  | 150    |       |    |        |

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050716\  
 Data File : BF086815.D  
 Acq On : 8 May 2016 11:03  
 Operator : UM/SJ  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SSTDCCC040EC

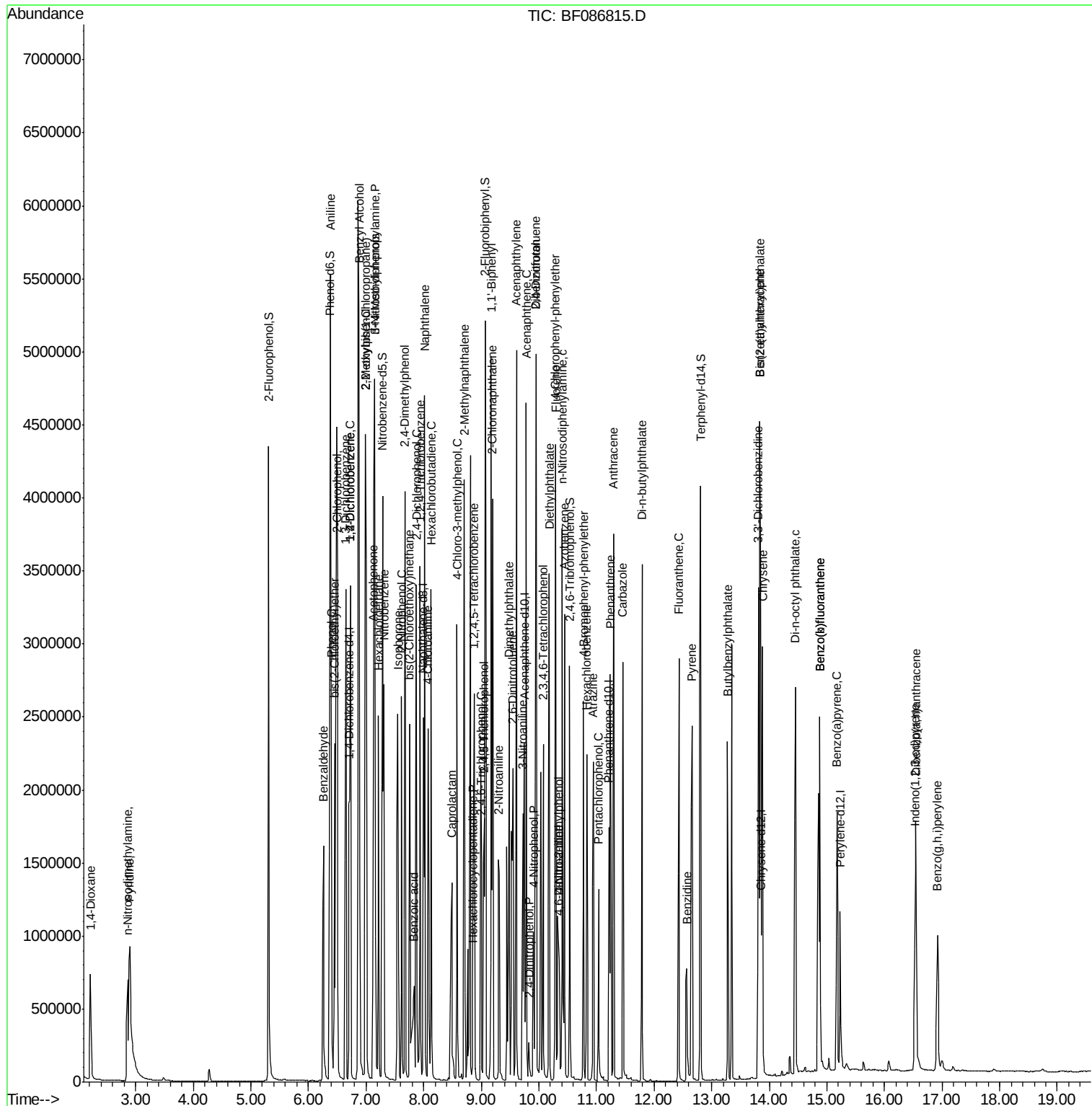
Quant Time: May 09 00:58:50 2016  
 Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF050416.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed May 04 15:12:53 2016  
 Response via : Initial Calibration

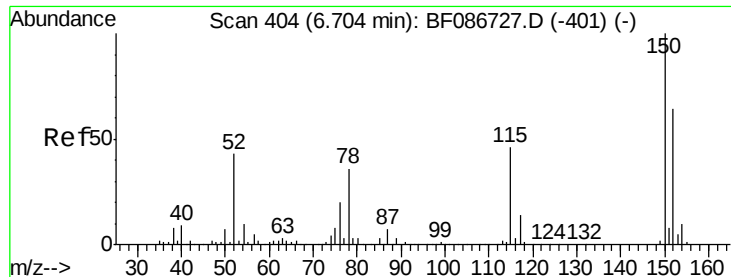
| Internal Standards             | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|--------------------------------|-------|------|----------|-------|-------|----------|
| 42) 2,4,6-Trichlorophenol      | 9.00  | 196  | 341318   | 44.54 | ng    | 92       |
| 43) 2,4,5-Trichlorophenol      | 9.05  | 196  | 342674   | 42.80 | ng    | 96       |
| 45) 1,1'-Biphenyl              | 9.17  | 154  | 1297125  | 38.11 | ng    | 95       |
| 46) 2-Chloronaphthalene        | 9.20  | 162  | 1015922  | 38.68 | ng    | 94       |
| 47) 2-Nitroaniline             | 9.30  | 65   | 399740   | 42.80 | ng    | # 70     |
| 48) Acenaphthylene             | 9.61  | 152  | 1453788  | 37.28 | ng    | 99       |
| 49) Dimethylphthalate          | 9.48  | 163  | 1293401  | 41.67 | ng    | 99       |
| 50) 2,6-Dinitrotoluene         | 9.55  | 165  | 308966   | 46.38 | ng    | 84       |
| 51) Acenaphthene               | 9.78  | 154  | 927124   | 36.04 | ng    | 99       |
| 52) 3-Nitroaniline             | 9.72  | 138  | 327469   | 43.24 | ng    | 98       |
| 53) 2,4-Dinitrophenol          | 9.82  | 184  | 40147    | 28.69 | ng    | 88       |
| 54) Dibenzofuran               | 9.95  | 168  | 1181542  | 37.06 | ng    | 95       |
| 55) 4-Nitrophenol              | 9.90  | 139  | 159871   | 36.21 | ng    | # 74     |
| 56) 2,4-Dinitrotoluene         | 9.95  | 165  | 328499   | 39.79 | ng    | # 88     |
| 57) Fluorene                   | 10.29 | 166  | 1038444  | 40.19 | ng    | 98       |
| 58) 2,3,4,6-Tetrachlorophenol  | 10.08 | 232  | 240748   | 40.94 | ng    | 97       |
| 59) Diethylphthalate           | 10.18 | 149  | 1239510  | 42.19 | ng    | 98       |
| 60) 4-Chlorophenyl-phenylether | 10.28 | 204  | 436045   | 35.50 | ng    | 96       |
| 61) 4-Nitroaniline             | 10.34 | 138  | 283513   | 38.97 | ng    | 85       |
| 62) Azobenzene                 | 10.44 | 77   | 1217816  | 37.41 | ng    | 85       |
| 64) 4,6-Dinitro-2-methylphenol | 10.35 | 198  | 80309    | 29.70 | ng    | # 10     |
| 65) n-Nitrosodiphenylamine     | 10.41 | 169  | 830683   | 40.97 | ng    | 99       |
| 66) 4-Bromophenyl-phenylether  | 10.77 | 248  | 300635   | 47.37 | ng    | 91       |
| 67) Hexachlorobenzene          | 10.83 | 284  | 302958   | 39.15 | ng    | # 67     |
| 68) Atrazine                   | 10.94 | 200  | 287538   | 46.27 | ng    | 95       |
| 69) Pentachlorophenol          | 11.04 | 266  | 148938   | 42.59 | ng    | 98       |
| 70) Phenanthrene               | 11.24 | 178  | 1343879  | 39.18 | ng    | 99       |
| 71) Anthracene                 |       |      |          |       |       |          |

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050716\  
 Data File : BF086815.D  
 Acq On : 8 May 2016 11:03  
 Operator : UM/SJ  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SSTDCCC040EC

Quant Time: May 09 00:58:50 2016  
 Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF050416.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed May 04 15:12:53 2016  
 Response via : Initial Calibration





#1

1,4-Dichlorobenzene-d4

Concen: 40.00 ng

RT: 6.70 min Scan# 404

Delta R.T. -0.03 min

Lab File: BF086815.D

Acq: 8 May 2016 11:03

Instrument :

BNA\_F

ClientSampleId :

SSTDCCC040EC

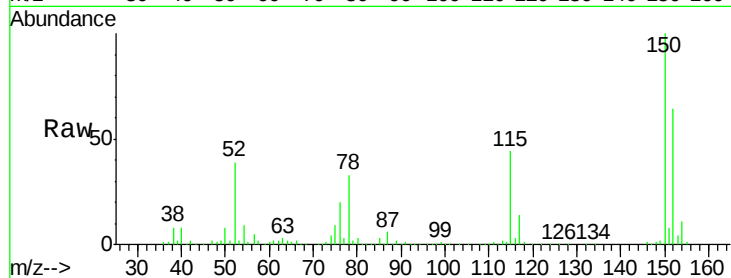
Tgt Ion:152 Resp: 248366

Ion Ratio Lower Upper

152 100

150 157.4 125.8 188.6

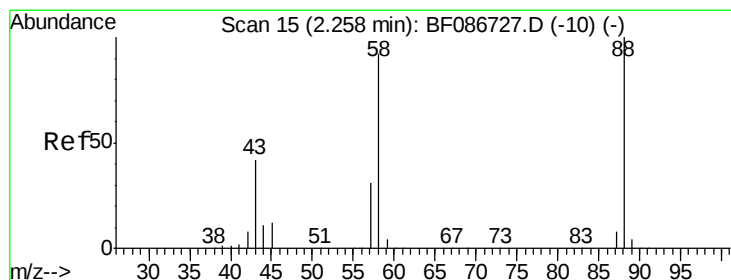
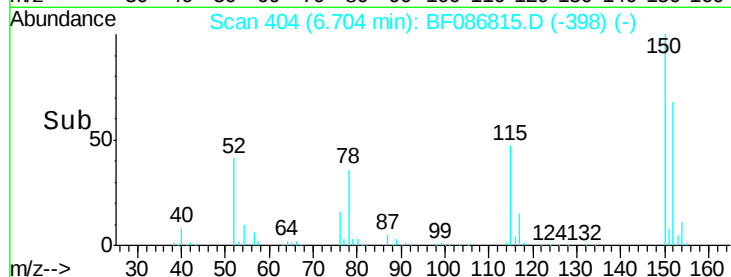
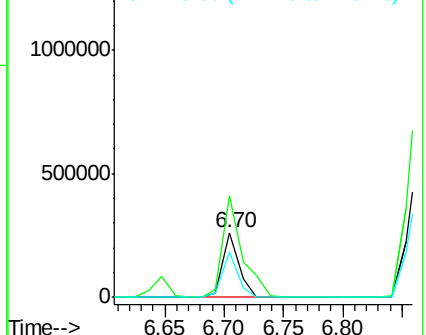
115 69.2 51.5 77.3



Abundance Ion 152.00 (151.70 to 152.70): E

Ion 150.00 (149.70 to 150.70): E

Ion 115.00 (114.70 to 115.70): E



#2

1,4-Dioxane

Concen: 38.73 ng

RT: 2.21 min Scan# 11

Delta R.T. -0.05 min

Lab File: BF086815.D

Acq: 8 May 2016 11:03

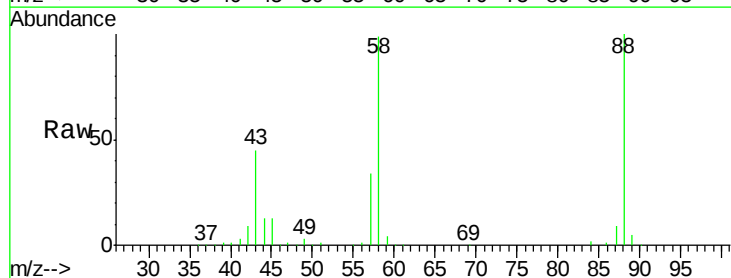
Tgt Ion: 88 Resp: 300354

Ion Ratio Lower Upper

88 100

58 97.3 59.6 89.4#

43 45.3 21.8 32.6#



Abundance Ion 88.00 (87.70 to 88.70): BF0

Ion 58.00 (57.70 to 58.70): BF0

Ion 43.00 (42.70 to 43.70): BF0

