

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF091216\  
 Data File : BF090062.D  
 Acq On : 12 Sep 2016 14:38  
 Operator : UM/SJ  
 Sample : H4716-02  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 EFF

Quant Time: Sep 12 16:11:41 2016  
 Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF083116.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Sep 12 15:34:24 2016  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response  | Conc    | Units | Dev(Min)  |
|--------------------------------|-------|------|-----------|---------|-------|-----------|
| 1) 1,4-Dichlorobenzene-d4      | 6.57  | 152  | 180493    | 20.00   | ng    | -0.01     |
| 21) Naphthalene-d8             | 7.86  | 136  | 658651    | 20.00   | ng    | 0.00      |
| 38) Acenaphthene-d10           | 9.61  | 164  | 351398    | 20.00   | ng    | 0.00      |
| 63) Phenanthrene-d10           | 11.08 | 188  | 662743    | 20.00   | ng    | 0.00      |
| 75) Chrysene-d12               | 13.70 | 240  | 227359    | 20.00   | ng    | 0.01      |
| 86) Perylene-d12               | 15.03 | 264  | 396137    | 20.00   | ng    | 0.00      |
| System Monitoring Compounds    |       |      |           |         |       |           |
| 5) 2-Fluorophenol              | 5.14  | 112  | 335115    | 29.72   | ng    | 0.00      |
| 7) Phenol-d6                   | 6.24  | 99   | 278900    | 20.36   | ng    | 0.01      |
| 23) Nitrobenzene-d5            | 7.14  | 82   | 861699    | 75.87   | ng    | -0.01     |
| 41) 2,4,6-Tribromophenol       | 10.40 | 330  | 247542    | 71.40   | ng    | 0.00      |
| 44) 2-Fluorobiphenyl           | 8.95  | 172  | 1790020   | 83.62   | ng    | 0.00      |
| 78) Terphenyl-d14              | 12.66 | 244  | 1173663   | 127.52  | ng    | 0.00      |
| Target Compounds               |       |      |           |         |       |           |
| 15) Benzyl Alcohol             | 6.76  | 79   | 200768    | 23.39   | ng    | Qvalue 97 |
| 22) Acetophenone               | 7.00  | 105  | 66238     | 4.48    | ng    | # 85      |
| 32) Benzoic acid               | 8.11  | 122  | 13517915m | 1879.57 | ng    |           |
| 49) Dimethylphthalate          | 9.35  | 163  | 318992    | 12.56   | ng    | 99        |
| 59) Diethylphthalate           | 10.04 | 149  | 76805     | 3.21    | ng    | 100       |
| 83) Bis(2-ethylhexyl)phthalate | 13.71 | 149  | 19280     | 2.16    | ng    | # 89      |

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

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