

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071716\  
 Data File : BF089078.D  
 Acq On : 17 Jul 2016 15:12  
 Operator : UM/SJ  
 Sample : H4046-11  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 C5-6

Quant Time: Jul 18 03:45:49 2016  
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jul 11 18:13:32 2016  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|-----------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 6.65  | 152  | 57044    | 20.00 | ng    | -0.06    |
| 21) Naphthalene-d8          | 7.94  | 136  | 236572   | 20.00 | ng    | -0.06    |
| 38) Acenaphthene-d10        | 9.69  | 164  | 98485    | 20.00 | ng    | -0.06    |
| 63) Phenanthrene-d10        | 11.16 | 188  | 153456   | 20.00 | ng    | -0.06    |
| 75) Chrysene-d12            | 13.79 | 240  | 114971   | 20.00 | ng    | -0.06    |
| 86) Perylene-d12            | 15.15 | 264  | 100095   | 20.00 | ng    | -0.06    |
| System Monitoring Compounds |       |      |          |       |       |          |
| 5) 2-Fluorophenol           | 5.24  | 112  | 327098   | 85.94 | ng    | -0.05    |
| 7) Phenol-d6                | 6.31  | 99   | 401190   | 81.57 | ng    | -0.05    |
| 23) Nitrobenzene-d5         | 7.22  | 82   | 255337   | 56.56 | ng    | -0.06    |
| 41) 2,4,6-Tribromophenol    | 10.48 | 330  | 71129    | 77.78 | ng    | -0.06    |
| 44) 2-Fluorobiphenyl        | 9.02  | 172  | 465784   | 74.71 | ng    | -0.06    |
| 78) Terphenyl-d14           | 12.74 | 244  | 301240   | 58.36 | ng    | -0.06    |
| Target Compounds            |       |      |          |       |       | Qvalue   |
| 48) Acenaphthylene          | 9.55  | 152  | 42101    | 4.13  | ng    | 99       |
| 49) Dimethylphthalate       | 9.42  | 163  | 91172    | 12.99 | ng    | 99       |
| 70) Phenanthrene            | 11.19 | 178  | 27667    | 3.30  | ng    | 99       |
| 71) Anthracene              | 11.23 | 178  | 36038    | 4.32  | ng    | 99       |
| 74) Fluoranthene            | 12.37 | 202  | 124942   | 14.69 | ng    | 99       |
| 77) Pyrene                  | 12.59 | 202  | 141896   | 14.56 | ng    | 98       |
| 80) Benzo(a)anthracene      | 13.78 | 228  | 91226    | 12.32 | ng    | # 87     |
| 82) Chrysene                | 13.82 | 228  | 80630m   | 11.01 | ng    |          |
| 85) Indeno(1,2,3-cd)pyrene  | 16.42 | 276  | 51980    | 9.22  | ng    | # 100    |
| 87) Benzo(b)fluoranthene    | 14.79 | 252  | 177608m  | 28.29 | ng    |          |
| 88) Benzo(k)fluoranthene    | 14.80 | 252  | 24416m   | 4.47  | ng    |          |
| 89) Benzo(a)pyrene          | 15.10 | 252  | 73728    | 13.87 | ng    | 96       |
| 90) Dibenzo(a,h)anthracene  | 16.42 | 278  | 12344m   | 2.78  | ng    |          |
| 91) Benzo(g,h,i)perylene    | 16.80 | 276  | 51450    | 11.20 | ng    | 96       |

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

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