

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071416\  
 Data File : BF089000.D  
 Acq On : 14 Jul 2016 22:21  
 Operator : UM/SJ  
 Sample : H3996-15  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 C8-B1(0-6)C

## Manual Integrations

APPROVED  
 Sohil  
 7/15/2016 9:16:30 AM

Quant Time: Jul 15 03:29:57 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      | 6.68  | 152  | 61000    | 20.00 | ng    | -0.02     |
| 21) Naphthalene-d8             | 7.98  | 136  | 222410   | 20.00 | ng    | -0.02     |
| 38) Acenaphthene-d10           | 9.71  | 164  | 86890    | 20.00 | ng    | -0.03     |
| 63) Phenanthrene-d10           | 11.19 | 188  | 138138   | 20.00 | ng    | -0.03     |
| 75) Chrysene-d12               | 13.82 | 240  | 127545   | 20.00 | ng    | -0.03     |
| 86) Perylene-d12               | 15.19 | 264  | 95898    | 20.00 | ng    | -0.02     |
| System Monitoring Compounds    |       |      |          |       |       |           |
| 5) 2-Fluorophenol              | 5.28  | 112  | 337708   | 82.97 | ng    | -0.01     |
| 7) Phenol-d6                   | 6.34  | 99   | 516706   | 98.25 | ng    | -0.01     |
| 23) Nitrobenzene-d5            | 7.26  | 82   | 325557   | 76.71 | ng    | -0.02     |
| 41) 2,4,6-Tribromophenol       | 10.50 | 330  | 62089    | 76.95 | ng    | -0.03     |
| 44) 2-Fluorobiphenyl           | 9.05  | 172  | 508986   | 92.53 | ng    | -0.02     |
| 78) Terphenyl-d14              | 12.78 | 244  | 328152   | 57.31 | ng    | -0.02     |
| Target Compounds               |       |      |          |       |       |           |
| 49) Dimethylphthalate          | 9.44  | 163  | 73130    | 11.81 | ng    | Qvalue 99 |
| 70) Phenanthrene               | 11.21 | 178  | 38923    | 5.15  | ng    | 99        |
| 74) Fluoranthene               | 12.40 | 202  | 72543    | 9.48  | ng    | 93        |
| 77) Pyrene                     | 12.62 | 202  | 69567    | 6.43  | ng    | 98        |
| 80) Benzo(a)anthracene         | 13.81 | 228  | 36397    | 4.43  | ng    | 96        |
| 82) Chrysene                   | 13.84 | 228  | 44366    | 5.46  | ng    | 96        |
| 83) Bis(2-ethylhexyl)phthalate | 13.82 | 149  | 16852    | 3.06  | ng    | 99        |
| 85) Indeno(1,2,3-cd)pyrene     | 16.46 | 276  | 20035    | 3.20  | ng    | # 100     |
| 87) Benzo(b)fluoranthene       | 14.81 | 252  | 53905m   | 8.96  | ng    |           |
| 88) Benzo(k)fluoranthene       | 14.82 | 252  | 13111m   | 2.50  | ng    |           |
| 89) Benzo(a)pyrene             | 15.13 | 252  | 29167    | 5.73  | ng    | 95        |
| 91) Benzo(a,h,i)perylene       | 16.83 | 276  | 19305    | 4.39  | ng    | 93        |

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

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