

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101518\  
 Data File : BF109890.D  
 Acq On : 15 Oct 2018 20:20  
 Operator : JU/SJ  
 Sample : J5193-01  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 BU-01-101118

Manual Integrations  
 APPROVED

Sohil  
 10/16/2018 9:57:33 AM

Quant Time: Oct 16 08:53:30 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Oct 15 18:09:48 2018  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 6.97  | 152  | 236155   | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.26  | 136  | 827017   | 20.00  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 10.02 | 164  | 312321   | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.50 | 188  | 538977   | 20.00  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.14 | 240  | 575694   | 20.00  | ng    | 0.00     |
| 87) Perylene-d12            | 15.67 | 264  | 430341   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |       |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.59  | 112  | 1391533  | 93.40  | ng    | 0.00     |
| 7) Phenol-d6                | 6.59  | 99   | 1683685  | 91.01  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.54  | 82   | 1021862  | 83.23  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.80 | 330  | 270973   | 100.21 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.33  | 172  | 1592395  | 81.36  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.09 | 244  | 1486231  | 54.21  | ng    | 0.00     |
| Target Compounds            |       |      |          |        |       |          |
| 10) Phenol                  | 6.60  | 94   | 49993    | 2.503  | ng    | 88       |
| 50) Dimethylphthalate       | 9.72  | 163  | 205733   | 9.303  | ng    | 100      |
| 71) Phenanthrene            | 11.53 | 178  | 112050   | 4.016  | ng    | 97       |
| 75) Fluoranthene            | 12.72 | 202  | 335208   | 12.025 | ng    | 100      |
| 78) Pyrene                  | 12.95 | 202  | 315026   | 7.450  | ng    | 97       |
| 81) Benzo(a)anthracene      | 14.14 | 228  | 161442   | 4.761  | ng    | 96       |
| 83) Chrysene                | 14.17 | 228  | 170166   | 5.272  | ng    | 98       |
| 86) Indeno(1,2,3-cd)pyrene  | 17.22 | 276  | 65070    | 2.504  | ng    | 96       |
| 88) Benzo(b)fluoranthene    | 15.22 | 252  | 180221m  | 7.264  | ng    |          |
| 89) Benzo(k)fluoranthene    | 15.24 | 252  | 63466m   | 2.595  | ng    |          |
| 90) Benzo(a)pyrene          | 15.60 | 252  | 114490   | 5.017  | ng    | # 92     |
| 92) Benzo(a,h,i)perylene    | 17.70 | 276  | 67970    | 3.327  | ng    | # 90     |

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

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