

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120216\  
 Data File : BF091412.D  
 Acq On : 3 Dec 2016 1:52  
 Operator : UM/SJ  
 Sample : H5825-04  
 Misc :  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 S12016SF-S-4

Manual Integrations  
 APPROVED

sohil  
 12/5/2016 11:21:10 AM

Quant Time: Dec 03 03:05:23 2016  
 Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF112916.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Nov 29 13:33:46 2016  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 6.86  | 152  | 161303   | 20.00  | ng    | -0.02    |
| 21) Naphthalene-d8          | 8.15  | 136  | 612721   | 20.00  | ng    | -0.02    |
| 38) Acenaphthene-d10        | 9.90  | 164  | 348443   | 20.00  | ng    | -0.02    |
| 63) Phenanthrene-d10        | 11.38 | 188  | 486546   | 20.00  | ng    | -0.02    |
| 75) Chrysene-d12            | 14.01 | 240  | 407523   | 20.00  | ng    | -0.03    |
| 86) Perylene-d12            | 15.44 | 264  | 313495   | 20.00  | ng    | -0.03    |
| System Monitoring Compounds |       |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.46  | 112  | 1013491  | 102.64 | ng    | 0.00     |
| 7) Phenol-d6                | 6.49  | 99   | 1320897  | 108.67 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.42  | 82   | 804214   | 73.56  | ng    | -0.03    |
| 41) 2,4,6-Tribromophenol    | 10.69 | 330  | 325286   | 98.61  | ng    | -0.02    |
| 44) 2-Fluorobiphenyl        | 9.22  | 172  | 1473292  | 76.56  | ng    | -0.02    |
| 78) Terphenyl-d14           | 12.97 | 244  | 1190030  | 63.47  | ng    | -0.02    |
| Target Compounds            |       |      |          |        |       | Qvalue   |
| 10) Phenol                  | 6.50  | 94   | 113443   | 7.93   | ng    | 87       |
| 49) Dimethylphthalate       | 9.61  | 163  | 162395   | 7.70   | ng    | 98       |
| 70) Phenanthrene            | 11.41 | 178  | 147511   | 5.83   | ng    | 97       |
| 74) Fluoranthene            | 12.59 | 202  | 245844   | 10.27  | ng    | 97       |
| 77) Pyrene                  | 12.81 | 202  | 236076   | 7.63   | ng    | 98       |
| 80) Benzo(a)anthracene      | 14.00 | 228  | 139727   | 5.86   | ng    | 100      |
| 82) Chrysene                | 14.04 | 228  | 91279    | 3.87   | ng    | 96       |
| 85) Indeno(1,2,3-cd)pyrene  | 16.84 | 276  | 45303    | 2.10   | ng    | # 90     |
| 87) Benzo(b)fluoranthene    | 15.03 | 252  | 138713m  | 7.12   | ng    |          |
| 89) Benzo(a)pyrene          | 15.39 | 252  | 87067m   | 4.98   | ng    |          |
| 91) Benzo(g,h,i)perylene    | 17.26 | 276  | 44685    | 2.68   | ng    | 99       |

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

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120216\  
Data File : BF091412.D  
Acq On : 3 Dec 2016 1:52  
Operator : UM/SJ  
Sample : H5825-04  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
S12016SF-S-4

Manual Integrations  
APPROVED

sohil  
12/5/2016 11:21:10 AM

Quant Time: Dec 03 03:05:23 2016  
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF112916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Tue Nov 29 13:33:46 2016  
Response via : Initial Calibration

