

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081719\  
 Data File : BF116376.D  
 Acq On : 18 Aug 2019 4:06  
 Operator : HP/JU  
 Sample : K4384-04DL 20X  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 B4(3-4)DL

Manual Integrations  
 APPROVED

Jagrut  
 8/19/2019 3:31:32 PM

Quant Time: Aug 19 17:53:01 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Aug 07 11:13:18 2019  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|-----------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 6.81  | 152  | 122644   | 20.00 | ng    | -0.04    |
| 21) Naphthalene-d8          | 8.09  | 136  | 491906   | 20.00 | ng    | -0.04    |
| 39) Acenaphthene-d10        | 9.84  | 164  | 243959   | 20.00 | ng    | -0.05    |
| 64) Phenanthrene-d10        | 11.33 | 188  | 367971   | 20.00 | ng    | -0.04    |
| 76) Chrysene-d12            | 13.97 | 240  | 234711   | 20.00 | ng    | -0.04    |
| 87) Perylene-d12            | 15.43 | 264  | 269672   | 20.00 | ng    | -0.05    |
| System Monitoring Compounds |       |      |          |       |       |          |
| 5) 2-Fluorophenol           | 5.48  | 112  | 30112    | 4.11  | ng    | 0.01     |
| 7) Phenol-d6                | 6.49  | 99   | 50380    | 5.45  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.40  | 82   | 29670    | 3.48  | ng    | -0.02    |
| 42) 2,4,6-Tribromophenol    | 10.65 | 330  | 7719     | 3.26  | ng    | -0.03    |
| 45) 2-Fluorobiphenyl        | 9.16  | 172  | 62044    | 4.76  | ng    | -0.04    |
| 79) Terphenyl-d14           | 12.90 | 244  | 49189    | 4.42  | ng    | -0.05    |
| Target Compounds            |       |      |          |       |       |          |
| 71) Phenanthrene            | 11.35 | 178  | 105369   | 5.601 | ng    | 97       |
| 75) Fluoranthene            | 12.54 | 202  | 127601   | 6.479 | ng    | 99       |
| 78) Pyrene                  | 12.77 | 202  | 118141   | 6.677 | ng    | 98       |
| 81) Benzo(a)anthracene      | 13.96 | 228  | 54474    | 3.631 | ng    | 97       |
| 83) Chrysene                | 14.00 | 228  | 49745    | 3.415 | ng    | 98       |
| 86) Indeno(1,2,3-cd)pyrene  | 16.90 | 276  | 30706    | 2.510 | ng    | 96       |
| 88) Benzo(b)fluoranthene    | 15.01 | 252  | 61899m   | 3.970 | ng    |          |
| 90) Benzo(a)pyrene          | 15.37 | 252  | 48156    | 3.455 | ng    | 96       |
| 92) Benzo(g,h,i)perylene    | 17.34 | 276  | 29226    | 2.466 | ng    | 97       |

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081719\  
Data File : BF116376.D  
Acq On : 18 Aug 2019 4:06  
Operator : HP/JU  
Sample : K4384-04DL 20X  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B4(3-4)DL

Manual Integrations  
APPROVED

Jagrut  
8/19/2019 3:31:32 PM

Quant Time: Aug 19 17:53:01 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Wed Aug 07 11:13:18 2019  
Response via : Initial Calibration

