

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF042219\  
 Data File : BF113897.D  
 Acq On : 23 Apr 2019 00:02  
 Operator : JU/SJ  
 Sample : K2202-01 2X  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SA-06-041819-A

Manual Integrations  
 APPROVED

Sohil  
 4/23/2019 4:08:12 PM

Quant Time: Apr 23 05:58:22 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF042219.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Apr 23 03:32:35 2019  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc  | Units | Dev(Min)   |
|-----------------------------|-------|------|----------|-------|-------|------------|
| 1) 1,4-Dichlorobenzene-d4   | 6.90  | 152  | 194490   | 20.00 | ng    | 0.00       |
| 21) Naphthalene-d8          | 8.18  | 136  | 658756   | 20.00 | ng    | 0.00       |
| 39) Acenaphthene-d10        | 9.93  | 164  | 312243   | 20.00 | ng    | 0.00       |
| 64) Phenanthrene-d10        | 11.42 | 188  | 646741   | 20.00 | ng    | 0.00       |
| 76) Chrysene-d12            | 14.06 | 240  | 642259   | 20.00 | ng    | -0.01      |
| 87) Perylene-d12            | 15.54 | 264  | 532731   | 20.00 | ng    | -0.03      |
| System Monitoring Compounds |       |      |          |       |       |            |
| 5) 2-Fluorophenol           | 5.52  | 112  | 590188   | 51.90 | ng    | 0.00       |
| 7) Phenol-d6                | 6.52  | 99   | 729372   | 51.12 | ng    | 0.00       |
| 23) Nitrobenzene-d5         | 7.45  | 82   | 407779   | 38.22 | ng    | 0.00       |
| 42) 2,4,6-Tribromophenol    | 10.72 | 330  | 216542   | 56.93 | ng    | 0.00       |
| 45) 2-Fluorobiphenyl        | 9.25  | 172  | 745984   | 37.37 | ng    | 0.00       |
| 79) Terphenyl-d14           | 13.00 | 244  | 991852   | 31.66 | ng    | 0.00       |
| Target Compounds            |       |      |          |       |       |            |
| 50) Dimethylphthalate       | 9.64  | 163  | 124436   | 5.510 | ng    | Qvalue 100 |
| 71) Phenanthrene            | 11.44 | 178  | 137100   | 4.218 | ng    | 98         |
| 75) Fluoranthene            | 12.63 | 202  | 337423   | 9.515 | ng    | 98         |
| 78) Pyrene                  | 12.86 | 202  | 330975   | 7.369 | ng    | 98         |
| 81) Benzo(a)anthracene      | 14.04 | 228  | 180279   | 4.764 | ng    | 95         |
| 83) Chrysene                | 14.08 | 228  | 158230   | 4.294 | ng    | 97         |
| 86) Indeno(1,2,3-cd)pyrene  | 17.02 | 276  | 69860m   | 2.001 | ng    |            |
| 88) Benzo(b)fluoranthene    | 15.10 | 252  | 202667m  | 6.013 | ng    |            |
| 90) Benzo(a)pyrene          | 15.47 | 252  | 127739   | 4.382 | ng    | # 94       |
| 92) Benzo(g,h,i)perylene    | 17.47 | 276  | 67679    | 2.451 | ng    | 94         |

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

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