

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072219\  
 Data File : BF115712.D  
 Acq On : 22 Jul 2019 16:33  
 Operator : HP/JU  
 Sample : K3622-09 2X  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 OR-02-071919-A

Manual Integrations  
 APPROVED

mohammad  
 7/23/2019 4:24:30 PM

Quant Time: Jul 23 03:33:49 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Jul 12 14:03:52 2019  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 6.88  | 152  | 174050   | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.16  | 136  | 672443   | 20.00  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.92  | 164  | 335403   | 20.00  | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.40 | 188  | 513783   | 20.00  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.04 | 240  | 413219   | 20.00  | ng    | 0.00     |
| 87) Perylene-d12            | 15.53 | 264  | 408624   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |       |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.50  | 112  | 411234   | 38.65  | ng    | 0.00     |
| 7) Phenol-d6                | 6.50  | 99   | 547406   | 40.54  | ng    | -0.01    |
| 23) Nitrobenzene-d5         | 7.43  | 82   | 324084   | 28.02  | ng    | -0.02    |
| 42) 2,4,6-Tribromophenol    | 10.70 | 330  | 125206   | 31.98  | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 9.23  | 172  | 573765   | 29.94  | ng    | -0.02    |
| 79) Terphenyl-d14           | 12.99 | 244  | 497065   | 22.20  | ng    | -0.01    |
| Target Compounds            |       |      |          |        |       |          |
| 50) Dimethylphthalate       | 9.63  | 163  | 75860    | 3.006  | ng    | 98       |
| 71) Phenanthrene            | 11.43 | 178  | 272205   | 10.336 | ng    | 97       |
| 72) Anthracene              | 11.47 | 178  | 82023    | 3.014  | ng    | 99       |
| 75) Fluoranthene            | 12.62 | 202  | 487479   | 17.516 | ng    | 98       |
| 78) Pyrene                  | 12.84 | 202  | 473351   | 13.521 | ng    | 97       |
| 81) Benzo(a)anthracene      | 14.03 | 228  | 226059   | 8.304  | ng    | 99       |
| 83) Chrysene                | 14.07 | 228  | 187410   | 6.858  | ng    | 95       |
| 86) Indeno(1,2,3-cd)pyrene  | 17.01 | 276  | 70041    | 3.017  | ng    | 95       |
| 88) Benzo(b)fluoranthene    | 15.09 | 252  | 211796m  | 8.049  | ng    |          |
| 89) Benzo(k)fluoranthene    | 15.12 | 252  | 59121m   | 2.520  | ng    |          |
| 90) Benzo(a)pyrene          | 15.46 | 252  | 150106   | 6.637  | ng    | # 92     |
| 92) Benzo(a,h,i)perylene    | 17.46 | 276  | 66503    | 3.359  | ng    | # 90     |

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

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