

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122018\  
 Data File : BF111653.D  
 Acq On : 20 Dec 2018 21:55  
 Operator : JU/SJ  
 Sample : J6357-09  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-09B

Manual Integrations  
 APPROVED

Sohil  
 12/21/2018 2:46:53 PM

Quant Time: Dec 21 07:12:46 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Dec 19 17:39:39 2018  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 6.90  | 152  | 222062   | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.17  | 136  | 811903   | 20.00  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.93  | 164  | 378657   | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.42 | 188  | 579546   | 20.00  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.05 | 240  | 503807   | 20.00  | ng    | 0.00     |
| 87) Perylene-d12            | 15.52 | 264  | 445329   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |       |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.53  | 112  | 1617818  | 124.18 | ng    | 0.01     |
| 7) Phenol-d6                | 6.53  | 99   | 2082736  | 125.62 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.46  | 82   | 1324666  | 94.97  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.72 | 330  | 441772   | 98.74  | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.25  | 172  | 2072988  | 79.80  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.00 | 244  | 1450688  | 54.61  | ng    | 0.00     |
| Target Compounds            |       |      |          |        |       |          |
| 10) Phenol                  | 6.54  | 94   | 135384   | 7.237  | ng    | 94       |
| 50) Dimethylphthalate       | 9.64  | 163  | 158777   | 5.734  | ng    | 96       |
| 71) Phenanthrene            | 11.44 | 178  | 417936   | 13.598 | ng    | 94       |
| 72) Anthracene              | 11.49 | 178  | 72949    | 2.365  | ng    | 96       |
| 75) Fluoranthene            | 12.62 | 202  | 414232   | 13.309 | ng    | 96       |
| 78) Pyrene                  | 12.85 | 202  | 470106   | 13.214 | ng    | 96       |
| 81) Benzo(a)anthracene      | 14.03 | 228  | 220493   | 7.669  | ng    | 99       |
| 83) Chrysene                | 14.07 | 228  | 221825   | 7.850  | ng    | 97       |
| 86) Indeno(1,2,3-cd)pyrene  | 16.99 | 276  | 66345    | 2.762  | ng    | # 83     |
| 88) Benzo(b)fluoranthene    | 15.09 | 252  | 217813m  | 8.369  | ng    |          |
| 90) Benzo(a)pyrene          | 15.46 | 252  | 152439m  | 6.447  | ng    |          |
| 92) Benzo(a,h,i)perylene    | 17.44 | 276  | 63829    | 3.157  | ng    | # 88     |

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

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