

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040620\  
 Data File : BF119813.D  
 Acq On : 7 Apr 2020 11:15  
 Operator : MA/SJ  
 Sample : L2147-03  
 Misc :  
 ALS Vial : 51 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 FK-GF-02-056-B

Manual Integrations  
 APPROVED

mohammad  
 4/8/2020 8:12:49 AM

Quant Time: Apr 07 11:46:33 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Apr 06 11:15:16 2020  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 6.78  | 152  | 175564   | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.07  | 136  | 640702   | 20.00  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.83  | 164  | 305856   | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.31 | 188  | 549311   | 20.00  | ng    | 0.00     |
| 76) Chrysene-d12            | 13.95 | 240  | 379222   | 20.00  | ng    | 0.00     |
| 87) Perylene-d12            | 15.40 | 264  | 382552   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |       |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.39  | 112  | 1334575  | 121.01 | ng    | 0.00     |
| 7) Phenol-d6                | 6.42  | 99   | 1721651  | 122.21 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.35  | 82   | 1052643  | 89.29  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.62 | 330  | 425126   | 134.73 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.15  | 172  | 1792250  | 86.81  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.90 | 244  | 1767479  | 91.32  | ng    | 0.00     |
| Target Compounds            |       |      |          |        |       |          |
| 50) Dimethylphthalate       | 9.54  | 163  | 251050   | 11.556 | ng    | 99       |
| 71) Phenanthrene            | 11.33 | 178  | 264689   | 9.293  | ng    | 96       |
| 75) Fluoranthene            | 12.53 | 202  | 428691   | 13.907 | ng    | 97       |
| 78) Pyrene                  | 12.76 | 202  | 396401   | 12.737 | ng    | 97       |
| 81) Benzo(a)anthracene      | 13.94 | 228  | 177588   | 7.201  | ng    | 97       |
| 83) Chrysene                | 13.98 | 228  | 173925   | 6.834  | ng    | 97       |
| 86) Indeno(1,2,3-cd)pyrene  | 16.82 | 276  | 88327    | 3.685  | ng    | 95       |
| 88) Benzo(b)fluoranthene    | 14.98 | 252  | 223107m  | 8.731  | ng    |          |
| 89) Benzo(k)fluoranthene    | 15.00 | 252  | 76258m   | 3.134  | ng    |          |
| 90) Benzo(a)pyrene          | 15.34 | 252  | 154086   | 6.635  | ng    | 98       |
| 92) Benzo(g,h,i)perylene    | 17.25 | 276  | 86206    | 4.224  | ng    | 99       |

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

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