

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100520\  
 Data File : BF121830.D  
 Acq On : 5 Oct 2020 21:53  
 Operator : JU/CG  
 Sample : L4198-31 2X  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 RO-4

Manual Integrations  
 APPROVED

mohammad  
 10/7/2020 10:36:26 AM

Quant Time: Oct 06 01:28:17 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Oct 01 14:29:14 2020  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc  | Units | Dev(Min)   |
|-----------------------------|-------|------|----------|-------|-------|------------|
| 1) 1,4-Dichlorobenzene-d4   | 6.76  | 152  | 101251   | 20.00 | ng    | 0.00       |
| 21) Naphthalene-d8          | 8.05  | 136  | 376246   | 20.00 | ng    | -0.01      |
| 39) Acenaphthene-d10        | 9.80  | 164  | 167454   | 20.00 | ng    | -0.01      |
| 64) Phenanthrene-d10        | 11.28 | 188  | 271347   | 20.00 | ng    | -0.01      |
| 76) Chrysene-d12            | 13.92 | 240  | 265388   | 20.00 | ng    | -0.01      |
| 86) Perylene-d12            | 15.36 | 264  | 262801   | 20.00 | ng    | 0.00       |
| System Monitoring Compounds |       |      |          |       |       |            |
| 5) 2-Fluorophenol           | 5.38  | 112  | 265373   | 38.95 | ng    | 0.00       |
| 7) Phenol-d6                | 6.40  | 99   | 323736   | 36.22 | ng    | -0.01      |
| 23) Nitrobenzene-d5         | 7.32  | 82   | 207758   | 29.65 | ng    | -0.02      |
| 42) 2,4,6-Tribromophenol    | 10.59 | 330  | 65976    | 31.70 | ng    | -0.02      |
| 45) 2-Fluorobiphenyl        | 9.12  | 172  | 354282   | 32.04 | ng    | -0.01      |
| 79) Terphenyl-d14           | 12.86 | 244  | 280528   | 21.41 | ng    | -0.02      |
| Target Compounds            |       |      |          |       |       |            |
| 50) Dimethylphthalate       | 9.52  | 163  | 32077    | 2.643 | ng    | Qvalue 100 |
| 75) Fluoranthene            | 12.49 | 202  | 66370    | 4.205 | ng    | 98         |
| 78) Pyrene                  | 12.72 | 202  | 69655    | 3.592 | ng    | 98         |
| 81) Benzo(a)anthracene      | 13.91 | 228  | 55235    | 3.290 | ng    | 99         |
| 83) Chrysene                | 13.94 | 228  | 44756    | 2.632 | ng    | 96         |
| 88) Benzo(b)fluoranthene    | 14.94 | 252  | 80372m   | 4.575 | ng    |            |
| 90) Benzo(a)pyrene          | 15.30 | 252  | 55182    | 3.695 | ng    | # 92       |
| 92) Benzo(g,h,i)perylene    | 17.20 | 276  | 29181    | 2.084 | ng    | 97         |

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

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