

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG041521\  
 Data File : BG048716.D  
 Acq On : 15 Apr 2021 19:46  
 Operator : CG/JU  
 Sample : M2017-01RE  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 FISHKILL-SP-1RE

Manual Integrations  
 APPROVED

mohammad  
 4/16/2021 11:28:29 AM

Quant Time: Apr 16 00:34:56 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG041421.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Apr 15 10:59:37 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units | Dev(Min) |
|-----------------------------|--------|------|----------|-------|-------|----------|
| -----                       |        |      |          |       |       |          |
| Internal Standards          |        |      |          |       |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.087  | 152  | 26611    | 20.00 | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.913 | 136  | 104290   | 20.00 | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.744 | 164  | 57053    | 20.00 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.499 | 188  | 120931   | 20.00 | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.800 | 240  | 124030   | 20.00 | ng    | # 0.00   |
| 86) Perylene-d12            | 25.155 | 264  | 122456   | 20.00 | ng    | 0.02     |
| System Monitoring Compounds |        |      |          |       |       |          |
| 5) 2-Fluorophenol           | 5.643  | 112  | 138558   | 87.08 | ng    | 0.01     |
| 7) Phenol-d6                | 7.247  | 99   | 178677   | 79.39 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.256  | 82   | 137310   | 58.95 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.236 | 330  | 59726    | 78.08 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.363 | 172  | 267461   | 67.70 | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.108 | 244  | 365893   | 50.77 | ng    | 0.00     |
| Target Compounds            |        |      |          |       |       |          |
|                             |        |      |          |       |       | Qvalue   |
| 50) Dimethylphthalate       | 14.186 | 163  | 12455    | 2.91  | ng    | # 93     |
| 71) Phenanthrene            | 17.540 | 178  | 25917    | 4.13  | ng    | # 93     |
| 75) Fluoranthene            | 19.556 | 202  | 65431    | 8.24  | ng    | 98       |
| 78) Pyrene                  | 19.920 | 202  | 71055    | 8.28  | ng    | 99       |
| 81) Benzo(a)anthracene      | 21.783 | 228  | 43075    | 5.08  | ng    | # 90     |
| 83) Chrysene                | 21.847 | 228  | 45950m   | 5.69  | ng    |          |
| 88) Benzo(b)fluoranthene    | 24.097 | 252  | 52307    | 6.30  | ng    | # 95     |
| 89) Benzo(k)fluoranthene    | 24.144 | 252  | 19256m   | 2.46  | ng    |          |
| 90) Benzo(a)pyrene          | 24.991 | 252  | 38008    | 4.97  | ng    | # 90     |
| -----                       |        |      |          |       |       |          |

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

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