

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091620\  
 Data File : BG046645.D  
 Acq On : 16 Sep 2020 18:01  
 Operator : CG/JU  
 Sample : L4021-01  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 B14-1-3

Quant Time: Sep 16 18:43:59 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082520.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Sep 16 11:07:52 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.92  | 152  | 69268    | 20.00 | ng    | 0.00     |
| 21) Naphthalene-d8        | 10.71 | 136  | 280998   | 20.00 | ng    | 0.00     |
| 39) Acenaphthene-d10      | 14.55 | 164  | 189083   | 20.00 | ng    | 0.00     |
| 64) Phenanthrene-d10      | 17.29 | 188  | 425837   | 20.00 | ng    | 0.00     |
| 76) Chrysene-d12          | 21.54 | 240  | 428875   | 20.00 | ng    | 0.00     |
| 86) Perylene-d12          | 24.63 | 264  | 427721   | 20.00 | ng    | 0.00     |

|                             |       |     |         |       |    |      |
|-----------------------------|-------|-----|---------|-------|----|------|
| System Monitoring Compounds |       |     |         |       |    |      |
| 5) 2-Fluorophenol           | 5.51  | 112 | 91537   | 23.41 | ng | 0.00 |
| 7) Phenol-d6                | 7.10  | 99  | 358015  | 64.58 | ng | 0.00 |
| 23) Nitrobenzene-d5         | 9.08  | 82  | 293332  | 59.66 | ng | 0.00 |
| 42) 2,4,6-Tribromophenol    | 16.04 | 330 | 7402    | 2.89  | ng | 0.00 |
| 45) 2-Fluorobiphenyl        | 13.17 | 172 | 777804  | 62.80 | ng | 0.00 |
| 79) Terphenyl-d14           | 19.91 | 244 | 1139665 | 56.55 | ng | 0.00 |

|                          |       |     |        |        |    |      |
|--------------------------|-------|-----|--------|--------|----|------|
| Target Compounds         |       |     |        | Qvalue |    |      |
| 10) Phenol               | 7.12  | 94  | 16788  | 3.288  | ng | 91   |
| 50) Dimethylphthalate    | 14.00 | 163 | 73533  | 5.027  | ng | 98   |
| 71) Phenanthrene         | 17.34 | 178 | 80797  | 3.743  | ng | 96   |
| 75) Fluoranthene         | 19.35 | 202 | 138576 | 4.989  | ng | 97   |
| 78) Pyrene               | 19.71 | 202 | 152657 | 5.590  | ng | 95   |
| 81) Benzo(a)anthracene   | 21.52 | 228 | 77453  | 2.897  | ng | 96   |
| 83) Chrysene             | 21.58 | 228 | 81426  | 3.167  | ng | 99   |
| 88) Benzo(b)fluoranthene | 23.66 | 252 | 98803  | 3.699  | ng | # 97 |
| 90) Benzo(a)pyrene       | 24.49 | 252 | 69447  | 2.786  | ng | 95   |
| 92) Benzo(g,h,i)perylene | 29.22 | 276 | 52302  | 2.113  | ng | 98   |

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

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