

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021621\  
 Data File : BG048188.D  
 Acq On : 17 Feb 2021 13:15  
 Operator : CG/JU  
 Sample : M1379-17  
 Misc :  
 ALS Vial : 30 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 DBH01

Manual Integrations  
 APPROVED

mohammad  
 2/18/2021 1:38:33 PM

Quant Time: Feb 17 17:46:50 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG021321MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Feb 17 15:04:50 2021  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.13  | 152  | 51837    | 20.00 | ng/ul | 0.03     |
| 18) Naphthalene-d8        | 10.95 | 136  | 191390   | 20.00 | ng/ul | 0.03     |
| 36) Acenaphthene-d10      | 14.76 | 164  | 125922   | 20.00 | ng/ul | 0.03     |
| 62) Phenanthrene-d10      | 17.50 | 188  | 273011   | 20.00 | ng/ul | 0.03     |
| 78) Chrysene-d12          | 21.85 | 240  | 380241m  | 20.00 | ng/ul | 0.12     |
| 86) Perylene-d12          | 25.29 | 264  | 238060m  | 20.00 | ng/ul | 0.29     |

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| System Monitoring Compounds    |       |     |         |       |       |      |
| 3) 1,4-Dioxane-d8              | 3.54  | 96  | 3237    | 2.60  | ng/uL | 0.03 |
| 5) Phenol-d5                   | 7.34  | 99  | 79889   | 18.29 | ng/ul | 0.09 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.46  | 67  | 47604   | 18.30 | ng/ul | 0.04 |
| 9) 2-Chlorophenol-d4           | 7.68  | 132 | 58278   | 19.01 | ng/ul | 0.05 |
| 13) 4-Methylphenol-d8          | 8.88  | 113 | 62709   | 18.30 | ng/ul | 0.08 |
| 19) Nitrobenzene-d5            | 9.31  | 128 | 30489   | 21.33 | ng/ul | 0.05 |
| 22) 2-Nitrophenol-d4           | 10.03 | 143 | 35729   | 20.66 | ng/ul | 0.04 |
| 26) 2,4-Dichlorophenol-d3      | 10.60 | 165 | 65823m  | 19.43 | ng/ul | 0.07 |
| 29) 4-Chloroaniline-d4         | 11.11 | 131 | 5245m   | 1.57  | ng/ul | 0.07 |
| 44) Dimethylphthalate-d6       | 14.18 | 166 | 172055  | 18.45 | ng/ul | 0.06 |
| 47) Acenaphthylene-d8          | 14.45 | 160 | 220051  | 19.30 | ng/ul | 0.03 |
| 52) 4-Nitrophenol-d4           | 15.06 | 143 | 25990m  | 16.40 | ng/ul | 0.16 |
| 58) Fluorene-d10               | 15.74 | 176 | 151051  | 18.15 | ng/ul | 0.03 |
| 63) 4,6-Dinitro-2-methylphenol | 15.88 | 200 | 8707m   | 5.20  | ng/ul | 0.06 |
| 71) Anthracene-d10             | 17.60 | 188 | 241245  | 20.61 | ng/ul | 0.03 |
| 79) Pyrene-d10                 | 19.89 | 212 | 287827  | 15.16 | ng/ul | 0.04 |
| 90) Benzo(a)pyrene-d12         | 25.06 | 264 | 259858m | 21.40 | ng/ul | 0.28 |

|                       |       |     |        |        |        |    |
|-----------------------|-------|-----|--------|--------|--------|----|
| Target Compounds      |       |     |        | Ovalue |        |    |
| 4) Benzaldehyde       | 7.28  | 77  | 35751  | 13.923 | ng/ul  | 85 |
| 6) Phenol             | 7.36  | 94  | 16828  | 3.732  | ng/ul  | 92 |
| 14) Acetophenone      | 8.97  | 105 | 15713  | 2.816  | ng/ul# | 83 |
| 16) 4-Methylphenol    | 8.95  | 108 | 16459  | 4.675  | ng/ul  | 91 |
| 45) Dimethylphthalate | 14.23 | 163 | 12013m | 1.246  | ng/ul  |    |
| 53) 4-Nitrophenol     | 15.03 | 109 | 33462m | 18.701 | ng/ul  |    |

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

