

Data Path : Z:\HPCHEM1\BNA G\DATA\BG111715\
 Data File : BG019676.D
 Acq On : 18 Nov 2015 1:29
 Operator : UM/NP
 Sample : G4427-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampled :
 BC795

Manual Integrations
 APPROVED

mohammad
 11/19/2015 8:27:09 AM

Quant Time: Nov 18 03:41:31 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG111615.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 18 03:01:13 2015
 Response via : Initial Calibration

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.15 | 152 | 18644 | 20.00 | ng/ul | 0.00 |
| 18) Naphthalene-d8 | 10.98 | 136 | 87644 | 20.00 | ng/ul | 0.00 |
| 36) Acenaphthene-d10 | 14.79 | 164 | 93288 | 20.00 | ng/ul | 0.00 |
| 62) Phenanthrene-d10 | 17.52 | 188 | 274334 | 20.00 | ng/ul | 0.00 |
| 78) Chrysene-d12 | 21.79 | 240 | 339274 | 20.00 | ng/ul | 0.00 |
| 86) Perylene-d12 | 25.11 | 264 | 329646 | 20.00 | ng/ul | 0.00 |

System Monitoring Compounds

| | | | | | | |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8 | 3.44 | 96 | 3248 | 7.44 | ng/uL | 0.00 |
| 5) Phenol-d5 | 7.29 | 99 | 70073 | 36.69 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d | 7.46 | 67 | 48362 | 38.65 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4 | 7.67 | 132 | 44516 | 36.35 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8 | 8.85 | 113 | 59868 | 37.86 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5 | 9.32 | 128 | 24552 | 34.12 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4 | 10.05 | 143 | 28543 | 35.24 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3 | 10.59 | 165 | 61251 | 34.98 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4 | 11.10 | 131 | 67394 | 40.54 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6 | 14.18 | 166 | 241950 | 28.68 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8 | 14.48 | 160 | 232641 | 25.44 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4 | 14.97 | 143 | 40672 | 30.60 | ng/ul | 0.00 |
| 58) Fluorene-d10 | 15.77 | 176 | 183626 | 24.32 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.88 | 200 | 40387 | 22.79 | ng/ul | 0.00 |
| 71) Anthracene-d10 | 17.62 | 188 | 325868 | 24.47 | ng/ul | 0.00 |
| 79) Pyrene-d10 | 19.89 | 212 | 347708 | 22.49 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12 | 24.88 | 264 | 334570 | 19.44 | ng/ul | 0.00 |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Ovalue |
|----------------------------|-------|------|----------|-------|--------|--------|
| 6) Phenol | 7.32 | 94 | 4175 | 2.02 | ng/ul# | 92 |
| 45) Dimethylphthalate | 14.22 | 163 | 219653 | 26.26 | ng/ul | 100 |
| 70) Phenanthrene | 17.57 | 178 | 19117 | 1.30 | ng/ul | 98 |
| 77) Fluoranthene | 19.56 | 202 | 113147 | 6.05 | ng/ul | 99 |
| 80) Pyrene | 19.92 | 202 | 98129 | 4.95 | ng/ul | 98 |
| 83) Benzo(a)anthracene | 21.77 | 228 | 89071 | 4.33 | ng/ul | 94 |
| 85) Chrysene | 21.84 | 228 | 66111 | 3.41 | ng/ul | 98 |
| 88) Benzo(b)fluoranthene | 24.05 | 252 | 85236 | 4.20 | ng/ul | 98 |
| 89) Benzo(k)fluoranthene | 24.11 | 252 | 30613m | 1.56 | ng/ul | |
| 91) Benzo(a)pyrene | 24.95 | 252 | 69068 | 3.53 | ng/ul# | 95 |
| 92) Indeno(1,2,3-cd)pyrene | 28.89 | 276 | 38974m | 1.78 | ng/ul | |
| 94) Benzo(g,h,i)perylene | 30.09 | 276 | 31923m | 1.76 | ng/ul | |

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

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