

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE111816\
 Data File : BE092569.D
 Acq On : 18 Nov 2016 16:57
 Operator : UM/SJ
 Sample : PB94794BSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampled :
 PB94794BSD

Manual Integrations
 APPROVED

sohil
 11/21/2016 3:43:09 PM

Quant Time: Nov 18 23:38:06 2016
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE111616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 16 15:43:18 2016
 Response via : Initial Calibration

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------|-------|------|----------|------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.14 | 152 | 10026 | 5.00 | ng | -0.02 |
| 7) Naphthalene-d8 | 10.93 | 136 | 45697 | 5.00 | ng | 0.00 |
| 12) Acenaphthene-d10 | 14.80 | 164 | 29360 | 5.00 | ng | 0.00 |
| 18) Phenanthrene-d10 | 17.56 | 188 | 53078 | 5.00 | ng | 0.00 |
| 24) Chrysene-d12 | 21.72 | 240 | 68066 | 5.00 | ng | 0.00 |
| 31) Perylene-d12 | 24.08 | 264 | 59429 | 5.00 | ng | -0.01 |

System Monitoring Compounds

| | | | | | | |
|--------------------------|-------|-----|-------|-------|----|-------|
| 4) 2-Fluorophenol | 5.65 | 112 | 20257 | 11.96 | ng | -0.02 |
| 5) Phenol-d6 | 7.31 | 99 | 29680 | 10.47 | ng | -0.05 |
| 8) Nitrobenzene-d5 | 9.31 | 82 | 13601 | 4.59 | ng | -0.05 |
| 13) 2,4,6-Tribromophenol | 16.28 | 330 | 6842 | 10.39 | ng | -0.02 |
| 14) 2-Fluorobiphenyl | 13.41 | 172 | 30495 | 4.26 | ng | -0.02 |
| 26) Terphenyl-d14 | 20.16 | 244 | 35205 | 4.52 | ng | -0.02 |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|--------------------------------|-------|------|----------|------|-------|--------|
| 2) 1,4-Dioxane | 3.39 | 88 | 4311 | 3.64 | ng | # 52 |
| 3) n-Nitrosodimethylamine | 3.73 | 42 | 6909 | 4.15 | ng | # 97 |
| 6) bis(2-Chloroethyl)ether | 7.56 | 93 | 12475 | 4.09 | ng | 86 |
| 9) Naphthalene | 10.97 | 128 | 37023 | 3.93 | ng | # 95 |
| 10) Hexachlorobutadiene | 11.26 | 225 | 5448 | 3.80 | ng | 100 |
| 11) 2-Methylnaphthalene | 12.60 | 142 | 24680 | 4.10 | ng | 93 |
| 15) Acenaphthylene | 14.51 | 152 | 165471 | 4.05 | ng | 100 |
| 16) Acenaphthene | 14.85 | 154 | 24306 | 3.64 | ng | 92 |
| 17) Fluorene | 15.85 | 166 | 32363 | 3.92 | ng | 99 |
| 19) Hexachlorobenzene | 16.87 | 284 | 9732 | 4.64 | ng | # 65 |
| 20) Pentachlorophenol | 17.21 | 266 | 7243 | 7.36 | ng | # 85 |
| 21) Phenanthrene | 17.58 | 178 | 53391 | 4.50 | ng | # 94 |
| 22) Anthracene | 17.67 | 178 | 51378 | 4.59 | ng | 99 |
| 23) Fluoranthene | 19.58 | 202 | 217373 | 3.92 | ng | 98 |
| 25) Pyrene | 19.95 | 202 | 230871 | 4.27 | ng | 99 |
| 27) Benzo(a)anthracene | 21.70 | 228 | 263986m | 4.11 | ng | |
| 28) Chrysene | 21.75 | 228 | 239368m | 4.14 | ng | |
| 29) Bis(2-ethylhexyl)phthalate | 21.61 | 149 | 22725 | 5.14 | ng | # 75 |
| 30) Indeno(1,2,3-cd)pyrene | 26.55 | 276 | 298018 | 4.72 | ng | 98 |
| 32) Benzo(b)fluoranthene | 23.35 | 252 | 59507 | 4.13 | ng | # 97 |
| 33) Benzo(k)fluoranthene | 23.39 | 252 | 56610 | 3.75 | ng | 94 |
| 34) Benzo(a)pyrene | 23.96 | 252 | 57112 | 4.28 | ng | # 94 |
| 35) Dibenzo(a,h)anthracene | 26.56 | 278 | 61326 | 4.65 | ng | 96 |
| 36) Benzo(g,h,i)perylene | 27.31 | 276 | 258350 | 4.62 | ng | 97 |

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

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