Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA E\DATA\BE022318\
Data File : BE095675.D
Acq On : 23 Feb 2018 10:17
Operator : SJ/JU
Sample : SSTDICC0.8
Misc :
ALS Vial : 5 Sample Multiplier: 1

**Instrument :** BNA_E
**ClientSampleId :** SSTDICC0.8

Quant Time: Feb 23 11:06:40 2018
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE022318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Feb 23 10:57:33 2018
Response via : Initial Calibration

| Internal Standards | R.T. | QIon | Response | Conc | Units | | Dev(Min) |
|---|---|---|---|---|---|---|---|
| 1) 1,4-Dichlorobenzene-d4 | 8.43 | 152 | 1172 | 0.40 | ng | | 0.00 |
| 7) Naphthalene-d8 | 11.26 | 136 | 4687 | 0.40 | ng | | -0.02 |
| 13) Acenaphthene-d10 | 15.00 | 164 | 2644 | 0.40 | ng | | 0.00 |
| 19) Phenanthrene-d10 | 17.71 | 188 | 5930 | 0.40 | ng | | 0.01 |
| 27) Chrysene-d12 | 21.79 | 240 | 6247 | 0.40 | ng | | 0.00 |
| 34) Perylene-d12 | 24.42 | 264 | 5506 | 0.40 | ng | | 0.00 |
| **System Monitoring Compounds** | | | | | | | |
| 4) 2-Fluorophenol | 5.92 | 112 | 2413 | 0.68 | ng | | 0.00 |
| 5) Phenol-d6 | 7.55 | 99 | 3325 | 0.68 | ng | | 0.00 |
| 8) Nitrobenzene-d5 | 9.60 | 82 | 3492 | 0.70 | ng | | -0.02 |
| 11) 2-Methylnaphthalene-d10 | 12.80 | 152 | 6066 | 0.78 | ng | | 0.00 |
| 14) 2,4,6-Tribromophenol | 16.47 | 330 | 751 | 0.62 | ng | | 0.00 |
| 15) 2-Fluorobiphenyl | 13.64 | 172 | 9263 | 0.81 | ng | | 0.00 |
| 25) Fluoranthene-d10 | 19.69 | 212 | 59792 | 0.77 | ng | | 0.00 |
| 29) Terphenyl-d14 | 20.25 | 244 | 9802 | 0.74 | ng | | 0.00 |
| **Target Compounds** | | | | | | | Qvalue |
| 2) 1,4-Dioxane | 3.65 | 88 | 1411 | 0.73 | ng | # | 90 |
| 3) n-Nitrosodimethylamine | 3.99 | 42 | 1823 | 0.75 | ng | # | 85 |
| 6) bis(2-Chloroethyl)ether | 7.83 | 93 | 3380 | 0.75 | ng | | 96 |
| 9) Naphthalene | 11.31 | 128 | 41708 | 0.77 | ng | | 99 |
| 10) Hexachlorobutadiene | 11.57 | 225 | 2522 | 0.86 | ng | | 96 |
| 12) 2-Methylnaphthalene | 12.88 | 142 | 7222 | 0.78 | ng | | 98 |
| 16) Acenaphthylene | 14.73 | 152 | 11520 | 0.76 | ng | | 99 |
| 17) Acenaphthene | 15.06 | 154 | 7400 | 0.79 | ng | | 97 |
| 18) Fluorene | 16.02 | 166 | 9097 | 0.77 | ng | # | 78 |
| 20) 4-Bromophenyl-phenylether | 16.89 | 248 | 2866 | 0.80 | ng | # | 71 |
| 21) Hexachlorobenzene | 17.03 | 284 | 2906 | 0.79 | ng | # | 82 |
| 22) Pentachlorophenol | 17.35 | 266 | 694 | 0.68 | ng | # | 75 |
| 23) Phenanthrene | 17.74 | 178 | 14401 | 0.79 | ng | | 98 |
| 24) Anthracene | 17.83 | 178 | 13077 | 0.77 | ng | | 95 |
| 26) Fluoranthene | 19.72 | 202 | 17840 | 0.78 | ng | # | 97 |
| 28) Pyrene | 20.06 | 202 | 20694 | 0.79 | ng | | 98 |
| 30) Benzo(a)anthracene | 21.78 | 228 | 16089 | 0.77 | ng | | 97 |
| 31) Chrysene | 21.84 | 228 | 15797 | 0.77 | ng | | 99 |
| 32) Bis(2-ethylhexyl)phthalate | 21.65 | 149 | 26438 | 0.55 | ng | | 100 |
| 33) Indeno(1,2,3-cd)pyrene | 27.23 | 276 | 15473 | 0.78 | ng | # | 94 |
| 35) Benzo(b)fluoranthene | 23.61 | 252 | 15181 | 0.81 | ng | # | 88 |
| 36) Benzo(k)fluoranthene | 23.66 | 252 | 15434 | 0.80 | ng | | 95 |
| 37) Benzo(a)pyrene | 24.30 | 252 | 13884 | 0.79 | ng | # | 93 |
| 38) Dibenzo(a,h)anthracene | 27.25 | 278 | 12535 | 0.82 | ng | # | 96 |
| 39) Benzo(g,h,i)perylene | 28.10 | 276 | 13102 | 0.80 | ng | # | 92 |

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

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