

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030121\
 Data File : BM028956.D
 Acq On : 01 Mar 2021 09:46
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD.444

Quant Time: Mar 01 10:25:34 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM020521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 01 10:25:15 2021
 Response via : Initial Calibration

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------|-------|------|----------|------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.78 | 152 | 273 | 0.40 | ng/ul | 0.00 |
| 2) Naphthalene-d8 | 10.58 | 136 | 1138 | 0.40 | ng/ul | 0.00 |
| 6) Acenaphthene-d10 | 14.44 | 164 | 634 | 0.40 | ng/ul | 0.00 |
| 10) Phenanthrene-d10 | 17.18 | 188 | 1267 | 0.40 | ng/ul | 0.00 |
| 16) Chrysene-d12 | 21.35 | 240 | 1019 | 0.40 | ng/ul | 0.00 |
| 20) Perylene-d12 | 23.54 | 264 | 978 | 0.40 | ng/ul | 0.00 |

| System Monitoring Compounds | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------|-------|------|----------|------|-------|----------|
| 4) 2-Methylnaphthalene-d10 | 12.17 | 152 | 695 | 0.43 | ng/ul | 0.00 |
| 14) Fluoranthene-d10 | 19.20 | 212 | 1296 | 0.40 | ng/ul | 0.00 |

| Target Compounds | R.T. | QIon | Response | Conc | Units | Ovalue |
|----------------------------|-------|------|----------|-------|--------|--------|
| 3) Naphthalene | 10.63 | 128 | 1187 | 0.419 | ng/ul | 94 |
| 5) 2-Methylnaphthalene | 12.25 | 142 | 798 | 0.421 | ng/ul | 99 |
| 7) Acenaphthylene | 14.16 | 152 | 993 | 0.384 | ng/ul | 95 |
| 8) Acenaphthene | 14.50 | 153 | 844 | 0.411 | ng/ul | 99 |
| 9) Fluorene | 15.49 | 166 | 949 | 0.417 | ng/ul | 99 |
| 11) Pentachlorophenol | 16.85 | 266 | 81 | 0.245 | ng/ul | 97 |
| 12) Phenanthrene | 17.22 | 178 | 1435 | 0.405 | ng/ul | 97 |
| 13) Anthracene | 17.31 | 178 | 1314 | 0.389 | ng/ul | 97 |
| 15) Fluoranthene | 19.23 | 202 | 1511 | 0.375 | ng/ul | 96 |
| 17) Pyrene | 19.60 | 202 | 1537 | 0.415 | ng/ul | 98 |
| 18) Benzo(a)anthracene | 21.33 | 228 | 1307 | 0.376 | ng/ul | 98 |
| 19) Chrysene | 21.38 | 228 | 1361 | 0.394 | ng/ul | 98 |
| 21) Benzo(b)fluoranthene | 22.88 | 252 | 1328 | 0.397 | ng/ul | 94 |
| 22) Benzo(k)fluoranthene | 22.93 | 252 | 1366 | 0.409 | ng/ul# | 92 |
| 23) Benzo(a)pyrene | 23.45 | 252 | 1190 | 0.399 | ng/ul# | 92 |
| 24) Indeno(1,2,3-cd)pyrene | 25.75 | 276 | 1484 | 0.382 | ng/ul# | 100 |
| 25) Dibenzo(a,h)anthracene | 25.76 | 278 | 1241 | 0.381 | ng/ul | 94 |
| 26) Benzo(g,h,i)perylene | 26.43 | 276 | 1295 | 0.401 | ng/ul | 98 |

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

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