

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111324\
 Data File : BM048693.D
 Acq On : 15 Nov 2024 03:34
 Operator : RC/JU
 Sample : PB164923BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS923

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/15/2024
 Supervised By :mohammad ahmed 11/15/2024

Quant Time: Nov 15 04:16:18 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM110624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 13 11:38:55 2024
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| Internal Standards | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 7.775 | 152 | 2729m | 0.400 | ng/ul | 0.04 |
| 4) Naphthalene-d8 | 10.507 | 136 | 9195 | 0.400 | ng/ul | 0.01 |
| 9) Acenaphthene-d10 | 14.326 | 164 | 5618 | 0.400 | ng/ul | 0.00 |
| 13) Phenanthrene-d10 | 17.090 | 188 | 11359 | 0.400 | ng/ul | 0.00 |
| 17) Chrysene-d12 | 21.268 | 240 | 8967 | 0.400 | ng/ul | 0.00 |
| 23) Perylene-d12 | 23.487 | 264 | 9130 | 0.400 | ng/ul | 0.00 |
| System Monitoring Compounds | | | | | | |
| 3) 1,4-Dioxane-d8 | 3.122 | 96 | 2772 | 0.842 | ng/ul | 0.00 |
| 6) 2-Methylnaphthalene-d10 | 12.113 | 152 | 4756 | 0.400 | ng/ul | 0.01 |
| 18) Fluoranthene-d10 | 19.114 | 212 | 11106 | 0.441 | ng/ul | 0.00 |
| Target Compounds | | | | | | |
| | | | | | | Qvalue |
| 2) 1,4-Dioxane | 3.155 | 88 | 8031 | 2.122 | ng/ul# | 63 |
| 5) Naphthalene | 10.557 | 128 | 9200 | 0.399 | ng/ul | 95 |
| 7) 2-Methylnaphthalene | 12.190 | 142 | 5233 | 0.361 | ng/ul | 92 |
| 8) 1-Methylnaphthalene | 12.383 | 142 | 6278 | 0.423 | ng/ul | 100 |
| 10) Acenaphthylene | 14.053 | 152 | 8545 | 0.400 | ng/ul | 99 |
| 11) Acenaphthene | 14.391 | 153 | 6917 | 0.405 | ng/ul | 99 |
| 12) Fluorene | 15.404 | 166 | 7393 | 0.399 | ng/ul | 99 |
| 14) Pentachlorophenol | 16.753 | 266 | 2218 | 0.532 | ng/ul | 99 |
| 15) Phenanthrene | 17.133 | 178 | 10815 | 0.362 | ng/ul | 98 |
| 16) Anthracene | 17.251 | 178 | 9486 | 0.342 | ng/ul | 97 |
| 19) Fluoranthene | 19.146 | 202 | 14431 | 0.435 | ng/ul | 99 |
| 20) Pyrene | 19.504 | 202 | 15640 | 0.434 | ng/ul | 99 |
| 21) Benzo(a)anthracene | 21.262 | 228 | 9232 | 0.314 | ng/ul | 99 |
| 22) Chrysene | 21.303 | 228 | 17351 | 0.493 | ng/ul | 99 |
| 24) Benzo(b)fluoranthene | 22.820 | 252 | 12532 | 0.440 | ng/ul | 97 |
| 25) Benzo(k)fluoranthene | 22.864 | 252 | 16287 | 0.488 | ng/ul | 99 |
| 26) Benzo(a)pyrene | 23.402 | 252 | 12173 | 0.438 | ng/ul | 99 |
| 27) Indeno(1,2,3-cd)pyrene | 25.759 | 276 | 14658 | 0.378 | ng/ul# | 97 |
| 28) Dibenzo(a,h)anthracene | 25.779 | 278 | 10664 | 0.353 | ng/ul | 94 |
| 29) Benzo(g,h,i)perylene | 26.436 | 276 | 13476 | 0.414 | ng/ul | 99 |

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

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