

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061323\
 Data File : BN025837.D
 Acq On : 14 Jun 2023 11:46
 Operator : MA/JU
 Sample : 02950-09DL 5X
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EW986DL

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 06/14/2023
 Supervised By :mohammad ahmed 06/14/2023

Quant Time: Jun 14 18:58:59 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN060923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 14 18:25:53 2023
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------|--------|------|----------|-------|---------|----------|
| Internal Standards | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 8.001 | 152 | 8881 | 0.400 | ng/ul | 0.00 |
| 4) Naphthalene-d8 | 10.814 | 136 | 26943 | 0.400 | ng/ul # | 0.00 |
| 9) Acenaphthene-d10 | 14.636 | 164 | 15611 | 0.400 | ng/ul | 0.00 |
| 13) Phenanthrene-d10 | 17.381 | 188 | 31336 | 0.400 | ng/ul | 0.00 |
| 17) Chrysene-d12 | 21.562 | 240 | 17568 | 0.400 | ng/ul | 0.00 |
| 23) Perylene-d12 | 24.029 | 264 | 16600 | 0.400 | ng/ul | -0.02 |
| System Monitoring Compounds | | | | | | |
| 3) 1,4-Dioxane-d8 | 3.352 | 96 | 6617 | 0.758 | ng/ul | 0.00 |
| 6) 2-Methylnaphthalene-d10 | 12.403 | 152 | 1665 | 0.049 | ng/ul | 0.00 |
| 18) Fluoranthene-d10 | 19.401 | 212 | 3216 | 0.058 | ng/ul | 0.00 |
| Target Compounds | | | | | | |
| | | | | | | Qvalue |
| 5) Naphthalene | 10.864 | 128 | 1948 | 0.028 | ng/ul# | 82 |
| 7) 2-Methylnaphthalene | 12.475 | 142 | 962 | 0.023 | ng/ul | 100 |
| 8) 1-Methylnaphthalene | 12.689 | 142 | 9080 | 0.200 | ng/ul | 99 |
| 10) Acenaphthylene | 14.358 | 152 | 9803 | 0.159 | ng/ul# | 87 |
| 11) Acenaphthene | 14.701 | 153 | 21790 | 0.423 | ng/ul | 99 |
| 12) Fluorene | 15.681 | 166 | 23731 | 0.414 | ng/ul | 98 |
| 15) Phenanthrene | 17.419 | 178 | 187532 | 2.052 | ng/ul | 100 |
| 16) Anthracene | 17.512 | 178 | 40406 | 0.524 | ng/ul | 99 |
| 19) Fluoranthene | 19.434 | 202 | 127439 | 1.586 | ng/ul | 99 |
| 20) Pyrene | 19.796 | 202 | 188140 | 2.262 | ng/ul | 99 |
| 21) Benzo(a)anthracene | 21.544 | 228 | 56545m | 1.009 | ng/ul | |
| 22) Chrysene | 21.600 | 228 | 79974 | 1.248 | ng/ul | 96 |
| 24) Benzo(b)fluoranthene | 23.278 | 252 | 41731m | 0.615 | ng/ul | |
| 25) Benzo(k)fluoranthene | 23.322 | 252 | 14298m | 0.210 | ng/ul | |
| 26) Benzo(a)pyrene | 23.918 | 252 | 39022 | 0.678 | ng/ul | 92 |
| 27) Indeno(1,2,3-cd)pyrene | 26.605 | 276 | 16416 | 0.250 | ng/ul# | 93 |
| 28) Dibenzo(a,h)anthracene | 26.609 | 278 | 5672 | 0.114 | ng/ul# | 71 |
| 29) Benzo(g,h,i)perylene | 27.400 | 276 | 18317 | 0.311 | ng/ul | 96 |

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

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