

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072324\
 Data File : BM046776.D
 Acq On : 23 Jul 2024 13:55
 Operator : MA/JU
 Sample : P3238-18DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
ClientSampleId :
 A4CG8DL

Manual Integrations
APPROVED
 Reviewed By :Jagrut Upadhyay 07/24/2024
 Supervised By :mohammad ahmed 07/25/2024

Quant Time: Jul 23 14:51:14 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM071024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 19 11:38:04 2024
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|------------------------------------|--------|------|----------|-------|--------|----------|
| Internal Standards | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 7.481 | 152 | 2831 | 0.400 | ng/ul | 0.00 |
| 4) Naphthalene-d8 | 10.237 | 136 | 7680 | 0.400 | ng/ul | #-0.01 |
| 9) Acenaphthene-d10 | 14.122 | 164 | 5209 | 0.400 | ng/ul | 0.00 |
| 13) Phenanthrene-d10 | 16.878 | 188 | 11394 | 0.400 | ng/ul | 0.00 |
| 17) Chrysene-d12 | 21.088 | 240 | 8844 | 0.400 | ng/ul | # 0.00 |
| 23) Perylene-d12 | 23.218 | 264 | 8633 | 0.400 | ng/ul | -0.01 |
| System Monitoring Compounds | | | | | | |
| 3) 1,4-Dioxane-d8 | 3.105 | 96 | 1247 | 0.536 | ng/ul | 0.00 |
| 6) 2-Methylnaphthalene-d10 | 11.837 | 152 | 491 | 0.040 | ng/ul | -0.01 |
| 18) Fluoranthene-d10 | 18.917 | 212 | 1213 | 0.045 | ng/ul | 0.00 |
| Target Compounds | | | | | | |
| | | | | | | Qvalue |
| 5) Naphthalene | 10.287 | 128 | 917 | 0.047 | ng/ul# | 88 |
| 7) 2-Methylnaphthalene | 11.914 | 142 | 350 | 0.025 | ng/ul | 97 |
| 8) 1-Methylnaphthalene | 12.134 | 142 | 309 | 0.022 | ng/ul | 96 |
| 10) Acenaphthylene | 13.840 | 152 | 1766 | 0.074 | ng/ul | 94 |
| 11) Acenaphthene | 14.186 | 153 | 1371 | 0.086 | ng/ul | 97 |
| 12) Fluorene | 15.181 | 166 | 1728 | 0.088 | ng/ul | 98 |
| 15) Phenanthrene | 16.916 | 178 | 40043 | 1.232 | ng/ul | 98 |
| 16) Anthracene | 17.009 | 178 | 7090 | 0.223 | ng/ul | 99 |
| 19) Fluoranthene | 18.950 | 202 | 88325 | 2.431 | ng/ul | 98 |
| 20) Pyrene | 19.312 | 202 | 64948 | 1.651 | ng/ul | 98 |
| 21) Benzo(a)anthracene | 21.071 | 228 | 37496 | 0.958 | ng/ul | 98 |
| 22) Chrysene | 21.123 | 228 | 38465 | 1.016 | ng/ul | 99 |
| 24) Benzo(b)fluoranthene | 22.593 | 252 | 47732m | 1.384 | ng/ul | |
| 25) Benzo(k)fluoranthene | 22.628 | 252 | 17951m | 0.530 | ng/ul | |
| 26) Benzo(a)pyrene | 23.128 | 252 | 19206 | 0.677 | ng/ul | 96 |
| 27) Indeno(1,2,3-cd)pyrene | 25.327 | 276 | 17084 | 0.396 | ng/ul# | 100 |
| 28) Dibenzo(a,h)anthracene | 25.340 | 278 | 5358 | 0.157 | ng/ul# | 79 |
| 29) Benzo(g,h,i)perylene | 25.967 | 276 | 2898 | 0.084 | ng/ul# | 78 |

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

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