

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071224\
 Data File : BM046584.D
 Acq On : 13 Jul 2024 15:41
 Operator : MA/JU
 Sample : P3035-24
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_M
ClientSampleId :
 A4CC9

Manual Integrations
APPROVED
 Reviewed By :Jagrut Upadhyay 07/15/2024
 Supervised By :mohammad ahmed 07/16/2024

Quant Time: Jul 15 09:02:50 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 : Mon Jul 15 08:57:27 2024
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|------------------------------------|--------|------|----------|-------|---------|----------|
| Internal Standards | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 7.510 | 152 | 3795 | 0.400 | ng/ul | 0.00 |
| 4) Naphthalene-d8 | 10.270 | 136 | 10543 | 0.400 | ng/ul # | 0.00 |
| 9) Acenaphthene-d10 | 14.154 | 164 | 6537 | 0.400 | ng/ul | 0.00 |
| 13) Phenanthrene-d10 | 16.904 | 188 | 14004m | 0.400 | ng/ul | 0.00 |
| 17) Chrysene-d12 | 21.111 | 240 | 9764 | 0.400 | ng/ul # | 0.00 |
| 23) Perylene-d12 | 23.253 | 264 | 9598 | 0.400 | ng/ul | 0.01 |
| System Monitoring Compounds | | | | | | |
| 3) 1,4-Dioxane-d8 | 3.117 | 96 | 8158 | 2.616 | ng/ul | 0.00 |
| 6) 2-Methylnaphthalene-d10 | 11.870 | 152 | 3748 | 0.220 | ng/ul | 0.00 |
| 18) Fluoranthene-d10 | 18.945 | 212 | 8323 | 0.280 | ng/ul | 0.00 |
| Target Compounds | | | | | | |
| | | | | | | Qvalue |
| 10) Acenaphthylene | 13.867 | 152 | 1744 | 0.059 | ng/ul# | 91 |
| 11) Acenaphthene | 14.214 | 153 | 1011 | 0.051 | ng/ul | 98 |
| 12) Fluorene | 15.209 | 166 | 1318 | 0.053 | ng/ul# | 94 |
| 15) Phenanthrene | 16.946 | 178 | 29662 | 0.742 | ng/ul | 99 |
| 16) Anthracene | 17.039 | 178 | 6505 | 0.167 | ng/ul | 99 |
| 19) Fluoranthene | 18.973 | 202 | 66889 | 1.668 | ng/ul# | 98 |
| 20) Pyrene | 19.340 | 202 | 53433 | 1.231 | ng/ul | 98 |
| 21) Benzo(a)anthracene | 21.096 | 228 | 26662 | 0.617 | ng/ul | 98 |
| 22) Chrysene | 21.149 | 228 | 28773 | 0.688 | ng/ul | 98 |
| 24) Benzo(b)fluoranthene | 22.622 | 252 | 35221m | 0.919 | ng/ul | |
| 25) Benzo(k)fluoranthene | 22.660 | 252 | 12768m | 0.339 | ng/ul | |
| 26) Benzo(a)pyrene | 23.163 | 252 | 21263 | 0.674 | ng/ul | 96 |
| 27) Indeno(1,2,3-cd)pyrene | 25.373 | 276 | 16726 | 0.348 | ng/ul# | 100 |
| 28) Dibenzo(a,h)anthracene | 25.383 | 278 | 3949 | 0.104 | ng/ul# | 32 |
| 29) Benzo(g,h,i)perylene | 26.017 | 276 | 16765 | 0.438 | ng/ul# | 95 |

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

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