

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042224\
 Data File : BM045506.D
 Acq On : 23 Apr 2024 10:11
 Operator : MA/JU
 Sample : P2104-04
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AA6

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 04/25/2024
 Supervised By :mohammad ahmed 04/26/2024

Quant Time: Apr 24 04:49:18 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM042024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 24 04:47:50 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.555	152	1250	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.314	136	3903	0.400	ng/ul	#-0.03
9) Acenaphthene-d10	14.200	164	2199	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.963	188	4480m	0.400	ng/ul	-0.03
17) Chrysene-d12	21.199	240	2999	0.400	ng/ul	0.00
23) Perylene-d12	23.388	264	3129m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.163	96	5659	3.378	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.942	152	1260	0.245	ng/ul	-0.01
18) Fluoranthene-d10	19.010	212	2546	0.270	ng/ul	-0.01
Target Compounds						
15) Phenanthrene	17.005	178	815m	0.062	ng/ul	Qvalue
19) Fluoranthene	19.038	202	1362	0.098	ng/ul	95
20) Pyrene	19.401	202	1526	0.107	ng/ul#	87
21) Benzo(a)anthracene	21.190	228	695	0.073	ng/ul#	75
22) Chrysene	21.237	228	624	0.042	ng/ul#	74
24) Benzo(b)fluoranthene	22.727	252	798m	0.076	ng/ul	
27) Indeno(1,2,3-cd)pyrene	25.608	276	409	0.026	ng/ul#	87
29) Benzo(g,h,i)perylene	26.276	276	496	0.035	ng/ul#	11

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

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