

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042224\
 Data File : BM045510.D
 Acq On : 23 Apr 2024 12:37
 Operator : MA/JU
 Sample : P2104-11
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AB1

Manual Integrations
 APPROVED

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

Quant Time: Apr 24 04:50:16 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	7.551	152	1620	0.400	ng/ul	-0.02	
4) Naphthalene-d8	10.314	136	4477	0.400	ng/ul	#-0.03	
9) Acenaphthene-d10	14.205	164	2112	0.400	ng/ul	0.00	
13) Phenanthrene-d10	16.967	188	3985m	0.400	ng/ul	-0.02	
17) Chrysene-d12	21.190	240	3504	0.400	ng/ul	-0.01	
23) Perylene-d12	23.373	264	4517m	0.400	ng/ul	-0.02	
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.167	96	7601	3.501	ng/ul	0.00	
6) 2-Methylnaphthalene-d10	11.925	152	1432	0.243	ng/ul	-0.03	
18) Fluoranthene-d10	19.010	212	2269	0.206	ng/ul	-0.01	
Target Compounds							
							Qvalue
10) Acenaphthylene	13.923	152	485	0.047	ng/ul#		69
15) Phenanthrene	17.005	178	895m	0.077	ng/ul		
19) Fluoranthene	19.038	202	3131	0.194	ng/ul		97
20) Pyrene	19.401	202	2568m	0.155	ng/ul		
21) Benzo(a)anthracene	21.176	228	1680	0.150	ng/ul#		81
22) Chrysene	21.225	228	1553	0.090	ng/ul#		84
24) Benzo(b)fluoranthene	22.721	252	2972m	0.196	ng/ul		
25) Benzo(k)fluoranthene	22.759	252	1146m	0.060	ng/ul		
26) Benzo(a)pyrene	23.274	252	1951	0.124	ng/ul#		67
27) Indeno(1,2,3-cd)pyrene	25.568	276	2000	0.087	ng/ul#		95
29) Benzo(g,h,i)perylene	26.232	276	2016	0.098	ng/ul#		74

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

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