

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092823\
 Data File : BM042112.D
 Acq On : 28 Sep 2023 13:36
 Operator : MA/JU
 Sample : 04417-37
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EZYE4

Quant Time: Sep 28 16:02:19 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM092223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 22 13:50:58 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.941	152	2810	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.751	136	7586	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.583	164	3948	0.400	ng/ul	-0.02
13) Phenanthrene-d10	17.337	188	8961	0.400	ng/ul	-0.02
17) Chrysene-d12	21.512	240	6729	0.400	ng/ul	-0.01
23) Perylene-d12	23.912	264	7935	0.400	ng/ul	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	11117	2.903	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.335	152	2780	0.293	ng/ul	-0.02
18) Fluoranthene-d10	19.361	212	5837	0.211	ng/ul	-0.01
Target Compounds						
					Qvalue	
5) Naphthalene	10.806	128	661	0.026	ng/ul#	69
15) Phenanthrene	17.379	178	1846	0.044	ng/ul#	88
19) Fluoranthene	19.389	202	5632	0.074	ng/ul	98
20) Pyrene	19.757	202	6789	0.091	ng/ul	99
21) Benzo(a)anthracene	21.495	228	1858	0.036	ng/ul#	83
22) Chrysene	21.547	228	2837	0.054	ng/ul#	89
24) Benzo(b)fluoranthene	23.172	252	5775	0.095	ng/ul#	47
26) Benzo(a)pyrene	23.804	252	2527	0.050	ng/ul#	25
27) Indeno(1,2,3-cd)pyrene	26.391	276	4081	0.059	ng/ul#	88
29) Benzo(g,h,i)perylene	27.169	276	5856	0.101	ng/ul#	88

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

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