

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100323\
 Data File : BM042161.D
 Acq On : 03 Oct 2023 13:22
 Operator : MA/JU
 Sample : 04588-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BBJM6

Quant Time: Oct 03 13:52:45 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM100223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 02 19:28:41 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.933	152	3814	0.400	ng/ul	# 0.00
4) Naphthalene-d8	10.746	136	9095	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.583	164	4314	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	17.333	188	10013	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.510	240	5190	0.400	ng/ul	0.00
23) Perylene-d12	23.912	264	5887	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.360	96	21488	4.231	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.330	152	4211	0.338	ng/ul	0.00
18) Fluoranthene-d10	19.357	212	8463	0.434	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.397	88	39404	7.429	ng/ul#	80
5) Naphthalene	10.796	128	661715	21.808	ng/ul	98
7) 2-Methylnaphthalene	12.401	142	328751	18.096	ng/ul	99
8) 1-Methylnaphthalene	12.621	142	226601	12.382	ng/ul	99
10) Acenaphthylene	14.296	152	1884	0.069	ng/ul#	1
11) Acenaphthene	14.648	153	5922	0.287	ng/ul#	88
12) Fluorene	15.633	166	15344	0.660	ng/ul	98
15) Phenanthrene	17.375	178	27936	0.677	ng/ul	99
16) Anthracene	17.468	178	932	0.025	ng/ul#	66
19) Fluoranthene	19.389	202	1606	0.041	ng/ul#	74
20) Pyrene	19.757	202	3404	0.081	ng/ul#	87
21) Benzo(a)anthracene	21.492	228	1884	0.056	ng/ul#	80
22) Chrysene	21.545	228	2415	0.074	ng/ul#	80
24) Benzo(b)fluoranthene	23.173	252	982	0.028	ng/ul#	1
29) Benzo(g,h,i)perylene	27.179	276	775	0.022	ng/ul#	27

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

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