

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101022\
 Data File : BM036930.D
 Acq On : 11 Oct 2022 05:01
 Operator : CG/JU
 Sample : N4991-08
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BG0

Quant Time: Oct 11 05:31:10 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM100622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 10 00:54:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.951	152	7458	0.400	ng/ul	0.00
4) Naphthalene-d8	10.754	136	25787	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.580	164	14562	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.313	188	30061	0.400	ng/ul	0.00
17) Chrysene-d12	21.482	240	24329	0.400	ng/ul	0.00
23) Perylene-d12	23.876	264	21866	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.348	96	56091	5.333	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.338	152	17340	0.445	ng/ul	-0.01
18) Fluoranthene-d10	19.336	212	37315	0.513	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.803	128	6257	0.085	ng/ul	99
7) 2-Methylnaphthalene	12.415	142	5137	0.114	ng/ul	93
8) 1-Methylnaphthalene	12.629	142	4142	0.090	ng/ul	100
15) Phenanthrene	17.355	178	5051	0.053	ng/ul	95
19) Fluoranthene	19.363	202	4305	0.044	ng/ul#	85
21) Benzo(a)anthracene	21.465	228	2187	0.025	ng/ul#	84
22) Chrysene	21.517	228	3199	0.034	ng/ul#	89
24) Benzo(b)fluoranthene	23.145	252	3159	0.032	ng/ul#	1

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

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