

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101022\
 Data File : BM036913.D
 Acq On : 10 Oct 2022 18:26
 Operator : CG/JU
 Sample : N4951-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CB8T8

Quant Time: Oct 11 02:18:00 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	7.955	152	8017	0.400	ng/ul	0.00
4) Naphthalene-d8	10.759	136	29657	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.581	164	18494	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.318	188	37227	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.487	240	25950	0.400	ng/ul	# 0.00
23) Perylene-d12	23.890	264	20300	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.348	96	39971	3.535	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.343	152	12165	0.272	ng/ul	0.00
18) Fluoranthene-d10	19.341	212	27719	0.357	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.814	128	2568	0.030	ng/ul#	62
8) 1-Methylnaphthalene	12.635	142	1576	0.030	ng/ul#	65
10) Acenaphthylene	14.321	152	2316	0.030	ng/ul#	35
11) Acenaphthene	14.645	153	66397	0.995	ng/ul	98
12) Fluorene	15.626	166	5123	0.068	ng/ul#	81
15) Phenanthrene	17.361	178	4469	0.038	ng/ul#	45
16) Anthracene	17.453	178	4031	0.041	ng/ul#	39

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

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