

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM050322\
 Data File : BM034894.D
 Acq On : 03 May 2022 17:08
 Operator : CG/JU
 Sample : N2562-21
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW480

Quant Time: May 04 03:01:45 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM042222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 04 02:59:36 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	5133	0.400	ng/ul	0.00
4) Naphthalene-d8	10.732	136	13733	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.561	164	7952	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.300	188	15834	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.473	240	12361	0.400	ng/ul	# 0.00
23) Perylene-d12	23.861	264	13002	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.373	96	20660	3.639	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.321	152	4708	0.244	ng/ul	0.00
18) Fluoranthene-d10	19.326	212	10484	0.269	ng/ul	0.00
Target Compounds						
					Qvalue	
7) 2-Methylnaphthalene	12.393	142	722	0.028	ng/ul	95
8) 1-Methylnaphthalene	12.607	142	677	0.026	ng/ul	82
15) Phenanthrene	17.343	178	3824	0.073	ng/ul#	93
19) Fluoranthene	19.354	202	1212	0.020	ng/ul#	60
20) Pyrene	19.717	202	2556	0.043	ng/ul#	69
22) Chrysene	21.511	228	3744	0.073	ng/ul#	73
24) Benzo(b)fluoranthene	23.133	252	1766	0.030	ng/ul#	1
29) Benzo(g,h,i)perylene	27.084	276	1918	0.033	ng/ul#	61

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

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