

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100923\
 Data File : BM042244.D
 Acq On : 10 Oct 2023 03:24
 Operator : MA/JU
 Sample : 04654-18
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BBBW7

Quant Time: Oct 10 05:06:03 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 09 23:59:28 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.051	152	4744	0.400	ng/ul	0.00
4) Naphthalene-d8	10.873	136	12450	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.689	164	5071	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	17.443	188	10137	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.627	240	4234	0.400	ng/ul	# 0.00
23) Perylene-d12	24.120	264	4701	0.400	ng/ul	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.410	96	28281	4.477	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.456	152	5620	0.329	ng/ul	0.00
18) Fluoranthene-d10	19.464	212	7565	0.476	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.448	88	15543	2.356	ng/ul#	76
14) Pentachlorophenol	17.075	266	105	0.023	ng/ul	92
19) Fluoranthene	19.497	202	827	0.026	ng/ul#	65
24) Benzo(b)fluoranthene	23.351	252	560	0.020	ng/ul#	1

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

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