

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101023\
 Data File : BM042253.D
 Acq On : 10 Oct 2023 15:16
 Operator : MA/JU
 Sample : 04654-14
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BBBT9

Quant Time: Oct 10 16:40:26 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 : Tue Oct 10 14:01:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.026	152	3886	0.400	ng/ul	0.00
4) Naphthalene-d8	10.845	136	9814	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.666	164	5034	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.413	188	9266	0.400	ng/ul #	0.00
17) Chrysene-d12	21.600	240	4216	0.400	ng/ul #	0.00
23) Perylene-d12	24.079	264	4016	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	21245	4.105	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.429	152	3266	0.243	ng/ul	0.00
18) Fluoranthene-d10	19.441	212	5007	0.316	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.419	88	2956	0.547	ng/ul	96
14) Pentachlorophenol	17.046	266	30905	7.357	ng/ul	99
15) Phenanthrene	17.455	178	900	0.024	ng/ul#	41
20) Pyrene	19.808	202	1638	0.048	ng/ul#	1
24) Benzo(b)fluoranthene	23.307	252	515	0.022	ng/ul#	1
26) Benzo(a)pyrene	23.915	252	2546	0.125	ng/ul#	46

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

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