

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092223\
 Data File : BM041984.D
 Acq On : 22 Sep 2023 13:21
 Operator : MA/JU
 Sample : SSTD3.247
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.2041

Quant Time: Sep 22 13:49:34 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM092223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 22 13:22:12 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.954	152	5860	0.400	ng/ul	0.00
4) Naphthalene-d8	10.768	136	14290	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.601	164	6320	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.350	188	12153	0.400	ng/ul	0.00
17) Chrysene-d12	21.524	240	5749	0.400	ng/ul	0.00
23) Perylene-d12	23.936	264	5600	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	25643	3.214	ng/ul	0.00
6) 2-Methylnaphthalene-d10	0.000	152	0d	0.000	ng/ul	
18) Fluoranthene-d10	0.000	212	0d	0.000	ng/ul	
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.406	88	27031	3.222	ng/ul#	84
14) Pentachlorophenol	16.978	266	24260	3.951	ng/ul	100

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

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