

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101823\
 Data File : BM042348.D
 Acq On : 18 Oct 2023 14:00
 Operator : MA/JU
 Sample : SSTD3.259
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.2006

Quant Time: Oct 18 14:29:10 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM101823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 18 13:57:01 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.013	152	4407	0.400	ng/ul	0.00
4) Naphthalene-d8	10.834	136	12247	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.652	164	5259	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.400	188	9804	0.400	ng/ul	0.00
17) Chrysene-d12	21.580	240	4313	0.400	ng/ul	0.00
23) Perylene-d12	24.053	264	3782	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	21183	3.268	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	3.410	88	22561	3.275	ng/ul#	89
14) Pentachlorophenol	17.033	266	14818	4.191	ng/ul	99

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

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