

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062321\
 Data File : BM030563.D
 Acq On : 24 Jun 2021 00:52
 Operator : CG/JU
 Sample : SSTD3.206
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.2606

Quant Time: Jun 24 01:38:18 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 24 01:37:13 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.801	152	3660	0.400	ng/ul	0.00
4) Naphthalene-d8	10.585	136	13864	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.428	164	7784	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.165	188	14545	0.400	ng/ul	0.00
17) Chrysene-d12	21.328	240	12830	0.400	ng/ul	0.00
23) Perylene-d12	23.590	264	12672	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.279	96	12411	3.458	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	3.313	88	13528	3.492	ng/ul#	80
14) Pentachlorophenol	16.829	266	15131	3.337	ng/ul	95

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

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