

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082622\
 Data File : BM036424.D
 Acq On : 26 Aug 2022 13:52
 Operator : CG/JU
 Sample : SSTD3.261
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.2013

Quant Time: Aug 26 14:22:21 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM082622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 26 13:25:55 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.791	152	5507	0.400	ng/ul	0.00
4) Naphthalene-d8	10.584	136	21927	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.432	164	13384	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.174	188	29496	0.400	ng/ul	0.00
17) Chrysene-d12	21.360	240	49509	0.400	ng/ul	# 0.00
23) Perylene-d12	23.689	264	28850	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	21932	2.935	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	3.239	88	21287	2.880	ng/ul#	78
14) Pentachlorophenol	16.832	266	20927	4.493	ng/ul	98

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

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