

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112123\
 Data File : BM042941.D
 Acq On : 21 Nov 2023 19:25
 Operator : MA/JU
 Sample : SSTD3.271
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.2020

Quant Time: Nov 22 10:32:53 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM112123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 22 10:26:14 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.821	152	758	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.622	136	1976	0.400	ng/ul	#-0.02
9) Acenaphthene-d10	14.464	164	1058	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.220	188	2339	0.400	ng/ul	0.00
17) Chrysene-d12	21.406	240	1698	0.400	ng/ul	# 0.00
23) Perylene-d12	23.768	264	2010	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	3315	2.799	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	3.202	88	3759	2.834	ng/ul#	52
14) Pentachlorophenol	16.866	266	4605	3.533	ng/ul	95

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

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