

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062522\
 Data File : BM035687.D
 Acq On : 25 Jun 2022 15:01
 Operator : CG/JU
 Sample : SSTD3.255
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.2006

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 06/29/2022
 Supervised By :mohammad ahmed 07/01/2022

Quant Time: Jun 28 02:06:31 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 28 02:03:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.808	152	3461	0.400	ng/ul	-0.06
4) Naphthalene-d8	10.578	136	13623	0.400	ng/ul	-0.04
9) Acenaphthene-d10	14.418	164	8304m	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.170	188	16965	0.400	ng/ul	-0.03
17) Chrysene-d12	21.360	240	13299	0.400	ng/ul	-0.02
23) Perylene-d12	23.704	264	10608	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	13834	2.841	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						
2) 1,4-Dioxane	3.251	88	14430m	2.976	ng/ul	Qvalue
14) Pentachlorophenol	16.853	266	9961m	3.061	ng/ul	

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

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