

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM083022\
 Data File : BM036457.D
 Acq On : 29 Aug 2022 17:03
 Operator : CG/JU
 Sample : SSTD3.266
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.2020

Quant Time: Aug 29 17:39:29 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM083022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 29 16:16:27 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.746	152	3539	0.400	ng/ul	0.00
4) Naphthalene-d8	10.534	136	13789	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.385	164	7957	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.132	188	16375	0.400	ng/ul	0.00
17) Chrysene-d12	21.316	240	15195	0.400	ng/ul	0.00
23) Perylene-d12	23.619	264	12584	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.139	96	15026	3.153	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	3.177	88	14797	3.288	ng/ul#	78
14) Pentachlorophenol	16.790	266	10134	3.131	ng/ul	98

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

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