

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110921\
 Data File : BM032912.D
 Acq On : 09 Nov 2021 12:59
 Operator : CG/JU
 Sample : SSTD3.255
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.2006

Quant Time: Nov 09 13:29:07 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM110921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 09 11:53:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.470	152	2058	0.400	ng/ul	0.00
4) Naphthalene-d8	10.240	136	7622	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.126	164	4845	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.866	188	10146	0.400	ng/ul	0.00
17) Chrysene-d12	21.053	240	8738	0.400	ng/ul	0.00
23) Perylene-d12	23.164	264	6531	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.865	96	7449	2.777	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	2.899	88	8302	2.929	ng/ul#	88
14) Pentachlorophenol	16.537	266	9725	2.870	ng/ul	98

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

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