

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110624\
 Data File : BM048476.D
 Acq On : 06 Nov 2024 14:05
 Operator : RC/JU
 Sample : SSTD3.211
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.2039

Quant Time: Nov 06 14:38:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM110624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 06 14:05:20 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.682	152	3616	0.400	ng/ul	0.00
4) Naphthalene-d8	10.464	136	11315	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.322	164	6109	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.069	188	11601	0.400	ng/ul	-0.01
17) Chrysene-d12	21.254	240	9926	0.400	ng/ul	0.00
23) Perylene-d12	23.493	264	10689	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	13376	3.034	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	3.168	88	14561	2.838	ng/ul#	63
14) Pentachlorophenol	16.740	266	14149	3.355	ng/ul	98

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

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