

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042222\
 Data File : BM034770.D
 Acq On : 22 Apr 2022 16:14
 Operator : CG/JU
 Sample : SSTD3.234
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.2034

Quant Time: Apr 22 16:50:17 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM042222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 22 16:08:48 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.947	152	5968	0.400	ng/ul	0.00
4) Naphthalene-d8	10.748	136	14996	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.575	164	7923	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.313	188	14284	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.488	240	11445	0.400	ng/ul	0.00
23) Perylene-d12	23.885	264	10565	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.386	96	19426	2.325	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	3.420	88	19433	2.903	ng/ul#	79
14) Pentachlorophenol	16.963	266	14650	3.153	ng/ul	98

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

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