

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041222\
 Data File : BM034622.D
 Acq On : 12 Apr 2022 15:07
 Operator : CG/JU
 Sample : SSTD3.227
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.2027

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 04/13/2022
 Supervised By :Yogesh Patel 04/14/2022

Quant Time: Apr 12 15:37:35 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM041222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 12 14:58:22 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	1996	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.694	136	5718	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.529	164	2927	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.272	188	5275	0.400	ng/ul	-0.01
17) Chrysene-d12	21.449	240	4942	0.400	ng/ul	#-0.02
23) Perylene-d12	23.843	264	4686m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.256	96	9177	2.745	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	3.290	88	9293	2.787	ng/ul#	79
14) Pentachlorophenol	16.930	266	5846	2.614	ng/ul	98

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

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