

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092721\
 Data File : BM032203.D
 Acq On : 27 Sep 2021 13:02
 Operator : CG/JU
 Sample : SSTD3.234
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.2034

Quant Time: Sep 27 13:33:40 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM092721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 27 11:46:23 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.664	152	6557	0.400	ng/ul	0.00
4) Naphthalene-d8	10.456	136	20898	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.308	164	10850	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.033	188	21977	0.400	ng/ul	0.00
17) Chrysene-d12	21.196	240	20943	0.400	ng/ul	0.00
23) Perylene-d12	23.361	264	15893	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.989	96	21708	3.014	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	3.024	88	23577	3.351	ng/ul	89
14) Pentachlorophenol	16.700	266	20406	3.510	ng/ul	99

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

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