

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110921\
 Data File : BM032965.D
 Acq On : 10 Nov 2021 22:46
 Operator : CG/JU
 Sample : M4492-04
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG1S1

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 11/11/2021
 Supervised By :mohammad ahmed 11/11/2021

Quant Time: Nov 11 08:29:26 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 13:31:56 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.462	152	2077	0.400	ng/ul	0.00
4) Naphthalene-d8	10.235	136	7720	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.117	164	4712	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	16.862	188	9721	0.400	ng/ul	0.00
17) Chrysene-d12	21.052	240	7949	0.400	ng/ul	0.00
23) Perylene-d12	23.156	264	8299	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.857	96	9307m	3.509	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	11.836	152	3487	0.318	ng/ul	0.00
18) Fluoranthene-d10	18.892	212	10282	0.427	ng/ul	0.00

Target Compounds	Qvalue

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

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