

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062322\
 Data File : BM035648.D
 Acq On : 24 Jun 2022 05:32
 Operator : CG/JU
 Sample : N3359-13
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4S8

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 06/24/2022
 Supervised By :Yogesh Patel 06/28/2022

Quant Time: Jun 24 06:13:15 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM061822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 20 01:56:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.821	152	5854	0.400	ng/ul	-0.04
4) Naphthalene-d8	10.600	136	20466	0.400	ng/ul	-0.04
9) Acenaphthene-d10	14.446	164	11370m	0.400	ng/ul	-0.02
13) Phenanthrene-d10	17.195	188	20218m	0.400	ng/ul	-0.03
17) Chrysene-d12	21.389	240	15266	0.400	ng/ul	#-0.02
23) Perylene-d12	23.718	264	19287	0.400	ng/ul	#-0.05
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	31369	3.808	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.211	152	11452	0.377	ng/ul	-0.03
18) Fluoranthene-d10	19.229	212	19888	0.388	ng/ul	-0.02
Target Compounds				Qvalue		
2) 1,4-Dioxane	3.302	88	3107m	0.379	ng/ul	
19) Fluoranthene	19.261	202	1730	0.022	ng/ul#	1
21) Benzo(a)anthracene	21.377	228	1343	0.023	ng/ul#	17

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

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