

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062322\
 Data File : BM035622.D
 Acq On : 23 Jun 2022 13:02
 Operator : CG/JU
 Sample : N3358-10
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4S3

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 06/27/2022
 Supervised By :Yogesh Patel 06/28/2022

Quant Time: Jun 24 01:35:26 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	7.820	152	4767	0.400	ng/u1	-0.04
4) Naphthalene-d8	10.600	136	17479	0.400	ng/u1	-0.04
9) Acenaphthene-d10	14.446	164	11503m	0.400	ng/u1	-0.02
13) Phenanthrene-d10	17.195	188	25121m	0.400	ng/u1	-0.03
17) Chrysene-d12	21.401	240	18138	0.400	ng/u1	-0.01
23) Perylene-d12	23.724	264	11160m	0.400	ng/u1	-0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.260	96	28385	4.232	ng/u1	-0.01
6) 2-Methylnaphthalene-d10	12.211	152	10152	0.391	ng/u1	-0.03
18) Fluoranthene-d10	19.229	212	26302	0.432	ng/u1	-0.02
Target Compounds						
11) Acenaphthene	14.510	153	1284	0.027	ng/u1	Qvalue 93

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

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