

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062322\
 Data File : BM035639.D
 Acq On : 24 Jun 2022 00:07
 Operator : CG/JU
 Sample : N3358-18
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4T1

Manual Integrations
 APPROVED

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

Quant Time: Jun 24 01:36:35 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	7.808	152	6031	0.400	ng/ul	-0.05
4) Naphthalene-d8	10.622	136	33839882	0.400	ng/ul	#-0.02
9) Acenaphthene-d10	14.488	164	17769m	0.400	ng/ul	0.02
13) Phenanthrene-d10	17.200	188	21195m	0.400	ng/ul	-0.03
17) Chrysene-d12	21.382	240	16502m	0.400	ng/ul	-0.03
23) Perylene-d12	23.720	264	17634m	0.400	ng/ul	-0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.268	96	27347	3.223	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.261	152	207619	0.004	ng/ul	0.02
18) Fluoranthene-d10	19.225	212	25678m	0.463	ng/ul	-0.03
Target Compounds						Qvalue
10) Acenaphthylene	14.326	152	16479	0.165	ng/ul#	73
11) Acenaphthene	14.562	153	38677	0.522	ng/ul#	67
12) Fluorene	15.520	166	5946	0.071	ng/ul#	34
15) Phenanthrene	17.238	178	16290m	0.210	ng/ul	

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

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