

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062322\
 Data File : BM035626.D
 Acq On : 23 Jun 2022 15:27
 Operator : CG/JU
 Sample : N3358-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4R7

Manual Integrations
 APPROVED

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

Quant Time: Jun 24 01:35:41 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	4882	0.400	ng/u1	-0.04
4) Naphthalene-d8	10.600	136	20889	0.400	ng/u1	-0.04
9) Acenaphthene-d10	14.445	164	10106m	0.400	ng/u1	-0.02
13) Phenanthrene-d10	17.195	188	17076	0.400	ng/u1	#-0.03
17) Chrysene-d12	21.395	240	14165	0.400	ng/u1	-0.02
23) Perylene-d12	23.724	264	17157	0.400	ng/u1	#-0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	31310	4.558	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.206	152	9987	0.322	ng/u1	-0.04
18) Fluoranthene-d10	19.229	212	17476	0.367	ng/u1	-0.02
Target Compounds						
11) Acenaphthene	14.510	153	1214	0.029	ng/u1	Qvalue 90

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

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