

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062322\
 Data File : BM035621.D
 Acq On : 23 Jun 2022 12:25
 Operator : CG/JU
 Sample : N3358-06
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4S2

Manual Integrations
 APPROVED

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

Quant Time: Jun 24 01:35:22 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	7.816	152	5358	0.400	ng/ul	-0.04
4) Naphthalene-d8	10.600	136	19544	0.400	ng/ul	-0.04
9) Acenaphthene-d10	14.446	164	11779m	0.400	ng/ul	-0.02
13) Phenanthrene-d10	17.195	188	21913	0.400	ng/ul	-0.03
17) Chrysene-d12	21.398	240	14318	0.400	ng/ul	-0.01
23) Perylene-d12	23.730	264	17509	0.400	ng/ul	#-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.260	96	30540	4.051	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.217	152	9293	0.320	ng/ul	-0.03
18) Fluoranthene-d10	19.229	212	18807	0.391	ng/ul	-0.02
Target Compounds						
					Qvalue	
5) Naphthalene	10.650	128	2889	0.044	ng/ul#	89
11) Acenaphthene	14.510	153	1756	0.036	ng/ul	97
14) Pentachlorophenol	16.900	266	294	0.070	ng/ul	94
19) Fluoranthene	19.261	202	1913	0.026	ng/ul#	30
21) Benzo(a)anthracene	21.386	228	1109	0.020	ng/ul#	74
22) Chrysene	21.436	228	1399	0.022	ng/ul#	75

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

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