

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062322\
 Data File : BM035625.D
 Acq On : 23 Jun 2022 14:51
 Operator : CG/JU
 Sample : N3358-15
 Misc :
 ALS Vial : 8 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 01:35:37 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	5618	0.400	ng/ul	-0.04
4) Naphthalene-d8	10.600	136	20169	0.400	ng/ul	-0.04
9) Acenaphthene-d10	14.446	164	11018m	0.400	ng/ul	-0.02
13) Phenanthrene-d10	17.195	188	19602	0.400	ng/ul	#-0.03
17) Chrysene-d12	21.392	240	14880	0.400	ng/ul	#-0.02
23) Perylene-d12	23.724	264	18485	0.400	ng/ul	#-0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	29475	3.729	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.211	152	11570	0.386	ng/ul	-0.03
18) Fluoranthene-d10	19.229	212	19253	0.385	ng/ul	-0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.302	88	2825	0.359	ng/ul#	11
11) Acenaphthene	14.510	153	1404	0.031	ng/ul	95
15) Phenanthrene	17.237	178	1960	0.027	ng/ul#	55
19) Fluoranthene	19.266	202	1967	0.026	ng/ul#	1
21) Benzo(a)anthracene	21.380	228	1383	0.025	ng/ul#	31
22) Chrysene	21.430	228	1397	0.021	ng/ul#	33

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

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