

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062322\
 Data File : BM035632.D
 Acq On : 23 Jun 2022 19:16
 Operator : CG/JU
 Sample : N3358-04
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4S0

Manual Integrations
 APPROVED

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

Quant Time: Jun 24 01:36:06 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	7.842	152	6	0.400	ng/ul	#-0.02
4) Naphthalene-d8	10.556	136	6889m	0.400	ng/ul	-0.09
9) Acenaphthene-d10	14.450	164	1019m	0.400	ng/ul	-0.02
13) Phenanthrene-d10	17.199	188	1026	0.400	ng/ul	#-0.03
17) Chrysene-d12	21.409	240	834	0.400	ng/ul	# 0.00
23) Perylene-d12	23.721	264	4341m	0.400	ng/ul	-0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	31839	3771.410	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.211	152	11151	1.089	ng/ul	-0.03
18) Fluoranthene-d10	19.233	212	20315	7.251	ng/ul	-0.02
Target Compounds						
2) 1,4-Dioxane	3.302	88	277	32.958	ng/ul#	82
11) Acenaphthene	14.515	153	828m	0.195	ng/ul	
12) Fluorene	15.500	166	536m	0.112	ng/ul	
15) Phenanthrene	17.237	178	851	0.226	ng/ul#	53
16) Anthracene	17.326	178	695	0.211	ng/ul#	39
19) Fluoranthene	19.266	202	959	0.225	ng/ul#	39
21) Benzo(a)anthracene	21.401	228	586	0.186	ng/ul#	55
22) Chrysene	21.447	228	838	0.223	ng/ul#	55
25) Benzo(k)fluoranthene	23.075	252	468	0.024	ng/ul#	1

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

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