

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062022\
 Data File : BM035584.D
 Acq On : 21 Jun 2022 22:50
 Operator : CG/JU
 Sample : N3285-02RE
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4P6RE

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 06/22/2022
 Supervised By :Yogesh Patel 06/25/2022

Quant Time: Jun 22 01:33:32 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	7.812	152	5824	0.400	ng/ul	-0.05
4) Naphthalene-d8	10.605	136	20917	0.400	ng/ul	#-0.04
9) Acenaphthene-d10	14.451	164	16566m	0.400	ng/ul	-0.02
13) Phenanthrene-d10	17.196	188	27124m	0.400	ng/ul	-0.03
17) Chrysene-d12	21.388	240	18144	0.400	ng/ul	#-0.02
23) Perylene-d12	23.726	264	18493	0.400	ng/ul	#-0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	28414	3.467	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.200	152	10431	0.336	ng/ul	-0.04
18) Fluoranthene-d10	19.230	212	21663	0.355	ng/ul	-0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.298	88	8065m	0.989	ng/ul	
5) Naphthalene	10.655	128	3068	0.043	ng/ul#	67
7) 2-Methylnaphthalene	12.272	142	1203	0.030	ng/ul#	56
8) 1-Methylnaphthalene	12.492	142	1460	0.039	ng/ul#	56
15) Phenanthrene	17.238	178	2236	0.022	ng/ul#	1

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

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