

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062822\
 Data File : BM035730.D
 Acq On : 28 Jun 2022 18:47
 Operator : CG/JU
 Sample : N3401-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AG0

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 06/29/2022
 Supervised By :mohammad ahmed 07/01/2022

Quant Time: Jun 29 04:27:35 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 29 04:24:18 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.795	152	2441	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.578	136	7845	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.423	164	4351	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.170	188	8819m	0.400	ng/ul	-0.01
17) Chrysene-d12	21.376	240	8266	0.400	ng/ul	# 0.00
23) Perylene-d12	23.688	264	10504	0.400	ng/ul	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.213	96	14031	4.236	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.206	152	3580	0.291	ng/ul	0.00
18) Fluoranthene-d10	19.206	212	8338	0.299	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.627	128	534	0.022	ng/ul#	56
21) Benzo(a)anthracene	21.373	228	609	0.021	ng/ul#	25
22) Chrysene	21.423	228	977	0.027	ng/ul#	45
24) Benzo(b)fluoranthene	22.984	252	1354	0.030	ng/ul#	1
25) Benzo(k)fluoranthene	23.030	252	1331	0.029	ng/ul#	1

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

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