

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070821\
 Data File : BM030892.D
 Acq On : 08 Jul 2021 19:37
 Operator : CG/JU
 Sample : M2825-08
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGDR2

Manual Integrations
 APPROVED

mohammad
 7/9/2021 2:48:18 PM

Quant Time: Jul 09 02:52:53 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 09 02:52:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	1873	0.400	ng/ul	0.00
4) Naphthalene-d8	10.545	136	6829	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.394	164	3510	0.400	ng/ul	#-0.01
13) Phenanthrene-d10	17.131	188	5854m	0.400	ng/ul	-0.02
17) Chrysene-d12	21.311	240	4822	0.400	ng/ul	0.00
23) Perylene-d12	23.547	264	6508	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.262	96	9084	4.451	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.151	152	3518	0.328	ng/ul	0.00
18) Fluoranthene-d10	19.153	212	5408	0.391	ng/ul	-0.02
Target Compounds						
					Qvalue	
15) Phenanthrene	17.173	178	478	0.024	ng/ul#	66
19) Fluoranthene	19.188	202	545	0.026	ng/ul#	61
20) Pyrene	19.546	202	458	0.021	ng/ul#	36
21) Benzo(a)anthracene	21.299	228	436	0.023	ng/ul#	66
22) Chrysene	21.348	228	514	0.024	ng/ul#	75
24) Benzo(b)fluoranthene	22.873	252	739m	0.026	ng/ul	

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

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