

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080521\
 Data File : BM031577.D
 Acq On : 08 Aug 2021 08:04
 Operator : CG/JU
 Sample : M3163-15
 Misc :
 ALS Vial : 103 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGEZ4

Manual Integrations
 APPROVED

mohammad
 8/9/2021 12:14:37 PM

Quant Time: Aug 09 04:13:31 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 06 18:39:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.580	152	1002	0.400	ng/ul	0.00
4) Naphthalene-d8	10.338	136	3709	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.212	164	2117	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.961	188	3941	0.400	ng/ul	#-0.01
17) Chrysene-d12	21.159	240	3184	0.400	ng/ul	# 0.00
23) Perylene-d12	23.309	264	3432	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.186	96	2727	2.451	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.939	152	1230	0.197	ng/ul	0.00
18) Fluoranthene-d10	18.993	212	2423	0.228	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	17.006	178	286	0.023	ng/ul#	60
19) Fluoranthene	19.028	202	534	0.039	ng/ul#	62
20) Pyrene	19.386	202	519	0.038	ng/ul#	55
21) Benzo(a)anthracene	21.144	228	276	0.021	ng/ul#	61
22) Chrysene	21.193	228	305	0.023	ng/ul#	69
24) Benzo(b)fluoranthene	22.672	252	401m	0.026	ng/ul	
26) Benzo(a)pyrene	23.217	252	271	0.020	ng/ul#	1

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

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