

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090321\
 Data File : BM032090.D
 Acq On : 04 Sep 2021 23:24
 Operator : CG/JU
 Sample : M3517-06
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AB9

Manual Integrations
 APPROVED

mohammad
 9/7/2021 11:39:09 AM

Quant Time: Sep 06 04:18:07 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM090321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 03 16:21:27 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.448	152	1479	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.203	136	4796	0.400	ng/ul	#-0.02
9) Acenaphthene-d10	14.095	164	1847	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	16.851	188	3530	0.400	ng/ul	#-0.01
17) Chrysene-d12	21.064	240	2744	0.400	ng/ul	#-0.01
23) Perylene-d12	23.175	264	3975	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.061	96	7022	4.323	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	11.809	152	1609	0.217	ng/ul	-0.01
18) Fluoranthene-d10	18.888	212	3273	0.407	ng/ul	-0.02
Target Compounds						
					Qvalue	
15) Phenanthrene	16.893	178	291	0.027	ng/ul#	66
16) Anthracene	16.980	178	216	0.023	ng/ul#	40
19) Fluoranthene	18.922	202	298	0.025	ng/ul#	50
20) Pyrene	19.281	202	368m	0.030	ng/ul	
21) Benzo(a)anthracene	21.054	228	288	0.029	ng/ul#	66
22) Chrysene	21.100	228	334	0.029	ng/ul#	69
24) Benzo(b)fluoranthene	22.552	252	505	0.028	ng/ul#	10
26) Benzo(a)pyrene	23.081	252	361	0.023	ng/ul#	4
27) Indeno(1,2,3-cd)pyrene	25.252	276	493	0.023	ng/ul#	68
29) Benzo(g,h,i)perylene	25.879	276	494	0.026	ng/ul#	34

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

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