

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070821\
 Data File : BM030896.D
 Acq On : 08 Jul 2021 22:04
 Operator : CG/JU
 Sample : M2825-09
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGD3

Manual Integrations
 APPROVED

mohammad
 7/9/2021 2:48:34 PM

Quant Time: Jul 09 02:53:24 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	7.767	152	1992	0.400	ng/ul	0.00
4) Naphthalene-d8	10.577	136	7042	0.400	ng/ul #	0.02
9) Acenaphthene-d10	14.413	164	3249	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.155	188	5530	0.400	ng/ul	0.00
17) Chrysene-d12	21.316	240	4739	0.400	ng/ul #	0.00
23) Perylene-d12	23.564	264	4624	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	36	0.017	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.200	152	2065	0.187	ng/ul	0.04
18) Fluoranthene-d10	19.172	212	3242	0.239	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	17.200	178	474	0.025	ng/ul#	48
19) Fluoranthene	19.203	202	791	0.038	ng/ul#	70
20) Pyrene	19.565	202	703	0.033	ng/ul#	67
21) Benzo(a)anthracene	21.306	228	427	0.023	ng/ul#	54
22) Chrysene	21.355	228	455	0.022	ng/ul#	66
24) Benzo(b)fluoranthene	22.892	252	615m	0.030	ng/ul	
25) Benzo(k)fluoranthene	22.939	252	536m	0.025	ng/ul	

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

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