

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072621\
 Data File : BM031178.D
 Acq On : 27 Jul 2021 00:35
 Operator : CG/JU
 Sample : M3010-07
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGEE5

Manual Integrations
 APPROVED

mohammad
 7/27/2021 12:09:43 PM

Quant Time: Jul 27 06:23:15 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 22 01:52:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.632	152	1435	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.392	136	5930	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.258	164	4238	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.006	188	8903m	0.400	ng/ul	0.00
17) Chrysene-d12	21.188	240	8449	0.400	ng/ul	-0.01
23) Perylene-d12	23.360	264	8303	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.213	96	5824	3.864	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.989	152	3248	0.312	ng/ul	0.00
18) Fluoranthene-d10	19.031	212	9316	0.361	ng/ul	-0.01
Target Compounds						
					Qvalue	
15) Phenanthrene	17.044	178	651	0.023	ng/ul#	81
19) Fluoranthene	19.062	202	1714	0.050	ng/ul#	81
20) Pyrene	19.428	202	1356	0.039	ng/ul#	75
21) Benzo(a)anthracene	21.173	228	922	0.027	ng/ul#	80
22) Chrysene	21.226	228	749	0.021	ng/ul#	83
24) Benzo(b)fluoranthene	22.716	252	1182m	0.030	ng/ul	

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

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