

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072321\
 Data File : BM031112.D
 Acq On : 24 Jul 2021 05:47
 Operator : CG/JU
 Sample : M2970-03DL 5X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGE92DL

Manual Integrations
 APPROVED

mohammad
 7/26/2021 1:26:36 PM

Quant Time: Jul 24 06:23:05 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 : Sat Jul 24 04:20:53 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.639	152	1436	0.40	ng/ul	0.00
4) Naphthalene-d8	10.400	136	5512	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.265	164	3630	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.009	188	7880	0.40	ng/ul	0.00
17) Chrysene-d12	21.192	240	6950	0.40	ng/ul	0.00
23) Perylene-d12	23.365	264	6042	0.40	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	570	0.38	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.997	152	243	0.03	ng/ul	0.00
18) Fluoranthene-d10	19.035	212	633	0.03	ng/ul	0.00
Target Compounds						
						Qvalue
10) Acenaphthylene	13.985	152	784	0.05	ng/ul#	87
12) Fluorene	15.315	166	348	0.02	ng/ul#	92
15) Phenanthrene	17.051	178	7922	0.32	ng/ul	99
16) Anthracene	17.141	178	2732	0.12	ng/ul	97
19) Fluoranthene	19.069	202	14417	0.52	ng/ul	99
20) Pyrene	19.431	202	10142	0.36	ng/ul	99
21) Benzo(a)anthracene	21.175	228	6565	0.24	ng/ul	97
22) Chrysene	21.229	228	5514	0.18	ng/ul	96
24) Benzo(b)fluoranthene	22.718	252	5946m	0.21	ng/ul	
25) Benzo(k)fluoranthene	22.756	252	1923m	0.07	ng/ul	
26) Benzo(a)pyrene	23.268	252	2539	0.10	ng/ul#	74
27) Indeno(1,2,3-cd)pyrene	25.521	276	1511	0.05	ng/ul#	94
28) Dibenzo(a,h)anthracene	25.528	278	542	0.02	ng/ul#	22

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

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