

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080221\
 Data File : BM031389.D
 Acq On : 02 Aug 2021 23:29
 Operator : CG/JU
 Sample : M3123-03DL 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0066DL

Manual Integrations
 APPROVED

mohammad
 8/3/2021 11:50:58 AM

Quant Time: Aug 03 02:07:13 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 : Mon Aug 02 11:29:41 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.601	152	1670	0.40	ng/ul	0.00
4) Naphthalene-d8	10.361	136	6323	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.231	164	3673	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.978	188	6788	0.40	ng/ul	# 0.00
17) Chrysene-d12	21.171	240	5012	0.40	ng/ul	0.00
23) Perylene-d12	23.331	264	4800	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.196	96	499	0.27	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.953	152	217	0.02	ng/ul	0.00
18) Fluoranthene-d10	19.009	212	415	0.02	ng/ul	0.00
Target Compounds						
						Qvalue
5) Naphthalene	10.410	128	6622	0.35	ng/ul	100
7) 2-Methylnaphthalene	12.029	142	7321	0.57	ng/ul	100
8) 1-Methylnaphthalene	12.250	142	4836	0.38	ng/ul	99
12) Fluorene	15.285	166	399	0.03	ng/ul#	62
15) Phenanthrene	17.020	178	9186	0.43	ng/ul	99
19) Fluoranthene	19.043	202	1788	0.08	ng/ul#	84
20) Pyrene	19.405	202	1610	0.07	ng/ul#	85
21) Benzo(a)anthracene	21.152	228	663	0.03	ng/ul#	42
22) Chrysene	21.205	228	4025	0.19	ng/ul#	90
24) Benzo(b)fluoranthene	22.687	252	1470m	0.07	ng/ul	
26) Benzo(a)pyrene	23.237	252	403	0.02	ng/ul#	1

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

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