

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071723\
 Data File : BM040933.D
 Acq On : 17 Jul 2023 21:51
 Operator : MA/JU
 Sample : 03423-11
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A46T2

Quant Time: Jul 18 02:47:56 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 15 03:58:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.853	152	4739	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.653	136	14960	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.499	164	7090	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.253	188	12508	0.400	ng/ul	0.00
17) Chrysene-d12	21.448	240	8505	0.400	ng/ul	# 0.00
23) Perylene-d12	23.825	264	11149	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.280	96	27807	3.735	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.247	152	8237	0.393	ng/ul	-0.01
18) Fluoranthene-d10	19.287	212	12891	0.473	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.702	128	2571	0.062	ng/ul#	90
7) 2-Methylnaphthalene	12.319	142	949	0.039	ng/ul	98
8) 1-Methylnaphthalene	12.539	142	707	0.029	ng/ul	95
12) Fluorene	15.550	166	599	0.021	ng/ul#	97
15) Phenanthrene	17.295	178	866	0.022	ng/ul#	79
19) Fluoranthene	19.315	202	1126	0.031	ng/ul#	81
21) Benzo(a)anthracene	21.434	228	845	0.029	ng/ul#	87
22) Chrysene	21.486	228	957	0.029	ng/ul#	89
24) Benzo(b)fluoranthene	23.102	252	1230	0.029	ng/ul#	22
25) Benzo(k)fluoranthene	23.149	252	1031	0.024	ng/ul#	17
27) Indeno(1,2,3-cd)pyrene	26.294	276	1281	0.024	ng/ul#	92
28) Dibenzo(a,h)anthracene	26.317	278	886	0.021	ng/ul#	30
29) Benzo(g,h,i)perylene	27.052	276	1046	0.022	ng/ul#	64

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

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