

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM113023\
 Data File : BM043100.D
 Acq On : 01 Dec 2023 00:18
 Operator : MA/JU
 Sample : PB157409BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS409

Quant Time: Dec 01 01:35:33 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM112423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 01 00:42:42 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.847	152	919	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.629	136	2774	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.460	164	1545	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.226	188	3119	0.400	ng/ul	0.00
17) Chrysene-d12	21.409	240	2030	0.400	ng/ul	0.00
23) Perylene-d12	23.761	264	2695	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.160	96	837	0.594	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.240	152	1066	0.287	ng/ul	0.00
18) Fluoranthene-d10	19.248	212	2175	0.294	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.193	88	2275	1.502	ng/ul#	81
5) Naphthalene	10.678	128	2249	0.298	ng/ul	98
7) 2-Methylnaphthalene	12.317	142	1289	0.290	ng/ul	99
8) 1-Methylnaphthalene	12.509	142	1379	0.278	ng/ul	99
10) Acenaphthylene	14.192	152	2261	0.287	ng/ul	98
11) Acenaphthene	14.525	153	1752	0.293	ng/ul	99
12) Fluorene	15.529	166	1786	0.291	ng/ul	97
14) Pentachlorophenol	16.896	266	527	0.624	ng/ul	98
15) Phenanthrene	17.268	178	2622	0.295	ng/ul	97
16) Anthracene	17.378	178	2574	0.305	ng/ul	96
19) Fluoranthene	19.281	202	3397	0.295	ng/ul	99
20) Pyrene	19.648	202	3572	0.292	ng/ul	98
21) Benzo(a)anthracene	21.397	228	2421	0.368	ng/ul	98
22) Chrysene	21.447	228	3737	0.305	ng/ul	98
24) Benzo(b)fluoranthene	23.033	252	3055	0.336	ng/ul	98
25) Benzo(k)fluoranthene	23.089	252	4028	0.322	ng/ul	98
26) Benzo(a)pyrene	23.662	252	3192	0.322	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.218	276	4402	0.328	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.242	278	3407	0.328	ng/ul	94
29) Benzo(g,h,i)perylene	26.976	276	4094	0.313	ng/ul	100

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

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