

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071723\
 Data File : BM040922.D
 Acq On : 17 Jul 2023 14:40
 Operator : MA/JU
 Sample : PB153870BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS870

Quant Time: Jul 18 02:46:27 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	7.857	152	3634	0.400	ng/ul	0.00
4) Naphthalene-d8	10.658	136	11955	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.504	164	4908	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.261	188	7915	0.400	ng/ul	0.00
17) Chrysene-d12	21.454	240	5244	0.400	ng/ul	0.00
23) Perylene-d12	23.833	264	5535	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.280	96	3671	0.643	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.253	152	4707	0.281	ng/ul	0.00
18) Fluoranthene-d10	19.292	212	5368	0.319	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.313	88	9500	1.635	ng/ul	89
5) Naphthalene	10.708	128	9989	0.303	ng/ul	99
7) 2-Methylnaphthalene	12.330	142	5327	0.275	ng/ul	97
8) 1-Methylnaphthalene	12.544	142	5433	0.281	ng/ul	99
10) Acenaphthylene	14.226	152	6059	0.276	ng/ul	98
11) Acenaphthene	14.564	153	5342	0.289	ng/ul	97
12) Fluorene	15.559	166	5209	0.264	ng/ul	97
14) Pentachlorophenol	16.919	266	1015	0.666	ng/ul	91
15) Phenanthrene	17.299	178	7135	0.290	ng/ul	98
16) Anthracene	17.396	178	5639	0.273	ng/ul	96
19) Fluoranthene	19.320	202	6898	0.310	ng/ul	100
20) Pyrene	19.687	202	7383	0.309	ng/ul	98
21) Benzo(a)anthracene	21.436	228	5131	0.289	ng/ul	98
22) Chrysene	21.489	228	6075	0.300	ng/ul	99
24) Benzo(b)fluoranthene	23.108	252	6568	0.307	ng/ul	93
25) Benzo(k)fluoranthene	23.155	252	6413	0.304	ng/ul#	89
26) Benzo(a)pyrene	23.728	252	5786	0.300	ng/ul#	85
27) Indeno(1,2,3-cd)pyrene	26.293	276	8619	0.321	ng/ul	95
28) Dibenzo(a,h)anthracene	26.307	278	6864	0.326	ng/ul	95
29) Benzo(g,h,i)perylene	27.045	276	7709	0.327	ng/ul	98

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

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