

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122823\
 Data File : BM043651.D
 Acq On : 28 Dec 2023 22:39
 Operator : MA/JU
 Sample : PB158092BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS092

Quant Time: Dec 28 23:20:29 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM121123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 26 22:58:31 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.967	152	4447	0.400	ng/ul	0.00
4) Naphthalene-d8	10.774	136	12865	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.601	164	5146	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.350	188	10525	0.400	ng/ul #	0.00
17) Chrysene-d12	21.527	240	6926	0.400	ng/ul #	0.00
23) Perylene-d12	23.939	264	7769	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.377	96	3488	0.621	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.363	152	4663	0.256	ng/ul	0.00
18) Fluoranthene-d10	19.371	212	6729	0.272	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.410	88	11010	1.954	ng/ul#	83
5) Naphthalene	10.823	128	10468	0.271	ng/ul	98
7) 2-Methylnaphthalene	12.440	142	5945	0.238	ng/ul	97
8) 1-Methylnaphthalene	12.654	142	5698	0.228	ng/ul	100
10) Acenaphthylene	14.324	152	8121	0.287	ng/ul	98
11) Acenaphthene	14.666	153	5910	0.287	ng/ul	99
12) Fluorene	15.656	166	6595	0.282	ng/ul	97
14) Pentachlorophenol	16.995	266	1506	0.646	ng/ul	97
15) Phenanthrene	17.392	178	10081	0.287	ng/ul	98
16) Anthracene	17.485	178	9499	0.298	ng/ul	99
19) Fluoranthene	19.404	202	10413	0.267	ng/ul#	92
20) Pyrene	19.761	202	10732	0.277	ng/ul#	90
21) Benzo(a)anthracene	21.513	228	8695	0.286	ng/ul	98
22) Chrysene	21.565	228	9843	0.307	ng/ul	99
24) Benzo(b)fluoranthene	23.202	252	9781	0.277	ng/ul	88
25) Benzo(k)fluoranthene	23.255	252	10962	0.300	ng/ul#	87
26) Benzo(a)pyrene	23.833	252	8933	0.307	ng/ul#	82
27) Indeno(1,2,3-cd)pyrene	26.458	276	12336	0.372	ng/ul#	44
28) Dibenzo(a,h)anthracene	26.482	278	10033	0.382	ng/ul#	92
29) Benzo(g,h,i)perylene	27.223	276	10610	0.375	ng/ul	93

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

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