

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090623\
 Data File : BM041728.D
 Acq On : 08 Sep 2023 12:20
 Operator : MA/JU
 Sample : PB155206BS
 Misc :
 ALS Vial : 69 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS206

Quant Time: Sep 08 22:54:30 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM090523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 08 06:08:11 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.996	152	8483	0.400	ng/ul	0.00
4) Naphthalene-d8	10.812	136	20847	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.638	164	8672	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.388	188	16427	0.400	ng/ul	0.00
17) Chrysene-d12	21.556	240	5383	0.400	ng/ul	0.00
23) Perylene-d12	23.985	264	5098	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.406	96	7800	0.602	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.390	152	9600	0.358	ng/ul	0.00
18) Fluoranthene-d10	19.408	212	10555	0.404	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.439	88	18493	1.373	ng/ul	91
5) Naphthalene	10.861	128	21933	0.320	ng/ul	99
7) 2-Methylnaphthalene	12.467	142	12951	0.333	ng/ul	99
8) 1-Methylnaphthalene	12.687	142	13686	0.350	ng/ul	100
10) Acenaphthylene	14.365	152	17874	0.334	ng/ul	99
11) Acenaphthene	14.698	153	13907	0.334	ng/ul	98
12) Fluorene	15.684	166	15555	0.333	ng/ul	100
14) Pentachlorophenol	17.012	266	3924	0.597	ng/ul	99
15) Phenanthrene	17.430	178	23199	0.334	ng/ul	99
16) Anthracene	17.523	178	19973	0.336	ng/ul	99
19) Fluoranthene	19.436	202	23371	0.368	ng/ul	99
20) Pyrene	19.803	202	23260	0.369	ng/ul	99
21) Benzo(a)anthracene	21.539	228	14595	0.332	ng/ul	100
22) Chrysene	21.594	228	15380	0.325	ng/ul	100
24) Benzo(b)fluoranthene	23.234	252	14685	0.322	ng/ul	100
25) Benzo(k)fluoranthene	23.284	252	13710	0.307	ng/ul	99
26) Benzo(a)pyrene	23.874	252	11785	0.318	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	26.501	276	15597	0.307	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.522	278	10979	0.308	ng/ul	98
29) Benzo(g,h,i)perylene	27.279	276	13429	0.298	ng/ul	98

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

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