

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090623\
 Data File : BM041714.D
 Acq On : 08 Sep 2023 03:05
 Operator : MA/JU
 Sample : PB155166BS
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS166

Quant Time: Sep 08 03:52:28 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 : Wed Sep 06 02:20:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.000	152	6995	0.400	ng/ul	0.00
4) Naphthalene-d8	10.817	136	17318	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.638	164	7597	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.388	188	14869	0.400	ng/ul	-0.01
17) Chrysene-d12	21.556	240	5017	0.400	ng/ul	-0.01
23) Perylene-d12	23.982	264	4702	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.402	96	6612	0.619	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.396	152	8639	0.388	ng/ul	-0.01
18) Fluoranthene-d10	19.408	212	10384	0.427	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.439	88	15907	1.432	ng/ul	90
5) Naphthalene	10.867	128	19249	0.338	ng/ul	99
7) 2-Methylnaphthalene	12.467	142	11532	0.357	ng/ul	100
8) 1-Methylnaphthalene	12.687	142	12235	0.377	ng/ul	100
10) Acenaphthylene	14.365	152	16373	0.349	ng/ul	99
11) Acenaphthene	14.703	153	12778	0.350	ng/ul	98
12) Fluorene	15.688	166	14279	0.349	ng/ul	99
14) Pentachlorophenol	17.012	266	3288	0.553	ng/ul	99
15) Phenanthrene	17.430	178	22214	0.354	ng/ul	100
16) Anthracene	17.523	178	19210	0.357	ng/ul	99
19) Fluoranthene	19.440	202	22382	0.378	ng/ul	98
20) Pyrene	19.803	202	22534	0.383	ng/ul	98
21) Benzo(a)anthracene	21.539	228	14256	0.348	ng/ul	100
22) Chrysene	21.594	228	14685	0.333	ng/ul	100
24) Benzo(b)fluoranthene	23.234	252	13689	0.325	ng/ul	97
25) Benzo(k)fluoranthene	23.284	252	12841	0.312	ng/ul	96
26) Benzo(a)pyrene	23.874	252	10860	0.318	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.501	276	14109	0.301	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.525	278	10043	0.305	ng/ul	99
29) Benzo(g,h,i)perylene	27.283	276	12151	0.292	ng/ul	97

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

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