

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM021924\
 Data File : BM044210.D
 Acq On : 20 Feb 2024 17:21
 Operator : MA/JU
 Sample : PB159061BS
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS061

Quant Time: Feb 20 23:05:38 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM021524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 15 13:40:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.896	152	4983	0.400	ng/ul	0.00
4) Naphthalene-d8	10.693	136	14291	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.543	164	6332	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.296	188	12863	0.400	ng/ul	0.00
17) Chrysene-d12	21.500	240	10174	0.400	ng/ul	0.00
23) Perylene-d12	23.888	264	11586	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	5415	0.767	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.288	152	7451	0.372	ng/ul	-0.01
18) Fluoranthene-d10	19.331	212	13180	0.434	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.386	88	14362	2.036	ng/ul#	77
5) Naphthalene	10.748	128	15895	0.374	ng/ul	99
7) 2-Methylnaphthalene	12.365	142	9262	0.354	ng/ul	99
8) 1-Methylnaphthalene	12.580	142	9406	0.355	ng/ul	99
10) Acenaphthylene	14.260	152	13591	0.391	ng/ul	97
11) Acenaphthene	14.603	153	9668	0.383	ng/ul	100
12) Fluorene	15.597	166	10074	0.354	ng/ul	100
14) Pentachlorophenol	16.946	266	3315	0.720	ng/ul	99
15) Phenanthrene	17.339	178	15538	0.381	ng/ul	99
16) Anthracene	17.436	178	16334	0.405	ng/ul	99
19) Fluoranthene	19.359	202	19252	0.415	ng/ul	98
20) Pyrene	19.721	202	20319	0.417	ng/ul	97
21) Benzo(a)anthracene	21.485	228	16124	0.346	ng/ul	99
22) Chrysene	21.538	228	20413	0.406	ng/ul	99
24) Benzo(b)fluoranthene	23.163	252	18158	0.380	ng/ul	96
25) Benzo(k)fluoranthene	23.212	252	20888	0.402	ng/ul	96
26) Benzo(a)pyrene	23.785	252	16878	0.380	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.373	276	21001	0.371	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.417	278	16579	0.372	ng/ul	96
29) Benzo(g,h,i)perylene	27.124	276	18751	0.375	ng/ul	98

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

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