

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022124\
 Data File : BM044257.D
 Acq On : 23 Feb 2024 02:03
 Operator : MA/JU
 Sample : PB159101BS
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS101

Quant Time: Feb 23 02:37:06 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	4440	0.400	ng/ul	0.00
4) Naphthalene-d8	10.699	136	12931	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.543	164	5522	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.301	188	11604	0.400	ng/ul	0.00
17) Chrysene-d12	21.500	240	10137	0.400	ng/ul	0.00
23) Perylene-d12	23.897	264	11603	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	5802	0.923	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.288	152	7984	0.441	ng/ul	-0.01
18) Fluoranthene-d10	19.331	212	15248	0.504	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.382	88	18261	2.906	ng/ul#	73
5) Naphthalene	10.749	128	18030	0.469	ng/ul	99
7) 2-Methylnaphthalene	12.365	142	10363	0.437	ng/ul	98
8) 1-Methylnaphthalene	12.580	142	10304	0.430	ng/ul	99
10) Acenaphthylene	14.265	152	15152	0.499	ng/ul	99
11) Acenaphthene	14.607	153	10715	0.487	ng/ul	99
12) Fluorene	15.602	166	11256	0.454	ng/ul	99
14) Pentachlorophenol	16.946	266	4035	0.971	ng/ul	98
15) Phenanthrene	17.343	178	17904	0.487	ng/ul	99
16) Anthracene	17.436	178	18653	0.513	ng/ul	99
19) Fluoranthene	19.359	202	23176	0.501	ng/ul	96
20) Pyrene	19.721	202	24585	0.506	ng/ul	96
21) Benzo(a)anthracene	21.485	228	21198	0.457	ng/ul	100
22) Chrysene	21.538	228	25763	0.515	ng/ul	99
24) Benzo(b)fluoranthene	23.166	252	23802	0.497	ng/ul	98
25) Benzo(k)fluoranthene	23.216	252	27451	0.528	ng/ul	99
26) Benzo(a)pyrene	23.789	252	22227	0.499	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.390	276	26758	0.472	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.427	278	20853	0.467	ng/ul	96
29) Benzo(g,h,i)perylene	27.134	276	24342	0.486	ng/ul	98

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

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