

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062424\
 Data File : BM046307.D
 Acq On : 26 Jun 2024 00:37
 Operator : MA/JU
 Sample : PB161514BS
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS514

Quant Time: Jun 26 01:41:21 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM061824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 25 08:04:59 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.430	152	1988	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.222	136	6370	0.400	ng/ul	#-0.03
9) Acenaphthene-d10	14.067	164	3649	0.400	ng/ul	-0.02
13) Phenanthrene-d10	16.829	188	7229	0.400	ng/ul	-0.03
17) Chrysene-d12	21.012	240	6096	0.400	ng/ul	#-0.01
23) Perylene-d12	23.095	264	7930	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.025	96	1596	0.585	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.839	152	2823	0.330	ng/ul	-0.03
18) Fluoranthene-d10	18.849	212	5936	0.325	ng/ul	-0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.050	88	3995	1.306	ng/ul#	88
5) Naphthalene	10.260	128	5722	0.322	ng/ul#	91
7) 2-Methylnaphthalene	11.905	142	3202	0.320	ng/ul	99
8) 1-Methylnaphthalene	12.103	142	3297	0.301	ng/ul	100
10) Acenaphthylene	13.790	152	6354	0.367	ng/ul	97
11) Acenaphthene	14.128	153	4247	0.340	ng/ul	98
12) Fluorene	15.141	166	4539	0.339	ng/ul	99
14) Pentachlorophenol	16.508	266	1460	0.674	ng/ul	99
15) Phenanthrene	16.871	178	6603	0.325	ng/ul	96
16) Anthracene	16.968	178	5359	0.354	ng/ul	99
19) Fluoranthene	18.882	202	9412	0.357	ng/ul	98
20) Pyrene	19.239	202	9800	0.343	ng/ul	99
21) Benzo(a)anthracene	20.997	228	8233	0.350	ng/ul	93
22) Chrysene	21.050	228	10101	0.336	ng/ul	94
24) Benzo(b)fluoranthene	22.484	252	9386	0.329	ng/ul	85
25) Benzo(k)fluoranthene	22.525	252	11668	0.344	ng/ul#	88
26) Benzo(a)pyrene	23.010	252	9833	0.371	ng/ul#	83
27) Indeno(1,2,3-cd)pyrene	25.145	276	12296	0.284	ng/ul#	93
28) Dibenzo(a,h)anthracene	25.155	278	9976	0.309	ng/ul#	87
29) Benzo(g,h,i)perylene	25.766	276	10709	0.288	ng/ul	91

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

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