

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062424\
 Data File : BM046246.D
 Acq On : 24 Jun 2024 10:48
 Operator : MA/JU
 Sample : PB161516BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS516

Quant Time: Jun 24 11:21:20 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 : Mon Jun 24 02:25:34 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.464	152	1369	0.400	ng/ul	0.00
4) Naphthalene-d8	10.266	136	4424	0.400	ng/ul	# 0.01
9) Acenaphthene-d10	14.090	164	2654	0.400	ng/ul	-0.02
13) Phenanthrene-d10	16.854	188	4878	0.400	ng/ul	0.00
17) Chrysene-d12	21.032	240	3904	0.400	ng/ul	-0.01
23) Perylene-d12	23.118	264	5281	0.400	ng/ul	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.029	96	1483	0.790	ng/ul	-0.02
6) 2-Methylnaphthalene-d10	11.882	152	2421	0.408	ng/ul	0.02
18) Fluoranthene-d10	18.872	212	4939	0.422	ng/ul	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.059	88	4072	1.933	ng/ul#	80
5) Naphthalene	10.315	128	4874	0.395	ng/ul#	92
7) 2-Methylnaphthalene	11.959	142	2703	0.389	ng/ul	97
8) 1-Methylnaphthalene	12.141	142	3138	0.412	ng/ul	96
10) Acenaphthylene	13.822	152	5024	0.399	ng/ul	95
11) Acenaphthene	14.155	153	3779	0.416	ng/ul	99
12) Fluorene	15.178	166	3827	0.392	ng/ul	99
14) Pentachlorophenol	16.533	266	1087	0.744	ng/ul	98
15) Phenanthrene	16.896	178	5409	0.395	ng/ul	99
16) Anthracene	17.002	178	3495	0.342	ng/ul	97
19) Fluoranthene	18.905	202	7470	0.442	ng/ul	98
20) Pyrene	19.262	202	7879	0.431	ng/ul	99
21) Benzo(a)anthracene	21.017	228	5582	0.370	ng/ul	99
22) Chrysene	21.067	228	8454	0.439	ng/ul	99
24) Benzo(b)fluoranthene	22.504	252	7209	0.379	ng/ul	93
25) Benzo(k)fluoranthene	22.545	252	9956	0.440	ng/ul	97
26) Benzo(a)pyrene	23.037	252	7992	0.453	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.179	276	10849	0.376	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.189	278	8329	0.388	ng/ul	95
29) Benzo(g,h,i)perylene	25.796	276	9556	0.386	ng/ul	96

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

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