

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071124\
 Data File : BM046528.D
 Acq On : 11 Jul 2024 22:11
 Operator : MA/JU
 Sample : PB161864BS
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS864

Quant Time: Jul 12 05:10:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM071024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 05:09:19 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.514	152	1749	0.400	ng/ul	0.00
4) Naphthalene-d8	10.277	136	4318	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.155	164	2826	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.909	188	6249	0.400	ng/ul	0.00
17) Chrysene-d12	21.111	240	5016	0.400	ng/ul	# 0.00
23) Perylene-d12	23.253	264	5997	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.122	96	682	0.475	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.877	152	1871	0.269	ng/ul	0.00
18) Fluoranthene-d10	18.946	212	4238	0.278	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.151	88	1872	1.221	ng/ul#	73
5) Naphthalene	10.320	128	2737	0.249	ng/ul	99
7) 2-Methylnaphthalene	11.948	142	1969	0.255	ng/ul	100
8) 1-Methylnaphthalene	12.174	142	1984	0.252	ng/ul	97
10) Acenaphthylene	13.868	152	3132	0.243	ng/ul	97
11) Acenaphthene	14.220	153	2233	0.259	ng/ul	96
12) Fluorene	15.214	166	2702	0.253	ng/ul	100
14) Pentachlorophenol	16.563	266	1184	0.374	ng/ul	97
15) Phenanthrene	16.947	178	4377	0.245	ng/ul	97
16) Anthracene	17.040	178	4171	0.240	ng/ul	98
19) Fluoranthene	18.974	202	5375	0.261	ng/ul	98
20) Pyrene	19.337	202	5706	0.256	ng/ul	99
21) Benzo(a)anthracene	21.096	228	5415	0.244	ng/ul	99
22) Chrysene	21.146	228	5319	0.248	ng/ul	99
24) Benzo(b)fluoranthene	22.618	252	6146	0.257	ng/ul	97
25) Benzo(k)fluoranthene	22.659	252	6081	0.259	ng/ul	96
26) Benzo(a)pyrene	23.162	252	5434	0.276	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	25.380	276	7777	0.259	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.396	278	6111	0.257	ng/ul	96
29) Benzo(g,h,i)perylene	26.020	276	6114	0.255	ng/ul	98

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

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