

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062424\
 Data File : BM046262.D
 Acq On : 24 Jun 2024 20:33
 Operator : MA/JU
 Sample : PB161673BS
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS673

Quant Time: Jun 25 02:27:13 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	7.459	152	1269	0.400	ng/ul	0.00
4) Naphthalene-d8	10.266	136	4275	0.400	ng/ul	# 0.01
9) Acenaphthene-d10	14.090	164	2687	0.400	ng/ul	-0.02
13) Phenanthrene-d10	16.854	188	4346	0.400	ng/ul	0.00
17) Chrysene-d12	21.035	240	3643	0.400	ng/ul	0.00
23) Perylene-d12	23.115	264	5143	0.400	ng/ul	#-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.029	96	1530	0.879	ng/ul	-0.02
6) 2-Methylnaphthalene-d10	11.888	152	2350	0.409	ng/ul	0.02
18) Fluoranthene-d10	18.877	212	4563	0.418	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.054	88	3553	1.819	ng/ul#	78
5) Naphthalene	10.310	128	4441	0.373	ng/ul#	90
7) 2-Methylnaphthalene	11.954	142	2349	0.350	ng/ul	96
8) 1-Methylnaphthalene	12.146	142	2894	0.393	ng/ul	99
10) Acenaphthylene	13.817	152	4555	0.357	ng/ul	94
11) Acenaphthene	14.155	153	3327	0.362	ng/ul	97
12) Fluorene	15.173	166	3343	0.339	ng/ul	99
14) Pentachlorophenol	16.550	266	800	0.614	ng/ul	90
15) Phenanthrene	16.896	178	4688	0.384	ng/ul	96
16) Anthracene	16.998	178	3177	0.349	ng/ul	94
19) Fluoranthene	18.909	202	6328	0.401	ng/ul	96
20) Pyrene	19.262	202	6775	0.397	ng/ul	97
21) Benzo(a)anthracene	21.023	228	4548	0.323	ng/ul	99
22) Chrysene	21.073	228	6835	0.381	ng/ul	98
24) Benzo(b)fluoranthene	22.504	252	6448	0.348	ng/ul	93
25) Benzo(k)fluoranthene	22.545	252	7908	0.359	ng/ul	96
26) Benzo(a)pyrene	23.034	252	6732	0.392	ng/ul	93
27) Indeno(1,2,3-cd)pyrene	25.185	276	9129	0.325	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.199	278	6982	0.334	ng/ul	92
29) Benzo(g,h,i)perylene	25.799	276	8056	0.334	ng/ul	97

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

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