

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM113023\
 Data File : BM043163.D
 Acq On : 02 Dec 2023 17:02
 Operator : MA/JU
 Sample : PB157370BS
 Misc :
 ALS Vial : 82 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS370

Quant Time: Dec 02 23:55:10 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM112423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 02 23:54:55 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.859	152	843	0.400	ng/ul	0.00
4) Naphthalene-d8	10.628	136	2951	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.456	164	1677	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.230	188	3649	0.400	ng/ul	0.00
17) Chrysene-d12	21.411	240	2200	0.400	ng/ul	0.00
23) Perylene-d12	23.764	264	2362	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.160	96	886	0.685	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.245	152	1465	0.371	ng/ul	0.00
18) Fluoranthene-d10	19.253	212	3274	0.408	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.193	88	2553	1.838	ng/ul#	85
5) Naphthalene	10.683	128	2758	0.344	ng/ul	97
7) 2-Methylnaphthalene	12.311	142	1576	0.333	ng/ul	97
8) 1-Methylnaphthalene	12.504	142	1662	0.315	ng/ul	93
10) Acenaphthylene	14.192	152	2911	0.340	ng/ul	99
11) Acenaphthene	14.516	153	2285	0.352	ng/ul	94
12) Fluorene	15.529	166	2370	0.356	ng/ul	98
14) Pentachlorophenol	16.884	266	624	0.632	ng/ul	96
15) Phenanthrene	17.268	178	3503	0.336	ng/ul	100
16) Anthracene	17.382	178	3327	0.337	ng/ul	97
19) Fluoranthene	19.281	202	4612	0.370	ng/ul	98
20) Pyrene	19.648	202	4906	0.370	ng/ul	98
21) Benzo(a)anthracene	21.400	228	2795	0.392	ng/ul	98
22) Chrysene	21.447	228	4708	0.354	ng/ul	98
24) Benzo(b)fluoranthene	23.036	252	3105	0.390	ng/ul	98
25) Benzo(k)fluoranthene	23.089	252	4357	0.397	ng/ul	99
26) Benzo(a)pyrene	23.665	252	3320	0.382	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.231	276	4244	0.361	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.252	278	3276	0.360	ng/ul	99
29) Benzo(g,h,i)perylene	26.979	276	4014	0.350	ng/ul	96

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

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