

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM113023\
 Data File : BM043164.D
 Acq On : 02 Dec 2023 17:38
 Operator : MA/JU
 Sample : PB157359BS
 Misc :
 ALS Vial : 83 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS359

Quant Time: Dec 02 23:55:18 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	7.847	152	860	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.633	136	2683	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.454	164	1487	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.228	188	3087	0.400	ng/ul	0.00
17) Chrysene-d12	21.412	240	1863	0.400	ng/ul	0.00
23) Perylene-d12	23.765	264	2315	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.160	96	857	0.649	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.233	152	1327	0.370	ng/ul	-0.01
18) Fluoranthene-d10	19.252	212	2674	0.394	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.193	88	2476	1.747	ng/ul#	84
5) Naphthalene	10.677	128	2572	0.353	ng/ul	96
7) 2-Methylnaphthalene	12.310	142	1490	0.346	ng/ul	99
8) 1-Methylnaphthalene	12.508	142	1581	0.330	ng/ul	96
10) Acenaphthylene	14.191	152	2596	0.342	ng/ul	97
11) Acenaphthene	14.515	153	2013	0.350	ng/ul	94
12) Fluorene	15.528	166	2080	0.352	ng/ul	97
14) Pentachlorophenol	16.882	266	491	0.588	ng/ul	97
15) Phenanthrene	17.271	178	3004	0.341	ng/ul	100
16) Anthracene	17.380	178	2860	0.343	ng/ul	98
19) Fluoranthene	19.284	202	3895	0.369	ng/ul	97
20) Pyrene	19.647	202	4118	0.367	ng/ul	99
21) Benzo(a)anthracene	21.400	228	2280	0.377	ng/ul	98
22) Chrysene	21.447	228	4130	0.367	ng/ul	98
24) Benzo(b)fluoranthene	23.037	252	3163	0.405	ng/ul	97
25) Benzo(k)fluoranthene	23.086	252	4205	0.391	ng/ul	97
26) Benzo(a)pyrene	23.671	252	3122	0.366	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.229	276	4420	0.383	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.259	278	3415	0.383	ng/ul	95
29) Benzo(g,h,i)perylene	26.986	276	4239	0.378	ng/ul	97

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

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