

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM113023\
 Data File : BM043151.D
 Acq On : 02 Dec 2023 09:09
 Operator : MA/JU
 Sample : PB157357BS
 Misc :
 ALS Vial : 69 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS357

Quant Time: Dec 02 22:41:03 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 01:12:12 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.876	152	551	0.400	ng/ul	0.05
4) Naphthalene-d8	10.634	136	1971	0.400	ng/ul	0.02
9) Acenaphthene-d10	14.456	164	1210	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.230	188	2524	0.400	ng/ul	0.01
17) Chrysene-d12	21.414	240	1507	0.400	ng/ul	0.01
23) Perylene-d12	23.767	264	1650	0.400	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.160	96	680	0.804	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.240	152	1075	0.408	ng/ul	0.02
18) Fluoranthene-d10	19.253	212	2310	0.421	ng/ul	0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.193	88	1919	2.114	ng/ul#	85
5) Naphthalene	10.683	128	2005	0.374	ng/ul	98
7) 2-Methylnaphthalene	12.317	142	1181	0.374	ng/ul	99
8) 1-Methylnaphthalene	12.504	142	1276	0.362	ng/ul	99
10) Acenaphthylene	14.192	152	2109	0.342	ng/ul	99
11) Acenaphthene	14.521	153	1631	0.349	ng/ul	98
12) Fluorene	15.529	166	1717	0.357	ng/ul	96
14) Pentachlorophenol	16.896	266	388	0.568	ng/ul	95
15) Phenanthrene	17.272	178	2556	0.355	ng/ul	98
16) Anthracene	17.386	178	2601	0.381	ng/ul	98
19) Fluoranthene	19.286	202	3390	0.397	ng/ul	98
20) Pyrene	19.648	202	3547	0.390	ng/ul	99
21) Benzo(a)anthracene	21.403	228	1894	0.387	ng/ul	98
22) Chrysene	21.449	228	3386	0.372	ng/ul	99
24) Benzo(b)fluoranthene	23.039	252	2354	0.423	ng/ul	99
25) Benzo(k)fluoranthene	23.089	252	3224	0.420	ng/ul	95
26) Benzo(a)pyrene	23.671	252	2381	0.392	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	26.235	276	3252	0.395	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.262	278	2436	0.383	ng/ul	99
29) Benzo(g,h,i)perylene	26.986	276	3229	0.404	ng/ul	98

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

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