

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082923\
 Data File : BM041577.D
 Acq On : 30 Aug 2023 13:19
 Operator : MA/JU
 Sample : PB155102BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS102

Quant Time: Aug 31 02:17:08 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM082923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 30 01:50:06 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.114	152	5439	0.400	ng/ul	0.00
4) Naphthalene-d8	10.944	136	12261	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.749	164	4806	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.502	188	8822	0.400	ng/ul	0.00
17) Chrysene-d12	21.691	240	3038	0.400	ng/ul	0.00
23) Perylene-d12	24.254	264	3190	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.452	96	5690	0.709	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.517	152	5682	0.384	ng/ul	0.00
18) Fluoranthene-d10	19.524	212	5634	0.376	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.490	88	14016	1.637	ng/ul#	85
5) Naphthalene	10.994	128	13364	0.341	ng/ul	99
7) 2-Methylnaphthalene	12.594	142	7479	0.336	ng/ul	99
8) 1-Methylnaphthalene	12.808	142	7844	0.349	ng/ul	99
10) Acenaphthylene	14.481	152	11361	0.339	ng/ul	99
11) Acenaphthene	14.814	153	7831	0.339	ng/ul	97
12) Fluorene	15.800	166	8500	0.338	ng/ul	97
14) Pentachlorophenol	17.126	266	1967	0.580	ng/ul	98
15) Phenanthrene	17.544	178	12846	0.333	ng/ul	99
16) Anthracene	17.637	178	11258	0.336	ng/ul	98
19) Fluoranthene	19.552	202	12980	0.326	ng/ul	97
20) Pyrene	19.919	202	12822	0.320	ng/ul	97
21) Benzo(a)anthracene	21.670	228	8549	0.315	ng/ul	99
22) Chrysene	21.726	228	9501	0.326	ng/ul	99
24) Benzo(b)fluoranthene	23.456	252	9355	0.316	ng/ul	97
25) Benzo(k)fluoranthene	23.509	252	9284	0.330	ng/ul	97
26) Benzo(a)pyrene	24.137	252	8161	0.309	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.961	276	10841	0.307	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.998	278	7348	0.310	ng/ul	96
29) Benzo(g,h,i)perylene	27.803	276	9725	0.329	ng/ul	99

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

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