

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082923\
 Data File : BM041589.D
 Acq On : 30 Aug 2023 21:52
 Operator : MA/JU
 Sample : PB155073BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS073

Quant Time: Aug 31 02:18:52 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	6620	0.400	ng/ul	0.00
4) Naphthalene-d8	10.944	136	14855	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.749	164	5748	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.502	188	10363	0.400	ng/ul	0.00
17) Chrysene-d12	21.682	240	3394	0.400	ng/ul	0.00
23) Perylene-d12	24.243	264	3838	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.452	96	6496	0.665	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.517	152	6511	0.363	ng/ul	0.00
18) Fluoranthene-d10	19.520	212	6231	0.372	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.486	88	15885	1.524	ng/ul#	83
5) Naphthalene	10.993	128	15413	0.325	ng/ul	100
7) 2-Methylnaphthalene	12.594	142	8639	0.321	ng/ul	100
8) 1-Methylnaphthalene	12.808	142	9058	0.333	ng/ul	99
10) Acenaphthylene	14.476	152	12654	0.316	ng/ul	99
11) Acenaphthene	14.814	153	8924	0.323	ng/ul	99
12) Fluorene	15.795	166	9536	0.317	ng/ul	98
14) Pentachlorophenol	17.122	266	2468	0.620	ng/ul	98
15) Phenanthrene	17.544	178	14258	0.315	ng/ul	99
16) Anthracene	17.637	178	12350	0.314	ng/ul	99
19) Fluoranthene	19.552	202	14137	0.318	ng/ul	100
20) Pyrene	19.915	202	13805	0.308	ng/ul	95
21) Benzo(a)anthracene	21.664	228	9351	0.309	ng/ul	100
22) Chrysene	21.723	228	10332	0.318	ng/ul	100
24) Benzo(b)fluoranthene	23.447	252	11100	0.311	ng/ul	98
25) Benzo(k)fluoranthene	23.500	252	10895	0.322	ng/ul	99
26) Benzo(a)pyrene	24.129	252	9829	0.309	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.954	276	13599	0.321	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.981	278	9239	0.324	ng/ul	99
29) Benzo(g,h,i)perylene	27.793	276	11675	0.328	ng/ul	100

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

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