

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030923\
 Data File : BM038975.D
 Acq On : 10 Mar 2023 11:19
 Operator : CG/JU
 Sample : PB151313BS
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS313

Quant Time: Mar 10 23:15:02 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM030323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 09 13:36:03 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.991	152	7401	0.400	ng/ul	0.00
4) Naphthalene-d8	10.806	136	24489	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.628	164	13773	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.370	188	29331	0.400	ng/ul	0.00
17) Chrysene-d12	21.550	240	24149	0.400	ng/ul	0.00
23) Perylene-d12	23.994	264	18323	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	3294	0.393	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.389	152	6373	0.215	ng/ul	0.00
18) Fluoranthene-d10	19.394	212	13092	0.228	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.409	88	8826	1.041	ng/ul	94
5) Naphthalene	10.855	128	12105	0.207	ng/ul	99
7) 2-Methylnaphthalene	12.461	142	7375	0.208	ng/ul	100
8) 1-Methylnaphthalene	12.681	142	7727	0.207	ng/ul	99
10) Acenaphthylene	14.346	152	9362	0.187	ng/ul	98
11) Acenaphthene	14.688	153	8129	0.197	ng/ul	99
12) Fluorene	15.674	166	9131	0.196	ng/ul	99
14) Pentachlorophenol	17.016	266	2094	0.297	ng/ul	99
15) Phenanthrene	17.413	178	15477	0.200	ng/ul	100
16) Anthracene	17.505	178	12699	0.199	ng/ul	99
19) Fluoranthene	19.426	202	17354	0.212	ng/ul	99
20) Pyrene	19.789	202	18092	0.215	ng/ul	98
21) Benzo(a)anthracene	21.532	228	15452	0.226	ng/ul	100
22) Chrysene	21.588	228	17104	0.208	ng/ul	99
24) Benzo(b)fluoranthene	23.248	252	15786	0.220	ng/ul	90
25) Benzo(k)fluoranthene	23.298	252	15142	0.211	ng/ul#	91
26) Benzo(a)pyrene	23.885	252	11954	0.209	ng/ul#	83
27) Indeno(1,2,3-cd)pyrene	26.534	276	14193	0.197	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.558	278	11679	0.212	ng/ul	97
29) Benzo(g,h,i)perylene	27.319	276	12842	0.191	ng/ul	98

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

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