

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031124\
 Data File : BM044573.D
 Acq On : 11 Mar 2024 11:43
 Operator : MA/JU
 Sample : SSTD0.411
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4017

Quant Time: Mar 11 12:24:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM031124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 11 12:23:17 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.724	152	5865	0.400	ng/ul	0.00
4) Naphthalene-d8	10.507	136	19080	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.367	164	10924	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.119	188	24518	0.400	ng/ul	0.00
17) Chrysene-d12	21.316	240	15552	0.400	ng/ul	0.00
23) Perylene-d12	23.575	264	14990	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.285	96	3091	0.387	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.101	152	10842	0.409	ng/ul	0.00
18) Fluoranthene-d10	19.155	212	22168	0.463	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.315	88	3055	0.378	ng/ul	100
5) Naphthalene	10.556	128	20902	0.379	ng/ul	100
7) 2-Methylnaphthalene	12.178	142	13357	0.390	ng/ul	100
8) 1-Methylnaphthalene	12.393	142	13438	0.388	ng/ul	100
10) Acenaphthylene	14.085	152	21466	0.369	ng/ul	100
11) Acenaphthene	14.427	153	15627	0.371	ng/ul	100
12) Fluorene	15.422	166	18038	0.379	ng/ul	100
14) Pentachlorophenol	16.777	266	1924	0.241	ng/ul	100
15) Phenanthrene	17.161	178	28640	0.377	ng/ul	100
16) Anthracene	17.259	178	26789	0.361	ng/ul	100
19) Fluoranthene	19.187	202	30882	0.431	ng/ul	100
20) Pyrene	19.550	202	31378	0.420	ng/ul	100
21) Benzo(a)anthracene	21.301	228	22660	0.334	ng/ul	100
22) Chrysene	21.351	228	26433	0.359	ng/ul	100
24) Benzo(b)fluoranthene	22.891	252	22368	0.374	ng/ul	100
25) Benzo(k)fluoranthene	22.941	252	24225	0.373	ng/ul	100
26) Benzo(a)pyrene	23.479	252	19722	0.358	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.863	276	22737	0.328	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.890	278	17435	0.320	ng/ul	100
29) Benzo(g,h,i)perylene	26.554	276	20542	0.334	ng/ul	100

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

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