

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070323\
 Data File : BM040582.D
 Acq On : 03 Jul 2023 19:22
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4148

Quant Time: Jul 04 01:29:36 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 03 13:40:38 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.923	152	5263	0.400	ng/ul	0.00
4) Naphthalene-d8	10.733	136	20018	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.566	164	9433	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.318	188	16566	0.400	ng/ul	0.00
17) Chrysene-d12	21.503	240	9921	0.400	ng/ul	0.00
23) Perylene-d12	23.912	264	10325	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.321	96	2792	0.338	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.322	152	10988	0.392	ng/ul	0.00
18) Fluoranthene-d10	19.346	212	13784	0.433	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.354	88	2875	0.342	ng/ul#	72
5) Naphthalene	10.782	128	21095	0.382	ng/ul	99
7) 2-Methylnaphthalene	12.393	142	12396	0.382	ng/ul	98
8) 1-Methylnaphthalene	12.613	142	12082	0.373	ng/ul	99
10) Acenaphthylene	14.289	152	16725	0.397	ng/ul	98
11) Acenaphthene	14.631	153	13051	0.367	ng/ul	99
12) Fluorene	15.621	166	13628	0.359	ng/ul	100
14) Pentachlorophenol	16.976	266	869	0.272	ng/ul	94
15) Phenanthrene	17.360	178	18522	0.360	ng/ul	99
16) Anthracene	17.453	178	16694	0.386	ng/ul	98
19) Fluoranthene	19.378	202	17533	0.416	ng/ul	98
20) Pyrene	19.741	202	17990	0.398	ng/ul	100
21) Benzo(a)anthracene	21.486	228	13609	0.406	ng/ul	98
22) Chrysene	21.538	228	13967	0.365	ng/ul	98
24) Benzo(b)fluoranthene	23.178	252	13781	0.345	ng/ul	88
25) Benzo(k)fluoranthene	23.225	252	14152	0.359	ng/ul#	87
26) Benzo(a)pyrene	23.804	252	12436	0.346	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	26.417	276	17430	0.348	ng/ul	94
28) Dibenzo(a,h)anthracene	26.431	278	13769	0.351	ng/ul	93
29) Benzo(g,h,i)perylene	27.185	276	14751	0.336	ng/ul	96

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

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