

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071423\
 Data File : BM040882.D
 Acq On : 14 Jul 2023 16:56
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4178

Quant Time: Jul 15 03:58:46 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 : Sat Jul 15 03:58:02 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.866	152	3642	0.400	ng/ul	0.00
4) Naphthalene-d8	10.664	136	13734	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.509	164	6822	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.261	188	12917	0.400	ng/ul	0.00
17) Chrysene-d12	21.451	240	8387	0.400	ng/ul	# 0.00
23) Perylene-d12	23.830	264	6762	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.280	96	2214	0.387	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.258	152	7857	0.408	ng/ul	0.00
18) Fluoranthene-d10	19.292	212	12008	0.447	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.313	88	2355	0.404	ng/ul#	89
5) Naphthalene	10.713	128	15019	0.396	ng/ul	98
7) 2-Methylnaphthalene	12.330	142	8926	0.401	ng/ul	99
8) 1-Methylnaphthalene	12.550	142	9087	0.409	ng/ul	98
10) Acenaphthylene	14.231	152	11379	0.373	ng/ul	99
11) Acenaphthene	14.569	153	9834	0.382	ng/ul	97
12) Fluorene	15.559	166	10296	0.375	ng/ul	98
14) Pentachlorophenol	16.919	266	1039	0.418	ng/ul	93
15) Phenanthrene	17.303	178	15095	0.376	ng/ul	100
16) Anthracene	17.396	178	12420	0.368	ng/ul	100
19) Fluoranthene	19.324	202	15139	0.425	ng/ul	96
20) Pyrene	19.687	202	16357	0.428	ng/ul	98
21) Benzo(a)anthracene	21.434	228	11320	0.399	ng/ul	99
22) Chrysene	21.486	228	12090	0.374	ng/ul	99
24) Benzo(b)fluoranthene	23.102	252	10354	0.396	ng/ul	98
25) Benzo(k)fluoranthene	23.149	252	10288	0.399	ng/ul	99
26) Benzo(a)pyrene	23.722	252	8777	0.372	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.290	276	10978	0.334	ng/ul	95
28) Dibenzo(a,h)anthracene	26.307	278	8848	0.344	ng/ul	98
29) Benzo(g,h,i)perylene	27.045	276	9457	0.329	ng/ul	98

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

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