

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112020\
 Data File : BM027903.D
 Acq On : 22 Nov 2020 05:37
 Operator : JU/CG
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.438

Quant Time: Nov 23 05:40:47 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 23 05:39:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	1229	0.40	ng/ul	0.00
2) Naphthalene-d8	11.17	136	4819	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	2679	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	4704	0.40	ng/ul	0.00
16) Chrysene-d12	21.76	240	2454	0.40	ng/ul	0.00
20) Perylene-d12	24.18	264	2294	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	3068	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	4102	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.22	128	5773	0.400	ng/ul	98
5) 2-Methylnaphthalene	12.80	142	3982	0.405	ng/ul	99
7) Acenaphthylene	14.67	152	5434	0.416	ng/ul	99
8) Acenaphthene	15.01	153	3927	0.401	ng/ul	100
9) Fluorene	15.98	166	4362	0.384	ng/ul	99
11) Pentachlorophenol	17.31	266	551	0.344	ng/ul	100
12) Phenanthrene	17.70	178	6040	0.406	ng/ul	100
13) Anthracene	17.79	178	5816	0.406	ng/ul	99
15) Fluoranthene	19.68	202	5590	0.339	ng/ul	97
17) Pyrene	20.04	202	5434	0.436	ng/ul	95
18) Benzo(a)anthracene	21.74	228	3989	0.403	ng/ul	96
19) Chrysene	21.80	228	3901	0.404	ng/ul	97
21) Benzo(b)fluoranthene	23.44	252	3716	0.385	ng/ul	94
22) Benzo(k)fluoranthene	23.49	252	3625	0.384	ng/ul#	93
23) Benzo(a)pyrene	24.07	252	3299	0.391	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	26.67	276	3975	0.436	ng/ul#	98
25) Dibenzo(a,h)anthracene	26.69	278	3279	0.431	ng/ul#	95
26) Benzo(g,h,i)perylene	27.45	276	3409	0.445	ng/ul	98

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

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