

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040620\
 Data File : BM025949.D
 Acq On : 06 Apr 2020 16:52
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.441

Quant Time: Apr 06 18:21:18 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 02 10:49:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	2157	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	9979	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	6324	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	16021	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	16132	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	16229	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	7035	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	20423	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	10568	0.360	ng/ul	99
5) 2-Methylnaphthalene	12.00	142	7960	0.388	ng/ul	99
7) Acenaphthylene	13.92	152	12280	0.454	ng/ul	100
8) Acenaphthene	14.26	153	8977	0.400	ng/ul	99
9) Fluorene	15.25	166	10986	0.395	ng/ul	100
11) Pentachlorophenol	16.61	266	1485	0.378	ng/ul	98
12) Phenanthrene	16.99	178	18655	0.370	ng/ul	100
13) Anthracene	17.08	178	17635	0.395	ng/ul	100
15) Fluoranthene	19.01	202	23251	0.373	ng/ul	99
17) Pyrene	19.37	202	24014	0.416	ng/ul	97
18) Benzo(a)anthracene	21.12	228	23653	0.410	ng/ul	99
19) Chrysene	21.18	228	24377	0.387	ng/ul	99
21) Benzo(b)fluoranthene	22.65	252	26969	0.452	ng/ul	98
22) Benzo(k)fluoranthene	22.69	252	25919	0.405	ng/ul	98
23) Benzo(a)pyrene	23.19	252	24030	0.450	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.40	276	24289	0.340	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.42	278	20068	0.336	ng/ul	97
26) Benzo(g,h,i)perylene	26.05	276	19675	0.325	ng/ul	98

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

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