

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111020\
 Data File : BM027781.D
 Acq On : 10 Nov 2020 21:55
 Operator : JU/CG
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.484

Quant Time: Nov 11 01:42:38 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM111020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 11 01:20:37 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	984	0.40	ng/ul	0.00
2) Naphthalene-d8	11.22	136	3557	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.99	164	2073	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	4364	0.40	ng/ul	0.00
16) Chrysene-d12	21.81	240	3256	0.40	ng/ul	0.00
20) Perylene-d12	24.24	264	2608	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.78	152	2304	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	4500	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.27	128	4225	0.372	ng/ul	98
5) 2-Methylnaphthalene	12.85	142	2931	0.371	ng/ul	99
7) Acenaphthylene	14.72	152	3973	0.389	ng/ul	98
8) Acenaphthene	15.06	153	2990	0.383	ng/ul	99
9) Fluorene	16.03	166	3521	0.375	ng/ul	99
11) Pentachlorophenol	17.35	266	572	0.363	ng/ul	98
12) Phenanthrene	17.75	178	5391	0.379	ng/ul	99
13) Anthracene	17.84	178	5109	0.369	ng/ul	99
15) Fluoranthene	19.72	202	5969	0.367	ng/ul	98
17) Pyrene	20.08	202	5895	0.408	ng/ul	97
18) Benzo(a)anthracene	21.79	228	5078	0.390	ng/ul	99
19) Chrysene	21.84	228	5146	0.388	ng/ul	98
21) Benzo(b)fluoranthene	23.50	252	4743	0.397	ng/ul	98
22) Benzo(k)fluoranthene	23.55	252	4603	0.393	ng/ul#	96
23) Benzo(a)pyrene	24.14	252	3917	0.373	ng/ul#	96
24) Indeno(1,2,3-cd)pyrene	26.76	276	4424	0.370	ng/ul#	95
25) Dibenzo(a,h)anthracene	26.77	278	3648	0.365	ng/ul#	96
26) Benzo(g,h,i)perylene	27.54	276	3675	0.363	ng/ul	98

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

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