

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111020\
 Data File : BM027769.D
 Acq On : 10 Nov 2020 13:59
 Operator : JU/CG
 Sample : SSTD0.881
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.881

Quant Time: Nov 10 14:36:32 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 : Tue Nov 10 14:33:58 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.39	152	508	0.40	ng/ul	0.00
2) Naphthalene-d8	11.22	136	1775	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.00	164	1117	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	2431	0.40	ng/ul	0.00
16) Chrysene-d12	21.81	240	2060	0.40	ng/ul	0.00
20) Perylene-d12	24.25	264	1716	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	2624	0.80	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	5771	0.80	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.27	128	4363	0.769	ng/ul	97
5) 2-Methylnaphthalene	12.86	142	3071	0.785	ng/ul	100
7) Acenaphthylene	14.72	152	4349	0.783	ng/ul	98
8) Acenaphthene	15.06	153	3310	0.786	ng/ul	99
9) Fluorene	16.03	166	4001	0.804	ng/ul	100
11) Pentachlorophenol	17.35	266	674	0.682	ng/ul	99
12) Phenanthrene	17.75	178	6194	0.786	ng/ul	99
13) Anthracene	17.84	178	6069	0.796	ng/ul	99
15) Fluoranthene	19.73	202	7134	0.794	ng/ul	98
17) Pyrene	20.08	202	7057	0.785	ng/ul	99
18) Benzo(a)anthracene	21.79	228	6375	0.782	ng/ul	99
19) Chrysene	21.85	228	6502	0.780	ng/ul	99
21) Benzo(b)fluoranthene	23.50	252	6237	0.826	ng/ul	97
22) Benzo(k)fluoranthene	23.55	252	6138	0.817	ng/ul	97
23) Benzo(a)pyrene	24.14	252	5428	0.816	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.77	276	6314	0.801	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.79	278	5302	0.808	ng/ul	99
26) Benzo(g,h,i)perylene	27.55	276	5303	0.780	ng/ul	98

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

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