

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112020\
 Data File : BM027852.D
 Acq On : 20 Nov 2020 14:43
 Operator : JU/CG
 Sample : SSTD0.876
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.876

Quant Time: Nov 20 17:15:13 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 : Fri Nov 20 17:13:49 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	859	0.40	ng/ul	0.00
2) Naphthalene-d8	11.17	136	3312	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	2010	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.67	188	4210	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	2918	0.40	ng/ul	0.00
20) Perylene-d12	24.19	264	2230	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	4730	0.85	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	9470	0.84	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.22	128	8136	0.820	ng/ul	99
5) 2-Methylnaphthalene	12.81	142	5587	0.827	ng/ul	99
7) Acenaphthylene	14.67	152	8151	0.831	ng/ul	97
8) Acenaphthene	15.01	153	6117	0.832	ng/ul	99
9) Fluorene	15.98	166	7078	0.830	ng/ul	99
11) Pentachlorophenol	17.31	266	1187	0.828	ng/ul	98
12) Phenanthrene	17.71	178	10873	0.818	ng/ul	99
13) Anthracene	17.80	178	10510	0.819	ng/ul	100
15) Fluoranthene	19.68	202	12108	0.821	ng/ul	99
17) Pyrene	20.04	202	12039	0.812	ng/ul	98
18) Benzo(a)anthracene	21.75	228	9698	0.824	ng/ul	100
19) Chrysene	21.81	228	9479	0.827	ng/ul	99
21) Benzo(b)fluoranthene	23.45	252	7827	0.835	ng/ul	95
22) Benzo(k)fluoranthene	23.50	252	7616	0.830	ng/ul	97
23) Benzo(a)pyrene	24.08	252	6624	0.808	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	26.69	276	7256	0.818	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.70	278	6043	0.818	ng/ul	97
26) Benzo(g,h,i)perylene	27.46	276	6125	0.823	ng/ul	98

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

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