

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112020\
 Data File : BM027857.D
 Acq On : 20 Nov 2020 17:59
 Operator : JU/CG
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.483

Quant Time: Nov 20 18:31:22 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 17:20:22 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	788	0.40	ng/ul	0.00
2) Naphthalene-d8	11.17	136	3104	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1913	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.67	188	3994	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	2924	0.40	ng/ul	0.00
20) Perylene-d12	24.18	264	2273	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	2170	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	4439	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.22	128	3760	0.404	ng/ul	100
5) 2-Methylnaphthalene	12.80	142	2586	0.409	ng/ul	98
7) Acenaphthylene	14.67	152	3703	0.397	ng/ul	99
8) Acenaphthene	15.01	153	2824	0.404	ng/ul	97
9) Fluorene	15.98	166	3252	0.401	ng/ul	100
11) Pentachlorophenol	17.31	266	552	0.406	ng/ul	99
12) Phenanthrene	17.70	178	5073	0.402	ng/ul	100
13) Anthracene	17.79	178	4840	0.398	ng/ul	99
15) Fluoranthene	19.68	202	5688	0.407	ng/ul	99
17) Pyrene	20.04	202	5693	0.383	ng/ul	98
18) Benzo(a)anthracene	21.75	228	4700	0.398	ng/ul	99
19) Chrysene	21.80	228	4646	0.404	ng/ul	99
21) Benzo(b)fluoranthene	23.45	252	3942	0.413	ng/ul	98
22) Benzo(k)fluoranthene	23.50	252	3826	0.409	ng/ul	97
23) Benzo(a)pyrene	24.08	252	3335	0.399	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	26.68	276	3749	0.415	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.70	278	3136	0.416	ng/ul	98
26) Benzo(g,h,i)perylene	27.46	276	3164	0.417	ng/ul	100

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

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