

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112020\
 Data File : BM027851.D
 Acq On : 20 Nov 2020 14:07
 Operator : JU/CG
 Sample : SSTD0.481
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.481

Quant Time: Nov 20 17:15:10 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	8.34	152	767	0.40	ng/ul	0.00
2) Naphthalene-d8	11.17	136	2985	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1838	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.67	188	3827	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	2669	0.40	ng/ul	0.00
20) Perylene-d12	24.19	264	2020	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	3066	0.61	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	6279	0.61	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.22	128	5269	0.589	ng/ul	100
5) 2-Methylnaphthalene	12.81	142	3636	0.597	ng/ul	100
7) Acenaphthylene	14.67	152	5283	0.589	ng/ul	100
8) Acenaphthene	15.01	153	3986	0.593	ng/ul	100
9) Fluorene	15.98	166	4613	0.592	ng/ul	100
11) Pentachlorophenol	17.31	266	763	0.585	ng/ul	100
12) Phenanthrene	17.71	178	7223	0.597	ng/ul	100
13) Anthracene	17.80	178	6932	0.594	ng/ul	100
15) Fluoranthene	19.68	202	7993	0.596	ng/ul	100
17) Pyrene	20.04	202	8057	0.594	ng/ul	100
18) Benzo(a)anthracene	21.75	228	6407	0.595	ng/ul	100
19) Chrysene	21.81	228	6263	0.597	ng/ul	100
21) Benzo(b)fluoranthene	23.45	252	5040	0.594	ng/ul	100
22) Benzo(k)fluoranthene	23.50	252	5025	0.604	ng/ul	100
23) Benzo(a)pyrene	24.08	252	4363	0.588	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.69	276	4723	0.588	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.70	278	3931	0.587	ng/ul	100
26) Benzo(g,h,i)perylene	27.46	276	3949	0.586	ng/ul	100

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

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