

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112020\
 Data File : BM027848.D
 Acq On : 20 Nov 2020 11:42
 Operator : JU/CG
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC0.4

Quant Time: Nov 20 17:22:12 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	1218	0.40	ng/ul	0.00
2) Naphthalene-d8	11.17	136	4716	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	2671	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.67	188	4877	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	2834	0.40	ng/ul	0.00
20) Perylene-d12	24.19	264	2577	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	3235	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	4777	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.22	128	5872	0.415	ng/ul	98
5) 2-Methylnaphthalene	12.81	142	3885	0.404	ng/ul	99
7) Acenaphthylene	14.68	152	5396	0.414	ng/ul	98
8) Acenaphthene	15.01	153	4034	0.413	ng/ul	100
9) Fluorene	15.98	166	4452	0.393	ng/ul	100
11) Pentachlorophenol	17.31	266	615	0.370	ng/ul	98
12) Phenanthrene	17.71	178	6381	0.414	ng/ul	99
13) Anthracene	17.80	178	6069	0.408	ng/ul	99
15) Fluoranthene	19.68	202	6315	0.370	ng/ul	98
17) Pyrene	20.04	202	6325	0.439	ng/ul	97
18) Benzo(a)anthracene	21.75	228	4721	0.413	ng/ul	98
19) Chrysene	21.81	228	4635	0.416	ng/ul	98
21) Benzo(b)fluoranthene	23.45	252	4164	0.384	ng/ul	98
22) Benzo(k)fluoranthene	23.50	252	4212	0.397	ng/ul	98
23) Benzo(a)pyrene	24.08	252	3836	0.405	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	26.69	276	4575	0.446	ng/ul#	98
25) Dibenzo(a,h)anthracene	26.71	278	3774	0.442	ng/ul	96
26) Benzo(g,h,i)perylene	27.46	276	3873	0.450	ng/ul	100

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

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