

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071519\
 Data File : BN006853.D
 Acq On : 15 Jul 2019 17:39
 Operator : HP/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.428

Quant Time: Jul 15 18:17:55 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN062719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 06:39:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	2181	0.40	ng/ul	0.01
2) Naphthalene-d8	10.37	136	8338	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	4345	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	8969	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	7261	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	7995	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	4759	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	9921	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.42	128	9255	0.387	ng/ul	98
5) 2-Methylnaphthalene	12.06	142	5824	0.360	ng/ul	100
7) Acenaphthylene	12.97	152	8030	0.358	ng/ul	99
8) Acenaphthene	14.31	153	6603	0.382	ng/ul	98
9) Fluorene	15.31	166	6855	0.362	ng/ul	99
11) Pentachlorophenol	16.72	266	423	0.317	ng/ul	92
12) Phenanthrene	17.06	178	11131	0.386	ng/ul	99
13) Anthracene	17.16	178	8832	0.354	ng/ul	98
15) Fluoranthene	19.09	202	12810	0.359	ng/ul	98
17) Pyrene	19.46	202	13527	0.408	ng/ul	97
18) Benzo(a)anthracene	21.23	228	10095	0.338	ng/ul	99
19) Chrysene	21.28	228	13688	0.411	ng/ul	98
21) Benzo(b)fluoranthene	22.81	252	11724	0.358	ng/ul	97
22) Benzo(k)fluoranthene	22.86	252	13314	0.388	ng/ul	98
23) Benzo(a)pyrene	23.39	252	11747	0.366	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.71	276	13904	0.428	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.71	278	10799	0.418	ng/ul	97
26) Benzo(g,h,i)perylene	26.39	276	12403	0.457	ng/ul	97

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

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