

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM080919\
 Data File : BM022019.D
 Acq On : 11 Aug 2019 05:58
 Operator : HP/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.451

Quant Time: Aug 11 06:49:28 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM080319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 09 01:29:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	985	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.40	136	4269	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.27	164	2300	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	5127	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	4937	0.40	ng/ul	-0.01
20) Perylene-d12	23.44	264	5907	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	2651	0.38	ng/ul	-0.01
14) Fluoranthene-d10	19.07	212	6164	0.36	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.45	128	4312	0.358	ng/ul	98
5) 2-Methylnaphthalene	12.09	142	2952	0.367	ng/ul	99
7) Acenaphthylene	14.00	152	3854	0.359	ng/ul#	83
8) Acenaphthene	14.34	153	3230	0.368	ng/ul	96
9) Fluorene	15.34	166	3491	0.358	ng/ul	98
11) Pentachlorophenol	16.71	266	395	0.418	ng/ul	87
12) Phenanthrene	17.08	178	5420	0.363	ng/ul	99
13) Anthracene	17.17	178	5090	0.400	ng/ul	98
15) Fluoranthene	19.10	202	7055	0.353	ng/ul	98
17) Pyrene	19.46	202	7437	0.367	ng/ul	99
18) Benzo(a)anthracene	21.21	228	6866	0.389	ng/ul	98
19) Chrysene	21.27	228	7698	0.339	ng/ul	97
21) Benzo(b)fluoranthene	22.77	252	7015	0.352	ng/ul	97
22) Benzo(k)fluoranthene	22.82	252	7553	0.335	ng/ul	98
23) Benzo(a)pyrene	23.35	252	7223	0.357	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.65	276	8832	0.365	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.66	278	7066	0.361	ng/ul	97
26) Benzo(g,h,i)perylene	26.33	276	7527	0.346	ng/ul	97

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

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