

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080619\
 Data File : BM021889.D
 Acq On : 07 Aug 2019 10:29
 Operator : HP/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.438

Quant Time: Aug 07 11:06:03 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 : Wed Aug 07 08:28:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	777	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	3538	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.28	164	2066	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	4671	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	4424	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	5386	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	2366	0.41	ng/ul	-0.01
14) Fluoranthene-d10	19.08	212	5947	0.38	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.47	128	3967	0.397	ng/ul	97
5) 2-Methylnaphthalene	12.09	142	2731	0.410	ng/ul	99
7) Acenaphthylene	14.01	152	3933	0.408	ng/ul	98
8) Acenaphthene	14.35	153	3093	0.392	ng/ul	96
9) Fluorene	15.35	166	3360	0.384	ng/ul	99
11) Pentachlorophenol	16.71	266	383	0.444	ng/ul	98
12) Phenanthrene	17.09	178	5254	0.387	ng/ul	99
13) Anthracene	17.18	178	5109	0.441	ng/ul	99
15) Fluoranthene	19.11	202	6960	0.382	ng/ul	99
17) Pyrene	19.47	202	7105	0.391	ng/ul	96
18) Benzo(a)anthracene	21.23	228	6725	0.425	ng/ul	92
19) Chrysene	21.28	228	7290	0.358	ng/ul#	91
21) Benzo(b)fluoranthene	22.79	252	7080	0.390	ng/ul	93
22) Benzo(k)fluoranthene	22.83	252	7206	0.351	ng/ul#	92
23) Benzo(a)pyrene	23.36	252	7042	0.381	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.67	276	8590	0.389	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.68	278	6918	0.388	ng/ul#	95
26) Benzo(g,h,i)perylene	26.36	276	7396	0.373	ng/ul#	92

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

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