

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091019\
 Data File : BM022648.D
 Acq On : 11 Sep 2019 09:35
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.461

Quant Time: Sep 11 10:11:50 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM091019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 11 07:37:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	15633	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	62980	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	32934	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	74449	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	64748	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	60389	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	35077	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	69838	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.69	128	59522	0.332	ng/ul	100
5) 2-Methylnaphthalene	12.30	142	40794	0.332	ng/ul	100
7) Acenaphthylene	14.20	152	49045	0.322	ng/ul	100
8) Acenaphthene	14.54	153	41369	0.324	ng/ul	99
9) Fluorene	15.52	166	46850	0.320	ng/ul	99
11) Pentachlorophenol	16.86	266	4402	0.291	ng/ul	99
12) Phenanthrene	17.25	178	76611	0.323	ng/ul	100
13) Anthracene	17.34	178	65014	0.319	ng/ul	100
15) Fluoranthene	19.27	202	85765	0.335	ng/ul	99
17) Pyrene	19.63	202	86240	0.294	ng/ul	99
18) Benzo(a)anthracene	21.37	228	78752	0.318	ng/ul	100
19) Chrysene	21.43	228	83754	0.321	ng/ul	100
21) Benzo(b)fluoranthene	22.98	252	84139	0.321	ng/ul	98
22) Benzo(k)fluoranthene	23.03	252	86721	0.324	ng/ul	99
23) Benzo(a)pyrene	23.57	252	74660	0.325	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.98	276	83277	0.297	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.00	278	67761	0.299	ng/ul	98
26) Benzo(g,h,i)perylene	26.69	276	63349	0.284	ng/ul	100

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

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