

Data Path : Z:\HPCHEM1\BNA_E\Data\BE060917\
 Data File : BE093085.D
 Acq On : 9 Jun 2017 11:16
 Operator : SJ/MA
 Sample : SSTDC00.4
 Misc : 0.07 PPM
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.460

Quant Time: Jun 09 15:18:24 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE082516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 26 11:18:03 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	5797	0.40	ng/ul	0.02
2) Naphthalene-d8	10.52	136	25726	0.40	ng/ul	0.03
6) Acenaphthene-d10	14.37	164	13262	0.40	ng/ul	0.04
10) Phenanthrene-d10	17.10	188	35151	0.40	ng/ul	0.03
16) Chrysene-d12	21.26	240	44585	0.40	ng/ul	0.03
20) Perylene-d12	23.67	264	36395	0.40	ng/ul	0.04
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	15009	0.44	ng/ul	0.02
14) Fluoranthene-d10	19.13	212	39224	0.45	ng/ul	0.03
Target Compounds					Qvalue	
3) Naphthalene	10.57	128	26031	0.442	ng/ul	97
5) 2-Methylnaphthalene	12.19	142	17458	0.445	ng/ul	99
7) Acenaphthylene	14.09	152	21259	0.370	ng/ul	99
8) Acenaphthene	14.43	153	18393	0.458	ng/ul	98
9) Fluorene	15.42	166	21225	0.455	ng/ul	98
11) Pentachlorophenol	16.75	266	1587	0.383	ng/ul	98
12) Phenanthrene	17.14	178	39837	0.459	ng/ul	98
13) Anthracene	17.14	178	39793	0.505	ng/ul	93
15) Fluoranthene	19.16	202	48304	0.473	ng/ul	97
17) Pyrene	19.52	202	53237	0.509	ng/ul#	100
18) Benzo(a)anthracene	21.24	228	49531	0.518	ng/ul	100
19) Chrysene	21.30	228	49622	0.443	ng/ul	100
21) Benzo(b)fluoranthene	22.93	252	42682	0.476	ng/ul	98
22) Benzo(k)fluoranthene	22.98	252	46654	0.443	ng/ul	97
23) Benzo(a)pyrene	23.55	252	42194	0.445	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.18	276	47431	0.460	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.21	278	39859	0.485	ng/ul	96
26) Benzo(g,h,i)perylene	26.96	276	40216	0.440	ng/ul	95

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

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