

Data Path : Z:\HPCHEM1\BNA_E\Data\BE100717\
 Data File : BE094196.D
 Acq On : 7 Oct 2017 11:01
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.418

Quant Time: Oct 07 23:51:29 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 22 01:38:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1232	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	5885	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	3405	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.24	188	8364	0.40	ng/ul	-0.03
16) Chrysene-d12	21.40	240	8972	0.40	ng/ul	-0.01
20) Perylene-d12	23.90	264	8778	0.40	ng/ul	-0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	3667	0.40	ng/ul	-0.02
14) Fluoranthene-d10	19.26	212	9879	0.35	ng/ul	-0.02
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	5710	0.396	ng/ul#	93
5) 2-Methylnaphthalene	12.35	142	3876	0.414	ng/ul	98
7) Acenaphthylene	14.24	152	5774	0.435	ng/ul	98
8) Acenaphthene	14.57	153	4382	0.401	ng/ul	98
9) Fluorene	15.56	166	5108	0.399	ng/ul	93
11) Pentachlorophenol	16.93	266	312	0.318	ng/ul	96
12) Phenanthrene	17.27	178	8890	0.407	ng/ul#	86
13) Anthracene	17.38	178	8471	0.398	ng/ul#	93
15) Fluoranthene	19.30	202	10184	0.355	ng/ul	84
17) Pyrene	19.66	202	10601	0.425	ng/ul#	83
18) Benzo(a)anthracene	21.39	228	10085	0.424	ng/ul#	86
19) Chrysene	21.44	228	10437	0.380	ng/ul	99
21) Benzo(b)fluoranthene	23.13	252	10096	0.423	ng/ul	86
22) Benzo(k)fluoranthene	23.18	252	10211	0.370	ng/ul#	88
23) Benzo(a)pyrene	23.79	252	10020	0.404	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.55	276	10767	0.425	ng/ul#	83
25) Dibenzo(a,h)anthracene	26.56	278	9000	0.423	ng/ul#	92
26) Benzo(g,h,i)perylene	27.36	276	9085	0.416	ng/ul	98

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

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