

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102017\
 Data File : BE094493.D
 Acq On : 21 Oct 2017 2:24
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.446

Quant Time: Oct 22 02:49:39 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 18 17:51:02 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1338	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	7424	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	4338	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	9783	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	10227	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	9935	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	4725	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	11780	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.75	128	7177	0.395	ng/ul#	93
5) 2-Methylnaphthalene	12.36	142	5004	0.424	ng/ul	99
7) Acenaphthylene	14.24	152	7985	0.472	ng/ul	97
8) Acenaphthene	14.58	153	5648	0.406	ng/ul	98
9) Fluorene	15.57	166	6465	0.397	ng/ul#	94
11) Pentachlorophenol	16.98	266	400	0.348	ng/ul	95
12) Phenanthrene	17.29	178	10335	0.404	ng/ul#	86
13) Anthracene	17.39	178	10334	0.416	ng/ul#	93
15) Fluoranthene	19.31	202	12182	0.363	ng/ul	84
17) Pyrene	19.67	202	12792	0.450	ng/ul#	82
18) Benzo(a)anthracene	21.40	228	11898	0.439	ng/ul#	87
19) Chrysene	21.46	228	11755	0.375	ng/ul	98
21) Benzo(b)fluoranthene	23.16	252	11137	0.413	ng/ul	88
22) Benzo(k)fluoranthene	23.22	252	11724	0.376	ng/ul	91
23) Benzo(a)pyrene	23.83	252	11033	0.393	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	26.63	276	11855	0.413	ng/ul#	85
25) Dibenzo(a,h)anthracene	26.63	278	10064	0.418	ng/ul	93
26) Benzo(g,h,i)perylene	27.47	276	9415	0.380	ng/ul	99

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

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