

Data Path : Z:\HPCHEM1\BNA E\DATA\BE100917\
 Data File : BE094230.D
 Acq On : 9 Oct 2017 20:47
 Operator : SJ/JU
 Sample : SSTDC000.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.423

Quant Time: Oct 10 09:21:32 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 10 09:18:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1990	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	10328	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	5695	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12584	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	13327	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	12877	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	6475	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	15285	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	9931	0.393	ng/ul#	92
5) 2-Methylnaphthalene	12.35	142	6847	0.417	ng/ul	99
7) Acenaphthylene	14.24	152	10298	0.464	ng/ul	96
8) Acenaphthene	14.57	153	7374	0.403	ng/ul	96
9) Fluorene	15.56	166	8283	0.387	ng/ul#	91
11) Pentachlorophenol	16.93	266	334	0.226	ng/ul	96
12) Phenanthrene	17.29	178	13251	0.403	ng/ul#	93
13) Anthracene	17.38	178	12995	0.406	ng/ul#	92
15) Fluoranthene	19.30	202	15602	0.361	ng/ul	84
17) Pyrene	19.66	202	16538	0.446	ng/ul#	83
18) Benzo(a)anthracene	21.39	228	15750	0.445	ng/ul#	87
19) Chrysene	21.45	228	15173	0.372	ng/ul	99
21) Benzo(b)fluoranthene	23.14	252	14995	0.429	ng/ul	92
22) Benzo(k)fluoranthene	23.20	252	15043	0.372	ng/ul	94
23) Benzo(a)pyrene	23.81	252	14744	0.406	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	26.57	276	16052	0.432	ng/ul#	84
25) Dibenzo(a,h)anthracene	26.59	278	13466	0.432	ng/ul	93
26) Benzo(g,h,i)perylene	27.39	276	13497	0.421	ng/ul	96

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

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