

Data Path : Z:\HPCHEM1\BNA E\DATA\BE101017\
 Data File : BE094275.D
 Acq On : 11 Oct 2017 5:54
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.428

Quant Time: Oct 11 08:28:14 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 11 08:24:24 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1938	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	9886	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	5298	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	11484	0.40	ng/ul	-0.02
16) Chrysene-d12	21.41	240	11928	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	11620	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	6151	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	13963	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	9468	0.391	ng/ul#	92
5) 2-Methylnaphthalene	12.35	142	6475	0.412	ng/ul	99
7) Acenaphthylene	14.24	152	9669	0.468	ng/ul	97
8) Acenaphthene	14.57	153	6735	0.396	ng/ul	97
9) Fluorene	15.56	166	7561	0.380	ng/ul#	89
11) Pentachlorophenol	16.95	266	375	0.278	ng/ul	96
12) Phenanthrene	17.29	178	12029	0.401	ng/ul#	93
13) Anthracene	17.38	178	12185	0.417	ng/ul#	92
15) Fluoranthene	19.30	202	14151	0.359	ng/ul	84
17) Pyrene	19.66	202	14946	0.450	ng/ul#	84
18) Benzo(a)anthracene	21.39	228	14371	0.454	ng/ul#	86
19) Chrysene	21.45	228	13772	0.377	ng/ul	98
21) Benzo(b)fluoranthene	23.14	252	13811	0.438	ng/ul	91
22) Benzo(k)fluoranthene	23.19	252	13355	0.366	ng/ul	94
23) Benzo(a)pyrene	23.81	252	13259	0.404	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.58	276	14430	0.430	ng/ul#	84
25) Dibenzo(a,h)anthracene	26.59	278	12106	0.430	ng/ul	94
26) Benzo(g,h,i)perylene	27.39	276	11762	0.406	ng/ul	94

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

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