

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA E\Data\Be111017\
 Data File : BE094800.D
 Acq On : 9 Nov 2017 13:17
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.483

Quant Time: Nov 09 18:45:52 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 07 01:37:48 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.94	152	1181	0.40	ng/ul	0.02
2) Naphthalene-d8	10.72	136	7238	0.40	ng/ul	0.02
6) Acenaphthene-d10	14.53	164	3909	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	8412	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	8787	0.40	ng/ul	0.00
20) Perylene-d12	23.99	264	8050	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.31	152	4261	0.36	ng/ul	0.01
14) Fluoranthene-d10	19.30	212	9548	0.38	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.77	128	6464	0.365	ng/ul#	95
5) 2-Methylnaphthalene	12.38	142	4317	0.349	ng/ul	99
7) Acenaphthylene	14.26	152	6475	0.366	ng/ul	97
8) Acenaphthene	14.59	153	4626	0.364	ng/ul	97
9) Fluorene	15.59	166	5139	0.360	ng/ul	94
11) Pentachlorophenol	17.08	266	85	0.202	ng/ul#	90
12) Phenanthrene	17.31	178	8283	0.373	ng/ul#	88
13) Anthracene	17.41	178	8145	0.371	ng/ul#	94
15) Fluoranthene	19.33	202	9856	0.384	ng/ul	84
17) Pyrene	19.69	202	10500	0.387	ng/ul#	81
18) Benzo(a)anthracene	21.42	228	9787	0.390	ng/ul#	87
19) Chrysene	21.48	228	9787	0.377	ng/ul	98
21) Benzo(b)fluoranthene	23.20	252	8625	0.381	ng/ul	89
22) Benzo(k)fluoranthene	23.25	252	9576	0.395	ng/ul	92
23) Benzo(a)pyrene	23.88	252	8480	0.375	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	26.72	276	6615	0.282	ng/ul#	82
25) Dibenzo(a,h)anthracene	26.71	278	5836	0.296	ng/ul	95
26) Benzo(g,h,i)perylene	27.56	276	4627	0.246	ng/ul#	89

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

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