

Data Path : Z:\HPCHEM1\BNA_E\Data\BE110817\
 Data File : BE094782.D
 Acq On : 8 Nov 2017 13:07
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.481

Quant Time: Nov 08 15:25:33 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 08 15:01:00 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1183	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	7289	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3865	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	8278	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	8553	0.40	ng/ul	0.00
20) Perylene-d12	23.98	264	8003	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	4276	0.36	ng/ul	0.01
14) Fluoranthene-d10	19.29	212	9489	0.39	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.76	128	6452	0.362	ng/ul#	95
5) 2-Methylnaphthalene	12.38	142	4329	0.347	ng/ul	98
7) Acenaphthylene	14.26	152	6491	0.371	ng/ul	98
8) Acenaphthene	14.59	153	4604	0.366	ng/ul	97
9) Fluorene	15.58	166	5072	0.360	ng/ul	93
11) Pentachlorophenol	17.08	266	110	0.266	ng/ul#	83
12) Phenanthrene	17.31	178	7988	0.365	ng/ul#	89
13) Anthracene	17.41	178	8056	0.373	ng/ul#	93
15) Fluoranthene	19.33	202	9630	0.381	ng/ul	84
17) Pyrene	19.69	202	10204	0.386	ng/ul#	81
18) Benzo(a)anthracene	21.42	228	9498	0.388	ng/ul#	87
19) Chrysene	21.47	228	9428	0.373	ng/ul	98
21) Benzo(b)fluoranthene	23.20	252	8481	0.377	ng/ul	93
22) Benzo(k)fluoranthene	23.25	252	9245	0.383	ng/ul	95
23) Benzo(a)pyrene	23.87	252	8435	0.375	ng/ul	91
24) Indeno(1,2,3-cd)pyrene	26.71	276	7058	0.303	ng/ul#	81
25) Dibenzo(a,h)anthracene	26.70	278	6213	0.317	ng/ul#	95
26) Benzo(g,h,i)perylene	27.55	276	5064	0.271	ng/ul#	87

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

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