

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112217\
 Data File : BE094861.D
 Acq On : 23 Nov 2017 8:26
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.496

Quant Time: Nov 23 23:16:26 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE111417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 23 23:11:00 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.94	152	579	0.40	ng/ul	0.02
2) Naphthalene-d8	10.71	136	3958	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.51	164	2834	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	7182	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	7616	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	7312	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	2588	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	8288	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.76	128	3370	0.357	ng/ul#	94
5) 2-Methylnaphthalene	12.37	142	2611	0.398	ng/ul	99
7) Acenaphthylene	14.24	152	4564	0.352	ng/ul	99
8) Acenaphthene	14.57	153	3318	0.363	ng/ul	96
9) Fluorene	15.57	166	3983	0.360	ng/ul	91
11) Pentachlorophenol	17.01	266	140	0.370	ng/ul	93
12) Phenanthrene	17.29	178	6887	0.357	ng/ul#	87
13) Anthracene	17.39	178	6939	0.353	ng/ul#	92
15) Fluoranthene	19.31	202	8473	0.335	ng/ul	84
17) Pyrene	19.68	202	8854	0.371	ng/ul#	80
18) Benzo(a)anthracene	21.41	228	8464	0.366	ng/ul#	86
19) Chrysene	21.46	228	8471	0.366	ng/ul	99
21) Benzo(b)fluoranthene	23.18	252	8070	0.370	ng/ul	87
22) Benzo(k)fluoranthene	23.23	252	8352	0.360	ng/ul#	88
23) Benzo(a)pyrene	23.86	252	7979	0.368	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.68	276	8860	0.375	ng/ul#	78
25) Dibenzo(a,h)anthracene	26.66	278	7339	0.369	ng/ul	91
26) Benzo(g,h,i)perylene	27.52	276	7411	0.375	ng/ul	95

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

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