

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE111417\
 Data File : BE094805.D
 Acq On : 14 Nov 2017 13:18
 Operator : SJ/JU
 Sample : SSTD0.887
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.887

Quant Time: Nov 14 14:05:35 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE111417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 14 13:51:20 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	771	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.70	136	4460	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	2741	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	7052	0.40	ng/ul	-0.02
16) Chrysene-d12	21.43	240	8149	0.40	ng/ul	0.00
20) Perylene-d12	23.98	264	7653	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	5730	0.79	ng/ul	-0.01
14) Fluoranthene-d10	19.29	212	18750	0.90	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	8296	0.761	ng/ul#	93
5) 2-Methylnaphthalene	12.37	142	5927	0.777	ng/ul	99
7) Acenaphthylene	14.24	152	9902	0.799	ng/ul	97
8) Acenaphthene	14.58	153	7011	0.787	ng/ul	99
9) Fluorene	15.57	166	8591	0.859	ng/ul#	94
11) Pentachlorophenol	17.01	266	292	0.829	ng/ul	97
12) Phenanthrene	17.29	178	14730	0.791	ng/ul#	86
13) Anthracene	17.40	178	15234	0.829	ng/ul#	92
15) Fluoranthene	19.32	202	19125	0.888	ng/ul	84
17) Pyrene	19.69	202	20383	0.809	ng/ul#	80
18) Benzo(a)anthracene	21.41	228	19569	0.840	ng/ul#	85
19) Chrysene	21.47	228	19195	0.797	ng/ul	98
21) Benzo(b)fluoranthene	23.19	252	17511	0.813	ng/ul	86
22) Benzo(k)fluoranthene	23.24	252	19061	0.826	ng/ul#	87
23) Benzo(a)pyrene	23.87	252	17747	0.825	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.69	276	19424	0.872	ng/ul#	81
25) Dibenzo(a,h)anthracene	26.68	278	16336	0.871	ng/ul	91
26) Benzo(g,h,i)perylene	27.53	276	16129	0.902	ng/ul	96

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

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