

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102017\
 Data File : BE094467.D
 Acq On : 20 Oct 2017 10:15
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.444

Quant Time: Oct 22 02:46:04 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 18 17:51:02 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1465	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	7983	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	4674	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	10678	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	11060	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	10799	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	5118	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	12836	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.75	128	7773	0.398	ng/ul#	93
5) 2-Methylnaphthalene	12.36	142	5466	0.431	ng/ul	99
7) Acenaphthylene	14.24	152	8627	0.473	ng/ul	97
8) Acenaphthene	14.58	153	6088	0.406	ng/ul	99
9) Fluorene	15.57	166	7054	0.402	ng/ul#	93
11) Pentachlorophenol	16.98	266	479	0.382	ng/ul	98
12) Phenanthrene	17.29	178	11364	0.407	ng/ul#	85
13) Anthracene	17.39	178	11025	0.406	ng/ul#	93
15) Fluoranthene	19.31	202	13206	0.360	ng/ul	85
17) Pyrene	19.68	202	13913	0.452	ng/ul#	82
18) Benzo(a)anthracene	21.40	228	13052	0.445	ng/ul#	86
19) Chrysene	21.46	228	12528	0.370	ng/ul	97
21) Benzo(b)fluoranthene	23.17	252	12198	0.416	ng/ul	90
22) Benzo(k)fluoranthene	23.22	252	12529	0.369	ng/ul	92
23) Benzo(a)pyrene	23.84	252	11908	0.391	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	26.63	276	13001	0.417	ng/ul#	84
25) Dibenzo(a,h)anthracene	26.64	278	10964	0.419	ng/ul	93
26) Benzo(g,h,i)perylene	27.47	276	10464	0.389	ng/ul	97

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

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