

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094397.D
 Acq On : 15 Oct 2017 17:30
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.440

Quant Time: Oct 16 06:56:29 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 06:50:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1648	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	8693	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5263	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	11958	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	12217	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	11671	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	5649	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	14409	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	8340	0.392	ng/ul#	93
5) 2-Methylnaphthalene	12.35	142	6041	0.437	ng/ul	98
7) Acenaphthylene	14.24	152	9774	0.476	ng/ul	96
8) Acenaphthene	14.58	153	6777	0.401	ng/ul	99
9) Fluorene	15.57	166	7930	0.401	ng/ul#	91
11) Pentachlorophenol	16.96	266	709	0.505	ng/ul	96
12) Phenanthrene	17.29	178	12602	0.403	ng/ul#	87
13) Anthracene	17.39	178	12662	0.417	ng/ul	93
15) Fluoranthene	19.31	202	14765	0.360	ng/ul	84
17) Pyrene	19.67	202	15410	0.453	ng/ul#	82
18) Benzo(a)anthracene	21.40	228	14447	0.446	ng/ul#	88
19) Chrysene	21.46	228	13594	0.363	ng/ul	99
21) Benzo(b)fluoranthene	23.17	252	13283	0.419	ng/ul	93
22) Benzo(k)fluoranthene	23.22	252	13625	0.372	ng/ul	93
23) Benzo(a)pyrene	23.83	252	12740	0.387	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.62	276	14294	0.424	ng/ul#	85
25) Dibenzo(a,h)anthracene	26.63	278	12129	0.429	ng/ul	93
26) Benzo(g,h,i)perylene	27.45	276	11969	0.412	ng/ul	97

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

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