

Data Path : Z:\HPCHEM1\BNA E\DATA\BE100917\
 Data File : BE094247.D
 Acq On : 10 Oct 2017 7:20
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.424

Quant Time: Oct 10 08:54:19 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 09 04:05:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1860	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	9461	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5181	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	11365	0.40	ng/ul	-0.02
16) Chrysene-d12	21.41	240	11899	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	11402	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	5863	0.40	ng/ul	-0.01
14) Fluoranthene-d10	19.27	212	14174	0.37	ng/ul	-0.01
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	9111	0.393	ng/ul#	92
5) 2-Methylnaphthalene	12.35	142	6283	0.418	ng/ul	98
7) Acenaphthylene	14.24	152	9882	0.489	ng/ul	96
8) Acenaphthene	14.58	153	6691	0.402	ng/ul	98
9) Fluorene	15.57	166	7562	0.388	ng/ul#	93
11) Pentachlorophenol	16.93	266	460	0.345	ng/ul	98
12) Phenanthrene	17.29	178	12190	0.410	ng/ul#	91
13) Anthracene	17.38	178	12466	0.432	ng/ul#	91
15) Fluoranthene	19.30	202	14396	0.369	ng/ul	85
17) Pyrene	19.66	202	15009	0.453	ng/ul#	84
18) Benzo(a)anthracene	21.40	228	14215	0.450	ng/ul#	88
19) Chrysene	21.45	228	13270	0.364	ng/ul	92
21) Benzo(b)fluoranthene	23.15	252	14228	0.459	ng/ul	90
22) Benzo(k)fluoranthene	23.20	252	13063	0.365	ng/ul#	90
23) Benzo(a)pyrene	23.81	252	13332	0.414	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	26.58	276	14530	0.441	ng/ul#	85
25) Dibenzo(a,h)anthracene	26.59	278	12133	0.439	ng/ul#	90
26) Benzo(g,h,i)perylene	27.41	276	12280	0.432	ng/ul#	86

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

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