

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA E\Data\Be111017\
 Data File : BE094795.D
 Acq On : 9 Nov 2017 9:09
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.482

Quant Time: Nov 09 12:49:02 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 07 01:37:48 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.94	152	1205	0.40	ng/ul	0.01
2) Naphthalene-d8	10.72	136	7319	0.40	ng/ul	0.02
6) Acenaphthene-d10	14.53	164	3946	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	8458	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	8800	0.40	ng/ul	0.00
20) Perylene-d12	23.98	264	8167	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.31	152	4315	0.36	ng/ul	0.01
14) Fluoranthene-d10	19.30	212	9597	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.77	128	6567	0.367	ng/ul#	94
5) 2-Methylnaphthalene	12.38	142	4367	0.349	ng/ul	100
7) Acenaphthylene	14.26	152	6660	0.373	ng/ul	98
8) Acenaphthene	14.59	153	4764	0.371	ng/ul	96
9) Fluorene	15.59	166	5198	0.361	ng/ul#	92
11) Pentachlorophenol	17.07	266	105	0.249	ng/ul#	88
12) Phenanthrene	17.31	178	8073	0.361	ng/ul#	88
13) Anthracene	17.41	178	8087	0.367	ng/ul#	93
15) Fluoranthene	19.33	202	9794	0.379	ng/ul	83
17) Pyrene	19.69	202	10467	0.385	ng/ul#	81
18) Benzo(a)anthracene	21.42	228	9596	0.381	ng/ul#	87
19) Chrysene	21.48	228	9859	0.379	ng/ul	100
21) Benzo(b)fluoranthene	23.20	252	8583	0.374	ng/ul	91
22) Benzo(k)fluoranthene	23.25	252	9336	0.379	ng/ul	93
23) Benzo(a)pyrene	23.87	252	8482	0.369	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.71	276	6961	0.293	ng/ul#	81
25) Dibenzo(a,h)anthracene	26.70	278	6101	0.305	ng/ul	95
26) Benzo(g,h,i)perylene	27.55	276	4900	0.257	ng/ul#	89

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

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