

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101117\
 Data File : BE094293.D
 Acq On : 11 Oct 2017 23:32
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.431

Quant Time: Oct 12 07:37:12 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 12 07:31:23 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1859	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	9155	0.40	ng/ul	0.02
6) Acenaphthene-d10	14.53	164	4827	0.40	ng/ul	0.01
10) Phenanthrene-d10	17.26	188	10379	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	10653	0.40	ng/ul	0.02
20) Perylene-d12	23.96	264	10830	0.40	ng/ul	0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	5624	0.40	ng/ul	0.01
14) Fluoranthene-d10	19.29	212	12748	0.37	ng/ul	0.01
Target Compounds					Ovalue	
3) Naphthalene	10.75	128	8802	0.393	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	5916	0.407	ng/ul	99
7) Acenaphthylene	14.25	152	9227	0.490	ng/ul	97
8) Acenaphthene	14.59	153	6188	0.399	ng/ul	96
9) Fluorene	15.57	166	7161	0.395	ng/ul	92
11) Pentachlorophenol	16.94	266	690	0.567	ng/ul#	80
12) Phenanthrene	17.31	178	10964	0.404	ng/ul	97
13) Anthracene	17.39	178	11362	0.431	ng/ul	96
15) Fluoranthene	19.32	202	13051	0.366	ng/ul	84
17) Pyrene	19.68	202	13870	0.468	ng/ul#	84
18) Benzo(a)anthracene	21.41	228	12818	0.454	ng/ul#	68
19) Chrysene	21.46	228	12113	0.371	ng/ul#	64
21) Benzo(b)fluoranthene	23.17	252	12123	0.412	ng/ul#	46
22) Benzo(k)fluoranthene	23.22	252	12231	0.360	ng/ul#	45
23) Benzo(a)pyrene	23.84	252	12363	0.404	ng/ul#	50
24) Indeno(1,2,3-cd)pyrene	26.63	276	12426	0.397	ng/ul#	89
25) Dibenzo(a,h)anthracene	26.64	278	11191	0.427	ng/ul#	35
26) Benzo(g,h,i)perylene	27.47	276	8489	0.315	ng/ul#	35

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

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