

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101117\
 Data File : BE094279.D
 Acq On : 11 Oct 2017 14:56
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.430

Quant Time: Oct 12 07:23:55 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 07:23:36 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	2041	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	10504	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	5646	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12038	0.40	ng/ul	0.02
16) Chrysene-d12	21.41	240	12781	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	12223	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	6538	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	14529	0.36	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.74	128	10130	0.394	ng/ul#	92
5) 2-Methylnaphthalene	12.35	142	6913	0.414	ng/ul	99
7) Acenaphthylene	14.24	152	10211	0.464	ng/ul	96
8) Acenaphthene	14.58	153	7256	0.400	ng/ul	97
9) Fluorene	15.56	166	8116	0.382	ng/ul#	91
11) Pentachlorophenol	16.94	266	372	0.263	ng/ul	96
12) Phenanthrene	17.29	178	12716	0.404	ng/ul#	93
13) Anthracene	17.38	178	12815	0.419	ng/ul#	92
15) Fluoranthene	19.30	202	14687	0.355	ng/ul	84
17) Pyrene	19.66	202	15414	0.433	ng/ul#	81
18) Benzo(a)anthracene	21.39	228	15294	0.451	ng/ul#	86
19) Chrysene	21.45	228	14537	0.371	ng/ul	97
21) Benzo(b)fluoranthene	23.15	252	14968	0.451	ng/ul	89
22) Benzo(k)fluoranthene	23.19	252	14533	0.379	ng/ul#	91
23) Benzo(a)pyrene	23.81	252	14071	0.408	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	26.58	276	15096	0.428	ng/ul#	83
25) Dibenzo(a,h)anthracene	26.59	278	12811	0.433	ng/ul	94
26) Benzo(g,h,i)perylene	27.40	276	12246	0.402	ng/ul	98

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

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