

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101117\
 Data File : BE094277.D
 Acq On : 11 Oct 2017 12:25
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.429

Quant Time: Oct 12 07:20:59 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 11 08:24:24 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1867	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	9936	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	5404	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	11628	0.40	ng/ul	-0.02
16) Chrysene-d12	21.41	240	12185	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	11985	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	6223	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	14011	0.36	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.74	128	9578	0.394	ng/ul#	92
5) 2-Methylnaphthalene	12.35	142	6594	0.418	ng/ul	98
7) Acenaphthylene	14.24	152	9797	0.465	ng/ul	96
8) Acenaphthene	14.57	153	6916	0.399	ng/ul	98
9) Fluorene	15.56	166	7776	0.383	ng/ul#	91
11) Pentachlorophenol	16.94	266	320	0.235	ng/ul	97
12) Phenanthrene	17.29	178	12099	0.398	ng/ul#	93
13) Anthracene	17.38	178	12208	0.413	ng/ul#	92
15) Fluoranthene	19.30	202	14280	0.358	ng/ul	84
17) Pyrene	19.66	202	15145	0.447	ng/ul#	85
18) Benzo(a)anthracene	21.39	228	14570	0.451	ng/ul#	85
19) Chrysene	21.44	228	14050	0.376	ng/ul	97
21) Benzo(b)fluoranthene	23.15	252	14291	0.439	ng/ul	90
22) Benzo(k)fluoranthene	23.20	252	13659	0.363	ng/ul	92
23) Benzo(a)pyrene	23.81	252	13669	0.404	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	26.58	276	14659	0.424	ng/ul#	83
25) Dibenzo(a,h)anthracene	26.59	278	12352	0.426	ng/ul	93
26) Benzo(g,h,i)perylene	27.40	276	11909	0.399	ng/ul	97

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

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