

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094385.D
 Acq On : 15 Oct 2017 10:08
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.439

Quant Time: Oct 16 05:11:29 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 16 05:09:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1413	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	7731	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	4698	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	11002	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	11070	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	10477	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	4957	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	12852	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.75	128	7458	0.394	ng/ul#	95
5) 2-Methylnaphthalene	12.36	142	5378	0.438	ng/ul	98
7) Acenaphthylene	14.24	152	8882	0.485	ng/ul	97
8) Acenaphthene	14.58	153	6191	0.411	ng/ul	97
9) Fluorene	15.57	166	7169	0.406	ng/ul#	94
11) Pentachlorophenol	16.96	266	567	0.439	ng/ul	96
12) Phenanthrene	17.29	178	11492	0.400	ng/ul#	87
13) Anthracene	17.40	178	11845	0.424	ng/ul#	93
15) Fluoranthene	19.31	202	13641	0.361	ng/ul	84
17) Pyrene	19.67	202	14758	0.479	ng/ul#	84
18) Benzo(a)anthracene	21.40	228	13093	0.446	ng/ul#	90
19) Chrysene	21.46	228	12372	0.365	ng/ul	92
21) Benzo(b)fluoranthene	23.17	252	12457	0.438	ng/ul	94
22) Benzo(k)fluoranthene	23.22	252	11962	0.364	ng/ul#	94
23) Benzo(a)pyrene	23.83	252	11486	0.388	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	26.62	276	13148	0.435	ng/ul#	84
25) Dibenzo(a,h)anthracene	26.63	278	11076	0.436	ng/ul#	92
26) Benzo(g,h,i)perylene	27.45	276	11148	0.427	ng/ul#	91

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

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