

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101317\
 Data File : BE094352.D
 Acq On : 14 Oct 2017 9:55
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.436

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:23:31 PM

Quant Time: Oct 15 04:34:43 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 14 05:22:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1075	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	5129	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	2900	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	7722	0.40	ng/ul	-0.02
16) Chrysene-d12	21.43	240	9417	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	8875	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	3315	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	11199	0.43	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.75	128	5147	0.41	ng/ul#	95
5) 2-Methylnaphthalene	12.36	142	3534	0.43	ng/ul	97
7) Acenaphthylene	14.25	152	5602m	0.50	ng/ul	
8) Acenaphthene	14.59	153	3972	0.43	ng/ul	97
9) Fluorene	15.57	166	4895	0.45	ng/ul	93
11) Pentachlorophenol	16.96	266	312	0.34	ng/ul	91
12) Phenanthrene	17.31	178	8596	0.43	ng/ul	95
13) Anthracene	17.40	178	8935m	0.46	ng/ul	
15) Fluoranthene	19.32	202	11361	0.43	ng/ul	84
17) Pyrene	19.68	202	12094	0.46	ng/ul#	83
18) Benzo(a)anthracene	21.41	228	11793	0.47	ng/ul#	88
19) Chrysene	21.47	228	10517m	0.36	ng/ul	
21) Benzo(b)fluoranthene	23.18	252	10780	0.45	ng/ul#	84
22) Benzo(k)fluoranthene	23.23	252	10766	0.39	ng/ul#	82
23) Benzo(a)pyrene	23.85	252	10474	0.42	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	26.64	276	11655	0.45	ng/ul#	83
25) Dibenzo(a,h)anthracene	26.65	278	9929	0.46	ng/ul#	82
26) Benzo(g,h,i)perylene	27.48	276	9593	0.43	ng/ul#	82

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

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