

Data Path : Z:\HPCHEM1\BNA E\DATA\BE093017\
 Data File : BE094153.D
 Acq On : 30 Sep 2017 10:33
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.416

Quant Time: Sep 30 22:14:57 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 22 01:38:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1427	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	7731	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	4874	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.24	188	12618	0.40	ng/ul	-0.03
16) Chrysene-d12	21.40	240	13545	0.40	ng/ul	-0.01
20) Perylene-d12	23.91	264	13111	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	4969	0.41	ng/ul	-0.01
14) Fluoranthene-d10	19.27	212	14488	0.34	ng/ul	-0.02
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	8775	0.463	ng/ul#	92
5) 2-Methylnaphthalene	12.35	142	5904	0.481	ng/ul	99
7) Acenaphthylene	14.24	152	9820	0.517	ng/ul	96
8) Acenaphthene	14.58	153	7350	0.470	ng/ul	98
9) Fluorene	15.56	166	8481	0.463	ng/ul#	93
11) Pentachlorophenol	16.93	266	680	0.459	ng/ul	97
12) Phenanthrene	17.29	178	14758	0.448	ng/ul#	93
13) Anthracene	17.38	178	14028	0.437	ng/ul#	92
15) Fluoranthene	19.30	202	17548	0.405	ng/ul	84
17) Pyrene	19.66	202	18654	0.495	ng/ul#	81
18) Benzo(a)anthracene	21.39	228	17694	0.492	ng/ul#	85
19) Chrysene	21.44	228	18115	0.437	ng/ul	98
21) Benzo(b)fluoranthene	23.14	252	17198	0.483	ng/ul	86
22) Benzo(k)fluoranthene	23.19	252	17254	0.419	ng/ul#	88
23) Benzo(a)pyrene	23.80	252	16692	0.451	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	26.55	276	17857	0.472	ng/ul#	84
25) Dibenzo(a,h)anthracene	26.56	278	15111	0.476	ng/ul#	90
26) Benzo(g,h,i)perylene	27.37	276	15290	0.468	ng/ul	98

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

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