

Data Path : Z:\HPCHEM1\BNA_E\Data\BE062017\
 Data File : BE093276.D
 Acq On : 20 Jun 2017 11:31
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.416

Quant Time: Jun 21 01:48:19 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 20 08:49:11 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1918	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	9968	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5697	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	14385	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	15245	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	13074	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	6329	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	16197	0.41	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	10165	0.411	ng/ul#	92
5) 2-Methylnaphthalene	12.18	142	7280	0.430	ng/ul	98
7) Acenaphthylene	14.09	152	10953	0.448	ng/ul	95
8) Acenaphthene	14.42	153	8094	0.410	ng/ul	96
9) Fluorene	15.42	166	9244	0.400	ng/ul	93
11) Pentachlorophenol	16.76	266	986	0.462	ng/ul	95
12) Phenanthrene	17.14	178	16172	0.409	ng/ul#	89
13) Anthracene	17.22	178	15268	0.443	ng/ul#	89
15) Fluoranthene	19.16	202	19684	0.406	ng/ul	83
17) Pyrene	19.51	202	21118	0.470	ng/ul#	88
18) Benzo(a)anthracene	21.24	228	18680	0.459	ng/ul#	86
19) Chrysene	21.29	228	17331	0.411	ng/ul	96
21) Benzo(b)fluoranthene	22.93	252	16959	0.427	ng/ul	90
22) Benzo(k)fluoranthene	22.98	252	16369	0.388	ng/ul#	91
23) Benzo(a)pyrene	23.55	252	15955	0.425	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	26.18	276	18284	0.421	ng/ul#	93
25) Dibenzo(a,h)anthracene	26.20	278	15271	0.417	ng/ul#	88
26) Benzo(g,h,i)perylene	26.96	276	15051	0.404	ng/ul	96

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

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