

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA_E\Data\BE062317\
 Data File : BE093334.D
 Acq On : 23 Jun 2017 12:30
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.421

Quant Time: Jun 24 03:56:14 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 22 02:10:17 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	1934	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	9927	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5517	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	13168	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	14215	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	11889	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	5977	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	13789	0.39	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.56	128	9759	0.396	ng/ul#	92
5) 2-Methylnaphthalene	12.18	142	6897	0.410	ng/ul	99
7) Acenaphthylene	14.08	152	9292	0.392	ng/ul	97
8) Acenaphthene	14.42	153	7492	0.392	ng/ul	98
9) Fluorene	15.41	166	8307	0.371	ng/ul#	90
11) Pentachlorophenol	16.76	266	656	0.336	ng/ul	96
12) Phenanthrene	17.14	178	14673	0.405	ng/ul#	92
13) Anthracene	17.22	178	13290	0.421	ng/ul#	91
15) Fluoranthene	19.15	202	16798	0.378	ng/ul	84
17) Pyrene	19.51	202	18580	0.443	ng/ul#	85
18) Benzo(a)anthracene	21.24	228	16120	0.425	ng/ul#	86
19) Chrysene	21.29	228	15663	0.398	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	14536	0.403	ng/ul	88
22) Benzo(k)fluoranthene	22.97	252	14797	0.386	ng/ul#	89
23) Benzo(a)pyrene	23.55	252	13849	0.406	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.18	276	15769	0.399	ng/ul#	94
25) Dibenzo(a,h)anthracene	26.20	278	13170	0.396	ng/ul#	87
26) Benzo(g,h,i)perylene	26.96	276	12973	0.383	ng/ul	97

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

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