

Data Path : Z:\HPCHEM1\BNA_E\Data\BE061417\
 Data File : BE093164.D
 Acq On : 14 Jun 2017 11:42
 Operator : SJ/MA
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.404

Quant Time: Jun 15 02:25:00 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 14 08:39:04 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	3109	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	14190	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	7902	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	21234	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	23944	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	19505	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	8759	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	23506	0.41	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	14594	0.415	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	10116	0.420	ng/ul	98
7) Acenaphthylene	14.08	152	13782	0.406	ng/ul	96
8) Acenaphthene	14.43	153	11225	0.410	ng/ul	99
9) Fluorene	15.42	166	12886	0.402	ng/ul#	90
11) Pentachlorophenol	16.76	266	1464	0.465	ng/ul	98
12) Phenanthrene	17.14	178	24035	0.411	ng/ul#	88
13) Anthracene	17.22	178	21639	0.425	ng/ul#	88
15) Fluoranthene	19.15	202	28781	0.402	ng/ul	83
17) Pyrene	19.51	202	30578	0.433	ng/ul#	85
18) Benzo(a)anthracene	21.24	228	27483	0.430	ng/ul#	86
19) Chrysene	21.29	228	27520	0.415	ng/ul	97
21) Benzo(b)fluoranthene	22.92	252	25087	0.424	ng/ul	87
22) Benzo(k)fluoranthene	22.97	252	25023	0.398	ng/ul#	89
23) Benzo(a)pyrene	23.55	252	23806	0.425	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	26.17	276	26704	0.412	ng/ul#	91
25) Dibenzo(a,h)anthracene	26.19	278	22030	0.404	ng/ul#	91
26) Benzo(g,h,i)perylene	26.95	276	22451	0.404	ng/ul	98

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

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