

Data Path : Z:\HPCHEM1\BNA_E\Data\BE061317\
 Data File : BE093162.D
 Acq On : 14 Jun 2017 10:28
 Operator : SJ/MA
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.403

Quant Time: Jun 14 10:58:51 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 08 17:42:17 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	3333	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	15780	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	9054	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	23398	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	25128	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	20510	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	9930	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	25797	0.41	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	16126	0.412	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	11410	0.426	ng/ul	99
7) Acenaphthylene	14.09	152	16137	0.415	ng/ul	96
8) Acenaphthene	14.43	153	12849	0.410	ng/ul	99
9) Fluorene	15.42	166	14819	0.403	ng/ul#	92
11) Pentachlorophenol	16.76	266	1561	0.450	ng/ul	93
12) Phenanthrene	17.14	178	26927	0.418	ng/ul#	88
13) Anthracene	17.22	178	23812	0.425	ng/ul#	87
15) Fluoranthene	19.15	202	31488	0.399	ng/ul	84
17) Pyrene	19.52	202	33831	0.457	ng/ul#	82
18) Benzo(a)anthracene	21.24	228	29385	0.438	ng/ul#	86
19) Chrysene	21.29	228	28918	0.416	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	25663	0.412	ng/ul	88
22) Benzo(k)fluoranthene	22.98	252	26658	0.403	ng/ul#	90
23) Benzo(a)pyrene	23.55	252	25060	0.426	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	26.17	276	27873	0.409	ng/ul#	91
25) Dibenzo(a,h)anthracene	26.20	278	23234	0.405	ng/ul#	91
26) Benzo(g,h,i)perylene	26.95	276	23500	0.402	ng/ul	94

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

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