

Data Path : Z:\HPCHEM1\BNA_E\Data\BE061517\
 Data File : BE093207.D
 Acq On : 15 Jun 2017 14:21
 Operator : SJ/MA
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.408

Quant Time: Jun 16 04:36:18 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 15 07:58:47 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2259	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	12189	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	7221	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	17334	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	17518	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	14850	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	7966	0.45	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	19202	0.41	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	12491	0.413	ng/ul#	91
5) 2-Methylnaphthalene	12.19	142	9184	0.444	ng/ul	99
7) Acenaphthylene	14.08	152	14265	0.460	ng/ul	96
8) Acenaphthene	14.43	153	10293	0.412	ng/ul	98
9) Fluorene	15.42	166	11602	0.396	ng/ul	96
11) Pentachlorophenol	16.77	266	1302	0.506	ng/ul	98
12) Phenanthrene	17.14	178	19360	0.406	ng/ul#	86
13) Anthracene	17.22	178	19143	0.461	ng/ul#	87
15) Fluoranthene	19.16	202	23327	0.399	ng/ul	83
17) Pyrene	19.52	202	24687	0.478	ng/ul#	84
18) Benzo(a)anthracene	21.24	228	21473	0.460	ng/ul#	85
19) Chrysene	21.30	228	19830	0.409	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	19103	0.424	ng/ul	91
22) Benzo(k)fluoranthene	22.98	252	18193	0.380	ng/ul#	92
23) Benzo(a)pyrene	23.56	252	18190	0.427	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	26.19	276	20630	0.418	ng/ul#	92
25) Dibenzo(a,h)anthracene	26.21	278	17133	0.412	ng/ul#	90
26) Benzo(g,h,i)perylene	26.97	276	17152	0.405	ng/ul	96

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

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