

Data Path : Z:\HPCHEM1\BNA_E\Data\BE061317\
 Data File : BE093128.D
 Acq On : 13 Jun 2017 13:23
 Operator : SJ/MA
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.401

Quant Time: Jun 14 01:22:27 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	4830	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	22016	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	11929	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	30805	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	36796	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	31585	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	13349	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	36241	0.43	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	22498	0.412	ng/ul#	91
5) 2-Methylnaphthalene	12.19	142	15270	0.409	ng/ul	99
7) Acenaphthylene	14.09	152	20914	0.409	ng/ul	95
8) Acenaphthene	14.43	153	16948	0.410	ng/ul	98
9) Fluorene	15.42	166	19315	0.399	ng/ul#	89
11) Pentachlorophenol	16.76	266	2397	0.524	ng/ul	98
12) Phenanthrene	17.14	178	35716	0.421	ng/ul#	86
13) Anthracene	17.22	178	32792	0.444	ng/ul#	86
15) Fluoranthene	19.16	202	43800	0.421	ng/ul	83
17) Pyrene	19.52	202	46462	0.428	ng/ul#	82
18) Benzo(a)anthracene	21.24	228	44207	0.450	ng/ul#	86
19) Chrysene	21.29	228	42540	0.418	ng/ul	96
21) Benzo(b)fluoranthene	22.93	252	40273	0.420	ng/ul	90
22) Benzo(k)fluoranthene	22.98	252	39891	0.392	ng/ul#	90
23) Benzo(a)pyrene	23.55	252	38546	0.425	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	26.18	276	42716	0.407	ng/ul#	92
25) Dibenzo(a,h)anthracene	26.21	278	35504	0.402	ng/ul#	90
26) Benzo(g,h,i)perylene	26.95	276	35984	0.400	ng/ul	95

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

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