

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061417\
 Data File : BE093196.D
 Acq On : 15 Jun 2017 7:27
 Operator : SJ/MA
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.406

Quant Time: Jun 15 07:57:03 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 03:59:13 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2524	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	13155	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	7905	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	19875	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	20336	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	17262	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	8502	0.44	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	21851	0.40	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.57	128	13477	0.413	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	9888	0.443	ng/ul	98
7) Acenaphthylene	14.09	152	15244	0.449	ng/ul	97
8) Acenaphthene	14.43	153	11280	0.412	ng/ul	98
9) Fluorene	15.42	166	12807	0.399	ng/ul	90
11) Pentachlorophenol	16.75	266	1444	0.490	ng/ul	96
12) Phenanthrene	17.14	178	22184	0.406	ng/ul#	87
13) Anthracene	17.22	178	21351	0.448	ng/ul#	88
15) Fluoranthene	19.15	202	26196	0.391	ng/ul	84
17) Pyrene	19.51	202	28037	0.468	ng/ul#	86
18) Benzo(a)anthracene	21.24	228	24757	0.456	ng/ul#	86
19) Chrysene	21.30	228	23110	0.411	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	22570	0.431	ng/ul	90
22) Benzo(k)fluoranthene	22.98	252	21369	0.384	ng/ul#	91
23) Benzo(a)pyrene	23.56	252	21135	0.427	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	26.19	276	23781	0.415	ng/ul#	92
25) Dibenzo(a,h)anthracene	26.20	278	19730	0.409	ng/ul#	90
26) Benzo(g,h,i)perylene	26.96	276	19962	0.406	ng/ul	95

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

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