

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA_E\Data\BE062317\
 Data File : BE093340.D
 Acq On : 23 Jun 2017 16:56
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.422

Quant Time: Jun 24 03:59:16 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 24 03:58:33 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1811	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	9653	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5637	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	14052	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	14824	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	12351	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	5938	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	15217	0.40	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	9477	0.396	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	6878	0.420	ng/ul	98
7) Acenaphthylene	14.09	152	10100	0.418	ng/ul	97
8) Acenaphthene	14.43	153	7817	0.401	ng/ul	98
9) Fluorene	15.42	166	8903	0.389	ng/ul	94
11) Pentachlorophenol	16.75	266	932	0.447	ng/ul	97
12) Phenanthrene	17.14	178	15587	0.403	ng/ul#	90
13) Anthracene	17.22	178	15022	0.446	ng/ul#	90
15) Fluoranthene	19.15	202	18255	0.385	ng/ul	84
17) Pyrene	19.51	202	19609	0.449	ng/ul#	86
18) Benzo(a)anthracene	21.24	228	17249	0.436	ng/ul#	85
19) Chrysene	21.29	228	16037	0.391	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	15472	0.413	ng/ul	89
22) Benzo(k)fluoranthene	22.97	252	15390	0.387	ng/ul#	90
23) Benzo(a)pyrene	23.55	252	14719	0.415	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	26.18	276	16727	0.408	ng/ul#	93
25) Dibenzo(a,h)anthracene	26.20	278	13879	0.402	ng/ul#	87
26) Benzo(g,h,i)perylene	26.96	276	13457	0.382	ng/ul	96

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

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