

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA_E\Data\BE062417\
 Data File : BE093353.D
 Acq On : 24 Jun 2017 9:32
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.424

Quant Time: Jun 26 01:11:59 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 04:29:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1617	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	8172	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	4619	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	11968	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13860	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	11863	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	4981	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	13357	0.41	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	8055	0.397	ng/ul#	92
5) 2-Methylnaphthalene	12.18	142	5695	0.411	ng/ul	100
7) Acenaphthylene	14.08	152	8328	0.420	ng/ul	97
8) Acenaphthene	14.42	153	6374	0.399	ng/ul	98
9) Fluorene	15.41	166	7280	0.388	ng/ul	89
11) Pentachlorophenol	16.76	266	766	0.431	ng/ul	98
12) Phenanthrene	17.14	178	12979	0.394	ng/ul#	91
13) Anthracene	17.22	178	12396	0.432	ng/ul#	91
15) Fluoranthene	19.15	202	16226	0.402	ng/ul	83
17) Pyrene	19.51	202	17868	0.437	ng/ul#	84
18) Benzo(a)anthracene	21.24	228	15829	0.428	ng/ul#	85
19) Chrysene	21.29	228	15160	0.395	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	14445	0.401	ng/ul	90
22) Benzo(k)fluoranthene	22.97	252	14706	0.385	ng/ul#	90
23) Benzo(a)pyrene	23.55	252	13853	0.407	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.18	276	15992	0.406	ng/ul#	94
25) Dibenzo(a,h)anthracene	26.20	278	13179	0.397	ng/ul#	86
26) Benzo(g,h,i)perylene	26.96	276	12955	0.383	ng/ul	97

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

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