

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA_E\Data\BE062317\
 Data File : BE093351.D
 Acq On : 23 Jun 2017 23:51
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.423

Quant Time: Jun 24 04:30:32 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	7.73	152	1619	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	8880	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5160	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	12908	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13489	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	11419	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	5513	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	13964	0.40	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	8702	0.395	ng/ul#	92
5) 2-Methylnaphthalene	12.18	142	6309	0.419	ng/ul	99
7) Acenaphthylene	14.08	152	9744	0.440	ng/ul	97
8) Acenaphthene	14.42	153	7070	0.396	ng/ul	99
9) Fluorene	15.41	166	8148	0.389	ng/ul	93
11) Pentachlorophenol	16.75	266	963	0.503	ng/ul	98
12) Phenanthrene	17.13	178	13932	0.392	ng/ul#	91
13) Anthracene	17.22	178	14066	0.455	ng/ul#	91
15) Fluoranthene	19.15	202	16646	0.382	ng/ul	84
17) Pyrene	19.51	202	17793	0.447	ng/ul#	84
18) Benzo(a)anthracene	21.24	228	15771	0.438	ng/ul#	85
19) Chrysene	21.29	228	14532	0.389	ng/ul	96
21) Benzo(b)fluoranthene	22.93	252	14232	0.411	ng/ul	90
22) Benzo(k)fluoranthene	22.97	252	13939	0.379	ng/ul#	91
23) Benzo(a)pyrene	23.55	252	13759	0.420	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	26.18	276	15363	0.405	ng/ul#	94
25) Dibenzo(a,h)anthracene	26.19	278	12697	0.397	ng/ul#	84
26) Benzo(g,h,i)perylene	26.95	276	12491	0.384	ng/ul	95

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

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