

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061517\
 Data File : BE093232.D
 Acq On : 16 Jun 2017 5:46
 Operator : SJ/MA
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.410

Quant Time: Jun 16 06:39:10 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 16 05:36:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2398	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	11979	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	6648	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	16509	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	17012	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	14273	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	7508	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	17915	0.40	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	12233	0.412	ng/ul#	92
5) 2-Methylnaphthalene	12.19	142	8697	0.428	ng/ul	99
7) Acenaphthylene	14.09	152	12554	0.440	ng/ul	96
8) Acenaphthene	14.43	153	9453	0.411	ng/ul	98
9) Fluorene	15.42	166	10614	0.393	ng/ul	94
11) Pentachlorophenol	16.76	266	1051	0.429	ng/ul	98
12) Phenanthrene	17.14	178	18454	0.406	ng/ul#	89
13) Anthracene	17.22	178	17187	0.435	ng/ul#	89
15) Fluoranthene	19.15	202	21977	0.395	ng/ul	84
17) Pyrene	19.51	202	23605	0.471	ng/ul#	87
18) Benzo(a)anthracene	21.24	228	20701	0.456	ng/ul#	84
19) Chrysene	21.29	228	19649	0.417	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	18798	0.434	ng/ul	88
22) Benzo(k)fluoranthene	22.98	252	17691	0.385	ng/ul#	90
23) Benzo(a)pyrene	23.55	252	17425	0.425	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	26.19	276	19690	0.415	ng/ul#	93
25) Dibenzo(a,h)anthracene	26.20	278	16326	0.409	ng/ul#	88
26) Benzo(g,h,i)perylene	26.96	276	16389	0.403	ng/ul	95

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

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