

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061517\
 Data File : BE093236.D
 Acq On : 16 Jun 2017 10:39
 Operator : SJ/MA
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.411

Quant Time: Jun 16 14:39:51 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	2208	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	11314	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	6330	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	15768	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	16539	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	14073	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	7151	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	17400	0.41	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	11566	0.412	ng/ul#	91
5) 2-Methylnaphthalene	12.19	142	8306	0.433	ng/ul	98
7) Acenaphthylene	14.09	152	12189	0.449	ng/ul	96
8) Acenaphthene	14.43	153	9073	0.414	ng/ul	98
9) Fluorene	15.42	166	10172	0.396	ng/ul	93
11) Pentachlorophenol	16.77	266	1024	0.438	ng/ul	96
12) Phenanthrene	17.13	178	17514	0.404	ng/ul#	86
13) Anthracene	17.22	178	16469	0.436	ng/ul#	86
15) Fluoranthene	19.16	202	21333	0.401	ng/ul	83
17) Pyrene	19.52	202	22747	0.466	ng/ul#	84
18) Benzo(a)anthracene	21.24	228	20345	0.461	ng/ul#	85
19) Chrysene	21.29	228	19170	0.419	ng/ul	96
21) Benzo(b)fluoranthene	22.93	252	18484	0.433	ng/ul	88
22) Benzo(k)fluoranthene	22.98	252	17569	0.387	ng/ul#	90
23) Benzo(a)pyrene	23.55	252	17274	0.428	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	26.19	276	19588	0.419	ng/ul#	92
25) Dibenzo(a,h)anthracene	26.20	278	16288	0.414	ng/ul#	89
26) Benzo(g,h,i)perylene	26.96	276	16175	0.403	ng/ul	96

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

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