

Data Path : Z:\HPCHEM1\BNA E\DATA\BE060717\
 Data File : BE093065.D
 Acq On : 7 Jun 2017 12:05
 Operator : SJ/MA
 Sample : SSTD0.152
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.152

Quant Time: Jun 07 14:52:11 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 07 14:48:46 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	3660	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	16767	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	8841	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	25370	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	31613	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	26944	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	2377	0.10	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	6898	0.10	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	4206	0.10	ng/ul#	92
5) 2-Methylnaphthalene	12.19	142	2756	0.09	ng/ul	99
7) Acenaphthylene	14.09	152	3350	0.08	ng/ul	99
8) Acenaphthene	14.43	153	2985	0.09	ng/ul	96
9) Fluorene	15.42	166	3515	0.09	ng/ul	90
12) Phenanthrene	17.14	178	7356	0.09	ng/ul#	88
13) Anthracene	17.22	178	5696	0.07	ng/ul#	88
15) Fluoranthene	19.16	202	8571	0.08	ng/ul	84
17) Pyrene	19.52	202	9167	0.07	ng/ul#	86
18) Benzo(a)anthracene	21.24	228	8321	0.06	ng/ul#	86
19) Chrysene	21.29	228	8820	0.07	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	7983	0.07	ng/ul	90
22) Benzo(k)fluoranthene	22.98	252	8365	0.07	ng/ul	92
23) Benzo(a)pyrene	23.56	252	7414	0.07	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	26.18	276	8670	0.06	ng/ul#	94
25) Dibenzo(a,h)anthracene	26.20	278	7194	0.06	ng/ul#	89
26) Benzo(a,h,i)perylene	26.95	276	7419	0.06	ng/ul	95

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

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