

Data Path : Z:\HPCHEM1\BNA E\DATA\BE101915\
 Data File : BE091026.D
 Acq On : 19 Oct 2015 18:28
 Operator : UM/NP
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.482

Quant Time: Oct 20 01:02:10 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM02.2-EPA-SIM-BE101415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 20 01:00:24 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	5553	0.40	ng/ul	0.00
2) Naphthalene-d8	9.62	136	26206	0.40	ng/ul	0.00
6) Acenaphthene-d10	13.43	164	14429	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.16	188	36823	0.40	ng/ul	0.00
16) Chrysene-d12	20.31	240	45979	0.40	ng/ul	0.00
20) Perylene-d12	22.21	264	43186	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.12	152	11259	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.15	212	32269	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	9.66	128	22114	0.36	ng/ul	98
5) 2-Methylnaphthalene	11.19	142	12665	0.38	ng/ul	97
7) Acenaphthylene	13.18	152	21087	0.40	ng/ul	99
8) Acenaphthene	13.49	153	14580	0.38	ng/ul	96
9) Fluorene	14.48	166	18355	0.39	ng/ul	95
11) Pentachlorophenol	15.91	266	753	0.18	ng/ul	98
12) Phenanthrene	16.20	178	128216	0.39	ng/ul	94
13) Anthracene	16.27	178	121284	0.38	ng/ul	99
15) Fluoranthene	18.17	202	153509	0.37	ng/ul	98
17) Pyrene	18.56	202	152962	0.39	ng/ul	97
18) Benzo(a)anthracene	20.27	228	37008	0.37	ng/ul	99
19) Chrysene	20.34	228	40893	0.38	ng/ul	99
21) Benzo(b)fluoranthene	21.64	252	41616	0.41	ng/ul	83
22) Benzo(k)fluoranthene	21.67	252	39943	0.38	ng/ul#	84
23) Benzo(a)pyrene	22.11	252	38391	0.38	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	23.86	276	156644	0.39	ng/ul	99
25) Dibenzo(a,h)anthracene	23.81	278	39293	0.39	ng/ul	94
26) Benzo(g,h,i)perylene	24.43	276	167174	0.40	ng/ul#	92

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

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