

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094640.D
 Acq On : 30 Oct 2017 9:32
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.466

Quant Time: Oct 30 18:32:46 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 30 04:29:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1231	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	7338	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	3968	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	8599	0.40	ng/ul	0.02
16) Chrysene-d12	21.43	240	8952	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	8617	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	4632	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	10552	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.75	128	6966	0.388	ng/ul#	93
5) 2-Methylnaphthalene	12.36	142	4821	0.384	ng/ul	99
7) Acenaphthylene	14.25	152	7215	0.402	ng/ul	96
8) Acenaphthene	14.58	153	5080	0.394	ng/ul	96
9) Fluorene	15.57	166	5709	0.394	ng/ul	94
11) Pentachlorophenol	17.00	266	215	0.501	ng/ul	96
12) Phenanthrene	17.31	178	8831	0.389	ng/ul#	93
13) Anthracene	17.39	178	9122	0.407	ng/ul#	92
15) Fluoranthene	19.32	202	10681	0.407	ng/ul	84
17) Pyrene	19.68	202	11255	0.407	ng/ul#	83
18) Benzo(a)anthracene	21.42	228	10503	0.410	ng/ul#	88
19) Chrysene	21.47	228	10191	0.385	ng/ul	98
21) Benzo(b)fluoranthene	23.19	252	10076	0.416	ng/ul	95
22) Benzo(k)fluoranthene	23.24	252	9915	0.382	ng/ul	96
23) Benzo(a)pyrene	23.86	252	9570	0.395	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	26.68	276	9033	0.360	ng/ul#	81
25) Dibenzo(a,h)anthracene	26.68	278	7984	0.378	ng/ul#	94
26) Benzo(g,h,i)perylene	27.53	276	6724	0.334	ng/ul#	88

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

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