

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102517\
 Data File : BE094525.D
 Acq On : 25 Oct 2017 9:33
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.454

Quant Time: Oct 25 17:01:53 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 24 20:21:52 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1038	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	6125	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	3750	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	8574	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	9040	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	8795	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	4057	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	10103	0.40	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	6018	0.402	ng/ul#	94
5) 2-Methylnaphthalene	12.37	142	4284	0.409	ng/ul	100
7) Acenaphthylene	14.25	152	6730	0.397	ng/ul	97
8) Acenaphthene	14.59	153	4941	0.405	ng/ul	97
9) Fluorene	15.58	166	5518	0.403	ng/ul	96
11) Pentachlorophenol	17.03	266	120	0.280	ng/ul#	58
12) Phenanthrene	17.31	178	8940	0.395	ng/ul#	93
13) Anthracene	17.39	178	8796	0.394	ng/ul#	91
15) Fluoranthene	19.32	202	10363	0.396	ng/ul	85
17) Pyrene	19.68	202	10940	0.391	ng/ul#	84
18) Benzo(a)anthracene	21.41	228	10093	0.391	ng/ul#	86
19) Chrysene	21.46	228	10625	0.398	ng/ul	98
21) Benzo(b)fluoranthene	23.17	252	9787	0.396	ng/ul	88
22) Benzo(k)fluoranthene	23.22	252	10117	0.382	ng/ul#	90
23) Benzo(a)pyrene	23.84	252	9708	0.393	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	26.64	276	9724	0.380	ng/ul#	85
25) Dibenzo(a,h)anthracene	26.65	278	8311	0.385	ng/ul	94
26) Benzo(g,h,i)perylene	27.48	276	7742	0.377	ng/ul	98

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

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