

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE092017\
 Data File : BE094031.D
 Acq On : 20 Sep 2017 13:46
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.409

Quant Time: Sep 20 16:59:28 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 20 13:07:25 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1298	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	6096	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	3541	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	10024	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	12838	0.40	ng/ul	0.00
20) Perylene-d12	23.94	264	11476	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.32	152	3765	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.30	212	12763	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	5964	0.399	ng/ul#	94
5) 2-Methylnaphthalene	12.39	142	3844	0.397	ng/ul	100
7) Acenaphthylene	14.27	152	5277	0.382	ng/ul	99
8) Acenaphthene	14.60	153	4447	0.391	ng/ul	96
9) Fluorene	15.60	166	5330	0.400	ng/ul	94
11) Pentachlorophenol	16.96	266	386	0.328	ng/ul	96
12) Phenanthrene	17.31	178	10502	0.401	ng/ul#	88
13) Anthracene	17.41	178	9584	0.376	ng/ul#	93
15) Fluoranthene	19.32	202	13242	0.385	ng/ul	85
17) Pyrene	19.68	202	14188	0.397	ng/ul#	84
18) Benzo(a)anthracene	21.41	228	12926	0.380	ng/ul#	86
19) Chrysene	21.47	228	15979	0.406	ng/ul	98
21) Benzo(b)fluoranthene	23.17	252	12630	0.405	ng/ul	87
22) Benzo(k)fluoranthene	23.22	252	14190	0.394	ng/ul#	88
23) Benzo(a)pyrene	23.83	252	12774	0.394	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.60	276	13018	0.393	ng/ul#	84
25) Dibenzo(a,h)anthracene	26.63	278	10930	0.393	ng/ul	93
26) Benzo(g,h,i)perylene	27.42	276	11291	0.395	ng/ul	98

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

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