

Data Path : Z:\HPCHEM1\BNA E\DATA\BE101017\
 Data File : BE094256.D
 Acq On : 10 Oct 2017 18:12
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.426

Quant Time: Oct 11 06:38:12 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 Oct 11 06:34:16 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	2121	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	10741	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	5679	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12254	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	12706	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	12391	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	6687	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	14597	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	10369	0.394	ng/ul#	92
5) 2-Methylnaphthalene	12.35	142	7110	0.417	ng/ul	98
7) Acenaphthylene	14.24	152	10363	0.468	ng/ul	96
8) Acenaphthene	14.57	153	7384	0.405	ng/ul	97
9) Fluorene	15.56	166	8317	0.390	ng/ul#	90
11) Pentachlorophenol	16.95	266	388	0.270	ng/ul	96
12) Phenanthrene	17.29	178	12978	0.405	ng/ul#	93
13) Anthracene	17.38	178	12990	0.417	ng/ul#	92
15) Fluoranthene	19.30	202	14693	0.349	ng/ul	85
17) Pyrene	19.66	202	15529	0.439	ng/ul#	82
18) Benzo(a)anthracene	21.39	228	15825	0.469	ng/ul#	87
19) Chrysene	21.45	228	14578	0.375	ng/ul	98
21) Benzo(b)fluoranthene	23.15	252	15120	0.449	ng/ul	90
22) Benzo(k)fluoranthene	23.19	252	14661	0.377	ng/ul	92
23) Benzo(a)pyrene	23.81	252	14052	0.402	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	26.58	276	15347	0.429	ng/ul#	84
25) Dibenzo(a,h)anthracene	26.59	278	12812	0.427	ng/ul	93
26) Benzo(g,h,i)perylene	27.40	276	12547	0.407	ng/ul	98

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

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