

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE121817\
 Data File : BE095026.D
 Acq On : 19 Dec 2017 00:46
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.409

Quant Time: Dec 19 04:23:11 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE121817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 19 04:08:16 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	4056	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	13086	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.05	164	7519	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.75	188	18244	0.40	ng/ul	0.00
16) Chrysene-d12	21.84	240	18279	0.40	ng/ul	0.00
20) Perylene-d12	24.51	264	16054	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	9170	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.73	212	23846	0.42	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.34	128	14329	0.412	ng/ul#	88
5) 2-Methylnaphthalene	12.91	142	10226	0.380	ng/ul	99
7) Acenaphthylene	14.77	152	15118	0.440	ng/ul	95
8) Acenaphthene	15.11	153	11189	0.424	ng/ul	96
9) Fluorene	16.07	166	12231	0.417	ng/ul	91
11) Pentachlorophenol	17.41	266	1000	0.540	ng/ul#	60
12) Phenanthrene	17.78	178	19741	0.411	ng/ul#	81
13) Anthracene	17.87	178	18660	0.409	ng/ul#	87
15) Fluoranthene	19.76	202	26017	0.415	ng/ul	84
17) Pyrene	20.11	202	26379	0.415	ng/ul#	78
18) Benzo(a)anthracene	21.82	228	23326	0.444	ng/ul#	84
19) Chrysene	21.89	228	22786	0.409	ng/ul	98
21) Benzo(b)fluoranthene	23.68	252	21903	0.391	ng/ul	84
22) Benzo(k)fluoranthene	23.73	252	21152	0.414	ng/ul#	87
23) Benzo(a)pyrene	24.38	252	20533	0.426	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.35	276	21952	0.415	ng/ul#	81
25) Dibenzo(a,h)anthracene	27.37	278	18049	0.422	ng/ul#	88
26) Benzo(g,h,i)perylene	28.23	276	18706	0.408	ng/ul	94

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

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