

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081417\
 Data File : BE093800.D
 Acq On : 14 Aug 2017 23:11
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.405

Quant Time: Aug 15 11:53:03 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 15 10:51:12 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1945	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	10424	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	6078	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	13855	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	15514	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	13892	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	6561	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	16658	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.76	128	10124	0.399	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	7043	0.381	ng/ul	99
7) Acenaphthylene	14.24	152	10937	0.430	ng/ul	98
8) Acenaphthene	14.59	153	7850	0.392	ng/ul	97
9) Fluorene	15.57	166	9011	0.356	ng/ul#	95
11) Pentachlorophenol	16.91	266	1344	0.588	ng/ul	97
12) Phenanthrene	17.29	178	14794	0.392	ng/ul#	89
13) Anthracene	17.38	178	14913	0.406	ng/ul#	88
15) Fluoranthene	19.30	202	17361	0.368	ng/ul	84
17) Pyrene	19.66	202	18683	0.413	ng/ul#	83
18) Benzo(a)anthracene	21.39	228	17610	0.392	ng/ul#	86
19) Chrysene	21.45	228	16741	0.388	ng/ul	96
21) Benzo(b)fluoranthene	23.14	252	15835	0.368	ng/ul	94
22) Benzo(k)fluoranthene	23.19	252	16561	0.391	ng/ul	95
23) Benzo(a)pyrene	23.80	252	16007	0.392	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	26.56	276	16579	0.367	ng/ul#	86
25) Dibenzo(a,h)anthracene	26.58	278	14194	0.376	ng/ul	93
26) Benzo(g,h,i)perylene	27.38	276	13572	0.357	ng/ul	95

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

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