

Data Path : Z:\HPCHEM1\BNA_E\Data\BE081517\
 Data File : BE093802.D
 Acq On : 15 Aug 2017 11:29
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.406

Quant Time: Aug 16 05:31:04 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	2071	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	10920	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	5991	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	13785	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	15837	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	13670	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	6763	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	16097	0.37	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.75	128	10494	0.395	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	7160	0.370	ng/ul	99
7) Acenaphthylene	14.24	152	10307	0.411	ng/ul	97
8) Acenaphthene	14.59	153	7707	0.391	ng/ul	98
9) Fluorene	15.57	166	8867	0.356	ng/ul#	94
11) Pentachlorophenol	16.93	266	944	0.415	ng/ul	96
12) Phenanthrene	17.29	178	14538	0.387	ng/ul#	88
13) Anthracene	17.38	178	14236	0.389	ng/ul#	87
15) Fluoranthene	19.30	202	17038	0.363	ng/ul	85
17) Pyrene	19.66	202	17949	0.388	ng/ul#	84
18) Benzo(a)anthracene	21.39	228	17607	0.383	ng/ul#	85
19) Chrysene	21.45	228	17444	0.396	ng/ul	97
21) Benzo(b)fluoranthene	23.14	252	15706	0.371	ng/ul	88
22) Benzo(k)fluoranthene	23.20	252	16833	0.404	ng/ul#	89
23) Benzo(a)pyrene	23.81	252	15844	0.394	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	26.57	276	15657	0.353	ng/ul#	86
25) Dibenzo(a,h)anthracene	26.58	278	13482	0.363	ng/ul#	93
26) Benzo(g,h,i)perylene	27.38	276	12321	0.329	ng/ul	98

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

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