

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081617\
 Data File : BE093846.D
 Acq On : 16 Aug 2017 18:48
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.410

Quant Time: Aug 17 01:00:38 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 17 00:58:19 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2074	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	10912	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	6012	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	13519	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	14682	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	13212	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	6824	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	15473	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.75	128	10472	0.395	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	7248	0.374	ng/ul	98
7) Acenaphthylene	14.24	152	10486	0.417	ng/ul	96
8) Acenaphthene	14.58	153	7683	0.388	ng/ul	99
9) Fluorene	15.57	166	8813	0.352	ng/ul#	91
11) Pentachlorophenol	16.93	266	795	0.357	ng/ul	95
12) Phenanthrene	17.29	178	13924	0.378	ng/ul#	89
13) Anthracene	17.38	178	13954	0.389	ng/ul#	89
15) Fluoranthene	19.30	202	16440	0.357	ng/ul	84
17) Pyrene	19.66	202	17131	0.400	ng/ul#	84
18) Benzo(a)anthracene	21.40	228	16410	0.386	ng/ul#	86
19) Chrysene	21.45	228	16337	0.400	ng/ul	98
21) Benzo(b)fluoranthene	23.14	252	14875	0.364	ng/ul	88
22) Benzo(k)fluoranthene	23.19	252	15941	0.396	ng/ul#	90
23) Benzo(a)pyrene	23.80	252	15206	0.391	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	26.56	276	15860	0.370	ng/ul#	86
25) Dibenzo(a,h)anthracene	26.58	278	13142	0.366	ng/ul	93
26) Benzo(g,h,i)perylene	27.38	276	13164	0.364	ng/ul	99

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

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