

Data Path : Z:\HPCHEM1\BNA_E\Data\BE081117\
 Data File : BE093781.D
 Acq On : 11 Aug 2017 22:06
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.402

Quant Time: Aug 12 02:03:19 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 12 02:01:53 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2262	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	11899	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	6947	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	16149	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	16802	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	15507	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	7731	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	17910	0.35	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	11820	0.408	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	8324	0.394	ng/ul	99
7) Acenaphthylene	14.24	152	11933	0.411	ng/ul	97
8) Acenaphthene	14.59	153	9065	0.397	ng/ul	99
9) Fluorene	15.57	166	10587	0.366	ng/ul#	95
11) Pentachlorophenol	16.93	266	830	0.312	ng/ul	97
12) Phenanthrene	17.29	178	17359	0.395	ng/ul#	89
13) Anthracene	17.38	178	16295	0.380	ng/ul#	89
15) Fluoranthene	19.30	202	19093	0.347	ng/ul	84
17) Pyrene	19.66	202	19962	0.407	ng/ul#	81
18) Benzo(a)anthracene	21.39	228	19040	0.391	ng/ul#	86
19) Chrysene	21.45	228	18765	0.402	ng/ul	98
21) Benzo(b)fluoranthene	23.14	252	18334	0.382	ng/ul	87
22) Benzo(k)fluoranthene	23.19	252	18466	0.390	ng/ul#	87
23) Benzo(a)pyrene	23.80	252	18188	0.399	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	26.55	276	19240	0.382	ng/ul#	88
25) Dibenzo(a,h)anthracene	26.57	278	16159	0.384	ng/ul#	91
26) Benzo(g,h,i)perylene	27.36	276	15533	0.366	ng/ul	96

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

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