

Data Path : Z:\HPCHEM1\BNA_E\Data\BE082917\
 Data File : BE093895.D
 Acq On : 28 Aug 2017 13:42
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.414

Quant Time: Aug 28 15:01:08 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 19:06:13 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2082	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	10681	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5923	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12954	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	13377	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	11986	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	6754	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	14474	0.35	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.75	128	10478	0.403	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	7041	0.372	ng/ul	99
7) Acenaphthylene	14.24	152	10033	0.405	ng/ul	96
8) Acenaphthene	14.59	153	7657	0.393	ng/ul	98
9) Fluorene	15.57	166	8598	0.349	ng/ul#	95
11) Pentachlorophenol	16.93	266	574	0.269	ng/ul	96
12) Phenanthrene	17.29	178	13808	0.391	ng/ul#	88
13) Anthracene	17.38	178	13303	0.387	ng/ul#	87
15) Fluoranthene	19.30	202	15494	0.351	ng/ul	84
17) Pyrene	19.66	202	16806	0.430	ng/ul#	84
18) Benzo(a)anthracene	21.40	228	14380	0.371	ng/ul#	86
19) Chrysene	21.45	228	15396	0.414	ng/ul	98
21) Benzo(b)fluoranthene	23.15	252	13352	0.360	ng/ul	86
22) Benzo(k)fluoranthene	23.20	252	14595	0.399	ng/ul#	88
23) Benzo(a)pyrene	23.81	252	13528	0.384	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	26.57	276	11133	0.286	ng/ul#	79
25) Dibenzo(a,h)anthracene	26.60	278	8720	0.268	ng/ul	93
26) Benzo(g,h,i)perylene	27.38	276	9500	0.290	ng/ul	98

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

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