

Data Path : Z:\HPCHEM1\BNA_E\Data\BE081717\
 Data File : BE093860.D
 Acq On : 17 Aug 2017 9:31
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.412

Quant Time: Aug 17 19:04:41 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 02:52:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1942	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	10237	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	5626	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12504	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	13673	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	12518	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	6359	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	14526	0.37	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.75	128	9885	0.397	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	6702	0.369	ng/ul	99
7) Acenaphthylene	14.25	152	9882	0.420	ng/ul	95
8) Acenaphthene	14.59	153	7290	0.394	ng/ul	99
9) Fluorene	15.57	166	8145	0.348	ng/ul#	93
11) Pentachlorophenol	16.93	266	744	0.361	ng/ul	97
12) Phenanthrene	17.29	178	12775	0.375	ng/ul#	87
13) Anthracene	17.38	178	13117	0.395	ng/ul#	87
15) Fluoranthene	19.31	202	15352	0.360	ng/ul	85
17) Pyrene	19.67	202	16283	0.408	ng/ul#	84
18) Benzo(a)anthracene	21.40	228	15471	0.390	ng/ul#	87
19) Chrysene	21.45	228	15236	0.401	ng/ul	99
21) Benzo(b)fluoranthene	23.15	252	14484	0.374	ng/ul	90
22) Benzo(k)fluoranthene	23.20	252	14617	0.383	ng/ul	91
23) Benzo(a)pyrene	23.81	252	14485	0.393	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	26.57	276	15101	0.371	ng/ul#	86
25) Dibenzo(a,h)anthracene	26.59	278	12594	0.371	ng/ul	95
26) Benzo(g,h,i)perylene	27.38	276	12746	0.372	ng/ul	96

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

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