

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080817\
 Data File : BE093738.D
 Acq On : 9 Aug 2017 3:56
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.446

Manual Integrations
 APPROVED

Sohil
 8/9/2017 7:15:12 PM

Quant Time: Aug 09 07:07:08 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 09 05:31:57 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2482	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	12950	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	7465	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	17282	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	18062	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	16273m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	8358	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	19239	0.35	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	12904	0.410	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	9100	0.396	ng/ul	98
7) Acenaphthylene	14.24	152	13036	0.417	ng/ul	95
8) Acenaphthene	14.59	153	9795	0.399	ng/ul	98
9) Fluorene	15.57	166	11454	0.369	ng/ul#	95
11) Pentachlorophenol	16.91	266	1040	0.365	ng/ul	99
12) Phenanthrene	17.29	178	18463	0.392	ng/ul#	89
13) Anthracene	17.38	178	17653	0.385	ng/ul#	89
15) Fluoranthene	19.30	202	20577	0.350	ng/ul	84
17) Pyrene	19.66	202	21303	0.404	ng/ul#	82
18) Benzo(a)anthracene	21.39	228	20315	0.388	ng/ul#	86
19) Chrysene	21.45	228	19874	0.396	ng/ul	97
21) Benzo(b)fluoranthene	23.14	252	19135	0.380	ng/ul	87
22) Benzo(k)fluoranthene	23.19	252	19679m	0.396	ng/ul	
23) Benzo(a)pyrene	23.80	252	19005m	0.397	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.54	276	20328m	0.385	ng/ul	
25) Dibenzo(a,h)anthracene	26.56	278	16953m	0.384	ng/ul	
26) Benzo(g,h,i)perylene	27.36	276	16188m	0.363	ng/ul	

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

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