

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081017\
 Data File : BE093752.D
 Acq On : 10 Aug 2017 18:21
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.448

Manual Integrations
 APPROVED

sohil
 8/14/2017 7:03:00 PM

Quant Time: Aug 11 18:17:32 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 11 06:07:04 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2377	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	12429	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	7345	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	17054	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	18530	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	16640m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	8133	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	19684	0.37	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.75	128	12373	0.409	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	8746	0.397	ng/ul	99
7) Acenaphthylene	14.24	152	12757	0.415	ng/ul	97
8) Acenaphthene	14.59	153	9591	0.397	ng/ul	98
9) Fluorene	15.57	166	11143	0.365	ng/ul#	91
11) Pentachlorophenol	16.91	266	992	0.353	ng/ul	96
12) Phenanthrene	17.29	178	18505	0.398	ng/ul#	89
13) Anthracene	17.38	178	17529	0.387	ng/ul#	88
15) Fluoranthene	19.30	202	20845	0.359	ng/ul	84
17) Pyrene	19.66	202	21926	0.405	ng/ul#	83
18) Benzo(a)anthracene	21.39	228	21468	0.400	ng/ul#	86
19) Chrysene	21.45	228	20724	0.402	ng/ul	98
21) Benzo(b)fluoranthene	23.14	252	19594	0.381	ng/ul	88
22) Benzo(k)fluoranthene	23.19	252	20406m	0.402	ng/ul	
23) Benzo(a)pyrene	23.80	252	19247m	0.393	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	20698m	0.383	ng/ul	
25) Dibenzo(a,h)anthracene	26.57	278	17247m	0.382	ng/ul	
26) Benzo(g,h,i)perylene	27.36	276	16838m	0.370	ng/ul	

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

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