

Data Path : Z:\HPCHEM1\BNA E\Data\BE080317\
 Data File : BE093649.D
 Acq On : 3 Aug 2017 8:53
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc : I4426-07
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Aug 03 09:45:05 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 02 10:58:31 2017
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	109	-0.01
2	1,4-Dioxane	0.400	0.404	-1.0	111	0.00
3	n-Nitrosodimethylamine	0.400	0.380	5.0	110	0.01
4 S	2-Fluorophenol	0.400	0.419	-4.7	119	0.00
5 S	Phenol-d6	0.400	0.408	-2.0	115	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.393	1.8	110	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	104	0.00
8 S	Nitrobenzene-d5	0.400	0.402	-0.5	112	0.00
9	Naphthalene	0.400	0.395	1.3	104	0.00
10	Hexachlorobutadiene	0.400	0.399	0.3	106	-0.02
11 SURR	2-Methylnaphthalene-d10	0.400	0.397	0.8	104	0.00
12	2-Methylnaphthalene	0.400	0.392	2.0	103	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	102	0.00
14 S	2,4,6-Tribromophenol	0.400	0.422	-5.5	127	0.00
15 S	2-Fluorobiphenyl	0.400	0.402	-0.5	102	-0.02
16	Acenaphthylene	0.400	0.401	-0.3	106	0.00
17	Acenaphthene	0.400	0.397	0.8	103	0.00
18	Fluorene	0.400	0.396	1.0	100	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	123	0.00
20	4-Bromophenyl-phenylether	0.400	0.332	17.0	92	0.00
21	Hexachlorobenzene	0.400	0.365	8.8	100	-0.03
22	Pentachlorophenol	0.400	0.473	-18.2	147	0.00
23	Phenanthrene	0.400	0.377	5.8	112	0.00
24	Anthracene	0.400	0.426	-6.5	123	0.00
25 SURR	Fluoranthene-d10	0.400	0.377	5.8	106	0.00
26	Fluoranthene	0.400	0.372	7.0	105	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	115	-0.01
28	Pvrene	0.400	0.369	7.8	106	0.00
29 S	Terphenyl-d14	0.400	0.371	7.3	105	0.00
30	Benzo(a)anthracene	0.400	0.410	-2.5	122	0.00
31	Chrysene	0.400	0.390	2.5	112	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.470	-17.5	157	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.426	-6.5	121	0.00
34 I	Perylene-d12	0.400	0.400	0.0	124	-0.01
35	Benzo(b)fluoranthene	0.400	0.368	8.0	117	0.00
36	Benzo(k)fluoranthene	0.400	0.378	5.5	121	0.00
37 C	Benzo(a)pyrene	0.400	0.383	4.3	123	0.00
38	Dibenzo(a,h)anthracene	0.400	0.382	4.5	120	0.00
39	Benzo(a,h,i)perylene	0.400	0.367	8.3	116	0.00

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(#) = Out of Range	SPCC's out = 0 CCC's out = 0				