

Data Path : Z:\HPCHEM1\BNA_E\Data\BE071317\
 Data File : BE093461.D
 Acq On : 13 Jul 2017 16:41
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 13 17:18:19 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE070217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 03 06:40:23 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	139	0.00
2	1,4-Dioxane	0.400	0.368	8.0	128	0.00
3	n-Nitrosodimethylamine	0.400	0.424	-6.0	183	0.00
4 S	2-Fluorophenol	0.400	0.369	7.8	132	0.00
5 S	Phenol-d6	0.400	0.396	1.0	147	0.00
6	bis(2-Chloroethyl)ether	0.400	0.395	1.3	141	-0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	156	0.00
8 S	Nitrobenzene-d5	0.400	0.440	-10.0	186	0.00
9	Naphthalene	0.400	0.384	4.0	154	0.00
10	Hexachlorobutadiene	0.400	0.381	4.8	151	-0.02
11 SURR	2-Methylnaphthalene-d10	0.400	0.398	0.5	161	0.00
12	2-Methylnaphthalene	0.400	0.397	0.8	161	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	153	0.00
14 S	2,4,6-Tribromophenol	0.400	0.269	32.8#	113	0.00
15 S	2-Fluorobiphenyl	0.400	0.396	1.0	157	0.00
16	Acenaphthylene	0.400	0.380	5.0	154	0.00
17	Acenaphthene	0.400	0.384	4.0	153	0.00
18	Fluorene	0.400	0.359	10.3	143	-0.02
19 I	Phenanthrene-d10	0.400	0.400	0.0	97	0.00
20	4-Bromophenyl-phenylether	0.400	0.583	-45.7#	153	0.00
21	Hexachlorobenzene	0.400	0.507	-26.7#	137	0.00
22	Pentachlorophenol	0.400	0.603	-50.7#	157	0.00
23	Phenanthrene	0.400	0.420	-5.0	111	0.00
24	Anthracene	0.400	0.396	1.0	102	0.00
25 SURR	Fluoranthene-d10	0.400	0.431	-7.7	114	0.00
26	Fluoranthene	0.400	0.430	-7.5	115	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	105	-0.01
28	Pyrene	0.400	0.420	-5.0	113	0.00
29 S	Terphenyl-d14	0.400	0.415	-3.7	111	0.00
30	Benzo(a)anthracene	0.400	0.385	3.8	106	0.00
31	Chrysene	0.400	0.399	0.3	106	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.371	7.3	119	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.424	-6.0	114	-0.02
34 I	Perylene-d12	0.400	0.400	0.0	110	0.00
35	Benzo(b)fluoranthene	0.400	0.383	4.3	108	0.00
36	Benzo(k)fluoranthene	0.400	0.409	-2.2	118	-0.01
37 C	Benzo(a)pyrene	0.400	0.388	3.0	113	0.00
38	Dibenzo(a,h)anthracene	0.400	0.398	0.5	114	-0.01
39	Benzo(g,h,i)perylene	0.400	0.387	3.3	112	0.00

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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