

Data Path : Z:\HPCHEM1\BNA_E\Data\BE070717\
 Data File : BE093448.D
 Acq On : 7 Jul 2017 18:12
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 08 05:35:15 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	126	0.00
2	1,4-Dioxane	0.400	0.343	14.2	108	0.00
3	n-Nitrosodimethylamine	0.400	0.397	0.8	151	0.00
4 S	2-Fluorophenol	0.400	0.383	4.3	124	0.00
5 S	Phenol-d6	0.400	0.405	-1.3	136	0.00
6	bis(2-Chloroethyl)ether	0.400	0.391	2.3	126	-0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	140	0.00
8 S	Nitrobenzene-d5	0.400	0.447	-11.7	169	0.00
9	Naphthalene	0.400	0.388	3.0	140	0.00
10	Hexachlorobutadiene	0.400	0.373	6.8	133	-0.02
11 SURR	2-Methylnaphthalene-d10	0.400	0.409	-2.2	148	0.00
12	2-Methylnaphthalene	0.400	0.406	-1.5	148	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	141	0.00
14 S	2,4,6-Tribromophenol	0.400	0.269	32.8#	105	0.00
15 S	2-Fluorobiphenyl	0.400	0.397	0.8	145	0.00
16	Acenaphthylene	0.400	0.388	3.0	145	0.00
17	Acenaphthene	0.400	0.392	2.0	144	0.00
18	Fluorene	0.400	0.361	9.8	132	-0.02
19 I	Phenanthrene-d10	0.400	0.400	0.0	87	0.00
20	4-Bromophenyl-phenylether	0.400	0.606	-51.5#	142	0.00
21	Hexachlorobenzene	0.400	0.522	-30.5#	126	0.00
22	Pentachlorophenol	0.400	0.698	-74.5#	164	0.00
23	Phenanthrene	0.400	0.435	-8.7	103	0.00
24	Anthracene	0.400	0.398	0.5	91	0.00
25 SURR	Fluoranthene-d10	0.400	0.455	-13.7	108	0.00
26	Fluoranthene	0.400	0.455	-13.7	108	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	101	0.00
28	Pyrene	0.400	0.412	-3.0	106	0.00
29 S	Terphenyl-d14	0.400	0.404	-1.0	104	0.00
30	Benzo(a)anthracene	0.400	0.382	4.5	101	0.00
31	Chrysene	0.400	0.395	1.3	101	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.389	2.8	120	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.417	-4.2	108	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	106	0.00
35	Benzo(b)fluoranthene	0.400	0.376	6.0	102	0.00
36	Benzo(k)fluoranthene	0.400	0.387	3.3	108	0.00
37 C	Benzo(a)pyrene	0.400	0.386	3.5	108	0.00
38	Dibenzo(a,h)anthracene	0.400	0.390	2.5	107	-0.01
39	Benzo(g,h,i)perylene	0.400	0.382	4.5	106	0.00

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

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