Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080315\
Data File : BE090335.D
Acq On : 4 Aug 2015 7:37
Operator : TP/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 04 08:13:38 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 04 02:31:31 2015
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	5.000	5.000	0.0	46	0.00
2		1,4-Dioxane	1.000	1.207	-20.7	46	0.00
3		n-Nitrosodimethylamine	1.000	1.058	-5.8	48	0.00
4	S	2-Fluorophenol	1.000	1.058	-5.8	50	0.00
5	S	Phenol-d6	1.000	0.893	10.7	44	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	45	0.00
7	S	Nitrobenzene-d5	1.000	0.974	2.6	46	0.00
8		Nitrobenzene	1.000	0.895	10.5	43	0.00
9		Naphthalene	1.000	1.002	-0.2	46	0.00
10		Hexachlorobutadiene	1.000	0.992	0.8	44	0.00
11		2-Methylnaphthalene	1.000	0.879	12.1	40	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	38	0.00
13	S	2,4,6-Tribromophenol	1.000	0.942	5.8	42	0.00
14	S	2-Fluorobiphenyl	1.000	1.066	-6.6	40	0.00
15		Acenaphthylene	1.000	1.021	-2.1	39	0.00
16		Acenaphthene	1.000	0.981	1.9	37	0.00
17		Fluorene	1.000	0.942	5.8	38	0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	34	0.02
19		4-Bromophenyl-phenylether	1.000	1.140	-14.0	37	0.00
20		Hexachlorobenzene	1.000	1.163	-16.3	38	0.00
21		Pentachlorophenol	1.000	7.459	-645.9#	613	0.00
22		Phenanthrene	1.000	0.977	2.3	34	0.00
23		Anthracene	1.000	0.967	3.3	35	0.00
24		Fluoranthene	1.000	0.952	4.8	33	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	30	0.00
26		Pyrene	1.000	1.176	-17.6	34	0.00
27	S	Terphenyl-d14	1.000	1.134	-13.4	33	0.00
28		Benzo(a)anthracene	1.000	0.930	7.0	31	0.00
29		Chrysene	1.000	1.329	-32.9#	34	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.997	0.3	36	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.340	-34.0#	54	0.00
32	I	Perylene-d12	5.000	5.000	0.0	30	0.00
33		Benzo(b)fluoranthene	1.000	1.091	-9.1	39	0.00
34		Benzo(k)fluoranthene	1.000	1.191	-19.1	33	0.00
35	C	Benzo(a)pyrene	1.000	1.210	-21.0#	43	0.00
36		Dibenzo(a,h)anthracene	1.000	1.438	-43.8#	61	0.00
37		Benzo(g,h,i)perylene	1.000	1.331	-33.1#	44	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1