Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE091515\
Data File : BE090828.D
Acq On : 15 Sep 2015 10:56
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 15 13:48:03 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 02 11:40:50 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	52	0.00
2		1,4-Dioxane	1.000	1.031	-3.1	50	0.00
3		n-Nitrosodimethylamine	1.000	0.807	19.3	41	0.00
4	S	2-Fluorophenol	1.000	1.020	-2.0	52	0.00
5	S	Phenol-d6	1.000	0.820	18.0	45	0.03
6	I	Naphthalene-d8	5.000	5.000	0.0	54	0.00
7	S	Nitrobenzene-d5	1.000	0.800	20.0	43	0.00
8		Nitrobenzene	1.000	0.799	20.1	44	0.00
9		Naphthalene	1.000	0.949	5.1	51	0.00
10		Hexachlorobutadiene	1.000	0.980	2.0	51	0.00
11		2-Methylnaphthalene	1.000	0.923	7.7	51	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	56	0.00
13	S	2,4,6-Tribromophenol	1.000	0.801	19.9	45	0.02
14	S	2-Fluorobiphenyl	1.000	0.896	10.4	51	0.00
15		Acenaphthylene	1.000	0.919	8.1	53	0.00
16		Acenaphthene	1.000	0.987	1.3	55	0.00
17		Fluorene	1.000	0.938	6.2	53	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	61	0.00
19		4-Bromophenyl-phenylether	1.000	0.851	14.9	54	0.00
20		Hexachlorobenzene	1.000	0.946	5.4	57	0.00
21		Pentachlorophenol	1.000	0.853	14.7	37	0.00
22		Phenanthrene	1.000	0.941	5.9	57	-0.02
23		Anthracene	1.000	0.880	12.0	56	0.00
24		Fluoranthene	1.000	0.870	13.0	57	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	68	0.00
26		Pyrene	1.000	0.830	17.0	60	0.00
27	S	Terphenyl-d14	1.000	0.800	20.0	57	0.00
28		Benzo(a)anthracene	1.000	0.882	11.8	61	0.00
29		Chrysene	1.000	0.998	0.2	66	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.643	35.7#	46	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.910	9.0	67	0.00
32	I	Perylene-d12	5.000	5.000	0.0	70	-0.01
33		Benzo(b)fluoranthene	1.000	0.892	10.8	64	0.00
34		Benzo(k)fluoranthene	1.000	0.920	8.0	65	0.00
35	C	Benzo(a)pyrene	1.000	0.879	12.1	66	0.00
36		Dibenzo(a,h)anthracene	1.000	0.885	11.5	65	-0.01
37		Benzo(g,h,i)perylene	1.000	0.885	11.5	65	-0.01

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