Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE091415\
Data File : BE090810.D
Acq On : 14 Sep 2015 22:06
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 15 00:58:53 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	115	0.00
2		1,4-Dioxane	1.000	0.820	18.0	92	0.00
3		n-Nitrosodimethylamine	1.000	0.832	16.8	94	0.00
4	S	2-Fluorophenol	1.000	0.773	22.7	87	0.00
5	S	Phenol-d6	1.000	0.786	21.4	95	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	124	0.00
7	S	Nitrobenzene-d5	1.000	0.832	16.8	104	0.00
8		Nitrobenzene	1.000	0.858	14.2	110	0.00
9		Naphthalene	1.000	0.873	12.7	109	-0.01
10		Hexachlorobutadiene	1.000	0.889	11.1	107	0.00
11		2-Methylnaphthalene	1.000	0.886	11.4	114	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	137	0.00
13	S	2,4,6-Tribromophenol	1.000	0.711	28.9#	94	0.00
14	S	2-Fluorobiphenyl	1.000	0.832	16.8	115	0.00
15		Acenaphthylene	1.000	0.827	17.3	118	0.00
16		Acenaphthene	1.000	0.860	14.0	116	0.00
17		Fluorene	1.000	0.838	16.2	116	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	129	0.00
19		4-Bromophenyl-phenylether	1.000	0.850	15.0	114	-0.02
20		Hexachlorobenzene	1.000	0.864	13.6	111	0.00
21		Pentachlorophenol	1.000	0.916	8.4	92	0.00
22		Phenanthrene	1.000	0.889	11.1	115	-0.02
23		Anthracene	1.000	0.876	12.4	118	0.00
24		Fluoranthene	1.000	0.849	15.1	118	-0.01
25	I	Chrysene-d12	5.000	5.000	0.0	135	0.00
26		Pyrene	1.000	0.977	2.3	139	-0.01
27	S	Terphenyl-d14	1.000	0.898	10.2	127	-0.01
28		Benzo(a)anthracene	1.000	0.931	6.9	127	0.00
29		Chrysene	1.000	0.901	9.9	118	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.862	13.8	128	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.005	-0.5	147	-0.01
32	I	Perylene-d12	5.000	5.000	0.0	147	-0.01
33		Benzo(b)fluoranthene	1.000	0.916	8.4	139	0.00
34		Benzo(k)fluoranthene	1.000	0.882	11.8	132	0.00
35	C	Benzo(a)pyrene	1.000	0.909	9.1	143	0.00
36		Dibenzo(a,h)anthracene	1.000	0.927	7.3	144	-0.02
37		Benzo(g,h,i)perylene	1.000	0.859	14.1	134	-0.02

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