Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE091615\
Data File : BE090842.D
Acq On : 16 Sep 2015 11:32
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 17 00:54:42 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	110	0.00
2		1,4-Dioxane	1.000	0.839	16.1	90	0.00
3		n-Nitrosodimethylamine	1.000	0.839	16.1	91	0.00
4	S	2-Fluorophenol	1.000	0.792	20.8	85	0.00
5	S	Phenol-d6	1.000	0.904	9.6	105	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	126	0.00
7	S	Nitrobenzene-d5	1.000	0.859	14.1	108	0.00
8		Nitrobenzene	1.000	0.882	11.8	115	0.00
9		Naphthalene	1.000	0.875	12.5	110	-0.01
10		Hexachlorobutadiene	1.000	0.845	15.5	103	0.00
11		2-Methylnaphthalene	1.000	0.901	9.9	117	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	138	0.00
13	S	2,4,6-Tribromophenol	1.000	0.750	25.0	102	0.00
14	S	2-Fluorobiphenyl	1.000	0.817	18.3	115	-0.01
15		Acenaphthylene	1.000	0.844	15.6	122	0.00
16		Acenaphthene	1.000	0.864	13.6	118	0.00
17		Fluorene	1.000	0.852	14.8	119	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	133	0.00
19		4-Bromophenyl-phenylether	1.000	0.790	21.0	109	-0.02
20		Hexachlorobenzene	1.000	0.841	15.9	111	0.00
21		Pentachlorophenol	1.000	0.949	5.1	101	0.00
22		Phenanthrene	1.000	0.889	11.1	118	-0.02
23		Anthracene	1.000	0.881	11.9	122	0.00
24		Fluoranthene	1.000	0.870	13.0	124	-0.01
25	I	Chrysene-d12	5.000	5.000	0.0	145	0.00
26		Pyrene	1.000	0.966	3.4	148	-0.01
27	S	Terphenyl-d14	1.000	0.865	13.5	131	-0.01
28		Benzo(a)anthracene	1.000	0.927	7.3	136	0.00
29		Chrysene	1.000	0.919	8.1	130	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.853	14.7	136	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.028	-2.8	161	-0.02
32	I	Perylene-d12	5.000	5.000	0.0	160	-0.01
33		Benzo(b)fluoranthene	1.000	0.891	10.9	147	-0.01
34		Benzo(k)fluoranthene	1.000	0.923	7.7	150	-0.01
35	C	Benzo(a)pyrene	1.000	0.939	6.1	161	-0.01
36		Dibenzo(a,h)anthracene	1.000	0.937	6.3	159	-0.02
37		Benzo(g,h,i)perylene	1.000	0.859	14.1	146	-0.03

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