Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE092315\
Data File : BE090950.D
Acq On : 23 Sep 2015 15:23
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 24 03:58:09 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE091815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 18 18:17:34 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	113	0.00
2		1,4-Dioxane	1.000	1.223	-22.3	123	0.00
3		n-Nitrosodimethylamine	1.000	0.951	4.9	101	0.00
4	S	2-Fluorophenol	1.000	0.878	12.2	95	0.00
5	S	Phenol-d6	1.000	0.847	15.3	98	0.04
6		bis(2-Chloroethyl)ether	1.000	1.235	-23.5	141	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	115	0.00
8	S	Nitrobenzene-d5	1.000	0.799	20.1	93	0.01
9		Nitrobenzene	1.000	0.770	23.0	92	0.01
10		Naphthalene	1.000	1.047	-4.7	114	0.00
11		Hexachlorobutadiene	1.000	1.183	-18.3	127	0.00
12		2-Methylnaphthalene	1.000	0.978	2.2	115	0.02
13	I	Acenaphthene-d10	5.000	5.000	0.0	124	0.00
14	S	2,4,6-Tribromophenol	1.000	0.837	16.3	107	0.00
15	S	2-Fluorobiphenyl	1.000	0.974	2.6	120	0.00
16		Acenaphthylene	1.000	0.976	2.4	122	0.00
17		Acenaphthene	1.000	1.009	-0.9	127	0.00
18		Fluorene	1.000	1.009	-0.9	124	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	120	0.00
20		4-Bromophenyl-phenylether	1.000	0.985	1.5	120	0.00
21		Hexachlorobenzene	1.000	1.144	-14.4	129	0.00
22		Pentachlorophenol	1.000	0.863	13.7	99	0.04
23		Phenanthrene	1.000	0.992	0.8	115	0.00
24		Anthracene	1.000	1.023	-2.3	127	0.00
25		Fluoranthene	1.000	0.941	5.9	114	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	103	0.00
27		Pyrene	1.000	1.023	-2.3	113	0.00
28	S	Terphenyl-d14	1.000	1.066	-6.6	112	0.01
29		Benzo(a)anthracene	1.000	0.857	14.3	92	0.00
30		Chrysene	1.000	1.246	-24.6	119	0.00
31		Bis(2-ethylhexyl)phthalate	1.000	0.918	8.2	100	0.00
32		Indeno(1,2,3-cd)pyrene	1.000	0.855	14.5	87	0.00
33	I	Perylene-d12	5.000	5.000	0.0	93	0.00
34		Benzo(b)fluoranthene	1.000	0.892	10.8	86	0.00
35		Benzo(k)fluoranthene	1.000	1.223	-22.3	115	0.00
36	C	Benzo(a)pyrene	1.000	1.064	-6.4	103	0.00
37		Dibenzo(a,h)anthracene	1.000	0.900	10.0	85	0.00
38		Benzo(g,h,i)perylene	1.000	0.938	6.2	89	0.00

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