Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070115\
Data File : BE090042.D
Acq On : 1 Jul 2015 12:45
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 02 06:34:21 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 30 07:01:01 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	100	-0.02
2		1,4-Dioxane	1.000	0.858	14.2	89	-0.01
3		n-Nitrosodimethylamine	1.000	0.817	18.3	85	-0.02
4	S	2-Fluorophenol	1.000	0.905	9.5	91	-0.01
5	S	Phenol-d6	1.000	0.916	8.4	93	-0.01
6	I	Naphthalene-d8	5.000	5.000	0.0	100	-0.01
7	S	Nitrobenzene-d5	1.000	0.854	14.6	87	-0.01
8		Nitrobenzene	1.000	0.858	14.2	88	-0.01
9		Naphthalene	1.000	0.921	7.9	93	-0.01
10		Hexachlorobutadiene	1.000	0.950	5.0	96	-0.01
11		2-Methylnaphthalene	1.000	0.893	10.7	91	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	94	-0.01
13	S	2,4,6-Tribromophenol	1.000	0.777	22.3	75	0.00
14	S	2-Fluorobiphenyl	1.000	0.905	9.5	87	-0.01
15		Acenaphthylene	1.000	0.922	7.8	90	0.00
16		Acenaphthene	1.000	0.945	5.5	90	-0.01
17		Fluorene	1.000	0.927	7.3	90	-0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	89	-0.02
19		4-Bromophenyl-phenylether	1.000	0.895	10.5	81	-0.02
20		Hexachlorobenzene	1.000	0.915	8.5	84	-0.02
21		Pentachlorophenol	1.000	0.721	27.9#	59	0.00
22		Phenanthrene	1.000	0.888	11.2	81	-0.02
23		Anthracene	1.000	0.918	8.2	83	0.00
24		Fluoranthene	1.000	0.901	9.9	82	-0.01
25	I	Chrysene-d12	5.000	5.000	0.0	87	-0.01
26		Pyrene	1.000	0.928	7.2	82	-0.01
27	S	Terphenyl-d14	1.000	0.890	11.0	77	-0.01
28		Benzo(a)anthracene	1.000	0.913	8.7	81	-0.01
29		Chrysene	1.000	0.897	10.3	79	-0.01
30		Bis(2-ethylhexyl)phthalate	1.000	0.884	11.6	84	-0.01
31		Indeno(1,2,3-cd)pyrene	1.000	0.951	4.9	84	-0.03
32	I	Perylene-d12	5.000	5.000	0.0	84	-0.02
33		Benzo(b)fluoranthene	1.000	1.028	-2.8	87	-0.01
34		Benzo(k)fluoranthene	1.000	0.989	1.1	83	-0.01
35	C	Benzo(a)pyrene	1.000	1.012	-1.2	85	-0.02
36		Dibenzo(a,h)anthracene	1.000	0.972	2.8	83	-0.03
37		Benzo(g,h,i)perylene	1.000	0.969	3.1	82	-0.03

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