Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062615\
Data File : BE089955.D
Acq On : 27 Jun 2015 00:24
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jun 27 01:48:05 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 25 16:51:27 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	85	0.00
2		1,4-Dioxane	1.000	0.948	5.2	85	0.00
3		n-Nitrosodimethylamine	1.000	0.856	14.4	80	0.00
4	S	2-Fluorophenol	1.000	0.853	14.7	77	0.00
5	S	Phenol-d6	1.000	0.896	10.4	82	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	83	0.00
7	S	Nitrobenzene-d5	1.000	0.854	14.6	79	0.00
8		Nitrobenzene	1.000	0.841	15.9	79	0.00
9		Naphthalene	1.000	0.928	7.2	83	0.00
10		Hexachlorobutadiene	1.000	0.944	5.6	85	0.00
11		2-Methylnaphthalene	1.000	0.899	10.1	82	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	83	0.00
13	S	2,4,6-Tribromophenol	1.000	0.909	9.1	85	0.00
14	S	2-Fluorobiphenyl	1.000	0.901	9.9	83	0.00
15		Acenaphthylene	1.000	0.925	7.5	83	0.00
16		Acenaphthene	1.000	0.942	5.8	83	0.00
17		Fluorene	1.000	1.015	-1.5	84	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	81	0.00
19		4-Bromophenyl-phenylether	1.000	0.895	10.5	83	-0.02
20		Hexachlorobenzene	1.000	0.944	5.6	85	0.00
21		Pentachlorophenol	1.000	2.648	-164.8#	364	-0.02
22		Phenanthrene	1.000	0.926	7.4	84	-0.02
23		Anthracene	1.000	0.905	9.5	81	0.00
24		Fluoranthene	1.000	0.935	6.5	84	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	89	0.00
26		Pyrene	1.000	0.898	10.2	84	0.00
27	S	Terphenyl-d14	1.000	0.865	13.5	85	0.00
28		Benzo(a)anthracene	1.000	0.913	8.7	89	0.00
29		Chrysene	1.000	0.884	11.6	88	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.947	5.3	100	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.941	5.9	89	0.00
32	I	Perylene-d12	5.000	5.000	0.0	88	0.00
33		Benzo(b)fluoranthene	1.000	0.939	6.1	87	0.00
34		Benzo(k)fluoranthene	1.000	0.965	3.5	90	0.00
35	C	Benzo(a)pyrene	1.000	0.923	7.7	87	0.00
36		Dibenzo(a,h)anthracene	1.000	0.952	4.8	90	0.00
37		Benzo(g,h,i)perylene	1.000	0.935	6.5	88	0.00

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