Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE091715\
Data File : BE090874.D
Acq On : 17 Sep 2015 20:56
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 18 02:29:53 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	68	0.00
2		1,4-Dioxane	1.000	0.949	5.1	61	0.00
3		n-Nitrosodimethylamine	1.000	0.827	17.3	55	0.00
4	S	2-Fluorophenol	1.000	0.959	4.1	64	0.00
5	S	Phenol-d6	1.000	0.938	6.2	67	0.01
6	I	Naphthalene-d8	5.000	5.000	0.0	78	0.00
7	S	Nitrobenzene-d5	1.000	0.885	11.5	69	0.00
8		Nitrobenzene	1.000	0.860	14.0	69	0.00
9		Naphthalene	1.000	0.946	5.4	74	-0.01
10		Hexachlorobutadiene	1.000	0.913	8.7	69	0.00
11		2-Methylnaphthalene	1.000	1.012	-1.2	82	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	86	0.00
13	S	2,4,6-Tribromophenol	1.000	0.841	15.9	73	0.00
14	S	2-Fluorobiphenyl	1.000	0.903	9.7	79	0.00
15		Acenaphthylene	1.000	0.947	5.3	85	0.00
16		Acenaphthene	1.000	0.963	3.7	82	0.00
17		Fluorene	1.000	0.913	8.7	79	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	82	0.00
19		4-Bromophenyl-phenylether	1.000	0.936	6.4	80	0.00
20		Hexachlorobenzene	1.000	0.983	1.7	79	0.00
21		Pentachlorophenol	1.000	0.918	8.2	59	0.02
22		Phenanthrene	1.000	0.965	3.5	80	-0.02
23		Anthracene	1.000	0.927	7.3	79	0.00
24		Fluoranthene	1.000	0.881	11.9	78	-0.01
25	I	Chrysene-d12	5.000	5.000	0.0	74	0.00
26		Pyrene	1.000	1.013	-1.3	80	0.00
27	S	Terphenyl-d14	1.000	0.989	1.1	77	0.00
28		Benzo(a)anthracene	1.000	0.962	3.8	73	0.00
29		Chrysene	1.000	1.000	0.0	72	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.830	17.0	68	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.847	15.3	68	0.00
32	I	Perylene-d12	5.000	5.000	0.0	73	-0.01
33		Benzo(b)fluoranthene	1.000	0.940	6.0	71	0.00
34		Benzo(k)fluoranthene	1.000	0.947	5.3	71	0.00
35	C	Benzo(a)pyrene	1.000	0.921	7.9	72	-0.01
36		Dibenzo(a,h)anthracene	1.000	0.885	11.5	68	0.00
37		Benzo(g,h,i)perylene	1.000	0.884	11.6	69	-0.02

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