Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071615\
Data File : BE090137.D
Acq On : 17 Jul 2015 2:20
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 17 13:16:53 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE071615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 17 12:59:41 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	79	0.00
2		1,4-Dioxane	1.000	1.102	-10.2	85	0.00
3		n-Nitrosodimethylamine	1.000	0.905	9.5	75	0.00
4	S	2-Fluorophenol	1.000	1.026	-2.6	86	0.00
5	S	Phenol-d6	1.000	1.193	-19.3	99	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	85	0.00
7	S	Nitrobenzene-d5	1.000	1.177	-17.7	105	0.00
8		Nitrobenzene	1.000	0.849	15.1	75	0.00
9		Naphthalene	1.000	0.866	13.4	77	0.00
10		Hexachlorobutadiene	1.000	0.877	12.3	78	0.00
11		2-Methylnaphthalene	1.000	0.869	13.1	77	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	91	0.00
13	S	2,4,6-Tribromophenol	1.000	1.203	-20.3	162	0.00
14	S	2-Fluorobiphenyl	1.000	1.369	-36.9#	128	0.00
15		Acenaphthylene	1.000	0.822	17.8	78	0.00
16		Acenaphthene	1.000	0.823	17.7	78	0.00
17		Fluorene	1.000	0.805	19.5	76	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	94	0.00
19		4-Bromophenyl-phenylether	1.000	0.791	20.9	77	0.00
20		Hexachlorobenzene	1.000	0.792	20.8	75	0.00
21		Pentachlorophenol	1.000	0.962	3.8	84	0.00
22		Phenanthrene	1.000	0.814	18.6	79	0.00
23		Anthracene	1.000	0.828	17.2	80	0.00
24		Fluoranthene	1.000	0.822	17.8	79	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	94	0.00
26		Pyrene	1.000	0.811	18.9	78	0.00
27	S	Terphenyl-d14	1.000	1.686	-68.6#	158	0.00
28		Benzo(a)anthracene	1.000	0.863	13.7	87	0.00
29		Chrysene	1.000	0.846	15.4	78	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.845	15.5	84	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.929	7.1	92	0.00
32	I	Perylene-d12	5.000	5.000	0.0	95	0.00
33		Benzo(b)fluoranthene	1.000	0.989	1.1	95	0.00
34		Benzo(k)fluoranthene	1.000	0.735	26.5#	73	0.00
35	C	Benzo(a)pyrene	1.000	0.810	19.0	91	0.00
36		Dibenzo(a,h)anthracene	1.000	0.981	1.9	94	0.00
37		Benzo(g,h,i)perylene	1.000	0.825	17.5	90	0.00

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