Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070815\
Data File : BE090074.D
Acq On : 9 Jul 2015 00:34
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 09 01:21:53 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	147	0.00
2		1,4-Dioxane	1.000	0.761	23.9	116	0.00
3		n-Nitrosodimethylamine	1.000	0.811	18.9	124	0.00
4	S	2-Fluorophenol	1.000	1.016	-1.6	149	0.00
5	S	Phenol-d6	1.000	1.077	-7.7	161	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	154	-0.01
7	S	Nitrobenzene-d5	1.000	0.911	8.9	143	0.00
8		Nitrobenzene	1.000	0.885	11.5	140	0.00
9		Naphthalene	1.000	0.921	7.9	144	-0.01
10		Hexachlorobutadiene	1.000	0.938	6.2	146	0.00
11		2-Methylnaphthalene	1.000	0.890	11.0	139	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	139	0.00
13	S	2,4,6-Tribromophenol	1.000	1.083	-8.3	153	0.00
14	S	2-Fluorobiphenyl	1.000	0.923	7.7	131	0.00
15		Acenaphthylene	1.000	0.950	5.0	136	0.00
16		Acenaphthene	1.000	0.944	5.6	133	-0.01
17		Fluorene	1.000	0.892	10.8	127	-0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	123	-0.02
19		4-Bromophenyl-phenylether	1.000	0.907	9.3	115	-0.02
20		Hexachlorobenzene	1.000	0.894	10.6	115	-0.02
21		Pentachlorophenol	1.000	0.984	1.6	129	0.00
22		Phenanthrene	1.000	0.909	9.1	116	-0.02
23		Anthracene	1.000	0.928	7.2	117	0.00
24		Fluoranthene	1.000	0.915	8.5	115	-0.01
25	I	Chrysene-d12	5.000	5.000	0.0	114	-0.01
26		Pyrene	1.000	1.001	-0.1	116	0.00
27	S	Terphenyl-d14	1.000	0.960	4.0	110	0.00
28		Benzo(a)anthracene	1.000	0.959	4.1	112	-0.01
29		Chrysene	1.000	0.918	8.2	106	-0.01
30		Bis(2-ethylhexyl)phthalate	1.000	1.290	-29.0#	188	-0.01
31		Indeno(1,2,3-cd)pyrene	1.000	1.006	-0.6	117	-0.04
32	I	Perylene-d12	5.000	5.000	0.0	110	-0.02
33		Benzo(b)fluoranthene	1.000	1.074	-7.4	119	-0.02
34		Benzo(k)fluoranthene	1.000	1.020	-2.0	113	-0.02
35	C	Benzo(a)pyrene	1.000	1.081	-8.1	119	-0.02
36		Dibenzo(a,h)anthracene	1.000	1.017	-1.7	114	-0.03
37		Benzo(g,h,i)perylene	1.000	1.029	-2.9	115	-0.04

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