Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090815\
Data File : BE090747.D
Acq On : 8 Sep 2015 9:48
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 08 23:03:45 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	129	0.00
2		1,4-Dioxane	1.000	0.945	5.5	116	0.00
3		n-Nitrosodimethylamine	1.000	0.949	5.1	119	0.00
4	S	2-Fluorophenol	1.000	1.045	-4.5	133	0.00
5	S	Phenol-d6	1.000	1.135	-13.5	154	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	143	0.00
7	S	Nitrobenzene-d5	1.000	1.041	-4.1	149	0.00
8		Nitrobenzene	1.000	1.021	-2.1	155	0.00
9		Naphthalene	1.000	1.000	0.0	144	0.00
10		Hexachlorobutadiene	1.000	0.937	6.3	130	0.00
11		2-Methylnaphthalene	1.000	1.138	-13.8	168	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	161	0.00
13	S	2,4,6-Tribromophenol	1.000	1.115	-11.5	191	0.00
14	S	2-Fluorobiphenyl	1.000	0.987	1.3	161	0.00
15		Acenaphthylene	1.000	1.055	-5.5	177	0.00
16		Acenaphthene	1.000	1.007	-0.7	160	0.00
17		Fluorene	1.000	1.025	-2.5	166	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	153	0.00
19		4-Bromophenyl-phenylether	1.000	1.055	-5.5	168	0.00
20		Hexachlorobenzene	1.000	0.993	0.7	149	0.00
21		Pentachlorophenol	1.000	1.185	-18.5	173	0.00
22		Phenanthrene	1.000	1.037	-3.7	159	-0.02
23		Anthracene	1.000	1.119	-11.9	178	0.00
24		Fluoranthene	1.000	1.101	-10.1	182	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	159	0.00
26		Pyrene	1.000	1.162	-16.2	194	0.00
27	S	Terphenyl-d14	1.000	1.107	-10.7	184	0.00
28		Benzo(a)anthracene	1.000	1.097	-9.7	176	0.00
29		Chrysene	1.000	1.101	-10.1	168	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.160	-16.0	211	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.174	-17.4	202	0.00
32	I	Perylene-d12	5.000	5.000	0.0	173	0.00
33		Benzo(b)fluoranthene	1.000	1.041	-4.1	185	0.00
34		Benzo(k)fluoranthene	1.000	1.068	-6.8	187	0.00
35	C	Benzo(a)pyrene	1.000	1.114	-11.4	206	0.00
36		Dibenzo(a,h)anthracene	1.000	1.077	-7.7	201	-0.01
37		Benzo(g,h,i)perylene	1.000	1.066	-6.6	195	-0.01

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