Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081715\
Data File : BE090548.D
Acq On : 18 Aug 2015 00:24
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 18 06:12:00 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE081415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 14 16:52:42 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	112	0.00
2		1,4-Dioxane	1.000	1.063	-6.3	103	0.00
3		n-Nitrosodimethylamine	1.000	0.948	5.2	108	0.00
4	S	2-Fluorophenol	1.000	1.068	-6.8	119	0.00
5	S	Phenol-d6	1.000	1.054	-5.4	123	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	115	0.00
7	S	Nitrobenzene-d5	1.000	0.999	0.1	121	0.00
8		Nitrobenzene	1.000	0.991	0.9	122	0.00
9		Naphthalene	1.000	0.981	1.9	114	0.00
10		Hexachlorobutadiene	1.000	0.977	2.3	112	0.00
11		2-Methylnaphthalene	1.000	1.019	-1.9	120	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	119	0.00
13	S	2,4,6-Tribromophenol	1.000	1.026	-2.6	134	0.00
14	S	2-Fluorobiphenyl	1.000	1.002	-0.2	119	0.00
15		Acenaphthylene	1.000	1.003	-0.3	121	0.00
16		Acenaphthene	1.000	0.998	0.2	120	0.00
17		Fluorene	1.000	0.987	1.3	118	0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	116	0.00
19		4-Bromophenyl-phenylether	1.000	1.024	-2.4	122	0.00
20		Hexachlorobenzene	1.000	1.007	-0.7	117	0.00
21		Pentachlorophenol	1.000	0.959	4.1	148	0.00
22		Phenanthrene	1.000	0.974	2.6	115	0.00
23		Anthracene	1.000	1.004	-0.4	120	0.00
24		Fluoranthene	1.000	0.982	1.8	117	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	114	0.00
26		Pyrene	1.000	1.009	-0.9	118	0.00
27	S	Terphenyl-d14	1.000	1.023	-2.3	118	0.00
28		Benzo(a)anthracene	1.000	0.979	2.1	117	0.00
29		Chrysene	1.000	1.000	0.0	115	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.147	-14.7	155	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.999	0.1	123	0.00
32	I	Perylene-d12	5.000	5.000	0.0	118	0.00
33		Benzo(b)fluoranthene	1.000	1.014	-1.4	123	0.00
34		Benzo(k)fluoranthene	1.000	1.019	-1.9	119	0.00
35	C	Benzo(a)pyrene	1.000	1.015	-1.5	124	0.00
36		Dibenzo(a,h)anthracene	1.000	1.002	-0.2	125	0.00
37		Benzo(g,h,i)perylene	1.000	1.026	-2.6	123	0.00

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