Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081715\
Data File : BE090546.D
Acq On : 17 Aug 2015 23:09
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 18 04:36:55 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	115	0.00
2		1,4-Dioxane	1.000	1.077	-7.7	107	0.00
3		n-Nitrosodimethylamine	1.000	0.945	5.5	110	0.00
4	S	2-Fluorophenol	1.000	1.063	-6.3	121	0.00
5	S	Phenol-d6	1.000	1.032	-3.2	124	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	118	0.00
7	S	Nitrobenzene-d5	1.000	0.994	0.6	123	0.00
8		Nitrobenzene	1.000	0.990	1.0	125	0.00
9		Naphthalene	1.000	0.982	1.8	117	0.00
10		Hexachlorobutadiene	1.000	0.981	1.9	116	0.00
11		2-Methylnaphthalene	1.000	1.016	-1.6	123	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	121	0.00
13	S	2,4,6-Tribromophenol	1.000	1.055	-5.5	141	0.00
14	S	2-Fluorobiphenyl	1.000	1.008	-0.8	123	0.00
15		Acenaphthylene	1.000	1.008	-0.8	125	0.00
16		Acenaphthene	1.000	0.995	0.5	122	0.00
17		Fluorene	1.000	0.991	0.9	121	0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	119	0.00
19		4-Bromophenyl-phenylether	1.000	1.019	-1.9	124	0.00
20		Hexachlorobenzene	1.000	1.001	-0.1	119	0.00
21		Pentachlorophenol	1.000	0.954	4.6	150	0.00
22		Phenanthrene	1.000	0.966	3.4	117	0.00
23		Anthracene	1.000	0.994	0.6	122	0.00
24		Fluoranthene	1.000	0.976	2.4	119	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	116	0.00
26		Pyrene	1.000	1.011	-1.1	120	0.00
27	S	Terphenyl-d14	1.000	1.032	-3.2	122	0.00
28		Benzo(a)anthracene	1.000	0.980	2.0	120	0.00
29		Chrysene	1.000	0.999	0.1	117	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.155	-15.5	160	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.010	-1.0	127	0.00
32	I	Perylene-d12	5.000	5.000	0.0	122	0.00
33		Benzo(b)fluoranthene	1.000	1.013	-1.3	126	0.00
34		Benzo(k)fluoranthene	1.000	1.018	-1.8	123	0.00
35	C	Benzo(a)pyrene	1.000	1.015	-1.5	127	0.00
36		Dibenzo(a,h)anthracene	1.000	1.004	-0.4	129	0.00
37		Benzo(g,h,i)perylene	1.000	1.023	-2.3	126	0.00

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