Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081415\
Data File : BE090514.D
Acq On : 14 Aug 2015 10:06
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 14 15:26:29 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 15:17:14 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	116	0.00
2		1,4-Dioxane	1.000	0.951	4.9	115	0.00
3		n-Nitrosodimethylamine	1.000	1.013	-1.3	124	0.00
4	S	2-Fluorophenol	1.000	0.959	4.1	118	0.00
5	S	Phenol-d6	1.000	0.994	0.6	123	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	117	0.00
7	S	Nitrobenzene-d5	1.000	1.011	-1.1	123	0.00
8		Nitrobenzene	1.000	1.021	-2.1	125	0.00
9		Naphthalene	1.000	0.938	6.2	117	0.00
10		Hexachlorobutadiene	1.000	0.934	6.6	116	0.00
11		2-Methylnaphthalene	1.000	0.981	1.9	123	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	121	0.00
13	S	2,4,6-Tribromophenol	1.000	1.025	-2.5	133	-0.02
14	S	2-Fluorobiphenyl	1.000	0.960	4.0	122	0.00
15		Acenaphthylene	1.000	0.967	3.3	124	0.00
16		Acenaphthene	1.000	0.943	5.7	122	0.00
17		Fluorene	1.000	0.957	4.3	123	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	122	0.00
19		4-Bromophenyl-phenylether	1.000	0.963	3.7	127	0.00
20		Hexachlorobenzene	1.000	0.922	7.8	120	0.00
21		Pentachlorophenol	1.000	1.109	-10.9	152	0.00
22		Phenanthrene	1.000	0.915	8.5	121	0.00
23		Anthracene	1.000	0.972	2.8	127	0.00
24		Fluoranthene	1.000	0.978	2.2	128	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	126	0.00
26		Pyrene	1.000	0.940	6.0	128	0.00
27	S	Terphenyl-d14	1.000	0.954	4.6	128	0.00
28		Benzo(a)anthracene	1.000	0.924	7.6	129	0.00
29		Chrysene	1.000	0.947	5.3	128	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.058	-5.8	159	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.973	2.7	135	0.00
32	I	Perylene-d12	5.000	5.000	0.0	129	0.00
33		Benzo(b)fluoranthene	1.000	0.978	2.2	135	0.00
34		Benzo(k)fluoranthene	1.000	0.963	3.7	129	0.00
35	C	Benzo(a)pyrene	1.000	0.986	1.4	134	0.00
36		Dibenzo(a,h)anthracene	1.000	0.983	1.7	135	0.00
37		Benzo(g,h,i)perylene	1.000	0.977	2.3	131	0.00

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