Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE082015\
Data File : BE090583.D
Acq On : 20 Aug 2015 20:35
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 21 01:33:01 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	86	0.00
2		1,4-Dioxane	1.000	1.248	-24.8	91	0.00
3		n-Nitrosodimethylamine	1.000	1.111	-11.1	97	0.00
4	S	2-Fluorophenol	1.000	0.920	8.0	78	0.00
5	S	Phenol-d6	1.000	0.840	16.0	75	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	80	0.00
7	S	Nitrobenzene-d5	1.000	1.007	-0.7	85	0.00
8		Nitrobenzene	1.000	0.968	3.2	83	0.00
9		Naphthalene	1.000	0.978	2.2	79	0.00
10		Hexachlorobutadiene	1.000	0.990	1.0	79	0.00
11		2-Methylnaphthalene	1.000	1.024	-2.4	84	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	90	0.00
13	S	2,4,6-Tribromophenol	1.000	1.204	-20.4	119	0.00
14	S	2-Fluorobiphenyl	1.000	0.973	2.7	87	0.00
15		Acenaphthylene	1.000	0.987	1.3	90	0.00
16		Acenaphthene	1.000	1.000	0.0	91	0.00
17		Fluorene	1.000	1.055	-5.5	95	0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	102	0.00
19		4-Bromophenyl-phenylether	1.000	0.965	3.5	101	0.00
20		Hexachlorobenzene	1.000	0.947	5.3	97	0.00
21		Pentachlorophenol	1.000	1.218	-21.8	182	0.00
22		Phenanthrene	1.000	0.964	3.6	100	0.00
23		Anthracene	1.000	1.002	-0.2	106	0.00
24		Fluoranthene	1.000	1.072	-7.2	112	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	119	0.00
26		Pyrene	1.000	0.939	6.1	115	0.00
27	S	Terphenyl-d14	1.000	0.966	3.4	117	0.00
28		Benzo(a)anthracene	1.000	0.972	2.8	122	0.00
29		Chrysene	1.000	0.992	0.8	119	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.063	-6.3	149	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.960	4.0	124	0.00
32	I	Perylene-d12	5.000	5.000	0.0	122	0.00
33		Benzo(b)fluoranthene	1.000	0.999	0.1	125	0.00
34		Benzo(k)fluoranthene	1.000	1.011	-1.1	122	0.00
35	C	Benzo(a)pyrene	1.000	0.989	1.1	124	0.00
36		Dibenzo(a,h)anthracene	1.000	0.960	4.0	123	0.00
37		Benzo(g,h,i)perylene	1.000	0.986	1.4	122	0.00

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