Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080415\
Data File : BE090344.D
Acq On : 4 Aug 2015 15:34
Operator : TP/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 04 17:41:38 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 04 02:31:31 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	95	-0.01
2		1,4-Dioxane	0.796	0.799	-0.4	94	-0.02
3		n-Nitrosodimethylamine	0.845	0.985	-16.6	109	-0.02
4	S	2-Fluorophenol	0.956	1.086	-13.6	111	-0.01
5	S	Phenol-d6	1.175	1.234	-5.0	108	-0.01
6	I	Naphthalene-d8	1.000	1.000	0.0	108	-0.02
7	S	Nitrobenzene-d5	0.295	0.284	3.7	108	0.00
8		Nitrobenzene	0.305	0.294	3.6	110	0.00
9		Naphthalene	1.074	1.053	2.0	106	-0.02
10		Hexachlorobutadiene	0.144	0.132	8.3	96	0.00
11		2-Methylnaphthalene	0.665	0.649	2.4	106	-0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	107	-0.01
13	S	2,4,6-Tribromophenol	0.067	0.074	-10.4	164#	-0.02
14	S	2-Fluorobiphenyl	1.449	1.490	-2.8	109	-0.01
15		Acenaphthylene	8.730	8.557	2.0	105	-0.01
16		Acenaphthene	1.262	1.213	3.9	104	-0.01
17		Fluorene	1.533	1.509	1.6	112	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	111	0.00
19		4-Bromophenyl-phenylether	0.188	0.184	2.1	106	-0.02
20		Hexachlorobenzene	0.710	0.678	4.5	104	0.00
21		Pentachlorophenol	0.038	0.794	-1989.5#	3601#	-0.03
22		Phenanthrene	1.139	1.113	2.3	114	-0.02
23		Anthracene	1.018	0.988	2.9	115	-0.02
24		Fluoranthene	1.146	1.149	-0.3	117	-0.01
25	I	Chrysene-d12	1.000	1.000	0.0	142	-0.01
26		Pyrene	1.352	1.195	11.6	120	-0.01
27	S	Terphenyl-d14	0.936	0.915	2.2	133	-0.02
28		Benzo(a)anthracene	0.951	0.968	-1.8	162#	-0.01
29		Chrysene	1.320	1.142	13.5	118	-0.01
30		Bis(2-ethylhexyl)phthalate	0.335	0.385	-14.9	230#	0.00
31		Indeno(1,2,3-cd)pyrene	0.548	0.891	-62.6#	322#	-0.04
32	I	Perylene-d12	1.000	1.000	0.0	165#	-0.02
33		Benzo(b)fluoranthene	0.676	0.820	-21.3	234#	-0.02
34		Benzo(k)fluoranthene	1.357	1.172	13.6	133	-0.01
35	C	Benzo(a)pyrene	0.673	0.801	-19.0	223#	-0.02
36		Dibenzo(a,h)anthracene	0.431	0.705	-63.6#	385#	-0.05
37		Benzo(g,h,i)perylene	0.744	0.924	-24.2	212#	-0.04

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