Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081215\
Data File : BE090478.D
Acq On : 12 Aug 2015 21:45
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 13 01:22:01 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 06 12:28:08 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	108	0.00
2		1,4-Dioxane	1.000	1.073	-7.3	103	0.00
3		n-Nitrosodimethylamine	1.000	1.048	-4.8	114	0.00
4	S	2-Fluorophenol	1.000	1.439	-43.9#	167	0.00
5	S	Phenol-d6	1.000	1.164	-16.4	145	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	109	0.00
7	S	Nitrobenzene-d5	1.000	1.041	-4.1	123	0.00
8		Nitrobenzene	1.000	1.027	-2.7	129	0.00
9		Naphthalene	1.000	1.018	-1.8	112	0.00
10		Hexachlorobutadiene	1.000	1.093	-9.3	109	0.00
11		2-Methylnaphthalene	1.000	1.085	-8.5	123	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	119	0.00
13	S	2,4,6-Tribromophenol	1.000	1.551	-55.1#	224	0.00
14	S	2-Fluorobiphenyl	1.000	1.036	-3.6	121	0.00
15		Acenaphthylene	1.000	1.041	-4.1	125	0.00
16		Acenaphthene	1.000	1.009	-0.9	120	0.00
17		Fluorene	1.000	1.030	-3.0	121	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	119	0.00
19		4-Bromophenyl-phenylether	1.000	0.989	1.1	120	0.00
20		Hexachlorobenzene	1.000	0.939	6.1	111	0.00
21		Pentachlorophenol	1.000	1.432	-43.2#	225	0.00
22		Phenanthrene	1.000	1.020	-2.0	122	0.00
23		Anthracene	1.000	1.100	-10.0	134	-0.02
24		Fluoranthene	1.000	1.104	-10.4	136	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	125	0.00
26		Pyrene	1.000	1.065	-6.5	138	0.00
27	S	Terphenyl-d14	1.000	1.089	-8.9	137	-0.01
28		Benzo(a)anthracene	1.000	1.083	-8.3	150	0.00
29		Chrysene	1.000	0.977	2.3	119	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.281	-28.1#	192	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.115	-11.5	176	0.00
32	I	Perylene-d12	5.000	5.000	0.0	135	0.00
33		Benzo(b)fluoranthene	1.000	1.100	-10.0	166	0.00
34		Benzo(k)fluoranthene	1.000	0.899	10.1	117	0.00
35	C	Benzo(a)pyrene	1.000	1.074	-7.4	160	0.00
36		Dibenzo(a,h)anthracene	1.000	1.092	-9.2	182	-0.01
37		Benzo(g,h,i)perylene	1.000	1.091	-9.1	152	0.00

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