Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081215\
Data File : BE090497.D
Acq On : 13 Aug 2015 9:11
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 13 10:51:00 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	137	0.00
2		1,4-Dioxane	1.000	1.033	-3.3	125	0.00
3		n-Nitrosodimethylamine	1.000	1.104	-10.4	152	0.00
4	S	2-Fluorophenol	1.000	2.003	-100.3#	294	0.00
5	S	Phenol-d6	1.000	1.589	-58.9#	268	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	143	0.00
7	S	Nitrobenzene-d5	1.000	1.222	-22.2	197	-0.01
8		Nitrobenzene	1.000	1.202	-20.2	206	0.00
9		Naphthalene	1.000	1.014	-1.4	146	0.00
10		Hexachlorobutadiene	1.000	1.039	-3.9	136	0.00
11		2-Methylnaphthalene	1.000	1.186	-18.6	177	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	151	0.00
13	S	2,4,6-Tribromophenol	1.000	2.464	-146.4#	502	-0.02
14	S	2-Fluorobiphenyl	1.000	1.084	-8.4	161	0.00
15		Acenaphthylene	1.000	1.141	-14.1	173	0.00
16		Acenaphthene	1.000	1.029	-2.9	155	0.00
17		Fluorene	1.000	1.063	-6.3	158	-0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	140	0.00
19		4-Bromophenyl-phenylether	1.000	1.127	-12.7	161	0.00
20		Hexachlorobenzene	1.000	0.921	7.9	128	-0.02
21		Pentachlorophenol	1.000	2.591	-159.1#	589	-0.01
22		Phenanthrene	1.000	1.045	-4.5	147	0.00
23		Anthracene	1.000	1.225	-22.5	176	-0.02
24		Fluoranthene	1.000	1.214	-21.4	176	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	137	0.00
26		Pyrene	1.000	1.261	-26.1#	179	0.00
27	S	Terphenyl-d14	1.000	1.302	-30.2#	180	-0.01
28		Benzo(a)anthracene	1.000	1.200	-20.0	183	0.00
29		Chrysene	1.000	1.012	-1.2	135	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	2.733	-173.3#	545	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.305	-30.5#	235	0.00
32	I	Perylene-d12	5.000	5.000	0.0	159	0.00
33		Benzo(b)fluoranthene	1.000	1.162	-16.2	210	0.00
34		Benzo(k)fluoranthene	1.000	0.958	4.2	148	0.00
35	C	Benzo(a)pyrene	1.000	1.215	-21.5#	221	0.00
36		Dibenzo(a,h)anthracene	1.000	1.243	-24.3	256	-0.01
37		Benzo(g,h,i)perylene	1.000	1.254	-25.4#	207	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1