Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE081215\
Data File : BE090462.D
Acq On : 12 Aug 2015 11:51
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 13 01:22:15 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	118	0.00
2		1,4-Dioxane	1.000	1.002	-0.2	105	0.00
3		n-Nitrosodimethylamine	1.000	1.041	-4.1	123	0.00
4	S	2-Fluorophenol	1.000	1.367	-36.7#	173	0.00
5	S	Phenol-d6	1.000	1.199	-19.9	164	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	125	0.00
7	S	Nitrobenzene-d5	1.000	1.036	-3.6	140	0.00
8		Nitrobenzene	1.000	1.007	-0.7	144	0.00
9		Naphthalene	1.000	1.019	-1.9	128	0.00
10		Hexachlorobutadiene	1.000	1.086	-8.6	123	0.00
11		2-Methylnaphthalene	1.000	1.058	-5.8	138	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	134	0.00
13	S	2,4,6-Tribromophenol	1.000	1.346	-34.6#	212	0.00
14	S	2-Fluorobiphenyl	1.000	1.023	-2.3	134	0.00
15		Acenaphthylene	1.000	1.013	-1.3	136	0.00
16		Acenaphthene	1.000	0.990	1.0	132	0.00
17		Fluorene	1.000	1.010	-1.0	133	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	128	0.00
19		4-Bromophenyl-phenylether	1.000	1.007	-0.7	132	0.00
20		Hexachlorobenzene	1.000	0.929	7.1	118	0.00
21		Pentachlorophenol	1.000	1.212	-21.2	195	0.00
22		Phenanthrene	1.000	1.012	-1.2	131	0.00
23		Anthracene	1.000	1.075	-7.5	141	-0.02
24		Fluoranthene	1.000	1.036	-3.6	138	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	126	0.00
26		Pyrene	1.000	1.075	-7.5	140	0.00
27	S	Terphenyl-d14	1.000	1.076	-7.6	136	-0.01
28		Benzo(a)anthracene	1.000	1.042	-4.2	146	0.00
29		Chrysene	1.000	0.972	2.8	119	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.077	-7.7	157	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.036	-3.6	161	0.00
32	I	Perylene-d12	5.000	5.000	0.0	129	0.00
33		Benzo(b)fluoranthene	1.000	1.067	-6.7	153	0.00
34		Benzo(k)fluoranthene	1.000	0.901	9.9	113	0.00
35	C	Benzo(a)pyrene	1.000	1.035	-3.5	146	0.00
36		Dibenzo(a,h)anthracene	1.000	1.061	-6.1	168	0.00
37		Benzo(g,h,i)perylene	1.000	1.037	-3.7	138	-0.01

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