Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081115\
Data File : BE090443.D
Acq On : 12 Aug 2015 00:13
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 12 05:46:15 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	120	0.00
2		1,4-Dioxane	1.000	0.973	2.7	104	0.00
3		n-Nitrosodimethylamine	1.000	1.015	-1.5	122	0.00
4	S	2-Fluorophenol	1.000	1.402	-40.2#	180	0.00
5	S	Phenol-d6	1.000	1.204	-20.4	167	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	126	0.00
7	S	Nitrobenzene-d5	1.000	1.022	-2.2	139	0.00
8		Nitrobenzene	1.000	1.011	-1.1	146	0.00
9		Naphthalene	1.000	1.018	-1.8	129	0.00
10		Hexachlorobutadiene	1.000	1.091	-9.1	125	0.00
11		2-Methylnaphthalene	1.000	1.095	-9.5	144	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	134	0.00
13	S	2,4,6-Tribromophenol	1.000	1.465	-46.5#	236	0.00
14	S	2-Fluorobiphenyl	1.000	1.050	-5.0	139	0.00
15		Acenaphthylene	1.000	1.052	-5.2	142	0.00
16		Acenaphthene	1.000	1.013	-1.3	136	0.00
17		Fluorene	1.000	1.029	-2.9	137	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	129	0.00
19		4-Bromophenyl-phenylether	1.000	1.030	-3.0	136	0.00
20		Hexachlorobenzene	1.000	0.938	6.2	120	0.00
21		Pentachlorophenol	1.000	1.236	-23.6	201	0.00
22		Phenanthrene	1.000	1.032	-3.2	134	0.00
23		Anthracene	1.000	1.116	-11.6	147	-0.02
24		Fluoranthene	1.000	1.069	-6.9	143	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	127	0.00
26		Pyrene	1.000	1.098	-9.8	145	0.00
27	S	Terphenyl-d14	1.000	1.118	-11.8	144	-0.01
28		Benzo(a)anthracene	1.000	1.060	-6.0	150	0.00
29		Chrysene	1.000	1.006	-0.6	125	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.210	-21.0	184	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.049	-4.9	166	0.00
32	I	Perylene-d12	5.000	5.000	0.0	132	0.00
33		Benzo(b)fluoranthene	1.000	1.086	-8.6	160	0.00
34		Benzo(k)fluoranthene	1.000	0.946	5.4	121	0.00
35	C	Benzo(a)pyrene	1.000	1.053	-5.3	153	0.00
36		Dibenzo(a,h)anthracene	1.000	1.066	-6.6	173	0.00
37		Benzo(g,h,i)perylene	1.000	1.063	-6.3	145	0.00

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