Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE091815\
Data File : BE090892.D
Acq On : 19 Sep 2015 2:25
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 19 03:47:26 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE091815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 18 18:17:34 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	113	0.00
2		1,4-Dioxane	1.000	0.968	3.2	101	0.00
3		n-Nitrosodimethylamine	1.000	0.910	9.0	97	0.00
4	S	2-Fluorophenol	1.000	0.860	14.0	93	0.00
5	S	Phenol-d6	1.000	0.801	19.9	93	0.00
6		bis(2-Chloroethyl)ether	1.000	0.962	3.8	111	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	113	0.00
8	S	Nitrobenzene-d5	1.000	0.868	13.2	99	0.00
9		Nitrobenzene	1.000	0.897	10.3	105	0.00
10		Naphthalene	1.000	0.970	3.0	105	0.00
11		Hexachlorobutadiene	1.000	1.036	-3.6	111	0.00
12		2-Methylnaphthalene	1.000	0.900	10.0	104	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	118	0.00
14	S	2,4,6-Tribromophenol	1.000	0.769	23.1	93	-0.02
15	S	2-Fluorobiphenyl	1.000	0.921	7.9	107	0.00
16		Acenaphthylene	1.000	0.891	10.9	106	0.00
17		Acenaphthene	1.000	0.937	6.3	111	0.00
18		Fluorene	1.000	0.941	5.9	110	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	123	0.00
20		4-Bromophenyl-phenylether	1.000	0.877	12.3	109	-0.02
21		Hexachlorobenzene	1.000	0.966	3.4	113	0.00
22		Pentachlorophenol	1.000	0.761	23.9	86	0.00
23		Phenanthrene	1.000	0.939	6.1	112	0.00
24		Anthracene	1.000	0.854	14.6	108	0.00
25		Fluoranthene	1.000	0.870	13.0	108	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	113	0.00
27		Pyrene	1.000	0.892	10.8	108	0.00
28	S	Terphenyl-d14	1.000	0.898	10.2	105	0.00
29		Benzo(a)anthracene	1.000	0.909	9.1	108	0.00
30		Chrysene	1.000	0.968	3.2	105	0.00
31		Bis(2-ethylhexyl)phthalate	1.000	0.812	18.8	95	0.00
32		Indeno(1,2,3-cd)pyrene	1.000	0.864	13.6	97	0.00
33	I	Perylene-d12	5.000	5.000	0.0	110	0.00
34		Benzo(b)fluoranthene	1.000	0.906	9.4	102	0.00
35		Benzo(k)fluoranthene	1.000	0.901	9.9	99	0.00
36	C	Benzo(a)pyrene	1.000	0.873	12.7	99	0.00
37		Dibenzo(a,h)anthracene	1.000	0.875	12.5	97	0.00
38		Benzo(g,h,i)perylene	1.000	0.890	11.0	99	0.00

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