Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE092415\
Data File : BE090967.D
Acq On : 24 Sep 2015 5:54
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 24 08:36:38 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	133	0.00
2		1,4-Dioxane	1.000	1.087	-8.7	131	0.00
3		n-Nitrosodimethylamine	1.000	0.833	16.7	105	0.00
4	S	2-Fluorophenol	1.000	0.746	25.4#	96	0.00
5	S	Phenol-d6	1.000	0.662	33.8#	91	0.04
6		bis(2-Chloroethyl)ether	1.000	1.054	-5.4	142	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	141	0.00
8	S	Nitrobenzene-d5	1.000	0.812	18.8	115	0.00
9		Nitrobenzene	1.000	0.742	25.8#	109	0.00
10		Naphthalene	1.000	1.040	-4.0	139	0.00
11		Hexachlorobutadiene	1.000	1.196	-19.6	157	0.00
12		2-Methylnaphthalene	1.000	0.967	3.3	140	0.01
13	I	Acenaphthene-d10	5.000	5.000	0.0	151	0.00
14	S	2,4,6-Tribromophenol	1.000	0.671	32.9#	104	0.00
15	S	2-Fluorobiphenyl	1.000	0.957	4.3	143	0.00
16		Acenaphthylene	1.000	0.955	4.5	145	0.00
17		Acenaphthene	1.000	1.018	-1.8	155	0.00
18		Fluorene	1.000	0.971	2.9	145	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	137	0.00
20		4-Bromophenyl-phenylether	1.000	0.999	0.1	139	0.00
21		Hexachlorobenzene	1.000	1.206	-20.6	154	0.00
22		Pentachlorophenol	1.000	0.541	45.9#	60	0.04
23		Phenanthrene	1.000	1.004	-0.4	133	0.00
24		Anthracene	1.000	0.902	9.8	128	0.00
25		Fluoranthene	1.000	0.879	12.1	121	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	107	0.00
27		Pyrene	1.000	1.048	-4.8	120	0.00
28	S	Terphenyl-d14	1.000	1.025	-2.5	112	0.01
29		Benzo(a)anthracene	1.000	0.848	15.2	95	0.00
30		Chrysene	1.000	1.144	-14.4	115	0.00
31		Bis(2-ethylhexyl)phthalate	1.000	0.787	21.3	86	0.00
32		Indeno(1,2,3-cd)pyrene	1.000	0.831	16.9	88	0.00
33	I	Perylene-d12	5.000	5.000	0.0	97	0.00
34		Benzo(b)fluoranthene	1.000	0.882	11.8	88	0.00
35		Benzo(k)fluoranthene	1.000	1.090	-9.0	106	0.00
36	C	Benzo(a)pyrene	1.000	1.044	-4.4	105	0.00
37		Dibenzo(a,h)anthracene	1.000	0.860	14.0	84	0.00
38		Benzo(g,h,i)perylene	1.000	0.909	9.1	89	0.00

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