Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090215\
Data File : BE090710.D
Acq On : 2 Sep 2015 18:36
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 02 23:06:26 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 02 23:04:24 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	76	0.00
2		1,4-Dioxane	1.000	1.073	-7.3	75	0.00
3		n-Nitrosodimethylamine	1.000	0.969	3.1	71	0.00
4	S	2-Fluorophenol	1.000	1.219	-21.9	91	0.00
5	S	Phenol-d6	1.000	1.130	-13.0	90	0.01
6	I	Naphthalene-d8	5.000	5.000	0.0	80	0.00
7	S	Nitrobenzene-d5	1.000	1.007	-0.7	81	0.00
8		Nitrobenzene	1.000	0.982	1.8	83	0.00
9		Naphthalene	1.000	1.000	0.0	81	0.00
10		Hexachlorobutadiene	1.000	0.948	5.2	74	0.00
11		2-Methylnaphthalene	1.000	1.102	-10.2	91	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	85	0.00
13	S	2,4,6-Tribromophenol	1.000	1.296	-29.6#	120	0.00
14	S	2-Fluorobiphenyl	1.000	1.018	-1.8	88	0.00
15		Acenaphthylene	1.000	1.068	-6.8	94	0.00
16		Acenaphthene	1.000	1.007	-0.7	84	0.00
17		Fluorene	1.000	1.029	-2.9	88	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	85	0.00
19		4-Bromophenyl-phenylether	1.000	1.001	-0.1	88	0.00
20		Hexachlorobenzene	1.000	0.967	3.3	81	0.00
21		Pentachlorophenol	1.000	1.401	-40.1#	125	0.00
22		Phenanthrene	1.000	1.004	-0.4	86	0.00
23		Anthracene	1.000	1.061	-6.1	94	0.00
24		Fluoranthene	1.000	1.132	-13.2	104	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	92	0.00
26		Pyrene	1.000	1.140	-14.0	111	0.00
27	S	Terphenyl-d14	1.000	1.089	-8.9	105	0.00
28		Benzo(a)anthracene	1.000	1.067	-6.7	100	0.00
29		Chrysene	1.000	1.099	-9.9	97	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.387	-38.7#	150	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.136	-13.6	113	0.00
32	I	Perylene-d12	5.000	5.000	0.0	101	0.00
33		Benzo(b)fluoranthene	1.000	1.042	-4.2	107	0.00
34		Benzo(k)fluoranthene	1.000	1.085	-8.5	111	0.00
35	C	Benzo(a)pyrene	1.000	1.122	-12.2	121	0.00
36		Dibenzo(a,h)anthracene	1.000	1.049	-4.9	113	0.00
37		Benzo(g,h,i)perylene	1.000	1.059	-5.9	112	0.00

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