Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE090415\
Data File : BE090745.D
Acq On : 5 Sep 2015 00:29
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 07 01:03:25 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 11:40:50 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	78	0.00
2		1,4-Dioxane	1.000	1.038	-3.8	76	0.00
3		n-Nitrosodimethylamine	1.000	0.936	6.4	71	0.00
4	S	2-Fluorophenol	1.000	1.103	-10.3	85	0.00
5	S	Phenol-d6	1.000	1.067	-6.7	88	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	85	0.00
7	S	Nitrobenzene-d5	1.000	0.980	2.0	83	0.00
8		Nitrobenzene	1.000	0.941	5.9	84	0.00
9		Naphthalene	1.000	1.005	-0.5	86	0.00
10		Hexachlorobutadiene	1.000	0.979	2.1	80	0.00
11		2-Methylnaphthalene	1.000	1.128	-12.8	99	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	98	0.00
13	S	2,4,6-Tribromophenol	1.000	1.142	-14.2	120	0.00
14	S	2-Fluorobiphenyl	1.000	0.977	2.3	97	0.00
15		Acenaphthylene	1.000	1.053	-5.3	107	0.00
16		Acenaphthene	1.000	1.024	-2.4	99	0.00
17		Fluorene	1.000	1.039	-3.9	103	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	95	0.00
19		4-Bromophenyl-phenylether	1.000	1.039	-3.9	103	0.00
20		Hexachlorobenzene	1.000	1.021	-2.1	95	0.00
21		Pentachlorophenol	1.000	1.172	-17.2	106	0.00
22		Phenanthrene	1.000	1.028	-2.8	98	-0.02
23		Anthracene	1.000	1.046	-4.6	103	0.00
24		Fluoranthene	1.000	1.056	-5.6	108	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	96	0.00
26		Pyrene	1.000	1.119	-11.9	113	0.00
27	S	Terphenyl-d14	1.000	1.076	-7.6	108	0.00
28		Benzo(a)anthracene	1.000	1.066	-6.6	104	0.00
29		Chrysene	1.000	1.099	-9.9	101	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.154	-15.4	127	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.132	-13.2	118	0.00
32	I	Perylene-d12	5.000	5.000	0.0	102	0.00
33		Benzo(b)fluoranthene	1.000	1.059	-5.9	111	0.00
34		Benzo(k)fluoranthene	1.000	1.037	-3.7	108	0.00
35	C	Benzo(a)pyrene	1.000	1.089	-8.9	119	0.00
36		Dibenzo(a,h)anthracene	1.000	1.052	-5.2	116	-0.01
37		Benzo(g,h,i)perylene	1.000	1.048	-4.8	113	-0.02

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