Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_E\DATA\BE072015\
Data File : BE090175.D
Acq On : 20 Jul 2015 14:04
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 21 02:09:46 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jul 18 00:40:53 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	125	0.00
2		1,4-Dioxane	1.000	1.116	-11.6	130	0.00
3		n-Nitrosodimethylamine	1.000	0.992	0.8	128	0.00
4	S	2-Fluorophenol	1.000	0.907	9.3	115	-0.01
5	S	Phenol-d6	1.000	0.939	6.1	117	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	121	0.00
7	S	Nitrobenzene-d5	1.000	0.947	5.3	118	0.00
8		Nitrobenzene	1.000	0.980	2.0	118	0.00
9		Naphthalene	1.000	0.969	3.1	121	0.00
10		Hexachlorobutadiene	1.000	0.972	2.8	120	0.00
11		2-Methylnaphthalene	1.000	0.955	4.5	119	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	118	0.00
13	S	2,4,6-Tribromophenol	1.000	0.671	32.9#	81	-0.02
14	S	2-Fluorobiphenyl	1.000	0.976	2.4	117	0.00
15		Acenaphthylene	1.000	0.948	5.2	115	0.00
16		Acenaphthene	1.000	0.944	5.6	114	0.00
17		Fluorene	1.000	0.932	6.8	113	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	113	0.00
19		4-Bromophenyl-phenylether	1.000	0.991	0.9	116	0.00
20		Hexachlorobenzene	1.000	0.933	6.7	108	0.00
21		Pentachlorophenol	1.000	0.800	20.0	90	0.00
22		Phenanthrene	1.000	0.984	1.6	114	0.00
23		Anthracene	1.000	0.972	2.8	113	0.00
24		Fluoranthene	1.000	0.966	3.4	112	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	115	0.00
26		Pyrene	1.000	0.952	4.8	112	0.00
27	S	Terphenyl-d14	1.000	0.960	4.0	112	0.00
28		Benzo(a)anthracene	1.000	0.919	8.1	111	0.00
29		Chrysene	1.000	0.948	5.2	112	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.604	39.6#	71	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.817	18.3	96	0.00
32	I	Perylene-d12	5.000	5.000	0.0	109	0.00
33		Benzo(b)fluoranthene	1.000	0.912	8.8	102	0.00
34		Benzo(k)fluoranthene	1.000	0.953	4.7	106	0.00
35	C	Benzo(a)pyrene	1.000	0.895	10.5	100	0.00
36		Dibenzo(a,h)anthracene	1.000	0.847	15.3	95	0.00
37		Benzo(g,h,i)perylene	1.000	0.871	12.9	97	0.00

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