Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071615\
Data File : BE090139.D
Acq On : 17 Jul 2015 4:07
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 17 13:17:11 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE071615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 17 12:59:41 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	1.110	-11.0	82	0.00
3		n-Nitrosodimethylamine	1.000	0.925	7.5	74	0.00
4	S	2-Fluorophenol	1.000	0.851	14.9	69	0.00
5	S	Phenol-d6	1.000	0.887	11.3	71	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	72	0.00
7	S	Nitrobenzene-d5	1.000	0.910	9.0	69	0.00
8		Nitrobenzene	1.000	0.912	8.8	68	0.00
9		Naphthalene	1.000	0.957	4.3	72	0.00
10		Hexachlorobutadiene	1.000	0.999	0.1	75	0.00
11		2-Methylnaphthalene	1.000	0.946	5.4	71	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	70	0.00
13	S	2,4,6-Tribromophenol	1.000	0.847	15.3	72	0.00
14	S	2-Fluorobiphenyl	1.000	0.977	2.3	70	0.00
15		Acenaphthylene	1.000	0.964	3.6	70	0.00
16		Acenaphthene	1.000	0.966	3.4	70	0.00
17		Fluorene	1.000	0.946	5.4	68	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	71	0.00
19		4-Bromophenyl-phenylether	1.000	0.957	4.3	70	0.00
20		Hexachlorobenzene	1.000	0.970	3.0	69	0.00
21		Pentachlorophenol	1.000	4.318	-331.8#	725	0.00
22		Phenanthrene	1.000	0.937	6.3	69	0.00
23		Anthracene	1.000	0.937	6.3	68	0.00
24		Fluoranthene	1.000	0.957	4.3	69	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	72	0.00
26		Pyrene	1.000	0.951	4.9	70	0.00
27	S	Terphenyl-d14	1.000	0.994	0.6	71	0.00
28		Benzo(a)anthracene	1.000	1.053	-5.3	81	0.00
29		Chrysene	1.000	1.014	-1.4	71	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.963	3.7	74	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.308	-30.8#	103	0.00
32	I	Perylene-d12	5.000	5.000	0.0	76	0.00
33		Benzo(b)fluoranthene	1.000	1.302	-30.2#	100	0.00
34		Benzo(k)fluoranthene	1.000	0.890	11.0	74	0.00
35	C	Benzo(a)pyrene	1.000	0.940	6.0	96	0.00
36		Dibenzo(a,h)anthracene	1.000	1.272	-27.2#	106	0.00
37		Benzo(g,h,i)perylene	1.000	0.974	2.6	96	0.00

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