Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062915\
Data File : BE089997.D
Acq On : 30 Jun 2015 1:09
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jun 30 07:17:49 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 30 07:01:01 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	89	0.00
2		1,4-Dioxane	1.000	0.975	2.5	90	0.00
3		n-Nitrosodimethylamine	1.000	0.908	9.2	84	0.00
4	S	2-Fluorophenol	1.000	0.785	21.5	70	0.00
5	S	Phenol-d6	1.000	0.780	22.0	70	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	88	0.00
7	S	Nitrobenzene-d5	1.000	0.867	13.3	78	0.00
8		Nitrobenzene	1.000	0.874	12.6	79	0.00
9		Naphthalene	1.000	0.960	4.0	86	0.00
10		Hexachlorobutadiene	1.000	1.021	-2.1	91	0.00
11		2-Methylnaphthalene	1.000	0.949	5.1	85	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	87	0.00
13	S	2,4,6-Tribromophenol	1.000	0.601	39.9#	54	0.00
14	S	2-Fluorobiphenyl	1.000	0.973	2.7	87	0.00
15		Acenaphthylene	1.000	0.946	5.4	85	0.00
16		Acenaphthene	1.000	0.972	2.8	86	0.00
17		Fluorene	1.000	0.959	4.1	86	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	86	0.00
19		4-Bromophenyl-phenylether	1.000	0.967	3.3	85	0.00
20		Hexachlorobenzene	1.000	1.003	-0.3	90	0.00
21		Pentachlorophenol	1.000	0.663	33.7#	50	0.00
22		Phenanthrene	1.000	0.962	3.8	86	0.00
23		Anthracene	1.000	0.963	3.7	85	0.00
24		Fluoranthene	1.000	0.952	4.8	84	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	88	0.00
26		Pyrene	1.000	0.950	5.0	85	0.00
27	S	Terphenyl-d14	1.000	0.953	4.7	84	0.00
28		Benzo(a)anthracene	1.000	0.931	6.9	84	0.00
29		Chrysene	1.000	0.965	3.5	86	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.588	41.2#	41	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.844	15.6	75	0.00
32	I	Perylene-d12	5.000	5.000	0.0	84	0.00
33		Benzo(b)fluoranthene	1.000	0.966	3.4	81	0.00
34		Benzo(k)fluoranthene	1.000	0.966	3.4	81	0.00
35	C	Benzo(a)pyrene	1.000	0.913	8.7	76	0.00
36		Dibenzo(a,h)anthracene	1.000	0.886	11.4	75	0.00
37		Benzo(g,h,i)perylene	1.000	0.901	9.9	76	0.00

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