Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062615\
Data File : BE089979.D
Acq On : 27 Jun 2015 16:31
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jun 29 00:57:28 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 25 16:51:27 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	100	0.00
2		1,4-Dioxane	1.000	0.938	6.2	99	0.00
3		n-Nitrosodimethylamine	1.000	0.859	14.1	94	0.00
4	S	2-Fluorophenol	1.000	1.625	-62.5#	173	0.00
5	S	Phenol-d6	1.000	2.083	-108.3#	226	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	108	0.00
7	S	Nitrobenzene-d5	1.000	1.205	-20.5	146	0.00
8		Nitrobenzene	1.000	0.839	16.1	102	0.00
9		Naphthalene	1.000	0.838	16.2	98	0.00
10		Hexachlorobutadiene	1.000	0.819	18.1	95	0.00
11		2-Methylnaphthalene	1.000	0.799	20.1	95	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	103	0.00
13	S	2,4,6-Tribromophenol	1.000	2.369	-136.9#	362	0.00
14	S	2-Fluorobiphenyl	1.000	1.273	-27.3#	146	0.00
15		Acenaphthylene	1.000	0.826	17.4	91	0.00
16		Acenaphthene	1.000	0.811	18.9	88	0.00
17		Fluorene	1.000	0.826	17.4	85	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	93	0.00
19		4-Bromophenyl-phenylether	1.000	0.791	20.9	84	-0.02
20		Hexachlorobenzene	1.000	0.812	18.8	84	0.00
21		Pentachlorophenol	1.000	1.417	-41.7#	185	-0.02
22		Phenanthrene	1.000	0.821	17.9	85	-0.02
23		Anthracene	1.000	0.795	20.5	82	0.00
24		Fluoranthene	1.000	0.851	14.9	87	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	93	0.00
26		Pyrene	1.000	0.888	11.2	87	0.00
27	S	Terphenyl-d14	1.000	1.286	-28.6#	131	0.00
28		Benzo(a)anthracene	1.000	0.887	11.3	90	0.00
29		Chrysene	1.000	0.781	21.9	81	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	2.673	-167.3#	294	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.102	-10.2	108	0.00
32	I	Perylene-d12	5.000	5.000	0.0	98	0.00
33		Benzo(b)fluoranthene	1.000	0.973	2.7	101	0.00
34		Benzo(k)fluoranthene	1.000	0.871	12.9	91	0.00
35	C	Benzo(a)pyrene	1.000	0.987	1.3	104	0.00
36		Dibenzo(a,h)anthracene	1.000	0.988	1.2	105	0.00
37		Benzo(g,h,i)perylene	1.000	1.026	-2.6	108	0.00

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