Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE070615\
Data File : BE090060.D
Acq On : 6 Jul 2015 19:18
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 07 01:17:15 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	118	0.00
2		1,4-Dioxane	1.000	0.814	18.6	100	-0.01
3		n-Nitrosodimethylamine	1.000	0.815	18.5	100	-0.01
4	S	2-Fluorophenol	1.000	0.920	8.0	109	0.00
5	S	Phenol-d6	1.000	0.951	4.9	114	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	121	-0.01
7	S	Nitrobenzene-d5	1.000	0.858	14.2	106	0.00
8		Nitrobenzene	1.000	0.854	14.6	106	0.00
9		Naphthalene	1.000	0.921	7.9	113	-0.01
10		Hexachlorobutadiene	1.000	0.944	5.6	116	0.00
11		2-Methylnaphthalene	1.000	0.882	11.8	109	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	113	-0.01
13	S	2,4,6-Tribromophenol	1.000	0.789	21.1	91	0.00
14	S	2-Fluorobiphenyl	1.000	0.909	9.1	105	-0.01
15		Acenaphthylene	1.000	0.918	8.2	107	-0.01
16		Acenaphthene	1.000	0.932	6.8	106	-0.01
17		Fluorene	1.000	0.889	11.1	103	-0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	104	-0.02
19		4-Bromophenyl-phenylether	1.000	0.870	13.0	92	-0.02
20		Hexachlorobenzene	1.000	0.896	10.4	97	-0.02
21		Pentachlorophenol	1.000	0.727	27.3#	70	0.00
22		Phenanthrene	1.000	0.873	12.7	93	-0.02
23		Anthracene	1.000	0.900	10.0	95	0.00
24		Fluoranthene	1.000	0.880	12.0	93	-0.01
25	I	Chrysene-d12	5.000	5.000	0.0	96	-0.01
26		Pyrene	1.000	0.952	4.8	93	-0.01
27	S	Terphenyl-d14	1.000	0.922	7.8	89	-0.01
28		Benzo(a)anthracene	1.000	0.919	8.1	91	-0.01
29		Chrysene	1.000	0.902	9.8	88	-0.01
30		Bis(2-ethylhexyl)phthalate	1.000	0.870	13.0	91	-0.01
31		Indeno(1,2,3-cd)pyrene	1.000	0.915	8.5	90	-0.04
32	I	Perylene-d12	5.000	5.000	0.0	91	-0.02
33		Benzo(b)fluoranthene	1.000	1.020	-2.0	94	-0.02
34		Benzo(k)fluoranthene	1.000	0.981	1.9	90	-0.02
35	C	Benzo(a)pyrene	1.000	1.011	-1.1	92	-0.03
36		Dibenzo(a,h)anthracene	1.000	0.950	5.0	88	-0.04
37		Benzo(g,h,i)perylene	1.000	0.948	5.2	87	-0.04

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