Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070915\
Data File : BE090079.D
Acq On : 9 Jul 2015 18:46
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 10 02:19:48 2015
Quant Method : Z:\HPCHEM1\BNA_E\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	125	-0.01
2		1,4-Dioxane	1.000	0.795	20.5	103	0.00
3		n-Nitrosodimethylamine	1.000	0.778	22.2	101	-0.01
4	S	2-Fluorophenol	1.000	1.034	-3.4	129	0.00
5	S	Phenol-d6	1.000	1.213	-21.3	154	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	134	-0.01
7	S	Nitrobenzene-d5	1.000	0.899	10.1	123	-0.01
8		Nitrobenzene	1.000	0.813	18.7	113	0.00
9		Naphthalene	1.000	0.866	13.4	118	-0.01
10		Hexachlorobutadiene	1.000	0.890	11.0	121	0.00
11		2-Methylnaphthalene	1.000	0.832	16.8	114	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	125	-0.01
13	S	2,4,6-Tribromophenol	1.000	1.073	-7.3	137	0.00
14	S	2-Fluorobiphenyl	1.000	0.964	3.6	124	-0.01
15		Acenaphthylene	1.000	0.856	14.4	110	0.00
16		Acenaphthene	1.000	0.871	12.9	110	-0.01
17		Fluorene	1.000	0.801	19.9	103	-0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	109	-0.02
19		4-Bromophenyl-phenylether	1.000	0.844	15.6	94	-0.02
20		Hexachlorobenzene	1.000	0.872	12.8	99	-0.02
21		Pentachlorophenol	1.000	0.869	13.1	96	0.00
22		Phenanthrene	1.000	0.982	1.8	111	-0.02
23		Anthracene	1.000	0.863	13.7	96	0.00
24		Fluoranthene	1.000	1.093	-9.3	122	-0.02
25	I	Chrysene-d12	5.000	5.000	0.0	97	-0.01
26		Pyrene	1.000	1.178	-17.8	116	-0.01
27	S	Terphenyl-d14	1.000	1.006	-0.6	97	-0.01
28		Benzo(a)anthracene	1.000	1.014	-1.4	100	-0.01
29		Chrysene	1.000	0.957	4.3	94	-0.01
30		Bis(2-ethylhexyl)phthalate	1.000	1.048	-4.8	120	-0.01
31		Indeno(1,2,3-cd)pyrene	1.000	0.939	6.1	92	-0.04
32	I	Perylene-d12	5.000	5.000	0.0	92	-0.02
33		Benzo(b)fluoranthene	1.000	1.106	-10.6	102	-0.02
34		Benzo(k)fluoranthene	1.000	1.020	-2.0	94	-0.02
35	C	Benzo(a)pyrene	1.000	1.097	-9.7	101	-0.02
36		Dibenzo(a,h)anthracene	1.000	0.929	7.1	87	-0.03
37		Benzo(g,h,i)perylene	1.000	0.994	0.6	92	-0.04

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