Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081715\
Data File : BE090532.D
Acq On : 17 Aug 2015 13:42
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 18 04:34:14 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE081415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 14 16:52:42 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	107	0.00
2		1,4-Dioxane	1.000	1.085	-8.5	100	0.00
3		n-Nitrosodimethylamine	1.000	0.989	1.1	108	0.00
4	S	2-Fluorophenol	1.000	1.033	-3.3	110	0.00
5	S	Phenol-d6	1.000	1.040	-4.0	116	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	108	0.00
7	S	Nitrobenzene-d5	1.000	1.047	-4.7	119	0.00
8		Nitrobenzene	1.000	1.039	-3.9	120	0.00
9		Naphthalene	1.000	0.985	1.5	107	0.00
10		Hexachlorobutadiene	1.000	0.978	2.2	106	0.00
11		2-Methylnaphthalene	1.000	1.035	-3.5	115	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	115	0.00
13	S	2,4,6-Tribromophenol	1.000	1.103	-10.3	139	0.00
14	S	2-Fluorobiphenyl	1.000	0.995	0.5	114	0.00
15		Acenaphthylene	1.000	1.009	-0.9	118	0.00
16		Acenaphthene	1.000	0.994	0.6	115	0.00
17		Fluorene	1.000	1.017	-1.7	117	0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	116	0.00
19		4-Bromophenyl-phenylether	1.000	1.021	-2.1	122	0.00
20		Hexachlorobenzene	1.000	0.996	0.4	116	0.00
21		Pentachlorophenol	1.000	0.846	15.4	123	0.00
22		Phenanthrene	1.000	0.987	1.3	117	0.00
23		Anthracene	1.000	1.014	-1.4	122	0.00
24		Fluoranthene	1.000	1.019	-1.9	121	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	120	0.00
26		Pyrene	1.000	1.004	-0.4	124	0.00
27	S	Terphenyl-d14	1.000	1.034	-3.4	126	0.00
28		Benzo(a)anthracene	1.000	1.012	-1.2	128	0.00
29		Chrysene	1.000	1.013	-1.3	123	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.331	-33.1#	194	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.045	-4.5	136	0.00
32	I	Perylene-d12	5.000	5.000	0.0	126	0.00
33		Benzo(b)fluoranthene	1.000	1.033	-3.3	133	0.00
34		Benzo(k)fluoranthene	1.000	1.046	-4.6	131	0.00
35	C	Benzo(a)pyrene	1.000	1.061	-6.1	138	0.00
36		Dibenzo(a,h)anthracene	1.000	1.053	-5.3	140	0.00
37		Benzo(g,h,i)perylene	1.000	1.053	-5.3	135	0.00

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