Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE032816\
Data File : BE091543.D
Acq On : 28 Mar 2016 18:33
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 29 01:22:14 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE032516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 25 19:00:32 2016
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	86	0.00
2		1,4-Dioxane	0.400	0.410	-2.5	83	-0.02
3		n-Nitrosodimethylamine	0.400	0.405	-1.3	87	0.00
4	S	2-Fluorophenol	0.400	0.380	5.0	87	-0.02
5	S	Phenol-d6	0.400	0.363	9.3	86	-0.02
6	I	Naphthalene-d8	5.000	5.000	0.0	87	0.00
7	S	Nitrobenzene-d5	0.400	0.373	6.8	94	-0.01
8		Naphthalene	0.400	0.395	1.3	87	-0.02
9		2-Methylnaphthalene	0.400	0.382	4.5	88	0.00
10	I	Acenaphthene-d10	5.000	5.000	0.0	85	-0.01
11	S	2,4,6-Tribromophenol	0.400	0.390	2.5	97	-0.01
12	S	2-Fluorobiphenyl	0.400	0.396	1.0	86	-0.01
13		Acenaphthylene	0.400	0.368	8.0	88	0.00
14		Acenaphthene	0.400	0.402	-0.5	86	-0.01
15		Fluorene	0.400	0.381	4.8	86	-0.01
16	I	Phenanthrene-d10	5.000	5.000	0.0	87	0.00
17		Hexachlorobenzene	0.400	0.383	4.3	88	0.00
18		Pentachlorophenol	0.400	0.441	-10.2	93	-0.02
19		Phenanthrene	0.400	0.394	1.5	86	0.00
20		Anthracene	0.400	0.366	8.5	89	0.00
21		Fluoranthene	0.400	0.370	7.5	88	-0.02
22	I	Chrysene-d12	5.000	5.000	0.0	90	0.00
23		Pyrene	0.400	0.357	10.8	89	0.00
24	S	Terphenyl-d14	0.400	0.361	9.8	88	0.00
25		Benzo(a)anthracene	0.400	0.356	11.0	88	0.00
26		Chrysene	0.400	0.374	6.5	84	0.00
27		Indeno(1,2,3-cd)pyrene	0.400	0.355	11.3	90	-0.01
28	I	Perylene-d12	5.000	5.000	0.0	88	-0.01
29		Benzo(b)fluoranthene	0.400	0.397	0.8	99	0.00
30		Benzo(k)fluoranthene	0.400	0.345	13.8	80	0.00
31	C	Benzo(a)pyrene	0.400	0.487	-21.7#	92	-0.01
32		Dibenzo(a,h)anthracene	0.400	0.363	9.3	90	-0.01
33		Benzo(g,h,i)perylene	0.400	0.365	8.8	88	-0.01

(#) = Out of Range SPCC's out = 0 CCC's out = 1