Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\Data\BE042916\
Data File : BE091732.D
Acq On : 29 Apr 2016 12:57
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Apr 29 23:12:25 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE042616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 26 16:09:46 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	97	0.00
2		1,4-Dioxane	0.400	0.361	9.8	84	0.00
3		n-Nitrosodimethylamine	0.400	0.348	13.0	86	0.02
4	S	2-Fluorophenol	0.400	0.358	10.5	89	0.00
5	S	Phenol-d6	0.400	0.341	14.8	87	0.02
6		bis(2-Chloroethyl)ether	0.400	0.386	3.5	92	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	98	0.00
8	S	Nitrobenzene-d5	0.400	0.335	16.3	88	0.01
9		Nitrobenzene	0.400	0.327	18.3	87	0.00
10		Naphthalene	0.400	0.400	0.0	97	0.00
11		Hexachlorobutadiene	0.400	0.404	-1.0	101	0.00
12		2-Methylnaphthalene	0.400	0.382	4.5	94	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	95	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.406	-1.5	77	0.00
15	S	2-Fluorobiphenyl	0.400	0.402	-0.5	96	0.00
16		Acenaphthylene	0.400	0.388	3.0	104	0.00
17		Acenaphthene	0.400	0.389	2.8	92	0.00
18		Fluorene	0.400	0.388	3.0	92	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	94	0.00
20		4-Bromophenyl-phenylether	0.400	0.379	5.3	88	0.00
21		Hexachlorobenzene	0.400	0.406	-1.5	95	0.00
22		Pentachlorophenol	0.400	0.333	16.8	76	0.00
23		Phenanthrene	0.400	0.412	-3.0	94	0.00
24		Anthracene	0.400	0.379	5.3	90	0.00
25		Fluoranthene	0.400	0.389	2.8	90	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	85	0.00
27		Benzidine	0.400	0.433	-8.2	102	0.00
28		Pyrene	0.400	0.406	-1.5	90	0.00
29	S	Terphenyl-d14	0.400	0.400	0.0	87	0.00
30		Benzo(a)anthracene	0.400	0.390	2.5	85	0.00
31		3,3'-Dichlorobenzidine	0.400	0.361	9.8	89	-0.02
32		Chrysene	0.400	0.465	-16.3	94	0.00
33		Bis(2-ethylhexyl)phthalate	0.400	0.404	-1.0	112	-0.02
34		Indeno(1,2,3-cd)pyrene	0.400	0.452	-13.0	95	-0.01
35	I	Perylene-d12	5.000	5.000	0.0	95	0.00
36		Benzo(b)fluoranthene	0.400	0.421	-5.2	103	0.00
37		Benzo(k)fluoranthene	0.400	0.351	12.3	87	0.00
38	C	Benzo(a)pyrene	0.400	0.384	4.0	96	0.00
39		Dibenzo(a,h)anthracene	0.400	0.384	4.0	94	-0.02
40		Benzo(g,h,i)perylene	0.400	0.378	5.5	92	-0.02

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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