Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042916\
Data File : BE091739.D
Acq On : 29 Apr 2016 19:07
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Apr 29 23:55:12 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	93	0.00
2		1,4-Dioxane	0.400	0.347	13.3	78	0.02
3		n-Nitrosodimethylamine	0.400	0.328	18.0	78	0.02
4	S	2-Fluorophenol	0.400	0.348	13.0	83	0.00
5	S	Phenol-d6	0.400	0.328	18.0	81	0.02
6		bis(2-Chloroethyl)ether	0.400	0.376	6.0	86	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	91	0.00
8	S	Nitrobenzene-d5	0.400	0.327	18.3	80	0.01
9		Nitrobenzene	0.400	0.318	20.5	79	0.00
10		Naphthalene	0.400	0.407	-1.7	91	0.00
11		Hexachlorobutadiene	0.400	0.404	-1.0	94	-0.02
12		2-Methylnaphthalene	0.400	0.382	4.5	87	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	87	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.408	-2.0	71	0.00
15	S	2-Fluorobiphenyl	0.400	0.399	0.3	88	0.00
16		Acenaphthylene	0.400	0.382	4.5	94	0.00
17		Acenaphthene	0.400	0.384	4.0	84	0.00
18		Fluorene	0.400	0.391	2.3	85	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	88	0.00
20		4-Bromophenyl-phenylether	0.400	0.370	7.5	80	0.00
21		Hexachlorobenzene	0.400	0.400	0.0	88	0.00
22		Pentachlorophenol	0.400	0.310	22.5	64	0.00
23		Phenanthrene	0.400	0.406	-1.5	87	0.00
24		Anthracene	0.400	0.376	6.0	84	0.00
25		Fluoranthene	0.400	0.405	-1.3	88	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	78	0.00
27		Benzidine	0.400	0.374	6.5	72	0.00
28		Pyrene	0.400	0.432	-8.0	87	0.00
29	S	Terphenyl-d14	0.400	0.421	-5.2	84	0.00
30		Benzo(a)anthracene	0.400	0.393	1.8	78	0.00
31		3,3'-Dichlorobenzidine	0.400	0.337	15.8	75	-0.02
32		Chrysene	0.400	0.465	-16.3	85	0.00
33		Bis(2-ethylhexyl)phthalate	0.400	0.385	3.8	97	-0.02
34		Indeno(1,2,3-cd)pyrene	0.400	0.457	-14.2	87	-0.01
35	I	Perylene-d12	5.000	5.000	0.0	88	-0.01
36		Benzo(b)fluoranthene	0.400	0.421	-5.2	96	-0.01
37		Benzo(k)fluoranthene	0.400	0.348	13.0	80	0.00
38	C	Benzo(a)pyrene	0.400	0.376	6.0	87	0.00
39		Dibenzo(a,h)anthracene	0.400	0.381	4.8	86	-0.02
40		Benzo(g,h,i)perylene	0.400	0.376	6.0	86	-0.02

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

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