Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\Data\BE041316\
Data File : BE091651.D
Acq On : 13 Apr 2016 10:21
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Apr 14 03:57:40 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE040816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 08 16:48:14 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	88	0.00
2		1,4-Dioxane	0.400	0.378	5.5	85	0.00
3		n-Nitrosodimethylamine	0.400	0.362	9.5	86	0.00
4	S	2-Fluorophenol	0.400	0.340	15.0	79	0.00
5	S	Phenol-d6	0.400	0.329	17.8	79	0.00
6		bis(2-Chloroethyl)ether	0.400	0.375	6.3	87	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	87	-0.02
8	S	Nitrobenzene-d5	0.400	0.318	20.5	78	0.00
9		Nitrobenzene	0.400	0.426	-6.5	79	0.00
10		Naphthalene	0.400	0.396	1.0	88	0.00
11		Hexachlorobutadiene	0.400	0.399	0.3	88	0.00
12		2-Methylnaphthalene	0.400	0.377	5.8	85	-0.01
13	I	Acenaphthene-d10	5.000	5.000	0.0	82	0.00
14	S	2,4,6-Tribromophenol	0.400	0.452	-13.0	73	0.00
15	S	2-Fluorobiphenyl	0.400	0.409	-2.2	85	0.00
16		Acenaphthylene	0.400	0.365	8.8	92	0.00
17		Acenaphthene	0.400	0.415	-3.7	88	-0.01
18		Fluorene	0.400	0.384	4.0	81	-0.01
19	I	Phenanthrene-d10	5.000	5.000	0.0	88	0.00
20		4-Bromophenyl-phenylether	0.400	0.352	12.0	80	0.00
21		Hexachlorobenzene	0.400	0.376	6.0	83	0.00
22		Pentachlorophenol	0.400	0.368	8.0	66	0.00
23		Phenanthrene	0.400	0.375	6.3	83	0.00
24		Anthracene	0.400	0.340	15.0	79	0.00
25		Fluoranthene	0.400	0.332	17.0	79	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	80	0.00
27		Benzidine	0.400	0.479	-19.7	69	0.00
28		Pyrene	0.400	0.366	8.5	80	0.00
29	S	Terphenyl-d14	0.400	0.405	-1.3	83	0.00
30		Benzo(a)anthracene	0.400	0.368	8.0	77	0.00
31		3,3'-Dichlorobenzidine	0.400	0.323	19.3	55	0.00
32		Chrysene	0.400	0.464	-16.0	95	-0.02
33		Bis(2-ethylhexyl)phthalate	0.400	0.335	16.3	39	0.00
34		Indeno(1,2,3-cd)pyrene	0.400	0.331	17.3	69	-0.01
35	I	Perylene-d12	5.000	5.000	0.0	78	-0.01
36		Benzo(b)fluoranthene	0.400	0.359	10.3	73	-0.01
37		Benzo(k)fluoranthene	0.400	0.394	1.5	80	-0.01
38	C	Benzo(a)pyrene	0.400	0.326	18.5	68	0.00
39		Dibenzo(a,h)anthracene	0.400	0.339	15.3	70	-0.01
40		Benzo(g,h,i)perylene	0.400	0.361	9.8	73	-0.01

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

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