Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE061616\
Data File : BE091940.D
Acq On : 16 Jun 2016 14:27
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 16 18:20:21 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE061416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 15 12:00:38 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	81	0.00
2		1,4-Dioxane	0.400	0.392	2.0	77	0.00
3		n-Nitrosodimethylamine	0.400	0.344	14.0	69	0.04
4	S	2-Fluorophenol	0.400	0.405	-1.3	83	0.00
5	S	Phenol-d6	0.400	0.446	-11.5	84	0.02
6		bis(2-Chloroethyl)ether	0.400	0.351	12.3	71	0.02
7		n-Nitroso-di-n-propylamine	0.400	0.432	-8.0	77	0.02
8	I	Naphthalene-d8	5.000	5.000	0.0	84	0.03
9	S	Nitrobenzene-d5	0.400	0.387	3.3	81	0.02
10		Nitrobenzene	0.400	0.374	6.5	77	0.02
11		bis(2-Chloroethoxy)methane	0.400	0.408	-2.0	80	0.00
12		Naphthalene	0.400	0.392	2.0	77	0.00
13		Hexachlorobutadiene	0.400	0.382	4.5	75	0.00
14		2-Methylnaphthalene	0.400	0.389	2.8	79	0.01
15	I	Acenaphthene-d10	5.000	5.000	0.0	90	0.00
16	S	2,4,6-Tribromophenol	0.400	0.375	6.3	90	0.00
17	S	2-Fluorobiphenyl	0.400	0.374	6.5	78	0.00
18		Acenaphthylene	0.400	0.356	11.0	80	0.00
19		2,6-Dinitrotoluene	0.400	0.513	-28.2#	121	0.00
20		Acenaphthene	0.400	0.358	10.5	80	0.00
21		2,4-Dinitrotoluene	0.400	0.367	8.3	78	0.03
22		Fluorene	0.400	0.363	9.3	81	0.02
23		4-Chlorophenyl-phenylether	0.400	0.371	7.3	79	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	87	0.00
25		4-Bromophenyl-phenylether	0.400	0.373	6.8	80	0.02
26		Hexachlorobenzene	0.400	0.374	6.5	78	0.00
27		Pentachlorophenol	0.400	0.383	4.3	85	0.02
28		Phenanthrene	0.400	0.389	2.8	81	0.00
29		Anthracene	0.400	0.377	5.8	83	0.00
30		Fluoranthene	0.400	0.365	8.8	78	0.00
31	I	Chrysene-d12	5.000	5.000	0.0	87	0.00
32		Benzidine	0.400	0.375	6.3	98	0.00
33		Pyrene	0.400	0.360	10.0	80	0.00
34	S	Terphenyl-d14	0.400	0.357	10.8	75	0.00
35		Benzo(a)anthracene	0.400	0.481	-20.2	107	0.00
36		3,3'-Dichlorobenzidine	0.400	0.395	1.3	103	0.03
37		Chrysene	0.400	0.443	-10.7	87	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	0.571	-42.7#	119	-0.03
39		Indeno(1,2,3-cd)pyrene	0.400	0.362	9.5	80	0.00
40	I	Perylene-d12	5.000	5.000	0.0	88	-0.01
41		Benzo(b)fluoranthene	0.400	0.360	10.0	76	-0.01

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.406	-1.5	89	-0.01
43 C	Benzo(a)pyrene	0.400	0.386	3.5	85	0.00
44	Dibenzo(a,h)anthracene	0.400	0.367	8.3	80	-0.02
45	Benzo(g,h,i)perylene	0.400	0.368	8.0	80	-0.02

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