Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE052316\
Data File : BE091852.D
Acq On : 23 May 2016 18:06
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 24 05:23:40 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE051916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 19 15:36:49 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	114	-0.02
2		1,4-Dioxane	0.400	0.416	-4.0	100	-0.02
3		n-Nitrosodimethylamine	0.400	0.372	7.0	115	0.00
4	S	2-Fluorophenol	0.400	0.397	0.8	120	0.00
5	S	Phenol-d6	0.400	0.397	0.8	126	0.00
6		bis(2-Chloroethyl)ether	0.400	0.397	0.8	123	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	119	-0.02
8	S	Nitrobenzene-d5	0.400	0.407	-1.7	136	-0.01
9		Nitrobenzene	0.400	0.395	1.3	136	-0.01
10		Naphthalene	0.400	0.377	5.8	117	0.00
11		Hexachlorobutadiene	0.400	0.384	4.0	119	-0.02
12		2-Methylnaphthalene	0.400	0.384	4.0	121	-0.02
13	I	Acenaphthene-d10	5.000	5.000	0.0	114	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.372	7.0	122	-0.01
15	S	2-Fluorobiphenyl	0.400	0.387	3.3	116	-0.01
16		Acenaphthylene	0.400	0.403	-0.8	121	-0.01
17		Acenaphthene	0.400	0.374	6.5	113	-0.01
18		Fluorene	0.400	0.377	5.8	114	-0.02
19	I	Phenanthrene-d10	5.000	5.000	0.0	121	-0.01
20		4-Bromophenyl-phenylether	0.400	0.367	8.3	114	-0.02
21		Hexachlorobenzene	0.400	0.353	11.8	109	0.00
22		Pentachlorophenol	0.400	0.236	41.0#	80	0.00
23		Phenanthrene	0.400	0.313	21.8	101	-0.01
24		Anthracene	0.400	0.351	12.3	110	-0.01
25		Fluoranthene	0.400	0.333	16.8	101	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	100	-0.02
27		Benzidine	0.400	0.657	-64.3#	186	-0.02
28		Pyrene	0.400	0.389	2.8	102	-0.02
29	S	Terphenyl-d14	0.400	0.439	-9.7	118	-0.02
30		Benzo(a)anthracene	0.400	0.438	-9.5	118	-0.02
31		3,3'-Dichlorobenzidine	0.400	0.365	8.8	105	-0.02
32		Chrysene	0.400	0.425	-6.2	98	-0.02
33		Bis(2-ethylhexyl)phthalate	0.400	0.681	-70.3#	183	-0.02
34		Indeno(1,2,3-cd)pyrene	0.400	0.380	5.0	101	-0.05
35	I	Perylene-d12	5.000	5.000	0.0	96	-0.02
36		Benzo(b)fluoranthene	0.400	0.422	-5.5	107	-0.02
37		Benzo(k)fluoranthene	0.400	0.371	7.3	94	-0.02
38	C	Benzo(a)pyrene	0.400	0.406	-1.5	103	-0.02
39		Dibenzo(a,h)anthracene	0.400	0.403	-0.8	103	-0.06
40		Benzo(g,h,i)perylene	0.400	0.398	0.5	101	-0.06

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

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