Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102315\
Data File : BE091035.D
Acq On : 23 Oct 2015 15:55
Operator : UM/NP
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Oct 26 15:44:51 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 07 17:11:43 2015
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	51	-0.04
2		1,4-Dioxane	1.000	0.759	24.1	42	-0.13
3		n-Nitrosodimethylamine	1.000	1.290	-29.0#	66	-0.21
4	S	2-Fluorophenol	1.000	1.156	-15.6	62	-0.06
5	S	Phenol-d6	1.000	1.070	-7.0	61	-0.07
6		bis(2-Chloroethyl)ether	1.000	1.905	-90.5#	89	-0.09
7	I	Naphthalene-d8	5.000	5.000	0.0	51	-0.05
8	S	Nitrobenzene-d5	1.000	1.347	-34.7#	79	-0.06
9		Nitrobenzene	1.000	1.346	-34.6#	83	-0.06
10		Naphthalene	1.000	1.008	-0.8	51	-0.05
11		Hexachlorobutadiene	1.000	0.842	15.8	42	-0.04
12		2-Methylnaphthalene	1.000	1.041	-4.1	56	-0.05
13	I	Acenaphthene-d10	5.000	5.000	0.0	48	-0.04
14	S	2,4,6-Tribromophenol	1.000	0.932	6.8	55	-0.04
15	S	2-Fluorobiphenyl	1.000	0.986	1.4	51	-0.05
16		Acenaphthylene	1.000	0.977	2.3	52	-0.05
17		Acenaphthene	1.000	0.954	4.6	49	-0.04
18		Fluorene	1.000	0.999	0.1	52	-0.05
19	I	Phenanthrene-d10	5.000	5.000	0.0	51	-0.05
20		4-Bromophenyl-phenylether	1.000	0.887	11.3	49	-0.05
21		Hexachlorobenzene	1.000	0.119	88.1#	11	-0.05
22		Pentachlorophenol	1.000	0.865	13.5	53	-0.04
23		Phenanthrene	1.000	0.988	1.2	52	-0.05
24		Anthracene	1.000	1.017	-1.7	56	-0.05
25		Fluoranthene	1.000	1.023	-2.3	57	-0.05
26	I	Chrysene-d12	5.000	5.000	0.0	54	-0.04
27		Benzidine	1.000	1.891	-89.1#	119	-0.07
28		Pyrene	1.000	0.945	5.5	56	-0.05
29	S	Terphenyl-d14	1.000	1.117	-11.7	66	-0.05
30		Benzo(a)anthracene	1.000	1.011	-1.1	57	-0.04
31		3,3'-Dichlorobenzidine	1.000	0.000	100.0#	0	-20.84#
32		Chrysene	1.000	0.908	9.2	51	-0.04
33		Bis(2-ethylhexyl)phthalate	1.000	1.088	-8.8	73	-0.04
34		Indeno(1,2,3-cd)pyrene	1.000	1.051	-5.1	62	-0.09
35	I	Perylene-d12	5.000	5.000	0.0	61	-0.06
36		Benzo(b)fluoranthene	1.000	1.030	-3.0	70	-0.05
37		Benzo(k)fluoranthene	1.000	0.837	16.3	55	-0.05
38	C	Benzo(a)pyrene	1.000	1.039	-3.9	74	-0.05
39		Dibenzo(a,h)anthracene	1.000	1.127	-12.7	82	-0.09
40		Benzo(g,h,i)perylene	1.000	1.123	-12.3	79	-0.09

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

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