Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM010216\
Data File : BM003922.D
Acq On : 31 Dec 2015 19:52
Operator : SJ/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Jan 04 13:59:14 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM010216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 04 13:33:21 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	1.000	1.021	-2.1	95	0.00
3		n-Nitrosodimethylamine	1.000	0.973	2.7	91	0.00
4	S	2-Fluorophenol	1.000	0.890	11.0	88	0.00
5	S	Phenol-d6	1.000	0.927	7.3	88	0.00
6		bis(2-Chloroethyl)ether	1.000	0.993	0.7	91	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	92	0.00
8	S	Nitrobenzene-d5	1.000	0.931	6.9	87	0.00
9		Nitrobenzene	1.000	0.924	7.6	87	0.00
10		Naphthalene	1.000	0.989	1.1	92	-0.02
11		Hexachlorobutadiene	1.000	0.963	3.7	91	0.00
12		2-Methylnaphthalene	1.000	0.968	3.2	90	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	90	0.00
14	S	2,4,6-Tribromophenol	1.000	0.882	11.8	82	0.00
15	S	2-Fluorobiphenyl	1.000	1.004	-0.4	91	0.00
16		Acenaphthylene	1.000	0.975	2.5	87	0.00
17		Acenaphthene	1.000	0.977	2.3	93	0.00
18		Fluorene	1.000	1.017	-1.7	90	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	90	0.00
20		4-Bromophenyl-phenylether	1.000	0.893	10.7	86	0.00
21		Hexachlorobenzene	1.000	0.903	9.7	90	0.00
22		Pentachlorophenol	1.000	0.793	20.7	72	0.00
23		Phenanthrene	1.000	0.953	4.7	85	0.00
24		Anthracene	1.000	0.904	9.6	88	0.00
25		Fluoranthene	1.000	0.891	10.9	84	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	90	0.00
27		Benzidine	1.000	0.734	26.6#	60	0.00
28		Pyrene	1.000	0.934	6.6	83	0.00
29	S	Terphenyl-d14	1.000	0.916	8.4	83	-0.02
30		Benzo(a)anthracene	1.000	0.993	0.7	92	0.00
31		3,3'-Dichlorobenzidine	1.000	0.789	21.1	69	0.00
32		Chrysene	1.000	0.797	20.3	71	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	0.687	31.3#	55	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	0.976	2.4	88	0.00
35	I	Perylene-d12	5.000	5.000	0.0	85	0.00
36		Benzo(b)fluoranthene	1.000	0.958	4.2	88	0.00
37		Benzo(k)fluoranthene	1.000	0.898	10.2	79	0.00
38	C	Benzo(a)pyrene	1.000	0.916	8.4	83	0.00
39		Dibenzo(a,h)anthracene	1.000	1.023	-2.3	88	0.00
40		Benzo(g,h,i)perylene	1.000	1.023	-2.3	88	0.00

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

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