

Study Number: I14013B

Test Type: TOX

Route: Application

Species/Strain: Mouse/Taconic BALB/c

C Number:

Study Gender:

PWG Approval Date

I04G: Mean Body Weight Gain

Test Compound: 4-Methylcyclohexanemethanol Crude

CAS Number: CRUDEMCHM

I14013B

Female

See web page for date of PWG Approval

Date Report Requested: 09/05/2019

Time Report Requested: 08:36:59

Lab: NTP

Study Number: I14013B

Test Type: TOX

Route: Application

Species/Strain: Mouse/Taconic BALB/c

I04G: Mean Body Weight Gain

Test Compound: 4-Methylcyclohexanemethanol Crude

CAS Number: CRUDEMCHM

Date Report Requested: 09/05/2019

Time Report Requested: 08:36:59

Lab: NTP

Females

Treatment Groups (%)

Phase	Litter ID	Days	0		1		5		25		50	
			Wt Gain (g)	N	Wt Gain (g)	N	Wt Gain (g)	N	Wt Gain (g)	N	Wt Gain (g)	N
SD		1 - 6	0.2 ± 0.1 *	13	0.1 ± 0.1	13	0.5 ± 0.1	13	0.3 ± 0.1	13	0.4 ± 0.1	13

Study Number: I14013B

Test Type: TOX

Route: Application

Species/Strain: Mouse/Taconic BALB/c

I04G: Mean Body Weight Gain

Test Compound: 4-Methylcyclohexanemethanol Crude

CAS Number: CRUDEMCHM

Date Report Requested: 09/05/2019

Time Report Requested: 08:36:59

Lab: NTP

Females

Phase	Litter ID	Days	Treatment Groups (%)			
			75		0.15% DNFB	
			Wt Gain (g)	N	Wt Gain (g)	N
SD		1 - 6	0.5 ± 0.1	13	0.3 ± 0.1	13

Study Number: I14013B

Test Type: TOX

Route: Application

Species/Strain: Mouse/Taconic BALB/c

I04G: Mean Body Weight Gain

Test Compound: 4-Methylcyclohexanemethanol Crude

CAS Number: CRUDEMCHM

Date Report Requested: 09/05/2019

Time Report Requested: 08:36:59

Lab: NTP

LEGEND

Data are displayed as mean \pm SEM

SD – Study Day

Statistical analysis of weight data performed by Jonckheere (trend) and Williams or Dunnett (pairwise) tests.

Statistical significance for the control group indicates a significant trend test

Statistical significance for a treatment group indicates a significant pairwise test compared to the vehicle control group

* Statistically significant at $P \leq 0.05$

** Statistically significant at $P \leq 0.01$

DNFB = 1-Fluoro-2,4 -dinitrofluorobenzene

**** END OF REPORT ****