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straight_chain_alkane(1,[carb(h,h,h,h)]).

straight_chain_alkane(2,[carb(h,h,h,c),carb(c,h,h,h)]).

straight_chain_alkane(N,[H,carb(c,h,h,c) | T]):-

N>2,

N2 is N-1,

straight_chain_alkane(N2,[H | T]).

branched_alkane(4,[carb(h,h,h,c),carb(c,c1h3,h,c),carb(c,h,h,h)]).

branched_alkane(N,BA):-

N>4,

N1 is N - 3,

mc_generator(N1,1,Mid ,Com),

mid_chain(Mid ,A),

N2 is Mid+2,

sums(Com ,Com1),

constructer(A,Com1,B),

valid_check(B,N2),

append([carb(h,h,h,c)],B,BB),

append(BB,[carb(c,h,h,h)],BA).

constructer(Mid ,Nums,Ans):-

findall(A,num_seperate(Mid , Nums,A),A2),

dup_remover(A2,A3),

break_list(A3,Ans).

dup_remover([],[]).

dup_remover([Elm | T],A):-

member(Elm,T),

dup_remover(T,A),

!.

dup_remover([Elm|T],Ans):-

\+ member(Elm,T),

dup_remover(T,A),

Ans = [Elm|A],

!.

num_seperate([carb(c,h,h,c)],[1,1],[carb(c,c1h3,c1h3,c)]).

num_seperate(ML,[T],Ans):-

carb_attacher(ML,T,Ans).

num_seperate(Mid,[H|T],A):-

T\=[],

carb_attacher(Mid,H,Out),

num_seperate(Out,T,A).

carb_attacher(H,0,H).

carb_attacher([H],X,A):-

add_branch_to_carbon(H,X,A).

carb_attacher([H|T],X,[H|Z]):-

carb_attacher(T,X,L),

flatten(L,Z).

carb_attacher([H|T],X,Sol2):-

add_branch_to_carbon(H,X,M),

\+ member(M,T),

T\==[],

Sol = [M|T],

flatten(Sol,Sol2).

mid_chain(1,[carb(c,h,h,c)]).

mid_chain(N,[carb(c,h,h,c)|T]):-

N > 1,

N2 is N - 1,
mid_chain(N2 ,T).

valid_check(L,MN):-

valid(L,MN,1),
reverse(L,L1),
valid(L1,MN,1).

add_branch_to_carbon(A,0,A).

add_branch_to_carbon(carb(c,h,h,c),N,carb(c,A,h,c)):-

branch_name(N,A).

add_branch_to_carbon(carb(c,NH,h,c),N,carb(c,NH,A,c)):-

NH \= h,
branch_name(N,A),
N2 is N-1,
(add_branch_to_carbon(carb(c,NH,h,c),N2,carb(c,NH,A2,c)),A2
\= NH).

mc_generator(N,Acc,N,Acc):-

N1 is (N+1)*2,
Acc=<N1.
% Acc=<N*2.

mc_generator(N,Acc,N1,N2):-

NC is (N+1)*2,
Acc=<NC,
% Acc=<N*2,
NA is N -1,
NB is Acc +1,
mc_generator(NA,NB,N1,N2).

breakdown(0,[]).

breakdown(N,[H|T]) :-

 range(1,N,H),

 M is N - H,

 breakdown(M,T).

range(Low,High,_) :-

 Low > High,

 !,

 fail.

range(Low,_,Low).

range(Low,High,Out) :-

 Current is Low + 1,

 range(Current,High,Out).

branch_name(0,h).

branch_name(S,N):-

 S > 0,

 HS is S*2+1,

 atomic_list_concat([c,S,h,HS],N).

extractor(S,N):-

 sub_atom(S, 1, 1, 2, A),

 atom_number(A,N).

valid([],_,_).

valid([carb(c,h,h,c)|T],MN,N):-

 N1 is N+1,

 valid(T,MN,N1).

valid([carb(c,X,_c) | T],MN,N):-

 X\==h,
 extractor(X,N1),
 N2 is N+1,
 N3 is N1 +N2,
 MN>=N3,
 valid(T,MN,N2).

sums(A,B):-

 findall(X,breakdown(A,X),C),
 sums_help(C,[],Y),
 break_list(Y,B).

sums_help([],Y,Y).

sums_help([H | T],Y,B):-

 permutation(H,X),
 member(X,T),
 sums_help(T,Y,B),
 !.

sums_help([H | T],Y,B):-

 permutation(H,X),
 \+member(X,T),
 sums_help(T,[H | Y],B),
 !.

break_list([],_):=

 fail.

break_list([H | _],H).

break_list([_ | T],Ans):-

 break_list(T,Ans).

isomers(N,A):-

N<4,

straight_chain_alkane(N,A)

.

isomers(N,A):-

N>3,

findall(B,branched_alkane(N,B),C),

remove_ht(C,[],L),

isomer_help(L,[],X),

straight_chain_alkane(N,X2),

A=[X2|X].

remove_ht([],A,A).

remove_ht([[_|T]|T2],Acc,Out):-

last(T,Last),

delete(T,Last,NL),

remove_ht(T2,[NL|Acc],Out).

isomer_help([],A,A).

isomer_help([H|T],A,B):-

(\+member(H,T),

reverse(H,H1),

\+member(H1,T)),

append([carb(h,h,h,c)],H1,X),

append(X,[carb(c,h,h,h)],Y),

isomer_help(T,[Y|A],B).

isomer_help([H|T],A,B):-

(member(H,T);

reverse(H,H1),

member(H1,T)),

isomer_help(T,A,B).