



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2023 – 11:56 PM EDT

PDB ID : 7JGL
Title : Crystal Structure of the Zn-bound Human Heavy-chain variant 122H-delta C-star with 2,5-furandihydroxamate collected at 100K
Authors : Bailey, J.B.; Tezcan, F.A.
Deposited on : 2020-07-19
Resolution : 2.34 Å (reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.35.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35.1

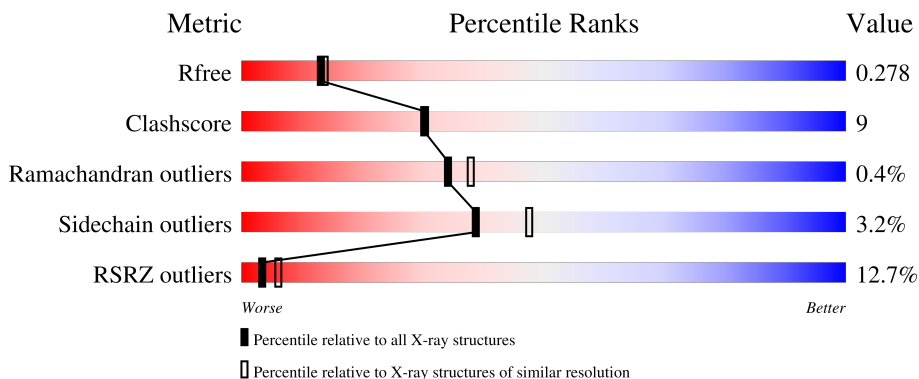
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.34 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2096 (2.36-2.32)
Clashscore	141614	2193 (2.36-2.32)
Ramachandran outliers	138981	2159 (2.36-2.32)
Sidechain outliers	138945	2160 (2.36-2.32)
RSRZ outliers	127900	2067 (2.36-2.32)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	182	
1	B	182	
1	C	182	
1	D	182	
1	E	182	

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Mol	Chain	Length	Quality of chain
1	F	182	2% 73% 21% . .
1	G	182	75% 21% .
1	H	182	73% 20% . .
1	I	182	74% 21% . .
1	J	182	81% 14% . .
1	K	182	79% 16% . .
1	L	182	73% 22% . .
1	M	182	81% 14% . .
1	N	182	76% 19% . .
1	O	182	76% 19% .
1	P	182	75% 20% .
1	Q	182	74% 20% . .
1	R	182	2% 70% 24% . .
1	S	182	75% 20% .
1	T	182	4% 68% 27% . .
1	U	182	71% 23% . .
1	V	182	81% 14% .
1	W	182	76% 18% . .
1	X	182	73% 23% . .
1	a	182	23% 96% .
1	b	182	24% 95% . .
1	c	182	23% 94% . .
1	d	182	23% 95% . .
1	e	182	28% 96% .
1	f	182	21% 95% . .

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Mol	Chain	Length	Quality of chain
1	g	182	25% 96% .
1	h	182	21% 95% ..
1	i	182	20% 95% ..
1	j	182	23% 95% ..
1	k	182	22% 95% ..
1	l	182	23% 95% ..
1	m	182	21% 96% .
1	n	182	21% 96% .
1	o	182	29% 95% ..
1	p	182	27% 94% ..
1	q	182	24% 96% .
1	r	182	25% 95% ..
1	s	182	20% 95% ..
1	t	182	27% 95% ..
1	u	182	29% 95% ..
1	v	182	24% 96% .
1	w	182	24% 95% ..
1	x	182	16% 95% ..

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 105214 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Ferritin heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	174	1431	896	252	279	4	0	0	0
1	B	174	1431	896	252	279	4	0	0	0
1	C	174	1431	896	252	279	4	0	0	0
1	D	174	1431	896	252	279	4	0	0	0
1	E	174	1431	896	252	279	4	0	0	0
1	F	174	1431	896	252	279	4	0	0	0
1	G	174	1431	896	252	279	4	0	0	0
1	H	174	1431	896	252	279	4	0	0	0
1	I	174	1431	896	252	279	4	0	0	0
1	J	174	1431	896	252	279	4	0	0	0
1	K	174	1431	896	252	279	4	0	0	0
1	L	174	1431	896	252	279	4	0	0	0
1	M	174	1431	896	252	279	4	0	0	0
1	N	174	1431	896	252	279	4	0	0	0
1	O	174	1431	896	252	279	4	0	0	0
1	P	174	1431	896	252	279	4	0	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	174	Total	C	N	O	S	0	0	0
			1431	896	252	279	4			
1	R	174	Total	C	N	O	S	0	0	0
			1431	896	252	279	4			
1	S	174	Total	C	N	O	S	0	0	0
			1431	896	252	279	4			
1	T	174	Total	C	N	O	S	0	0	0
			1431	896	252	279	4			
1	U	174	Total	C	N	O	S	0	0	0
			1431	896	252	279	4			
1	V	174	Total	C	N	O	S	0	0	0
			1431	896	252	279	4			
1	W	174	Total	C	N	O	S	0	0	0
			1431	896	252	279	4			
1	X	174	Total	C	N	O	S	0	0	0
			1431	896	252	279	4			
1	a	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	b	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	c	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	d	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	e	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	f	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	g	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	h	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	i	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	j	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	k	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	l	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	m	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	n	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	o	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	p	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	q	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	r	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	s	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	t	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	u	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	v	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	w	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			
1	x	174	Total	C	N	O	S	0	174	0
			2862	1792	504	558	8			

There are 240 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	86	GLN	LYS	engineered mutation	UNP P02794
A	90	GLU	CYS	engineered mutation	UNP P02794
A	102	ALA	CYS	engineered mutation	UNP P02794
A	122	HIS	THR	engineered mutation	UNP P02794
A	130	ALA	CYS	engineered mutation	UNP P02794
B	86	GLN	LYS	engineered mutation	UNP P02794
B	90	GLU	CYS	engineered mutation	UNP P02794
B	102	ALA	CYS	engineered mutation	UNP P02794
B	122	HIS	THR	engineered mutation	UNP P02794
B	130	ALA	CYS	engineered mutation	UNP P02794
C	86	GLN	LYS	engineered mutation	UNP P02794
C	90	GLU	CYS	engineered mutation	UNP P02794
C	102	ALA	CYS	engineered mutation	UNP P02794
C	122	HIS	THR	engineered mutation	UNP P02794
C	130	ALA	CYS	engineered mutation	UNP P02794
D	86	GLN	LYS	engineered mutation	UNP P02794
D	90	GLU	CYS	engineered mutation	UNP P02794

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Chain	Residue	Modelled	Actual	Comment	Reference
D	102	ALA	CYS	engineered mutation	UNP P02794
D	122	HIS	THR	engineered mutation	UNP P02794
D	130	ALA	CYS	engineered mutation	UNP P02794
E	86	GLN	LYS	engineered mutation	UNP P02794
E	90	GLU	CYS	engineered mutation	UNP P02794
E	102	ALA	CYS	engineered mutation	UNP P02794
E	122	HIS	THR	engineered mutation	UNP P02794
E	130	ALA	CYS	engineered mutation	UNP P02794
F	86	GLN	LYS	engineered mutation	UNP P02794
F	90	GLU	CYS	engineered mutation	UNP P02794
F	102	ALA	CYS	engineered mutation	UNP P02794
F	122	HIS	THR	engineered mutation	UNP P02794
F	130	ALA	CYS	engineered mutation	UNP P02794
G	86	GLN	LYS	engineered mutation	UNP P02794
G	90	GLU	CYS	engineered mutation	UNP P02794
G	102	ALA	CYS	engineered mutation	UNP P02794
G	122	HIS	THR	engineered mutation	UNP P02794
G	130	ALA	CYS	engineered mutation	UNP P02794
H	86	GLN	LYS	engineered mutation	UNP P02794
H	90	GLU	CYS	engineered mutation	UNP P02794
H	102	ALA	CYS	engineered mutation	UNP P02794
H	122	HIS	THR	engineered mutation	UNP P02794
H	130	ALA	CYS	engineered mutation	UNP P02794
I	86	GLN	LYS	engineered mutation	UNP P02794
I	90	GLU	CYS	engineered mutation	UNP P02794
I	102	ALA	CYS	engineered mutation	UNP P02794
I	122	HIS	THR	engineered mutation	UNP P02794
I	130	ALA	CYS	engineered mutation	UNP P02794
J	86	GLN	LYS	engineered mutation	UNP P02794
J	90	GLU	CYS	engineered mutation	UNP P02794
J	102	ALA	CYS	engineered mutation	UNP P02794
J	122	HIS	THR	engineered mutation	UNP P02794
J	130	ALA	CYS	engineered mutation	UNP P02794
K	86	GLN	LYS	engineered mutation	UNP P02794
K	90	GLU	CYS	engineered mutation	UNP P02794
K	102	ALA	CYS	engineered mutation	UNP P02794
K	122	HIS	THR	engineered mutation	UNP P02794
K	130	ALA	CYS	engineered mutation	UNP P02794
L	86	GLN	LYS	engineered mutation	UNP P02794
L	90	GLU	CYS	engineered mutation	UNP P02794
L	102	ALA	CYS	engineered mutation	UNP P02794
L	122	HIS	THR	engineered mutation	UNP P02794

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Chain	Residue	Modelled	Actual	Comment	Reference
L	130	ALA	CYS	engineered mutation	UNP P02794
M	86	GLN	LYS	engineered mutation	UNP P02794
M	90	GLU	CYS	engineered mutation	UNP P02794
M	102	ALA	CYS	engineered mutation	UNP P02794
M	122	HIS	THR	engineered mutation	UNP P02794
M	130	ALA	CYS	engineered mutation	UNP P02794
N	86	GLN	LYS	engineered mutation	UNP P02794
N	90	GLU	CYS	engineered mutation	UNP P02794
N	102	ALA	CYS	engineered mutation	UNP P02794
N	122	HIS	THR	engineered mutation	UNP P02794
N	130	ALA	CYS	engineered mutation	UNP P02794
O	86	GLN	LYS	engineered mutation	UNP P02794
O	90	GLU	CYS	engineered mutation	UNP P02794
O	102	ALA	CYS	engineered mutation	UNP P02794
O	122	HIS	THR	engineered mutation	UNP P02794
O	130	ALA	CYS	engineered mutation	UNP P02794
P	86	GLN	LYS	engineered mutation	UNP P02794
P	90	GLU	CYS	engineered mutation	UNP P02794
P	102	ALA	CYS	engineered mutation	UNP P02794
P	122	HIS	THR	engineered mutation	UNP P02794
P	130	ALA	CYS	engineered mutation	UNP P02794
Q	86	GLN	LYS	engineered mutation	UNP P02794
Q	90	GLU	CYS	engineered mutation	UNP P02794
Q	102	ALA	CYS	engineered mutation	UNP P02794
Q	122	HIS	THR	engineered mutation	UNP P02794
Q	130	ALA	CYS	engineered mutation	UNP P02794
R	86	GLN	LYS	engineered mutation	UNP P02794
R	90	GLU	CYS	engineered mutation	UNP P02794
R	102	ALA	CYS	engineered mutation	UNP P02794
R	122	HIS	THR	engineered mutation	UNP P02794
R	130	ALA	CYS	engineered mutation	UNP P02794
S	86	GLN	LYS	engineered mutation	UNP P02794
S	90	GLU	CYS	engineered mutation	UNP P02794
S	102	ALA	CYS	engineered mutation	UNP P02794
S	122	HIS	THR	engineered mutation	UNP P02794
S	130	ALA	CYS	engineered mutation	UNP P02794
T	86	GLN	LYS	engineered mutation	UNP P02794
T	90	GLU	CYS	engineered mutation	UNP P02794
T	102	ALA	CYS	engineered mutation	UNP P02794
T	122	HIS	THR	engineered mutation	UNP P02794
T	130	ALA	CYS	engineered mutation	UNP P02794
U	86	GLN	LYS	engineered mutation	UNP P02794

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Chain	Residue	Modelled	Actual	Comment	Reference
U	90	GLU	CYS	engineered mutation	UNP P02794
U	102	ALA	CYS	engineered mutation	UNP P02794
U	122	HIS	THR	engineered mutation	UNP P02794
U	130	ALA	CYS	engineered mutation	UNP P02794
V	86	GLN	LYS	engineered mutation	UNP P02794
V	90	GLU	CYS	engineered mutation	UNP P02794
V	102	ALA	CYS	engineered mutation	UNP P02794
V	122	HIS	THR	engineered mutation	UNP P02794
V	130	ALA	CYS	engineered mutation	UNP P02794
W	86	GLN	LYS	engineered mutation	UNP P02794
W	90	GLU	CYS	engineered mutation	UNP P02794
W	102	ALA	CYS	engineered mutation	UNP P02794
W	122	HIS	THR	engineered mutation	UNP P02794
W	130	ALA	CYS	engineered mutation	UNP P02794
X	86	GLN	LYS	engineered mutation	UNP P02794
X	90	GLU	CYS	engineered mutation	UNP P02794
X	102	ALA	CYS	engineered mutation	UNP P02794
X	122	HIS	THR	engineered mutation	UNP P02794
X	130	ALA	CYS	engineered mutation	UNP P02794
a	86	GLN	LYS	engineered mutation	UNP P02794
a	90	GLU	CYS	engineered mutation	UNP P02794
a	102	ALA	CYS	engineered mutation	UNP P02794
a	122	HIS	THR	engineered mutation	UNP P02794
a	130	ALA	CYS	engineered mutation	UNP P02794
b	86	GLN	LYS	engineered mutation	UNP P02794
b	90	GLU	CYS	engineered mutation	UNP P02794
b	102	ALA	CYS	engineered mutation	UNP P02794
b	122	HIS	THR	engineered mutation	UNP P02794
b	130	ALA	CYS	engineered mutation	UNP P02794
c	86	GLN	LYS	engineered mutation	UNP P02794
c	90	GLU	CYS	engineered mutation	UNP P02794
c	102	ALA	CYS	engineered mutation	UNP P02794
c	122	HIS	THR	engineered mutation	UNP P02794
c	130	ALA	CYS	engineered mutation	UNP P02794
d	86	GLN	LYS	engineered mutation	UNP P02794
d	90	GLU	CYS	engineered mutation	UNP P02794
d	102	ALA	CYS	engineered mutation	UNP P02794
d	122	HIS	THR	engineered mutation	UNP P02794
d	130	ALA	CYS	engineered mutation	UNP P02794
e	86	GLN	LYS	engineered mutation	UNP P02794
e	90	GLU	CYS	engineered mutation	UNP P02794
e	102	ALA	CYS	engineered mutation	UNP P02794

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Chain	Residue	Modelled	Actual	Comment	Reference
e	122	HIS	THR	engineered mutation	UNP P02794
e	130	ALA	CYS	engineered mutation	UNP P02794
f	86	GLN	LYS	engineered mutation	UNP P02794
f	90	GLU	CYS	engineered mutation	UNP P02794
f	102	ALA	CYS	engineered mutation	UNP P02794
f	122	HIS	THR	engineered mutation	UNP P02794
f	130	ALA	CYS	engineered mutation	UNP P02794
g	86	GLN	LYS	engineered mutation	UNP P02794
g	90	GLU	CYS	engineered mutation	UNP P02794
g	102	ALA	CYS	engineered mutation	UNP P02794
g	122	HIS	THR	engineered mutation	UNP P02794
g	130	ALA	CYS	engineered mutation	UNP P02794
h	86	GLN	LYS	engineered mutation	UNP P02794
h	90	GLU	CYS	engineered mutation	UNP P02794
h	102	ALA	CYS	engineered mutation	UNP P02794
h	122	HIS	THR	engineered mutation	UNP P02794
h	130	ALA	CYS	engineered mutation	UNP P02794
i	86	GLN	LYS	engineered mutation	UNP P02794
i	90	GLU	CYS	engineered mutation	UNP P02794
i	102	ALA	CYS	engineered mutation	UNP P02794
i	122	HIS	THR	engineered mutation	UNP P02794
i	130	ALA	CYS	engineered mutation	UNP P02794
j	86	GLN	LYS	engineered mutation	UNP P02794
j	90	GLU	CYS	engineered mutation	UNP P02794
j	102	ALA	CYS	engineered mutation	UNP P02794
j	122	HIS	THR	engineered mutation	UNP P02794
j	130	ALA	CYS	engineered mutation	UNP P02794
k	86	GLN	LYS	engineered mutation	UNP P02794
k	90	GLU	CYS	engineered mutation	UNP P02794
k	102	ALA	CYS	engineered mutation	UNP P02794
k	122	HIS	THR	engineered mutation	UNP P02794
k	130	ALA	CYS	engineered mutation	UNP P02794
l	86	GLN	LYS	engineered mutation	UNP P02794
l	90	GLU	CYS	engineered mutation	UNP P02794
l	102	ALA	CYS	engineered mutation	UNP P02794
l	122	HIS	THR	engineered mutation	UNP P02794
l	130	ALA	CYS	engineered mutation	UNP P02794
m	86	GLN	LYS	engineered mutation	UNP P02794
m	90	GLU	CYS	engineered mutation	UNP P02794
m	102	ALA	CYS	engineered mutation	UNP P02794
m	122	HIS	THR	engineered mutation	UNP P02794
m	130	ALA	CYS	engineered mutation	UNP P02794

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Chain	Residue	Modelled	Actual	Comment	Reference
n	86	GLN	LYS	engineered mutation	UNP P02794
n	90	GLU	CYS	engineered mutation	UNP P02794
n	102	ALA	CYS	engineered mutation	UNP P02794
n	122	HIS	THR	engineered mutation	UNP P02794
n	130	ALA	CYS	engineered mutation	UNP P02794
o	86	GLN	LYS	engineered mutation	UNP P02794
o	90	GLU	CYS	engineered mutation	UNP P02794
o	102	ALA	CYS	engineered mutation	UNP P02794
o	122	HIS	THR	engineered mutation	UNP P02794
o	130	ALA	CYS	engineered mutation	UNP P02794
p	86	GLN	LYS	engineered mutation	UNP P02794
p	90	GLU	CYS	engineered mutation	UNP P02794
p	102	ALA	CYS	engineered mutation	UNP P02794
p	122	HIS	THR	engineered mutation	UNP P02794
p	130	ALA	CYS	engineered mutation	UNP P02794
q	86	GLN	LYS	engineered mutation	UNP P02794
q	90	GLU	CYS	engineered mutation	UNP P02794
q	102	ALA	CYS	engineered mutation	UNP P02794
q	122	HIS	THR	engineered mutation	UNP P02794
q	130	ALA	CYS	engineered mutation	UNP P02794
r	86	GLN	LYS	engineered mutation	UNP P02794
r	90	GLU	CYS	engineered mutation	UNP P02794
r	102	ALA	CYS	engineered mutation	UNP P02794
r	122	HIS	THR	engineered mutation	UNP P02794
r	130	ALA	CYS	engineered mutation	UNP P02794
s	86	GLN	LYS	engineered mutation	UNP P02794
s	90	GLU	CYS	engineered mutation	UNP P02794
s	102	ALA	CYS	engineered mutation	UNP P02794
s	122	HIS	THR	engineered mutation	UNP P02794
s	130	ALA	CYS	engineered mutation	UNP P02794
t	86	GLN	LYS	engineered mutation	UNP P02794
t	90	GLU	CYS	engineered mutation	UNP P02794
t	102	ALA	CYS	engineered mutation	UNP P02794
t	122	HIS	THR	engineered mutation	UNP P02794
t	130	ALA	CYS	engineered mutation	UNP P02794
u	86	GLN	LYS	engineered mutation	UNP P02794
u	90	GLU	CYS	engineered mutation	UNP P02794
u	102	ALA	CYS	engineered mutation	UNP P02794
u	122	HIS	THR	engineered mutation	UNP P02794
u	130	ALA	CYS	engineered mutation	UNP P02794
v	86	GLN	LYS	engineered mutation	UNP P02794
v	90	GLU	CYS	engineered mutation	UNP P02794

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Chain	Residue	Modelled	Actual	Comment	Reference
v	102	ALA	CYS	engineered mutation	UNP P02794
v	122	HIS	THR	engineered mutation	UNP P02794
v	130	ALA	CYS	engineered mutation	UNP P02794
w	86	GLN	LYS	engineered mutation	UNP P02794
w	90	GLU	CYS	engineered mutation	UNP P02794
w	102	ALA	CYS	engineered mutation	UNP P02794
w	122	HIS	THR	engineered mutation	UNP P02794
w	130	ALA	CYS	engineered mutation	UNP P02794
x	86	GLN	LYS	engineered mutation	UNP P02794
x	90	GLU	CYS	engineered mutation	UNP P02794
x	102	ALA	CYS	engineered mutation	UNP P02794
x	122	HIS	THR	engineered mutation	UNP P02794
x	130	ALA	CYS	engineered mutation	UNP P02794

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	3	Total Zn 3 3	0	0
2	B	3	Total Zn 3 3	0	0
2	C	3	Total Zn 3 3	0	0
2	D	2	Total Zn 2 2	0	0
2	E	3	Total Zn 3 3	0	0
2	F	4	Total Zn 4 4	0	0
2	G	3	Total Zn 3 3	0	0
2	H	3	Total Zn 3 3	0	0
2	I	2	Total Zn 2 2	0	0
2	J	2	Total Zn 2 2	0	0
2	K	3	Total Zn 3 3	0	0
2	L	2	Total Zn 2 2	0	0
2	M	2	Total Zn 2 2	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	
2	N	3	Total 3	Zn 3	0	0
2	O	3	Total 3	Zn 3	0	0
2	P	3	Total 3	Zn 3	0	0
2	Q	2	Total 2	Zn 2	0	0
2	R	2	Total 2	Zn 2	0	0
2	S	2	Total 2	Zn 2	0	0
2	T	3	Total 3	Zn 3	0	0
2	U	2	Total 2	Zn 2	0	0
2	V	2	Total 2	Zn 2	0	0
2	W	3	Total 3	Zn 3	0	0
2	X	2	Total 2	Zn 2	0	0
2	a	3	Total 6	Zn 6	0	3
2	b	2	Total 4	Zn 4	0	2
2	c	2	Total 4	Zn 4	0	2
2	d	4	Total 8	Zn 8	0	4
2	e	3	Total 6	Zn 6	0	3
2	f	3	Total 6	Zn 6	0	3
2	g	3	Total 6	Zn 6	0	3
2	h	2	Total 4	Zn 4	0	2
2	i	2	Total 4	Zn 4	0	2
2	j	2	Total 4	Zn 4	0	2

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	k	2	Total 4	Zn 4	0	2
2	l	3	Total 6	Zn 6	0	3
2	m	2	Total 4	Zn 4	0	2
2	n	3	Total 6	Zn 6	0	3
2	o	2	Total 4	Zn 4	0	2
2	p	3	Total 6	Zn 6	0	3
2	q	2	Total 4	Zn 4	0	2
2	r	2	Total 4	Zn 4	0	2
2	s	3	Total 6	Zn 6	0	3
2	t	2	Total 4	Zn 4	0	2
2	u	2	Total 4	Zn 4	0	2
2	v	3	Total 6	Zn 6	0	3
2	w	4	Total 8	Zn 8	0	4
2	x	3	Total 6	Zn 6	0	3

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

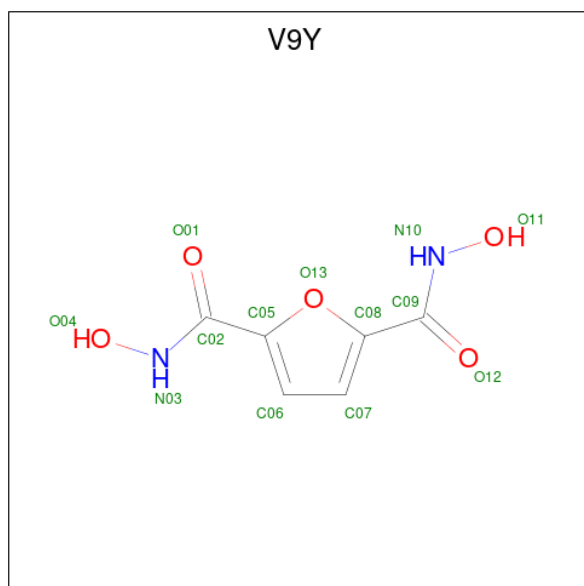
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Na 1	0	0
3	D	1	Total 1	Na 1	0	0
3	G	1	Total 1	Na 1	0	0
3	J	1	Total 1	Na 1	0	0
3	M	1	Total 1	Na 1	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	P	1	Total Na 1 1	0	0
3	S	1	Total Na 1 1	0	0
3	V	1	Total Na 1 1	0	0
3	a	1	Total Na 2 2	0	1
3	d	1	Total Na 2 2	0	1
3	g	1	Total Na 2 2	0	1
3	j	1	Total Na 2 2	0	1
3	n	1	Total Na 2 2	0	1
3	p	1	Total Na 2 2	0	1
3	s	1	Total Na 2 2	0	1
3	x	1	Total Na 2 2	0	1

- Molecule 4 is N 2 ,N 5 -dihydroxyfuran-2,5-dicarboxamide (three-letter code: V9Y) (formula: C₆H₆N₂O₅) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	1
			26	12	4	10		
4	D	1	Total	C	N	O	0	1
			26	12	4	10		
4	I	1	Total	C	N	O	0	1
			26	12	4	10		
4	L	1	Total	C	N	O	0	1
			26	12	4	10		
4	O	1	Total	C	N	O	0	1
			26	12	4	10		
4	R	1	Total	C	N	O	0	1
			26	12	4	10		
4	X	1	Total	C	N	O	0	1
			26	12	4	10		
4	t	1	Total	C	N	O	0	1
			26	12	4	10		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	92	Total	O	0	0
			92	92		
5	B	76	Total	O	0	0
			76	76		
5	C	86	Total	O	0	0
			86	86		
5	D	69	Total	O	0	0
			69	69		
5	E	78	Total	O	0	0
			78	78		
5	F	54	Total	O	0	0
			54	54		
5	G	66	Total	O	0	0
			66	66		
5	H	57	Total	O	0	0
			57	57		
5	I	59	Total	O	0	0
			59	59		
5	J	104	Total	O	0	0
			104	104		
5	K	83	Total	O	0	0
			83	83		
5	L	83	Total	O	0	0
			83	83		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	M	94	Total O 94 94	0	0
5	N	64	Total O 64 64	0	0
5	O	90	Total O 90 90	0	0
5	P	56	Total O 56 56	0	0
5	Q	54	Total O 54 54	0	0
5	R	58	Total O 58 58	0	0
5	S	67	Total O 67 67	0	0
5	T	61	Total O 61 61	0	0
5	U	58	Total O 58 58	0	0
5	V	83	Total O 83 83	0	0
5	W	70	Total O 70 70	0	0
5	X	91	Total O 91 91	0	0
5	c	1	Total O 1 1	0	0
5	d	1	Total O 1 1	0	0
5	j	1	Total O 1 1	0	0
5	l	2	Total O 2 2	0	0
5	m	1	Total O 1 1	0	0
5	n	1	Total O 1 1	0	0
5	q	1	Total O 1 1	0	0
5	r	1	Total O 1 1	0	0
5	u	1	Total O 1 1	0	0

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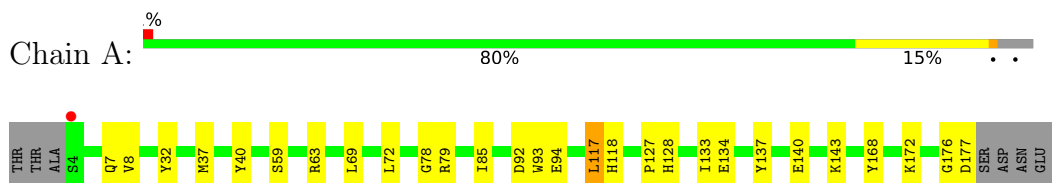
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	v	1	Total	O	0	0
			1	1		

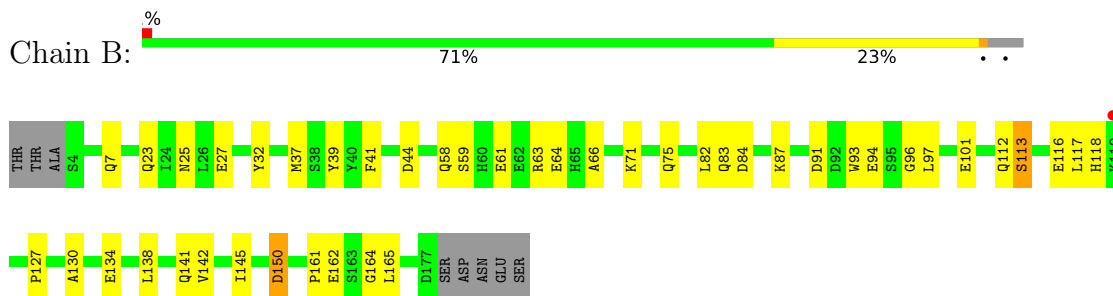
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

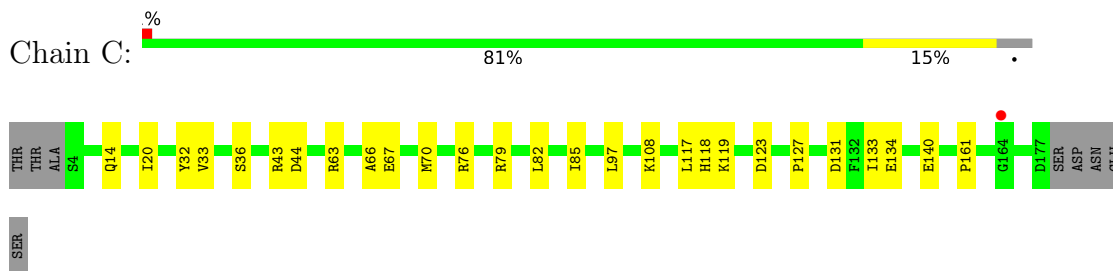
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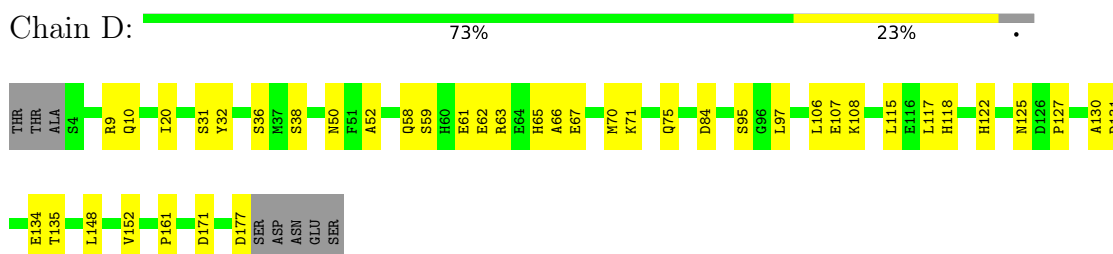
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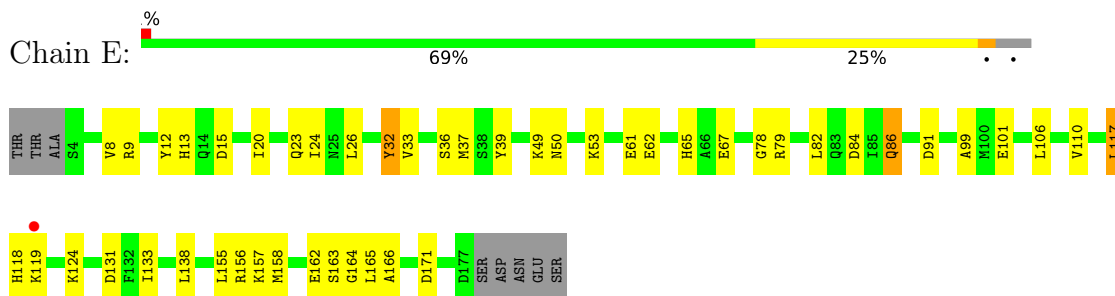
- Molecule 1: Ferritin heavy chain



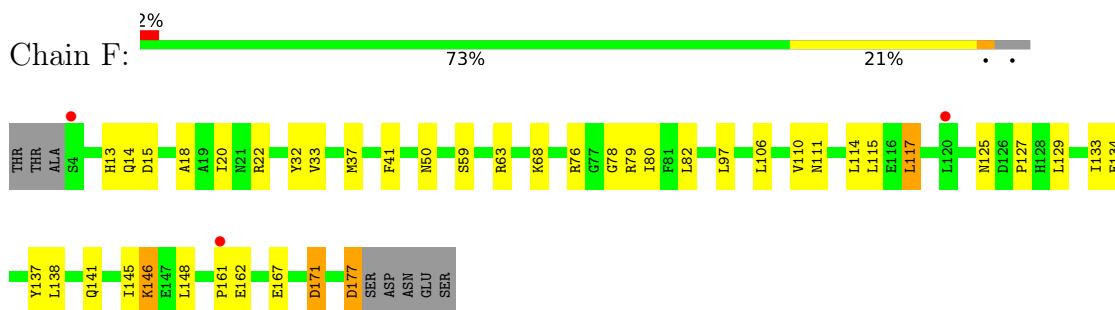
- Molecule 1: Ferritin heavy chain



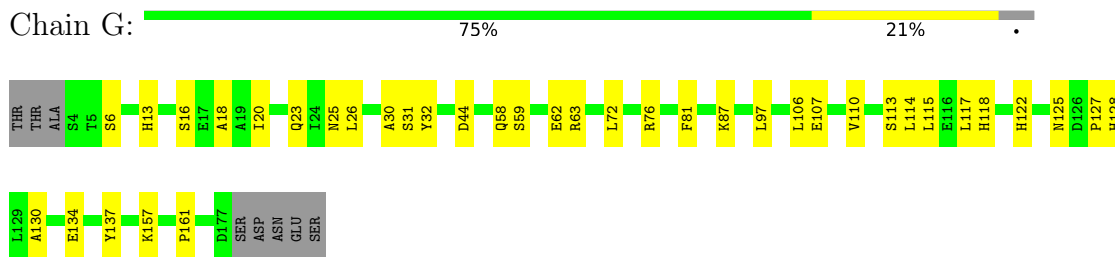
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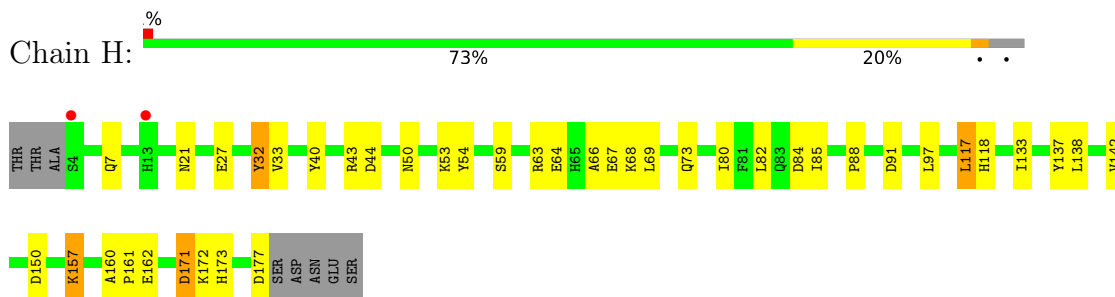
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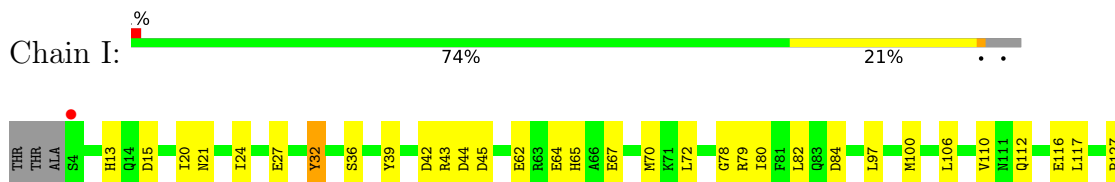
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- Molecule 1: Ferritin heavy chain

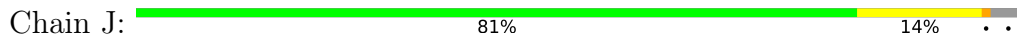


- Molecule 1: Ferritin heavy chain

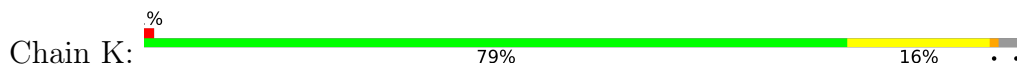




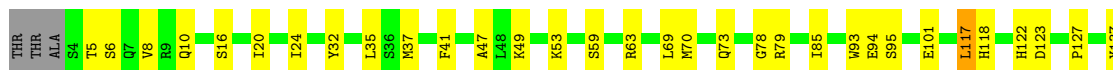
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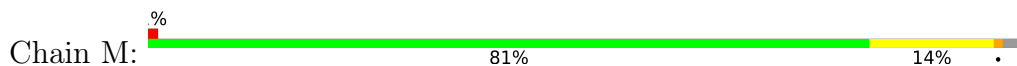
- Molecule 1: Ferritin heavy chain



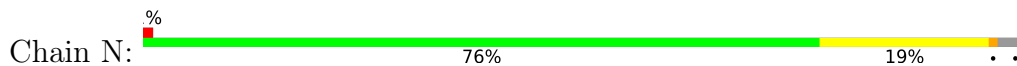
- Molecule 1: Ferritin heavy chain



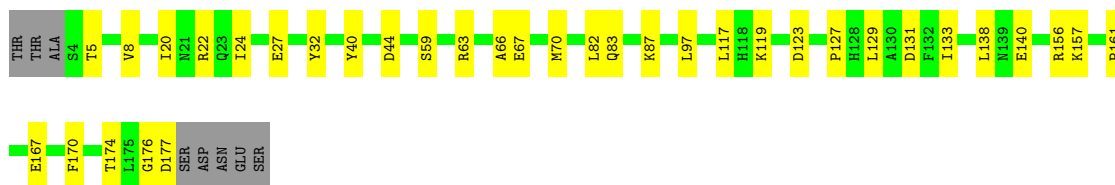
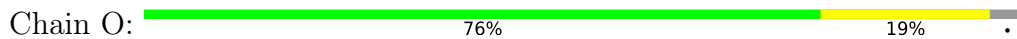
- Molecule 1: Ferritin heavy chain



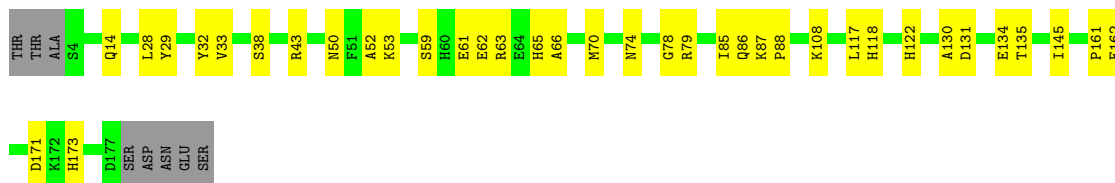
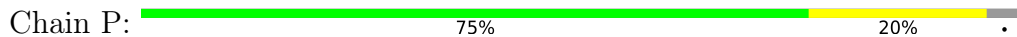
- Molecule 1: Ferritin heavy chain



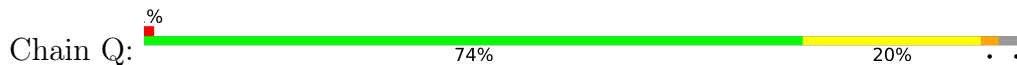
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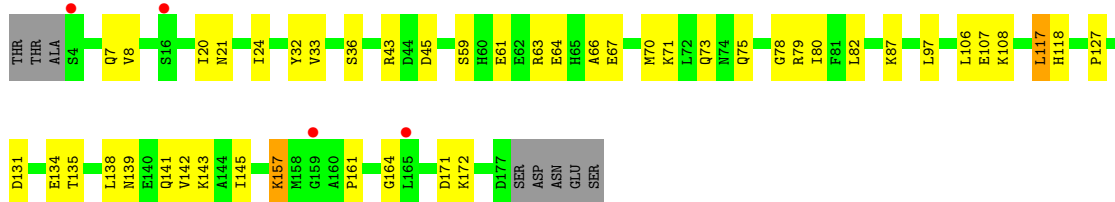
• Molecule 1: Ferritin heavy chain



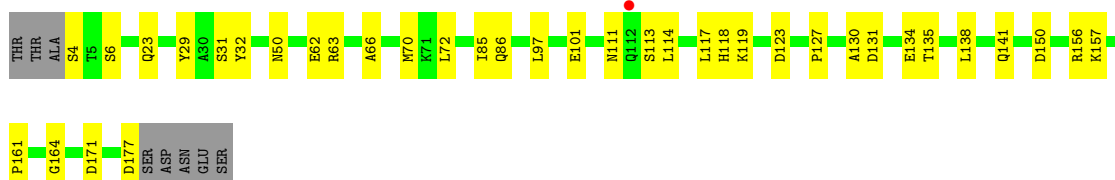
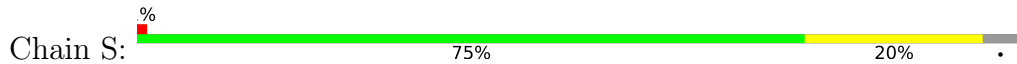
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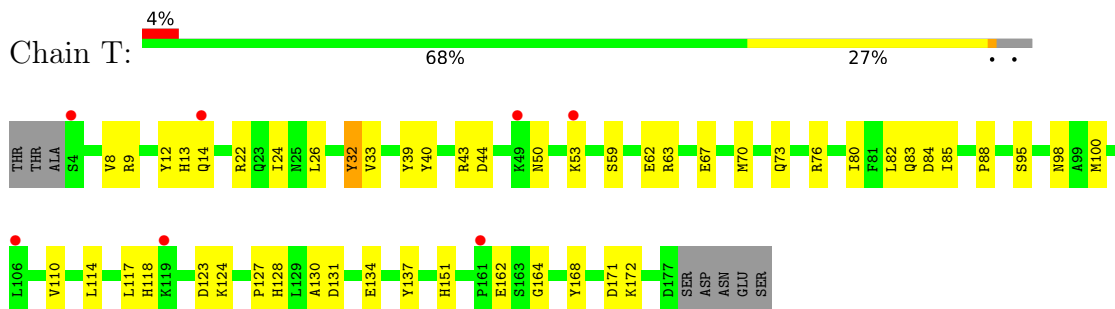
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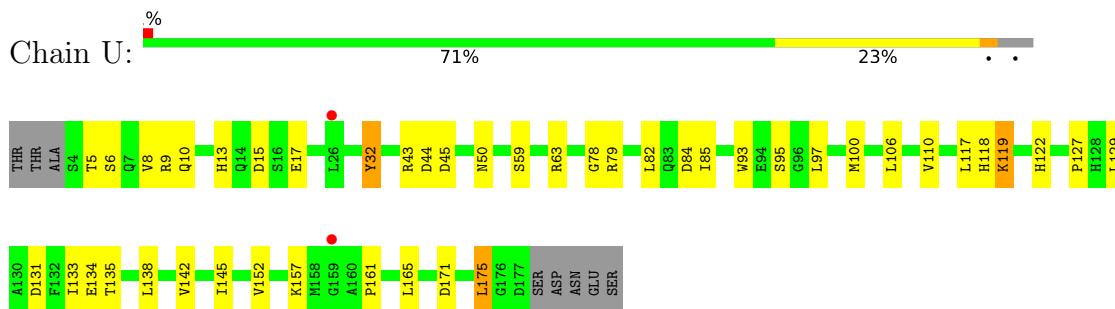
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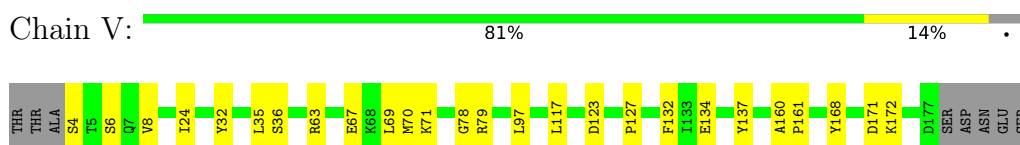
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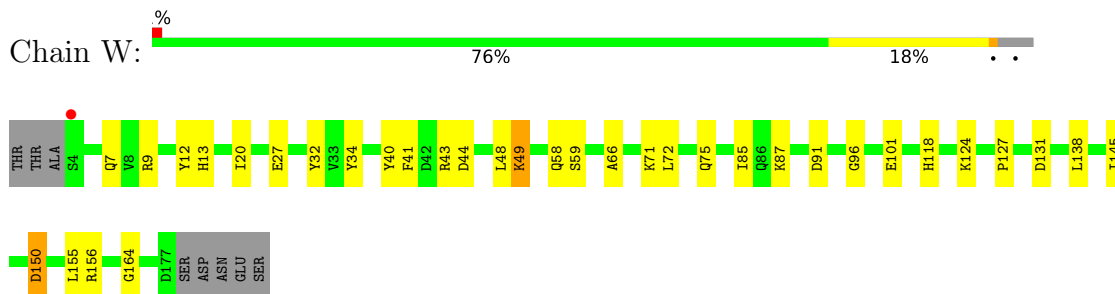
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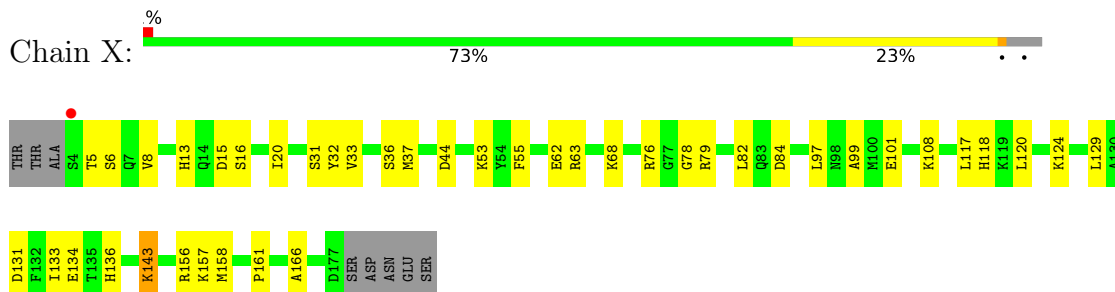
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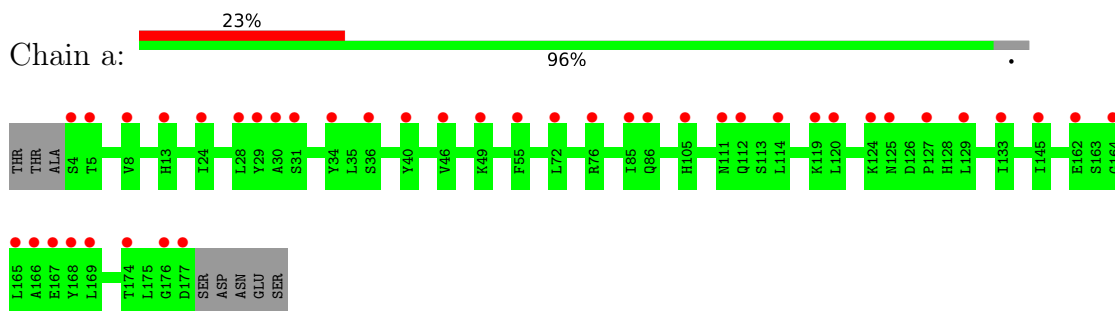
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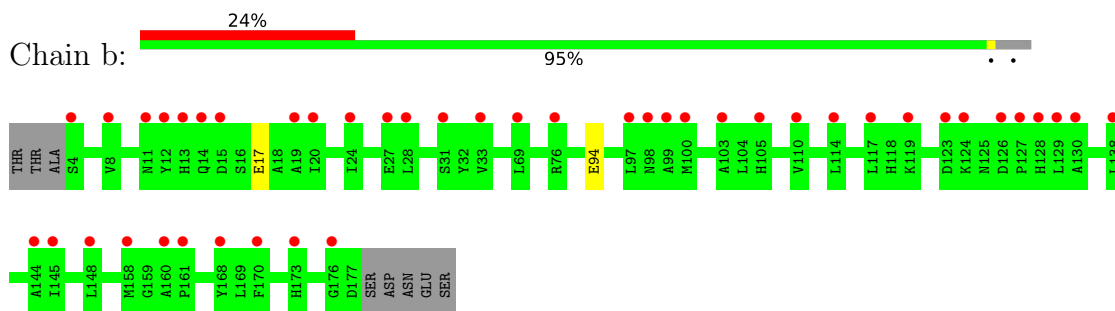
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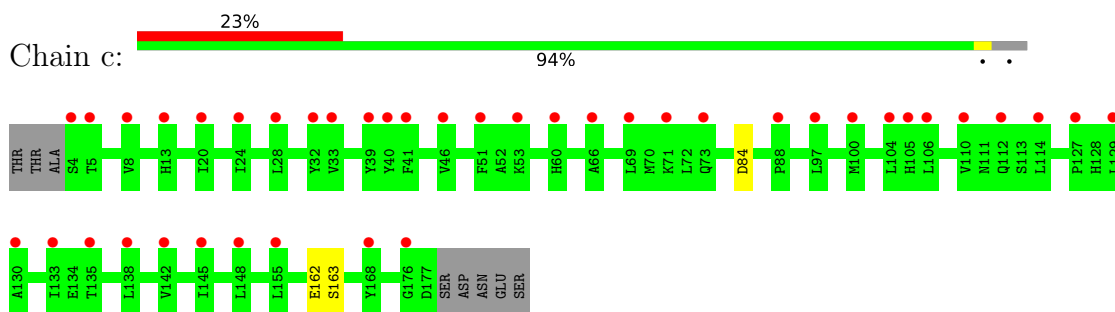
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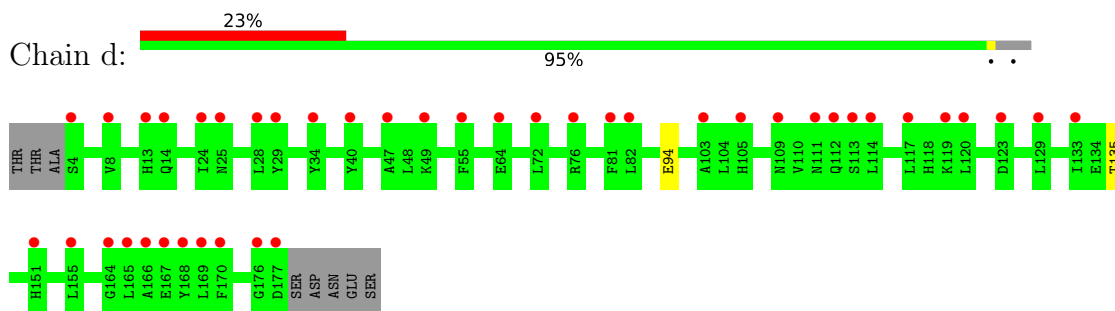
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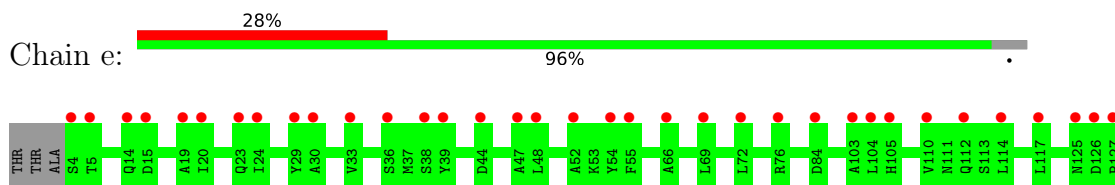
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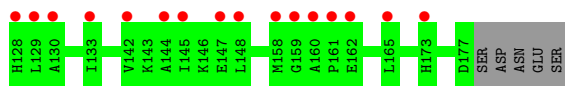


- Molecule 1: Ferritin heavy chain

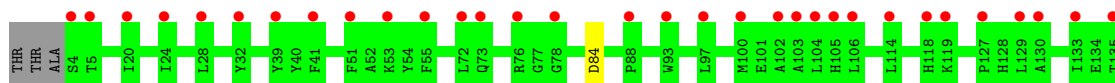
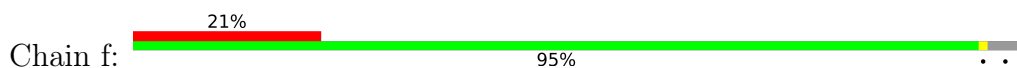


- Molecule 1: Ferritin heavy chain

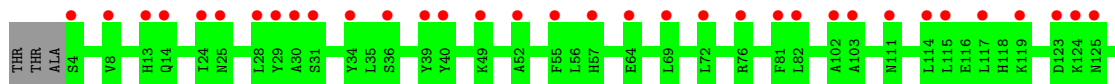




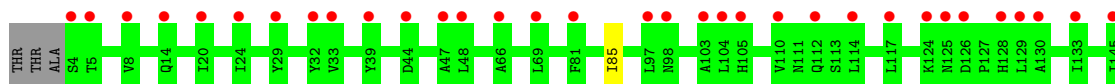
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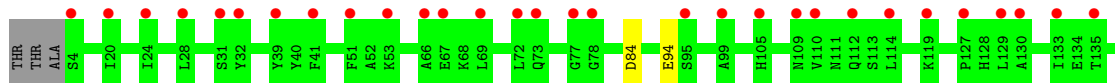
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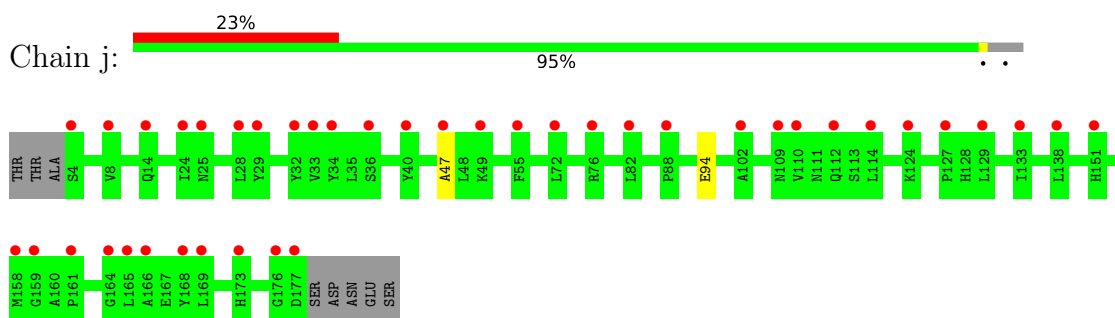
- Molecule 1: Ferritin heavy chain



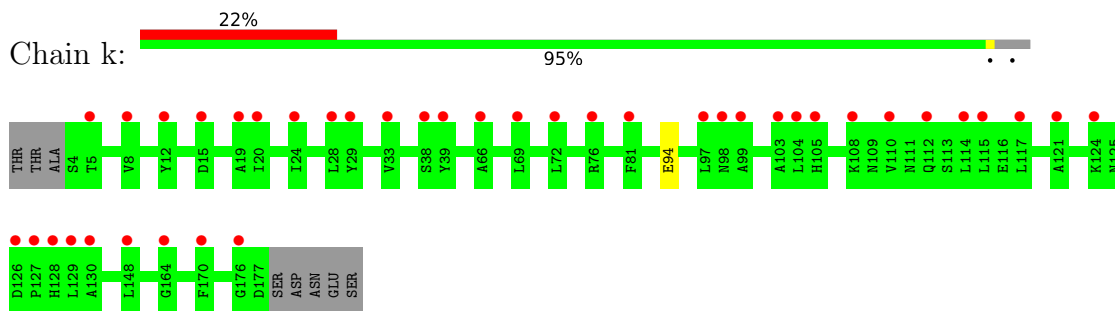
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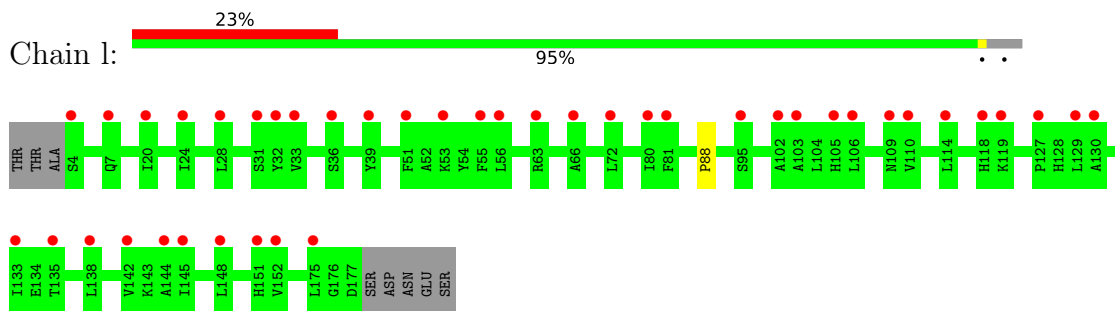
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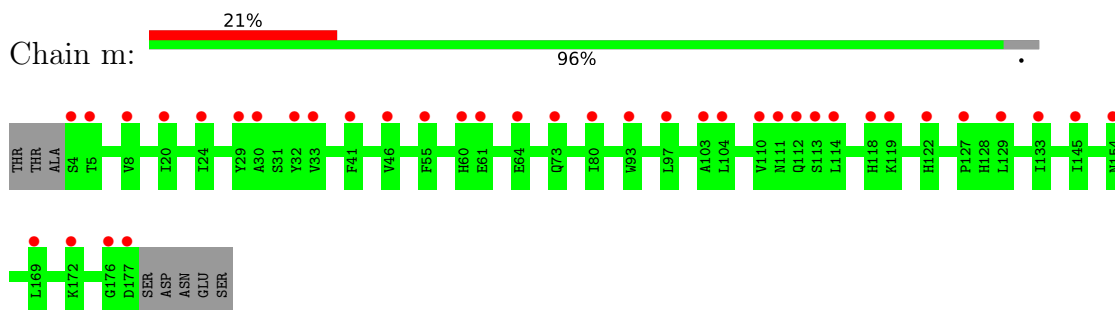
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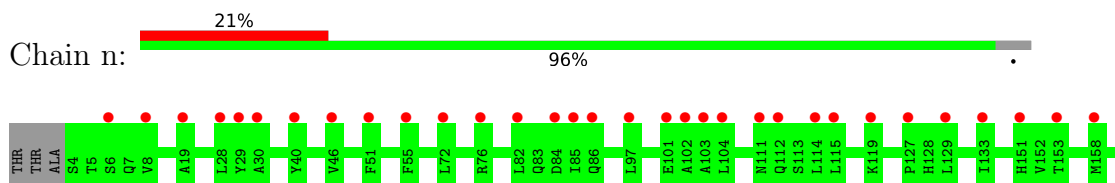
- Molecule 1: Ferritin heavy chain

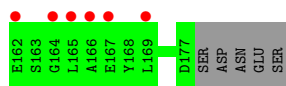


- Molecule 1: Ferritin heavy chain

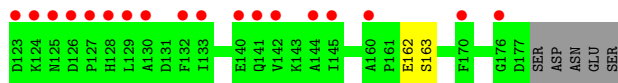
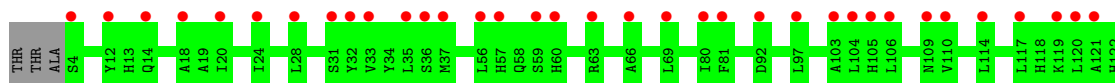
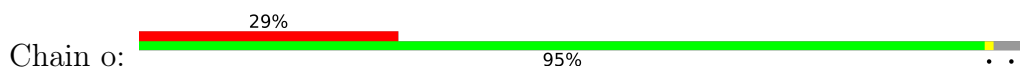


- Molecule 1: Ferritin heavy chain

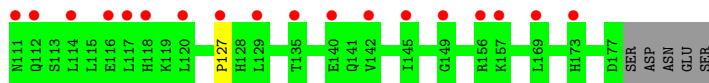
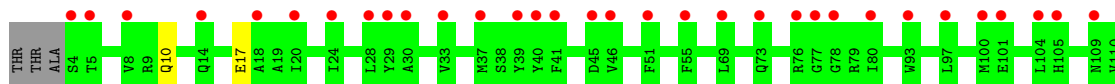
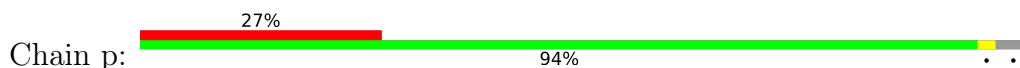




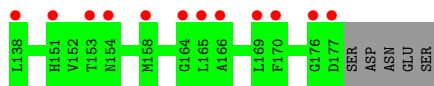
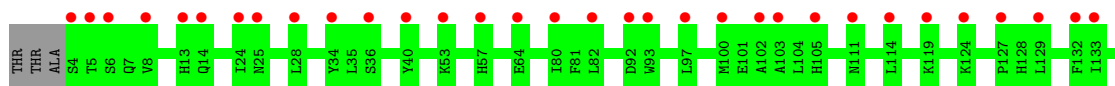
- Molecule 1: Ferritin heavy chain



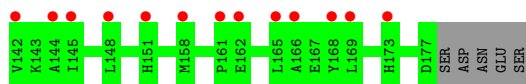
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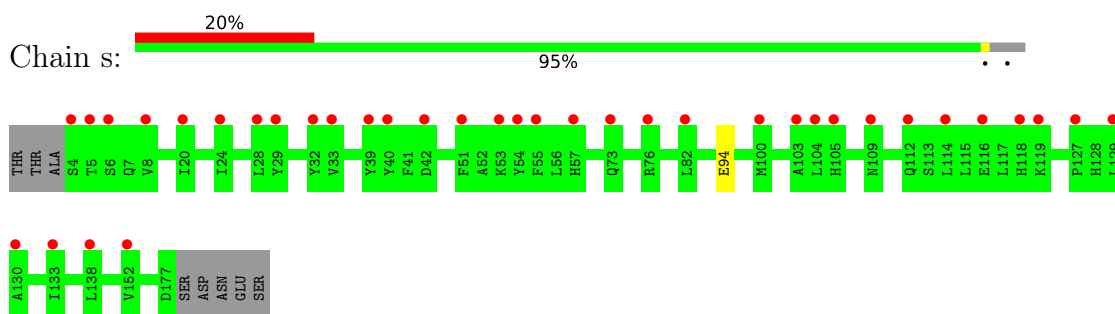
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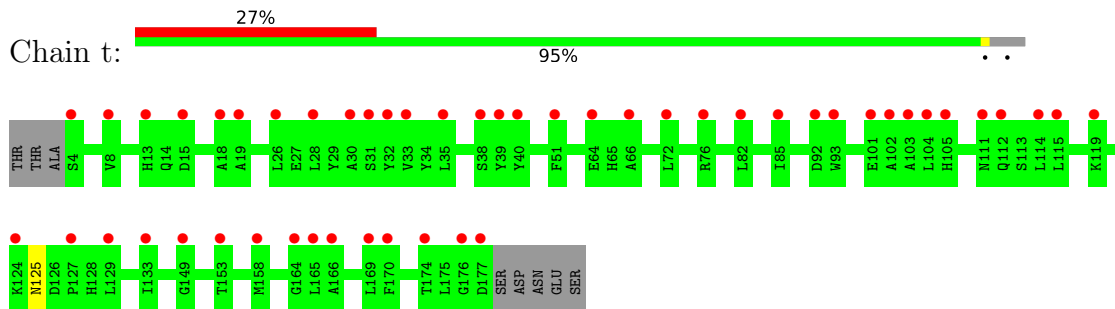
- Molecule 1: Ferritin heavy chain



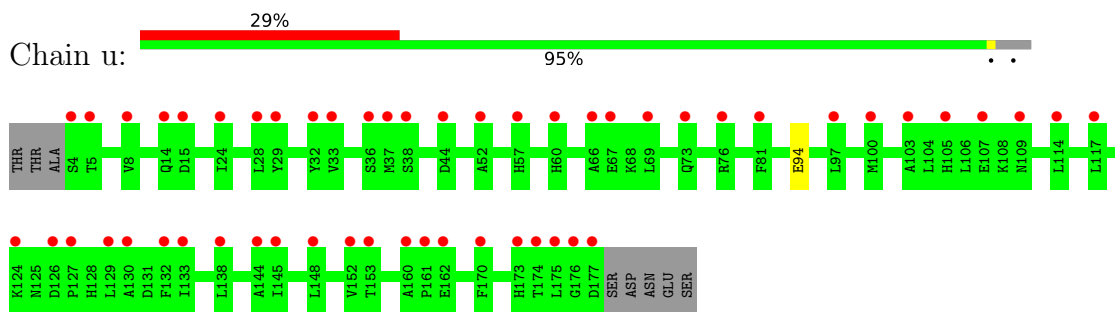
- Molecule 1: Ferritin heavy chain



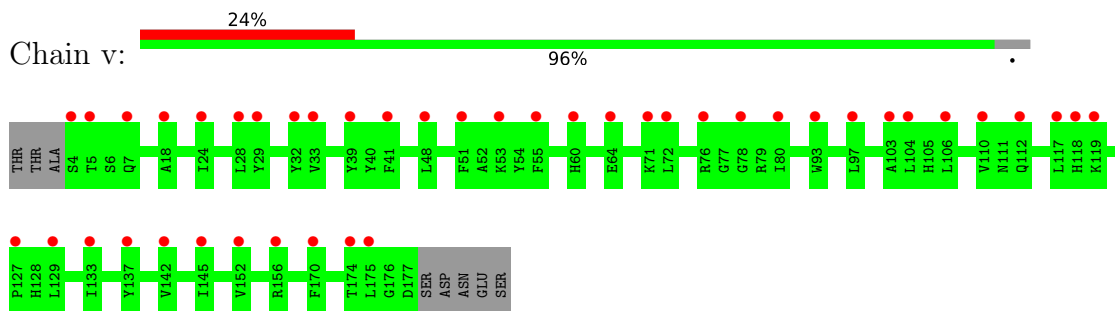
- Molecule 1: Ferritin heavy chain



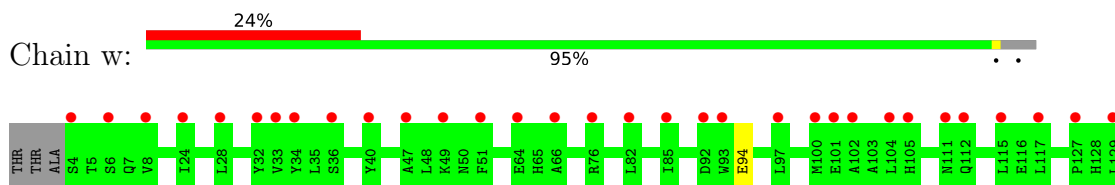
- Molecule 1: Ferritin heavy chain

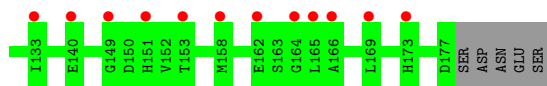


- Molecule 1: Ferritin heavy chain

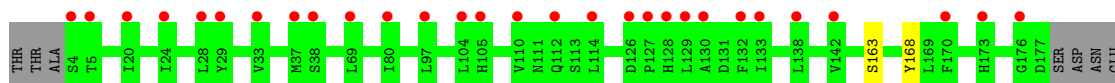
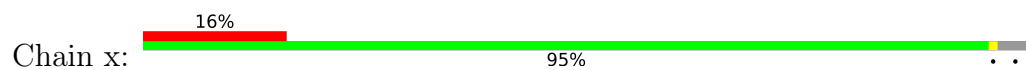


- Molecule 1: Ferritin heavy chain





- Molecule 1: Ferritin heavy chain



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4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	213.99Å 213.90Å 156.33Å 90.00° 90.14° 90.00°	Depositor
Resolution (Å)	38.49 – 2.34 39.19 – 2.34	Depositor EDS
% Data completeness (in resolution range)	94.7 (38.49-2.34) 93.7 (39.19-2.34)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.86 (at 2.34Å)	Xtrriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.207 , 0.278 0.207 , 0.278	Depositor DCC
R_{free} test set	55682 reflections (9.97%)	wwPDB-VP
Wilson B-factor (Å ²)	34.0	Xtrriage
Anisotropy	0.156	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 37.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.18$	Xtrriage
Estimated twinning fraction	0.427 for k,h,-l 0.440 for -k,-h,-l 0.429 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	105214	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.43% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: NA, ZN, V9Y

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.66	0/1461	0.72	0/1969
1	B	0.61	0/1461	0.68	1/1969 (0.1%)
1	C	0.65	0/1461	0.73	2/1969 (0.1%)
1	D	0.64	1/1461 (0.1%)	0.70	0/1969
1	E	0.60	0/1461	0.71	0/1969
1	F	0.59	0/1461	0.66	0/1969
1	G	0.64	1/1461 (0.1%)	0.71	2/1969 (0.1%)
1	H	0.58	0/1461	0.68	0/1969
1	I	0.58	1/1461 (0.1%)	0.66	0/1969
1	J	0.64	0/1461	0.70	0/1969
1	K	0.61	0/1461	0.71	0/1969
1	L	0.63	0/1461	0.72	1/1969 (0.1%)
1	M	0.63	0/1461	0.70	0/1969
1	N	0.57	0/1461	0.69	0/1969
1	O	0.63	0/1461	0.67	0/1969
1	P	0.62	0/1461	0.70	1/1969 (0.1%)
1	Q	0.56	0/1461	0.70	0/1969
1	R	0.57	0/1461	0.66	0/1969
1	S	0.65	0/1461	0.68	0/1969
1	T	0.58	0/1461	0.70	0/1969
1	U	0.58	0/1461	0.67	0/1969
1	V	0.65	0/1461	0.71	0/1969
1	W	0.59	0/1461	0.71	1/1969 (0.1%)
1	X	0.61	0/1461	0.69	0/1969
All	All	0.61	3/35064 (0.0%)	0.69	8/47256 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	107	GLU	CB-CG	6.58	1.64	1.52
1	G	107	GLU	CB-CG	6.09	1.63	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	27	GLU	CB-CG	6.05	1.63	1.52

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	123	ASP	CB-CG-OD1	6.90	124.51	118.30
1	G	72	LEU	CA-CB-CG	6.04	129.19	115.30
1	L	150	ASP	CB-CG-OD1	5.63	123.37	118.30
1	P	43	ARG	NE-CZ-NH1	5.43	123.02	120.30
1	B	150	ASP	CB-CG-OD1	5.38	123.14	118.30
1	W	150	ASP	CB-CG-OD1	5.13	122.92	118.30
1	G	115	LEU	CA-CB-CG	-5.12	103.54	115.30
1	C	43	ARG	NE-CZ-NH1	-5.09	117.75	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1431	0	1362	22	0
1	B	1431	0	1362	34	0
1	C	1431	0	1362	17	0
1	D	1431	0	1362	29	0
1	E	1431	0	1362	41	0
1	F	1431	0	1362	31	0
1	G	1431	0	1362	27	0
1	H	1431	0	1362	33	0
1	I	1431	0	1362	30	0
1	J	1431	0	1362	19	0
1	K	1431	0	1362	23	0
1	L	1431	0	1362	27	0
1	M	1431	0	1362	16	0
1	N	1431	0	1362	19	0
1	O	1431	0	1362	25	0
1	P	1431	0	1362	24	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	Q	1431	0	1362	27	0
1	R	1431	0	1362	35	0
1	S	1431	0	1362	24	0
1	T	1431	0	1362	38	0
1	U	1431	0	1362	36	0
1	V	1431	0	1362	16	0
1	W	1431	0	1362	26	0
1	X	1431	0	1362	29	0
1	a	2862	0	2710	0	0
1	b	2862	0	2700	0	0
1	c	2862	0	2703	0	0
1	d	2862	0	2710	0	0
1	e	2862	0	2702	0	0
1	f	2862	0	2704	0	0
1	g	2862	0	2705	0	0
1	h	2862	0	2706	0	0
1	i	2862	0	2701	0	0
1	j	2862	0	2707	0	0
1	k	2862	0	2703	0	0
1	l	2862	0	2701	0	0
1	m	2862	0	2705	0	0
1	n	2862	0	2710	0	0
1	o	2862	0	2708	0	0
1	p	2862	0	2709	0	0
1	q	2862	0	2713	0	0
1	r	2862	0	2705	0	0
1	s	2862	0	2702	0	0
1	t	2862	0	2703	0	0
1	u	2862	0	2702	0	0
1	v	2862	0	2706	0	0
1	w	2862	0	2707	0	0
1	x	2862	0	2703	0	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
2	C	3	0	0	0	0
2	D	2	0	0	0	0
2	E	3	0	0	0	0
2	F	4	0	0	0	0
2	G	3	0	0	0	0
2	H	3	0	0	0	0
2	I	2	0	0	0	0
2	J	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	K	3	0	0	0	0
2	L	2	0	0	0	0
2	M	2	0	0	0	0
2	N	3	0	0	0	0
2	O	3	0	0	0	0
2	P	3	0	0	0	0
2	Q	2	0	0	0	0
2	R	2	0	0	0	0
2	S	2	0	0	0	0
2	T	3	0	0	0	0
2	U	2	0	0	0	0
2	V	2	0	0	0	0
2	W	3	0	0	0	0
2	X	2	0	0	0	0
2	a	6	0	0	0	0
2	b	4	0	0	0	0
2	c	4	0	0	0	0
2	d	8	0	0	0	0
2	e	6	0	0	0	0
2	f	6	0	0	0	0
2	g	6	0	0	0	0
2	h	4	0	0	0	0
2	i	4	0	0	0	0
2	j	4	0	0	0	0
2	k	4	0	0	0	0
2	l	6	0	0	0	0
2	m	4	0	0	0	0
2	n	6	0	0	0	0
2	o	4	0	0	0	0
2	p	6	0	0	0	0
2	q	4	0	0	0	0
2	r	4	0	0	0	0
2	s	6	0	0	0	0
2	t	4	0	0	0	0
2	u	4	0	0	0	0
2	v	6	0	0	0	0
2	w	8	0	0	0	0
2	x	6	0	0	0	0
3	A	1	0	0	0	0
3	D	1	0	0	0	0
3	G	1	0	0	0	0
3	J	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	M	1	0	0	0	0
3	P	1	0	0	0	0
3	S	1	0	0	0	0
3	V	1	0	0	0	0
3	a	2	0	0	0	0
3	d	2	0	0	0	0
3	g	2	0	0	0	0
3	j	2	0	0	0	0
3	n	2	0	0	0	0
3	p	2	0	0	0	0
3	s	2	0	0	0	0
3	x	2	0	0	0	0
4	A	26	0	0	0	0
4	D	26	0	0	1	0
4	I	26	0	0	2	0
4	L	26	0	0	0	0
4	O	26	0	0	0	0
4	R	26	0	0	2	0
4	X	26	0	0	0	0
4	t	26	0	0	0	0
5	A	92	0	0	3	0
5	B	76	0	0	8	0
5	C	86	0	0	0	0
5	D	69	0	0	4	0
5	E	78	0	0	3	0
5	F	54	0	0	2	0
5	G	66	0	0	4	0
5	H	57	0	0	4	0
5	I	59	0	0	0	0
5	J	104	0	0	4	0
5	K	83	0	0	6	0
5	L	83	0	0	6	0
5	M	94	0	0	2	0
5	N	64	0	0	1	0
5	O	90	0	0	6	0
5	P	56	0	0	4	0
5	Q	54	0	0	2	0
5	R	58	0	0	1	0
5	S	67	0	0	3	0
5	T	61	0	0	4	0
5	U	58	0	0	1	0
5	V	83	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	W	70	0	0	4	0
5	X	91	0	0	2	0
5	c	1	0	0	0	0
5	d	1	0	0	0	0
5	j	1	0	0	0	0
5	l	2	0	0	0	0
5	m	1	0	0	0	0
5	n	1	0	0	0	0
5	q	1	0	0	0	0
5	r	1	0	0	0	0
5	u	1	0	0	0	0
5	v	1	0	0	0	0
All	All	105214	0	97613	537	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (537) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:101:GLU:OE2	1:L:156:ARG:NH2	1.99	0.95
1:I:158:MET:HE1	1:I:170:PHE:HB2	1.47	0.92
1:X:101:GLU:OE2	1:X:156:ARG:NH2	2.07	0.87
1:O:63:ARG:NE	1:O:67:GLU:OE2	2.08	0.86
1:G:122:HIS:HE1	4:I:203[C]:V9Y:O04	1.62	0.82
1:S:118:HIS:CE1	1:U:127:PRO:HB3	2.15	0.81
1:G:25:ASN:ND2	5:G:301:HOH:O	2.08	0.81
1:A:92:ASP:O	5:A:301:HOH:O	1.97	0.81
1:L:157:LYS:NZ	5:L:301:HOH:O	2.10	0.79
1:O:167:GLU:OE1	5:O:301:HOH:O	2.00	0.78
1:Q:101:GLU:OE2	1:Q:156:ARG:NH2	2.17	0.78
1:D:122:HIS:HE1	4:D:204[C]:V9Y:O04	1.66	0.77
1:M:134:GLU:HA	1:M:138:LEU:HD12	1.65	0.77
1:T:44:ASP:OD2	5:T:301:HOH:O	2.02	0.77
1:H:73:GLN:HG2	1:H:80:ILE:HG13	1.66	0.77
1:W:49:LYS:NZ	5:W:301:HOH:O	2.17	0.76
1:H:50:ASN:HA	1:H:53:LYS:HG2	1.69	0.75
1:M:58:GLN:HG2	5:M:311:HOH:O	1.87	0.75
1:P:14:GLN:O	5:P:301:HOH:O	2.04	0.75
1:B:63:ARG:NH2	1:D:59:SER:OG	2.21	0.74
1:D:97:LEU:HD23	1:D:161:PRO:HD3	1.71	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:63:ARG:NH2	1:U:59:SER:OG	2.22	0.73
1:Q:162:GLU:N	1:Q:162:GLU:OE2	2.21	0.73
1:K:90:GLU:OE1	5:K:301:HOH:O	2.07	0.73
1:J:140:GLU:OE1	5:J:302:HOH:O	2.07	0.71
1:B:87:LYS:NZ	5:B:303:HOH:O	2.22	0.71
1:V:71:LYS:NZ	5:V:301:HOH:O	2.22	0.70
1:S:127:PRO:HB3	1:T:118:HIS:CE1	2.27	0.70
1:P:173:HIS:HB3	1:T:172:LYS:HG2	1.74	0.70
1:R:59:SER:OG	1:T:63:ARG:NH2	2.25	0.70
1:R:134:GLU:HA	1:R:138:LEU:HD12	1.73	0.69
1:A:78:GLY:O	1:A:79:ARG:NH1	2.26	0.68
1:F:97:LEU:HD23	1:F:161:PRO:HD3	1.74	0.68
1:P:33:VAL:HG22	1:P:88:PRO:HB3	1.76	0.68
1:L:157:LYS:NZ	5:L:303:HOH:O	2.27	0.68
1:J:86:GLN:NE2	5:J:301:HOH:O	2.01	0.67
1:H:80:ILE:O	5:H:301:HOH:O	2.13	0.67
1:N:141:GLN:OE1	5:N:1501:HOH:O	2.13	0.66
1:M:134:GLU:HG2	1:O:131:ASP:HB2	1.76	0.65
1:X:68:LYS:HZ2	1:X:136:HIS:CE1	2.14	0.65
1:H:32:TYR:OH	1:U:82:LEU:HB3	1.95	0.65
1:A:63:ARG:HH21	1:L:35:LEU:HD13	1.61	0.64
1:F:59:SER:OG	1:Q:63:ARG:NH2	2.30	0.64
1:E:91:ASP:O	5:E:301:HOH:O	2.14	0.64
1:M:71:LYS:NZ	5:M:304:HOH:O	2.31	0.64
1:Q:97:LEU:HD23	1:Q:161:PRO:HD3	1.78	0.64
1:V:97:LEU:HD23	1:V:161:PRO:HD3	1.78	0.64
1:R:82:LEU:HD13	1:T:32:TYR:CE1	2.34	0.63
1:B:87:LYS:O	5:B:301:HOH:O	2.15	0.63
1:W:34:TYR:CD2	1:W:58:GLN:HB3	2.33	0.63
1:T:33:VAL:HG22	1:T:88:PRO:HB3	1.79	0.63
1:U:43:ARG:NH2	1:U:45:ASP:OD2	2.28	0.62
1:X:20:ILE:HD13	1:X:117:LEU:HD11	1.81	0.62
1:C:117:LEU:HG	1:C:133:ILE:HD11	1.80	0.62
1:O:83:GLN:NE2	5:O:302:HOH:O	2.08	0.62
1:O:20:ILE:HD13	1:O:117:LEU:HD21	1.82	0.62
1:T:22:ARG:NH2	5:T:302:HOH:O	2.07	0.62
1:J:121:ALA:HB2	1:J:129:LEU:HD23	1.80	0.61
1:I:97:LEU:HD23	1:I:161:PRO:HD3	1.83	0.61
1:U:131:ASP:O	1:U:135:THR:HG23	1.99	0.61
1:I:100:MET:HB3	1:I:152:VAL:HG23	1.83	0.61
1:Q:73:GLN:HG2	1:Q:80:ILE:HG13	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:63:ARG:NH2	1:P:59:SER:OG	2.34	0.61
1:D:118:HIS:CE1	1:F:127:PRO:HB3	2.36	0.60
1:F:82:LEU:HD13	1:Q:32:TYR:CE1	2.37	0.60
1:S:23:GLN:OE1	1:S:113:SER:OG	2.17	0.60
1:E:101:GLU:OE2	1:E:156:ARG:NH2	2.32	0.60
1:H:32:TYR:CE1	1:U:82:LEU:HD13	2.36	0.60
1:S:131:ASP:O	1:S:135:THR:HG23	2.02	0.60
1:D:127:PRO:HB3	1:E:118:HIS:CE1	2.37	0.60
1:N:13:HIS:CD2	1:N:124:LYS:HD2	2.36	0.60
1:R:97:LEU:HD23	1:R:161:PRO:HD3	1.83	0.60
1:U:171:ASP:HA	1:U:175:LEU:HD12	1.82	0.60
1:W:71:LYS:O	1:W:75:GLN:HG3	2.01	0.60
1:E:37:MET:CE	1:E:99:ALA:HB1	2.32	0.60
1:K:101:GLU:OE2	1:K:156:ARG:NH2	2.34	0.59
1:H:59:SER:OG	1:U:63:ARG:NH2	2.35	0.59
1:H:82:LEU:HD13	1:U:32:TYR:CE1	2.37	0.59
1:M:118:HIS:CE1	1:O:127:PRO:HB3	2.37	0.59
1:K:61:GLU:HG2	5:K:351:HOH:O	2.02	0.59
1:W:49:LYS:H	1:W:49:LYS:HD2	1.67	0.59
1:K:64:GLU:HB3	5:K:347:HOH:O	2.02	0.59
1:B:165:LEU:HD12	1:E:158:MET:HG2	1.85	0.59
1:H:33:VAL:HG22	1:H:88:PRO:HB3	1.85	0.59
1:W:164:GLY:N	5:W:305:HOH:O	2.35	0.58
1:B:39:TYR:OH	1:D:67:GLU:HB3	2.03	0.58
1:D:130:ALA:O	1:D:134:GLU:HG3	2.03	0.58
1:E:9:ARG:NE	1:E:12:TYR:O	2.31	0.58
1:E:49:LYS:NZ	5:E:304:HOH:O	2.35	0.58
1:L:177:ASP:HB2	5:L:315:HOH:O	2.04	0.57
1:G:157:LYS:HZ3	1:G:157:LYS:HB2	1.69	0.57
1:S:119:LYS:NZ	1:S:123:ASP:OD2	2.38	0.57
1:D:58:GLN:HG2	5:D:311:HOH:O	2.04	0.57
1:I:43:ARG:NH2	1:I:45:ASP:OD2	2.28	0.57
1:T:62:GLU:OE2	1:T:62:GLU:HA	2.05	0.57
1:X:37:MET:HE3	1:X:99:ALA:HB1	1.87	0.57
1:W:87:LYS:NZ	5:W:304:HOH:O	2.35	0.57
1:S:4:SER:N	5:S:304:HOH:O	2.38	0.56
1:E:13:HIS:HE1	1:E:15:ASP:OD2	1.88	0.56
1:S:101:GLU:OE2	1:S:156:ARG:NH2	2.38	0.56
1:I:70:MET:HE3	1:I:80:ILE:HG21	1.86	0.56
1:K:71:LYS:NZ	5:K:306:HOH:O	2.38	0.56
1:T:164:GLY:N	5:T:303:HOH:O	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:118:HIS:CE1	1:I:127:PRO:HB3	2.41	0.56
1:M:78:GLY:O	1:M:79:ARG:NH1	2.39	0.56
1:E:78:GLY:O	1:E:79:ARG:NH1	2.38	0.56
1:G:58:GLN:HG2	5:G:319:HOH:O	2.05	0.56
1:G:23:GLN:OE1	1:G:113:SER:OG	2.20	0.56
1:D:106:LEU:O	1:D:106:LEU:HD23	2.06	0.55
1:P:130:ALA:O	1:P:134:GLU:HG3	2.06	0.55
1:X:158:MET:HB3	1:X:166:ALA:HB1	1.89	0.55
1:G:130:ALA:O	1:G:134:GLU:HG3	2.05	0.55
1:E:106:LEU:O	1:E:110:VAL:HG23	2.06	0.55
1:L:20:ILE:HD13	1:L:117:LEU:HD11	1.88	0.55
1:P:161:PRO:HD2	1:P:162:GLU:OE2	2.07	0.55
1:D:177:ASP:HA	5:D:352:HOH:O	2.06	0.55
1:F:37:MET:O	1:F:41:PHE:HD2	1.88	0.55
1:R:172:LYS:HE2	5:R:316:HOH:O	2.07	0.55
1:X:97:LEU:HD23	1:X:161:PRO:HD3	1.89	0.55
1:A:118:HIS:CE1	1:C:127:PRO:HB3	2.41	0.55
1:H:117:LEU:HD22	1:H:133:ILE:HD11	1.89	0.55
1:Q:114:LEU:HD13	1:Q:137:TYR:HB3	1.89	0.55
1:E:32:TYR:CE1	1:I:82:LEU:HD13	2.42	0.55
1:Q:9:ARG:O	1:R:108:LYS:NZ	2.40	0.55
1:R:63:ARG:NH2	1:T:59:SER:OG	2.40	0.54
1:V:168:TYR:CE1	1:V:172:LYS:HE2	2.41	0.54
1:A:168:TYR:O	1:A:172:LYS:HG2	2.08	0.54
1:B:112:GLN:O	1:B:116:GLU:HG3	2.07	0.54
1:A:63:ARG:NH2	1:L:59:SER:OG	2.35	0.54
1:B:164:GLY:HA3	1:E:157:LYS:O	2.07	0.54
1:J:101:GLU:OE2	1:J:156:ARG:NH2	2.32	0.54
1:P:78:GLY:O	1:P:79:ARG:NH1	2.40	0.54
1:T:100:MET:HE2	1:T:151:HIS:HB2	1.89	0.54
1:U:97:LEU:HD23	1:U:161:PRO:HD3	1.89	0.54
1:R:78:GLY:O	1:R:79:ARG:NH1	2.39	0.54
1:O:59:SER:OG	1:V:63:ARG:NH2	2.41	0.54
1:R:82:LEU:HB3	1:T:32:TYR:OH	2.07	0.54
1:D:38:SER:OG	1:D:52:ALA:O	2.21	0.54
1:D:125:ASN:HB3	1:E:119:LYS:HE2	1.90	0.54
1:F:114:LEU:HD13	1:F:137:TYR:HB3	1.88	0.54
1:U:50:ASN:HB2	1:U:171:ASP:OD2	2.08	0.53
1:O:176:GLY:O	1:O:177:ASP:HB2	2.07	0.53
1:R:87:LYS:HD3	1:T:83:GLN:HA	1.90	0.53
1:I:21:ASN:HA	1:I:24:ILE:HD12	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:164:GLY:HA3	1:X:157:LYS:O	2.09	0.53
1:L:78:GLY:O	1:L:79:ARG:NH1	2.42	0.53
1:F:106:LEU:O	1:F:110:VAL:HG23	2.08	0.53
1:E:162:GLU:CD	1:E:162:GLU:H	2.12	0.53
1:Q:127:PRO:HB3	1:R:118:HIS:CE1	2.43	0.53
1:G:18:ALA:O	5:G:302:HOH:O	2.19	0.52
1:T:9:ARG:NH2	1:T:12:TYR:O	2.41	0.52
1:U:13:HIS:HE1	1:U:15:ASP:OD2	1.92	0.52
1:W:131:ASP:HB2	1:X:134:GLU:HG2	1.91	0.52
1:C:63:ARG:O	1:C:67:GLU:HG3	2.08	0.52
1:F:134:GLU:HA	1:F:138:LEU:HD12	1.90	0.52
1:I:43:ARG:HH21	1:I:45:ASP:CG	2.13	0.52
1:W:9:ARG:NH2	1:W:12:TYR:O	2.37	0.52
1:A:127:PRO:HB3	1:B:118:HIS:CE1	2.43	0.52
1:O:117:LEU:HG	1:O:133:ILE:HD11	1.91	0.52
1:G:59:SER:OG	1:K:63:ARG:NH2	2.41	0.52
1:G:127:PRO:HB3	1:H:118:HIS:CE1	2.45	0.52
1:K:130:ALA:O	1:K:134:GLU:HG3	2.10	0.52
1:U:119:LYS:HE3	1:U:122:HIS:HD1	1.73	0.52
1:Q:50:ASN:HB2	1:Q:171:ASP:OD2	2.10	0.52
1:F:18:ALA:O	1:F:22:ARG:HG3	2.10	0.51
1:O:63:ARG:NH2	1:V:35:LEU:HD13	2.25	0.51
1:V:134:GLU:HG2	1:X:131:ASP:HB2	1.91	0.51
1:D:31:SER:HB2	1:D:62:GLU:HB2	1.92	0.51
1:F:167:GLU:HB2	1:R:157:LYS:NZ	2.25	0.51
1:T:50:ASN:HB2	1:T:171:ASP:OD2	2.11	0.51
1:P:131:ASP:O	1:P:135:THR:HG23	2.11	0.51
1:U:13:HIS:CE1	1:U:15:ASP:OD2	2.64	0.51
1:D:148:LEU:O	1:D:152:VAL:HG23	2.11	0.51
1:Q:163:SER:HB3	1:Q:166:ALA:HB2	1.91	0.51
1:A:140:GLU:OE2	1:A:143:LYS:NZ	2.43	0.51
1:K:127:PRO:HD2	5:K:308:HOH:O	2.11	0.51
1:M:168:TYR:O	1:M:172:LYS:HG2	2.11	0.51
1:U:106:LEU:O	1:U:110:VAL:HG23	2.11	0.51
1:H:172:LYS:HE2	5:H:304:HOH:O	2.11	0.51
1:W:49:LYS:H	1:W:49:LYS:CD	2.24	0.51
1:B:59:SER:OG	1:D:63:ARG:NH2	2.44	0.50
1:Q:49:LYS:HB3	1:Q:49:LYS:NZ	2.26	0.50
1:V:127:PRO:HB3	1:W:118:HIS:CE1	2.46	0.50
1:G:26:LEU:HD21	1:G:110:VAL:HA	1.93	0.50
1:H:63:ARG:O	1:H:67:GLU:HG3	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:T:26:LEU:HD21	1:T:110:VAL:HA	1.93	0.50
1:A:69:LEU:HG	1:A:137:TYR:OH	2.12	0.50
1:R:20:ILE:HD13	1:R:117:LEU:HD11	1.93	0.50
1:D:131:ASP:O	1:D:135:THR:HG23	2.12	0.50
1:E:163:SER:HB3	1:E:166:ALA:CB	2.42	0.50
1:H:50:ASN:HB2	1:H:171:ASP:OD2	2.11	0.50
1:L:24:ILE:HD13	1:L:70:MET:HG2	1.93	0.50
1:T:127:PRO:HB3	1:U:118:HIS:CE1	2.47	0.50
1:J:32:TYR:OH	1:X:82:LEU:HB3	2.11	0.50
1:J:87:LYS:HG2	1:X:84:ASP:OD1	2.12	0.50
1:O:44:ASP:HA	1:W:150:ASP:OD1	2.12	0.50
5:H:311:HOH:O	1:S:157:LYS:HE2	2.12	0.49
1:I:13:HIS:CE1	1:I:15:ASP:OD2	2.65	0.49
1:P:122:HIS:CE1	4:R:203[D]:V9Y:C02	2.94	0.49
1:D:63:ARG:O	1:D:67:GLU:HG3	2.13	0.49
1:E:62:GLU:O	1:E:65:HIS:HB2	2.11	0.49
1:G:31:SER:HB2	1:G:62:GLU:HB2	1.93	0.49
1:S:31:SER:HB2	1:S:62:GLU:HB2	1.95	0.49
1:C:97:LEU:HD23	1:C:161:PRO:HD3	1.95	0.49
1:F:80:ILE:O	5:F:301:HOH:O	2.20	0.49
1:I:154:ASN:O	1:I:158:MET:HG3	2.13	0.49
1:Q:141:GLN:NE2	5:Q:303:HOH:O	2.33	0.49
1:R:67:GLU:HG2	1:T:39:TYR:OH	2.12	0.49
1:T:13:HIS:CE1	1:T:124:LYS:HE2	2.48	0.49
1:G:87:LYS:NZ	5:G:309:HOH:O	2.45	0.49
1:J:118:HIS:ND1	5:J:307:HOH:O	2.35	0.49
1:N:114:LEU:HD13	1:N:137:TYR:HB3	1.95	0.49
1:S:135:THR:HG21	5:S:348:HOH:O	2.11	0.49
1:U:9:ARG:NH1	1:U:17:GLU:OE1	2.31	0.49
1:A:59:SER:OG	1:L:63:ARG:NH2	2.42	0.49
1:N:92:ASP:OD1	1:N:94:GLU:N	2.41	0.48
1:S:50:ASN:HB2	1:S:171:ASP:OD2	2.12	0.48
1:U:50:ASN:HB3	1:U:175:LEU:HB2	1.95	0.48
1:G:97:LEU:HD23	1:G:161:PRO:HD3	1.93	0.48
1:L:123:ASP:OD1	5:L:302:HOH:O	2.20	0.48
1:G:63:ARG:NH2	1:K:59:SER:OG	2.46	0.48
1:V:63:ARG:NE	1:V:67:GLU:OE2	2.38	0.48
1:B:25:ASN:ND2	5:B:305:HOH:O	2.35	0.48
1:X:5:THR:O	5:X:301:HOH:O	2.20	0.48
1:I:20:ILE:HD13	1:I:117:LEU:HD11	1.94	0.48
1:T:73:GLN:HG2	1:T:80:ILE:HG13	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:6:SER:OG	1:X:8:VAL:HG22	2.14	0.48
1:X:37:MET:CE	1:X:99:ALA:HB1	2.43	0.48
1:T:128:HIS:CD2	1:U:142:VAL:HG21	2.48	0.48
1:W:127:PRO:HB3	1:X:118:HIS:CE1	2.47	0.48
1:A:85:ILE:HB	1:L:85:ILE:HB	1.96	0.48
1:A:117:LEU:HD22	1:A:133:ILE:HD11	1.95	0.48
1:B:25:ASN:OD1	1:B:83:GLN:HB2	2.13	0.48
1:V:69:LEU:HG	1:V:137:TYR:OH	2.14	0.48
1:E:163:SER:HB3	1:E:166:ALA:HB2	1.95	0.48
1:H:27:GLU:HB2	1:H:66:ALA:HB2	1.95	0.48
1:L:16:SER:O	1:L:20:ILE:HG12	2.14	0.48
1:Q:13:HIS:CE1	1:Q:124:LYS:HE3	2.48	0.48
1:V:168:TYR:O	1:V:172:LYS:HG2	2.13	0.48
1:F:63:ARG:NH2	1:Q:59:SER:OG	2.46	0.48
1:F:167:GLU:HB2	1:R:157:LYS:HZ1	1.79	0.48
1:F:177:ASP:HA	5:F:338:HOH:O	2.13	0.48
1:G:76:ARG:HD3	1:G:128:HIS:ND1	2.29	0.48
1:U:78:GLY:O	1:U:79:ARG:NH1	2.45	0.48
1:B:161:PRO:HD2	1:B:162:GLU:OE2	2.14	0.47
1:S:130:ALA:O	1:S:134:GLU:HG3	2.14	0.47
1:T:100:MET:HE2	1:T:151:HIS:CB	2.43	0.47
1:B:23:GLN:OE1	1:B:113:SER:OG	2.30	0.47
1:B:97:LEU:O	1:B:101:GLU:HG3	2.13	0.47
1:F:78:GLY:O	1:F:79:ARG:NH1	2.47	0.47
1:R:36:SER:HB2	1:T:82:LEU:CD1	2.44	0.47
1:U:9:ARG:HG2	5:U:323:HOH:O	2.13	0.47
1:R:43:ARG:NH2	1:R:45:ASP:OD2	2.38	0.47
1:C:20:ILE:HD13	1:C:117:LEU:HD21	1.96	0.47
1:H:82:LEU:HB3	1:U:32:TYR:OH	2.15	0.47
1:G:81:PHE:HD1	1:K:91:ASP:OD2	1.97	0.47
1:O:22:ARG:HD3	5:O:305:HOH:O	2.13	0.47
1:P:118:HIS:NE2	1:R:127:PRO:HB3	2.30	0.47
1:H:160:ALA:HB1	1:H:161:PRO:HA	1.97	0.47
1:I:112:GLN:O	1:I:116:GLU:HG3	2.14	0.47
1:M:97:LEU:HD23	1:M:161:PRO:HD3	1.95	0.47
1:N:33:VAL:HG22	1:N:88:PRO:HB3	1.96	0.47
1:N:85:ILE:HB	1:P:85:ILE:HB	1.97	0.47
1:W:41:PHE:CZ	1:W:96:GLY:HA2	2.49	0.47
1:X:129:LEU:O	1:X:133:ILE:HG12	2.15	0.47
1:D:50:ASN:HB2	1:D:171:ASP:OD2	2.14	0.47
1:R:107:GLU:OE2	1:R:141:GLN:NE2	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:V:78:GLY:O	1:V:79:ARG:NH1	2.47	0.47
1:B:44:ASP:OD2	5:B:302:HOH:O	2.20	0.47
1:L:122:HIS:CE1	5:L:339:HOH:O	2.68	0.47
1:R:21:ASN:HA	1:R:24:ILE:HD12	1.97	0.47
1:A:128:HIS:CE1	1:B:142:VAL:HG21	2.50	0.46
1:C:131:ASP:HA	1:C:134:GLU:HB2	1.96	0.46
1:E:36:SER:HB2	1:I:82:LEU:HD12	1.97	0.46
1:F:13:HIS:CE1	1:F:15:ASP:OD2	2.68	0.46
1:J:161:PRO:HD2	1:J:162:GLU:OE2	2.14	0.46
1:F:162:GLU:OE1	1:F:162:GLU:N	2.45	0.46
1:H:64:GLU:OE2	1:H:68:LYS:HE3	2.15	0.46
1:T:8:VAL:HB	1:U:145:ILE:HG22	1.96	0.46
1:T:130:ALA:O	1:T:134:GLU:HG3	2.16	0.46
1:L:79:ARG:HA	1:L:79:ARG:HD3	1.64	0.46
1:B:141:GLN:NE2	5:B:307:HOH:O	2.42	0.46
1:Q:128:HIS:CD2	1:R:142:VAL:HG21	2.50	0.46
1:C:79:ARG:HA	1:C:79:ARG:HD3	1.75	0.46
1:F:33:VAL:HG11	1:F:106:LEU:HD22	1.98	0.46
1:G:20:ILE:HD13	1:G:117:LEU:HD11	1.98	0.46
1:H:97:LEU:HD23	1:H:161:PRO:HG3	1.97	0.46
1:K:13:HIS:CD2	1:K:124:LYS:HD2	2.51	0.46
1:R:164:GLY:HA3	1:U:157:LYS:O	2.15	0.46
1:O:24:ILE:HD13	1:O:70:MET:HG2	1.97	0.46
1:H:173:HIS:HB3	1:K:172:LYS:HG2	1.98	0.46
1:P:66:ALA:O	1:P:70:MET:HG3	2.16	0.46
1:Q:44:ASP:OD2	5:Q:301:HOH:O	2.21	0.46
1:S:134:GLU:HA	1:S:138:LEU:HD12	1.98	0.46
1:D:20:ILE:HD13	1:D:117:LEU:HD11	1.98	0.46
1:E:131:ASP:HB2	1:F:134:GLU:HG2	1.98	0.46
1:O:119:LYS:HE2	1:O:123:ASP:OD2	2.16	0.46
1:P:108:LYS:HE3	1:R:7:GLN:O	2.15	0.46
1:B:71:LYS:O	1:B:75:GLN:HG3	2.15	0.46
1:E:23:GLN:NE2	1:E:26:LEU:HD23	2.31	0.46
1:E:155:LEU:HA	1:E:155:LEU:HD23	1.70	0.46
1:N:41:PHE:CZ	1:N:96:GLY:HA2	2.51	0.46
1:E:67:GLU:HB3	1:I:39:TYR:OH	2.16	0.45
1:X:143:LYS:HB3	1:X:143:LYS:HE2	1.65	0.45
1:C:140:GLU:N	1:C:140:GLU:OE2	2.47	0.45
1:D:66:ALA:O	1:D:70:MET:HG3	2.16	0.45
1:I:62:GLU:O	1:I:65:HIS:HB2	2.17	0.45
1:N:13:HIS:CG	1:N:124:LYS:HD2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:82:LEU:HD12	1:D:36:SER:HB2	1.97	0.45
1:X:78:GLY:O	1:X:79:ARG:NH1	2.48	0.45
1:E:82:LEU:HB3	1:I:32:TYR:OH	2.17	0.45
1:F:114:LEU:HD12	1:F:141:GLN:HG3	1.97	0.45
1:O:157:LYS:HE2	5:P:333:HOH:O	2.16	0.45
1:C:117:LEU:HG	1:C:133:ILE:CD1	2.45	0.45
1:E:23:GLN:HE22	1:E:26:LEU:HD23	1.81	0.45
1:P:38:SER:OG	1:P:52:ALA:O	2.27	0.45
1:R:43:ARG:HH21	1:R:45:ASP:CG	2.18	0.45
1:R:73:GLN:OE1	1:R:78:GLY:HA3	2.16	0.45
1:S:85:ILE:HB	1:W:85:ILE:HB	1.98	0.45
1:R:108:LYS:HE3	1:R:108:LYS:HB3	1.81	0.45
1:A:7:GLN:HG3	1:A:8:VAL:HG13	1.99	0.45
1:C:82:LEU:O	1:M:87:LYS:HD2	2.17	0.45
1:E:86:GLN:H	1:E:86:GLN:HG2	1.46	0.45
1:H:21:ASN:OD1	1:H:73:GLN:NE2	2.50	0.45
1:O:177:ASP:HA	5:O:327:HOH:O	2.15	0.45
1:R:139:ASN:ND2	1:R:143:LYS:HD2	2.31	0.45
1:I:106:LEU:O	1:I:110:VAL:HG23	2.16	0.45
1:T:162:GLU:H	1:T:162:GLU:CD	2.19	0.45
1:X:13:HIS:ND1	1:X:15:ASP:HB2	2.32	0.45
1:A:92:ASP:OD1	1:A:94:GLU:N	2.47	0.45
1:J:79:ARG:HD3	1:J:79:ARG:HA	1.84	0.45
1:S:63:ARG:NH2	1:W:59:SER:OG	2.49	0.45
1:U:50:ASN:HB2	1:U:171:ASP:CG	2.36	0.45
1:J:118:HIS:CE1	1:L:127:PRO:HB3	2.52	0.44
1:N:42:ASP:HB3	1:P:74:ASN:HD22	1.82	0.44
1:N:71:LYS:O	1:N:75:GLN:HG3	2.17	0.44
1:Q:8:VAL:HB	1:R:145:ILE:HG22	1.98	0.44
1:T:40:TYR:CD1	1:T:43:ARG:HD3	2.52	0.44
1:U:100:MET:HB3	1:U:152:VAL:HG23	1.99	0.44
1:B:63:ARG:HG2	5:B:371:HOH:O	2.16	0.44
1:H:44:ASP:HA	1:S:150:ASP:OD1	2.17	0.44
1:I:78:GLY:O	1:I:79:ARG:NH1	2.51	0.44
1:O:138:LEU:HD23	1:O:138:LEU:HA	1.77	0.44
1:Q:97:LEU:O	1:Q:101:GLU:HG3	2.17	0.44
1:T:9:ARG:NE	1:T:12:TYR:O	2.49	0.44
1:W:7:GLN:O	1:X:108:LYS:HE3	2.17	0.44
1:F:146:LYS:NZ	1:I:42:ASP:O	2.49	0.44
1:J:134:GLU:HA	1:J:138:LEU:HD12	1.98	0.44
1:D:9:ARG:HD2	5:D:338:HOH:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:13:HIS:O	1:G:16:SER:HB2	2.18	0.44
1:O:40:TYR:HA	5:O:372:HOH:O	2.18	0.44
1:S:97:LEU:HD23	1:S:161:PRO:HD3	2.00	0.44
1:W:138:LEU:HD23	1:W:138:LEU:HA	1.81	0.44
1:D:115:LEU:HD23	1:D:115:LEU:HA	1.79	0.44
1:F:125:ASN:O	1:F:127:PRO:HD3	2.18	0.44
1:J:63:ARG:NH1	5:J:310:HOH:O	2.50	0.44
1:K:155:LEU:HA	1:K:155:LEU:HD23	1.73	0.44
1:M:20:ILE:HD13	1:M:117:LEU:HD11	1.99	0.44
1:A:176:GLY:O	1:A:177:ASP:HB2	2.17	0.44
1:N:59:SER:OG	1:P:63:ARG:NH2	2.50	0.44
1:T:168:TYR:CE1	1:T:172:LYS:HE2	2.53	0.44
1:C:33:VAL:O	1:C:36:SER:HB3	2.18	0.44
1:K:40:TYR:CD1	1:K:43:ARG:HD3	2.53	0.44
1:P:108:LYS:NZ	5:P:304:HOH:O	2.36	0.44
1:U:93:TRP:O	1:U:95:SER:N	2.51	0.44
1:B:87:LYS:N	1:D:84:ASP:OD1	2.38	0.44
1:B:141:GLN:NE2	5:B:304:HOH:O	2.22	0.44
1:E:13:HIS:NE2	1:E:124:LYS:HE2	2.33	0.44
1:I:13:HIS:HE1	1:I:15:ASP:OD2	2.01	0.44
1:M:101:GLU:OE2	1:M:156:ARG:NH2	2.47	0.44
1:P:145:ILE:HG22	1:R:8:VAL:HB	1.99	0.44
1:Q:161:PRO:HD2	1:Q:162:GLU:OE2	2.18	0.44
1:U:129:LEU:O	1:U:133:ILE:HG12	2.18	0.44
1:A:8:VAL:HB	1:B:145:ILE:HG22	1.99	0.44
1:B:27:GLU:HB2	1:B:66:ALA:HB2	1.99	0.44
1:B:41:PHE:CZ	1:B:96:GLY:HA2	2.52	0.44
1:F:79:ARG:HD3	1:F:79:ARG:HA	1.83	0.44
1:K:37:MET:HA	1:K:93:TRP:CD1	2.52	0.44
1:V:160:ALA:HB1	1:V:161:PRO:HA	2.00	0.44
1:F:111:ASN:O	1:F:115:LEU:HD12	2.18	0.43
1:L:6:SER:OG	1:L:8:VAL:HG22	2.18	0.43
1:M:129:LEU:O	1:M:133:ILE:HG12	2.18	0.43
1:P:28:LEU:HA	1:P:28:LEU:HD23	1.80	0.43
1:P:50:ASN:HB2	1:P:171:ASP:OD2	2.18	0.43
1:S:111:ASN:ND2	1:U:10:GLN:OE1	2.38	0.43
1:P:62:GLU:O	1:P:65:HIS:HB2	2.18	0.43
1:R:66:ALA:O	1:R:70:MET:HG3	2.18	0.43
1:T:95:SER:HG	1:T:98:ASN:H	1.65	0.43
1:T:114:LEU:HD13	1:T:137:TYR:HB3	2.00	0.43
1:B:37:MET:HA	1:B:93:TRP:CD1	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:ASP:HB3	1:L:47:ALA:HB1	2.00	0.43
1:F:20:ILE:HD13	1:F:117:LEU:HD11	1.99	0.43
1:H:40:TYR:CD1	1:H:43:ARG:HD3	2.53	0.43
1:U:6:SER:OG	1:U:8:VAL:HG22	2.18	0.43
1:A:118:HIS:NE2	1:A:134:GLU:OE1	2.46	0.43
1:D:108:LYS:NZ	5:D:307:HOH:O	2.44	0.43
1:E:50:ASN:HB2	1:E:171:ASP:OD2	2.18	0.43
1:E:138:LEU:HD23	1:E:138:LEU:HA	1.73	0.43
1:G:114:LEU:HD13	1:G:137:TYR:HB3	2.01	0.43
1:H:32:TYR:HE1	1:U:82:LEU:HD13	1.83	0.43
1:K:164:GLY:N	5:K:310:HOH:O	2.50	0.43
1:O:27:GLU:HB2	1:O:66:ALA:HB2	2.00	0.43
1:E:117:LEU:HD22	1:E:133:ILE:HD11	2.01	0.43
1:T:85:ILE:HA	5:T:306:HOH:O	2.19	0.43
1:X:16:SER:O	1:X:20:ILE:HG12	2.18	0.43
1:X:31:SER:HB2	1:X:62:GLU:HB2	2.00	0.43
1:E:20:ILE:O	1:E:24:ILE:HG13	2.18	0.43
1:H:161:PRO:HD2	1:H:162:GLU:OE2	2.18	0.43
1:M:132:PHE:CE1	1:M:137:TYR:HE1	2.37	0.43
1:B:127:PRO:HB3	1:C:118:HIS:NE2	2.33	0.43
1:E:32:TYR:OH	1:I:82:LEU:HB3	2.19	0.43
1:G:81:PHE:CD1	1:K:91:ASP:OD2	2.72	0.43
1:K:106:LEU:O	1:K:109:ASN:HB2	2.18	0.43
1:N:40:TYR:CD1	1:N:43:ARG:HD3	2.54	0.43
1:S:141:GLN:NE2	5:S:309:HOH:O	2.49	0.43
1:W:48:LEU:HD23	1:W:48:LEU:HA	1.81	0.43
1:I:72:LEU:C	1:I:72:LEU:HD13	2.39	0.43
1:Q:163:SER:HB3	1:Q:166:ALA:CB	2.49	0.43
1:C:85:ILE:HB	1:M:85:ILE:HB	2.00	0.43
1:G:118:HIS:CE1	1:I:127:PRO:CB	3.02	0.43
1:H:150:ASP:OD1	1:K:44:ASP:HA	2.18	0.43
1:R:33:VAL:HG11	1:R:106:LEU:HD22	2.01	0.43
1:U:134:GLU:HA	1:U:138:LEU:HD12	2.00	0.43
1:W:155:LEU:HD23	1:W:155:LEU:HA	1.88	0.43
1:X:120:LEU:O	1:X:124:LYS:HG2	2.19	0.43
1:J:97:LEU:HD23	1:J:161:PRO:HD3	2.01	0.43
1:X:37:MET:HE1	1:X:55:PHE:CE1	2.54	0.42
1:H:138:LEU:HD23	1:H:138:LEU:HA	1.82	0.42
1:P:29:TYR:CD1	1:P:86:GLN:HB2	2.54	0.42
1:W:101:GLU:OE2	1:W:156:ARG:NH2	2.45	0.42
1:F:50:ASN:HB2	1:F:171:ASP:OD2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:37:MET:O	1:L:41:PHE:HD2	2.02	0.42
1:Q:155:LEU:HD23	1:Q:155:LEU:HA	1.57	0.42
1:C:66:ALA:O	1:C:70:MET:HG3	2.19	0.42
1:O:156:ARG:NH1	5:O:316:HOH:O	2.51	0.42
1:T:131:ASP:HB2	1:U:134:GLU:HG2	2.00	0.42
1:U:79:ARG:HD3	1:U:79:ARG:HA	1.75	0.42
1:E:164:GLY:N	5:E:310:HOH:O	2.51	0.42
1:G:128:HIS:CE1	1:H:142:VAL:HG21	2.53	0.42
1:Q:130:ALA:O	1:Q:134:GLU:HG3	2.19	0.42
1:B:61:GLU:O	1:B:64:GLU:HB2	2.19	0.42
1:E:37:MET:HE2	1:E:37:MET:HB3	1.75	0.42
1:N:138:LEU:HD23	1:N:138:LEU:HA	1.76	0.42
1:X:79:ARG:HA	1:X:79:ARG:HD3	1.77	0.42
1:H:157:LYS:O	1:K:164:GLY:HA3	2.20	0.42
1:I:158:MET:HB3	1:U:165:LEU:HD13	2.02	0.42
1:J:165:LEU:HA	1:J:165:LEU:HD12	1.82	0.42
1:L:69:LEU:HG	1:L:137:TYR:OH	2.20	0.42
1:Q:138:LEU:HD23	1:Q:138:LEU:HA	1.81	0.42
1:D:71:LYS:O	1:D:75:GLN:HG3	2.20	0.42
1:G:125:ASN:O	1:G:127:PRO:HD3	2.19	0.42
1:H:69:LEU:HG	1:H:137:TYR:OH	2.20	0.42
1:W:13:HIS:CD2	1:W:124:LYS:HD2	2.55	0.42
1:E:49:LYS:O	1:E:53:LYS:HG3	2.20	0.42
1:F:129:LEU:O	1:F:133:ILE:HG12	2.20	0.42
1:R:36:SER:HB2	1:T:82:LEU:HD12	2.02	0.42
1:S:157:LYS:HB3	1:S:157:LYS:HE3	1.67	0.42
1:T:63:ARG:O	1:T:67:GLU:HG3	2.19	0.42
1:W:40:TYR:CD1	1:W:43:ARG:HD3	2.55	0.42
1:A:127:PRO:HD2	5:A:318:HOH:O	2.18	0.41
1:B:7:GLN:O	1:C:108:LYS:HE3	2.19	0.41
1:G:30:ALA:HA	1:G:106:LEU:HD21	2.02	0.41
1:T:114:LEU:HA	1:T:114:LEU:HD23	1.82	0.41
1:V:24:ILE:HD13	1:V:70:MET:HG2	2.02	0.41
1:F:82:LEU:O	1:Q:87:LYS:HD2	2.20	0.41
1:F:148:LEU:HD23	1:F:148:LEU:HA	1.95	0.41
1:J:111:ASN:ND2	1:L:10:GLN:OE1	2.43	0.41
1:L:93:TRP:O	1:L:95:SER:N	2.50	0.41
1:C:76:ARG:HD2	1:C:76:ARG:HA	1.83	0.41
1:E:8:VAL:HB	1:F:145:ILE:HG22	2.02	0.41
1:E:39:TYR:OH	1:I:67:GLU:HG2	2.20	0.41
1:O:82:LEU:HD12	1:V:36:SER:HB2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:76:ARG:HD3	1:G:128:HIS:CE1	2.54	0.41
1:I:139:ASN:ND2	1:I:143:LYS:HD2	2.35	0.41
1:L:139:ASN:ND2	5:L:310:HOH:O	2.53	0.41
1:W:27:GLU:HB2	1:W:66:ALA:HB2	2.01	0.41
1:X:76:ARG:HD2	1:X:76:ARG:HA	1.76	0.41
1:A:37:MET:O	1:A:40:TYR:HB3	2.20	0.41
1:B:130:ALA:O	1:B:134:GLU:HG3	2.21	0.41
1:D:62:GLU:O	1:D:65:HIS:HB2	2.20	0.41
1:E:165:LEU:HA	1:E:165:LEU:HD12	1.78	0.41
1:H:85:ILE:HA	5:H:312:HOH:O	2.20	0.41
1:R:71:LYS:O	1:R:75:GLN:HG3	2.20	0.41
1:S:29:TYR:CD1	1:S:86:GLN:HB2	2.56	0.41
1:T:24:ILE:HD13	1:T:70:MET:HG2	2.03	0.41
1:T:76:ARG:HD2	1:T:76:ARG:HA	1.78	0.41
1:W:44:ASP:OD2	5:W:302:HOH:O	2.21	0.41
1:B:134:GLU:HA	1:B:138:LEU:HD12	2.03	0.41
1:H:85:ILE:HB	1:U:85:ILE:HB	2.02	0.41
1:K:127:PRO:HB3	1:L:118:HIS:CE1	2.55	0.41
1:N:24:ILE:HD13	1:N:70:MET:HG2	2.02	0.41
1:O:20:ILE:HD11	1:O:129:LEU:HD11	2.02	0.41
1:P:87:LYS:NZ	5:P:311:HOH:O	2.53	0.41
1:V:8:VAL:HB	1:W:145:ILE:HG22	2.01	0.41
1:I:79:ARG:HA	1:I:79:ARG:HD3	1.85	0.41
1:R:21:ASN:OD1	1:R:80:ILE:HA	2.20	0.41
1:S:66:ALA:O	1:S:70:MET:HG3	2.21	0.41
1:V:132:PHE:CE1	1:V:137:TYR:HE1	2.39	0.41
1:E:13:HIS:CE1	1:E:15:ASP:OD2	2.72	0.41
1:E:124:LYS:HA	1:E:124:LYS:HD2	1.86	0.41
1:F:76:ARG:HA	1:F:76:ARG:HD2	1.88	0.41
1:H:53:LYS:HG3	1:H:54:TYR:N	2.36	0.41
1:J:41:PHE:CZ	1:J:96:GLY:HA2	2.56	0.41
1:J:132:PHE:CE1	1:J:137:TYR:HE1	2.39	0.41
1:L:73:GLN:OE1	1:L:78:GLY:HA3	2.21	0.41
1:L:170:PHE:CZ	1:L:174:THR:HG21	2.55	0.41
1:N:139:ASN:OD1	1:N:143:LYS:HD2	2.21	0.41
1:O:170:PHE:CE1	1:O:174:THR:HG21	2.56	0.41
1:P:122:HIS:HE1	4:R:203[D]:V9Y:N03	2.19	0.41
1:Q:63:ARG:O	1:Q:67:GLU:HG3	2.21	0.41
1:R:131:ASP:O	1:R:135:THR:HG23	2.21	0.41
1:X:63:ARG:NH1	5:X:304:HOH:O	2.32	0.41
1:A:72:LEU:C	1:A:72:LEU:HD13	2.41	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:106:LEU:HD23	1:D:106:LEU:C	2.42	0.41
1:E:33:VAL:O	1:E:37:MET:HG3	2.21	0.41
1:M:145:ILE:HG22	1:O:8:VAL:HB	2.03	0.41
1:B:93:TRP:O	1:B:94:GLU:HB2	2.20	0.40
1:J:8:VAL:HB	1:K:145:ILE:HG22	2.02	0.40
1:K:49:LYS:HD2	1:K:49:LYS:H	1.86	0.40
1:L:20:ILE:HD13	1:L:20:ILE:HA	1.80	0.40
1:N:118:HIS:NE2	1:N:130:ALA:HB1	2.37	0.40
1:S:23:GLN:NE2	1:S:114:LEU:HG	2.36	0.40
1:A:93:TRP:HA	5:A:301:HOH:O	2.21	0.40
1:E:82:LEU:CD1	1:I:36:SER:HB2	2.51	0.40
1:J:69:LEU:HG	1:J:137:TYR:OH	2.20	0.40
1:M:60:HIS:O	1:M:64:GLU:HG3	2.22	0.40
1:N:4:SER:HB2	1:N:5:THR:H	1.54	0.40
1:Q:26:LEU:O	1:Q:29:TYR:HB3	2.21	0.40
1:X:20:ILE:HD11	1:X:129:LEU:HD11	2.02	0.40
1:I:97:LEU:CD2	1:I:161:PRO:HD3	2.51	0.40
1:O:97:LEU:HD23	1:O:161:PRO:HD3	2.03	0.40
1:W:20:ILE:HD12	1:W:72:LEU:HD12	2.04	0.40
1:B:63:ARG:NH1	5:B:314:HOH:O	2.51	0.40
1:C:44:ASP:HA	1:N:150:ASP:OD1	2.22	0.40
1:D:117:LEU:O	1:D:117:LEU:HD23	2.21	0.40
1:G:122:HIS:HE1	4:I:203[C]:V9Y:N03	2.19	0.40
1:X:33:VAL:O	1:X:36:SER:HB3	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	172/182 (94%)	168 (98%)	4 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	172/182 (94%)	166 (96%)	5 (3%)	1 (1%)	25	26
1	C	172/182 (94%)	166 (96%)	6 (4%)	0	100	100
1	D	172/182 (94%)	163 (95%)	8 (5%)	1 (1%)	25	26
1	E	172/182 (94%)	161 (94%)	10 (6%)	1 (1%)	25	26
1	F	172/182 (94%)	163 (95%)	9 (5%)	0	100	100
1	G	172/182 (94%)	164 (95%)	8 (5%)	0	100	100
1	H	172/182 (94%)	161 (94%)	10 (6%)	1 (1%)	25	26
1	I	172/182 (94%)	162 (94%)	9 (5%)	1 (1%)	25	26
1	J	172/182 (94%)	166 (96%)	6 (4%)	0	100	100
1	K	172/182 (94%)	164 (95%)	8 (5%)	0	100	100
1	L	172/182 (94%)	166 (96%)	6 (4%)	0	100	100
1	M	172/182 (94%)	169 (98%)	3 (2%)	0	100	100
1	N	172/182 (94%)	165 (96%)	6 (4%)	1 (1%)	25	26
1	O	172/182 (94%)	167 (97%)	5 (3%)	0	100	100
1	P	172/182 (94%)	166 (96%)	6 (4%)	0	100	100
1	Q	172/182 (94%)	158 (92%)	13 (8%)	1 (1%)	25	26
1	R	172/182 (94%)	163 (95%)	9 (5%)	0	100	100
1	S	172/182 (94%)	163 (95%)	9 (5%)	0	100	100
1	T	172/182 (94%)	164 (95%)	7 (4%)	1 (1%)	25	26
1	U	172/182 (94%)	163 (95%)	8 (5%)	1 (1%)	25	26
1	V	172/182 (94%)	170 (99%)	2 (1%)	0	100	100
1	W	172/182 (94%)	167 (97%)	5 (3%)	0	100	100
1	X	172/182 (94%)	165 (96%)	7 (4%)	0	100	100
1	a	344/182 (189%)	320 (93%)	24 (7%)	0	100	100
1	b	344/182 (189%)	320 (93%)	20 (6%)	4 (1%)	13	11
1	c	344/182 (189%)	328 (95%)	10 (3%)	6 (2%)	9	6
1	d	344/182 (189%)	324 (94%)	16 (5%)	4 (1%)	13	11
1	e	344/182 (189%)	320 (93%)	24 (7%)	0	100	100
1	f	344/182 (189%)	322 (94%)	20 (6%)	2 (1%)	25	26
1	g	344/182 (189%)	330 (96%)	14 (4%)	0	100	100
1	h	344/182 (189%)	312 (91%)	30 (9%)	2 (1%)	25	26

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	i	344/182 (189%)	328 (95%)	12 (4%)	4 (1%)	13	11
1	j	344/182 (189%)	326 (95%)	14 (4%)	4 (1%)	13	11
1	k	344/182 (189%)	320 (93%)	22 (6%)	2 (1%)	25	26
1	l	344/182 (189%)	318 (92%)	24 (7%)	2 (1%)	25	26
1	m	344/182 (189%)	326 (95%)	18 (5%)	0	100	100
1	n	344/182 (189%)	320 (93%)	24 (7%)	0	100	100
1	o	344/182 (189%)	318 (92%)	22 (6%)	4 (1%)	13	11
1	p	344/182 (189%)	318 (92%)	20 (6%)	6 (2%)	9	6
1	q	344/182 (189%)	328 (95%)	16 (5%)	0	100	100
1	r	344/182 (189%)	332 (96%)	10 (3%)	2 (1%)	25	26
1	s	344/182 (189%)	334 (97%)	8 (2%)	2 (1%)	25	26
1	t	344/182 (189%)	328 (95%)	14 (4%)	2 (1%)	25	26
1	u	344/182 (189%)	328 (95%)	14 (4%)	2 (1%)	25	26
1	v	344/182 (189%)	322 (94%)	22 (6%)	0	100	100
1	w	344/182 (189%)	326 (95%)	16 (5%)	2 (1%)	25	26
1	x	344/182 (189%)	322 (94%)	18 (5%)	4 (1%)	13	11
All	All	12384/8736 (142%)	11720 (95%)	601 (5%)	63 (0%)	34	31

All (63) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	f	84[C]	ASP
1	f	84[D]	ASP
1	i	94[C]	GLU
1	i	94[D]	GLU
1	o	163[C]	SER
1	o	163[D]	SER
1	Q	84	ASP
1	T	84	ASP
1	c	162[C]	GLU
1	c	162[D]	GLU
1	c	163[C]	SER
1	c	163[D]	SER
1	d	135[C]	THR
1	d	135[D]	THR
1	k	94[C]	GLU

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Mol	Chain	Res	Type
1	k	94[D]	GLU
1	H	84	ASP
1	I	84	ASP
1	N	84	ASP
1	U	84	ASP
1	c	84[C]	ASP
1	c	84[D]	ASP
1	o	162[C]	GLU
1	o	162[D]	GLU
1	r	94[C]	GLU
1	r	94[D]	GLU
1	x	168[C]	TYR
1	x	168[D]	TYR
1	B	84	ASP
1	b	94[C]	GLU
1	b	94[D]	GLU
1	d	94[C]	GLU
1	d	94[D]	GLU
1	i	84[C]	ASP
1	i	84[D]	ASP
1	p	17[C]	GLU
1	p	17[D]	GLU
1	s	94[C]	GLU
1	s	94[D]	GLU
1	u	94[C]	GLU
1	u	94[D]	GLU
1	D	10	GLN
1	j	47[C]	ALA
1	j	47[D]	ALA
1	p	10[C]	GLN
1	p	10[D]	GLN
1	p	127[C]	PRO
1	p	127[D]	PRO
1	t	125[C]	ASN
1	t	125[D]	ASN
1	E	84	ASP
1	b	17[C]	GLU
1	b	17[D]	GLU
1	h	85[C]	ILE
1	h	85[D]	ILE
1	j	94[C]	GLU
1	j	94[D]	GLU

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Mol	Chain	Res	Type
1	l	88[C]	PRO
1	l	88[D]	PRO
1	w	94[C]	GLU
1	w	94[D]	GLU
1	x	163[C]	SER
1	x	163[D]	SER

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	153/160 (96%)	151 (99%)	2 (1%)	69 79
1	B	153/160 (96%)	148 (97%)	5 (3%)	38 46
1	C	153/160 (96%)	150 (98%)	3 (2%)	55 66
1	D	153/160 (96%)	150 (98%)	3 (2%)	55 66
1	E	153/160 (96%)	149 (97%)	4 (3%)	46 56
1	F	153/160 (96%)	146 (95%)	7 (5%)	27 33
1	G	153/160 (96%)	150 (98%)	3 (2%)	55 66
1	H	153/160 (96%)	146 (95%)	7 (5%)	27 33
1	I	153/160 (96%)	149 (97%)	4 (3%)	46 56
1	J	153/160 (96%)	150 (98%)	3 (2%)	55 66
1	K	153/160 (96%)	147 (96%)	6 (4%)	32 41
1	L	153/160 (96%)	145 (95%)	8 (5%)	23 28
1	M	153/160 (96%)	149 (97%)	4 (3%)	46 56
1	N	153/160 (96%)	147 (96%)	6 (4%)	32 41
1	O	153/160 (96%)	149 (97%)	4 (3%)	46 56
1	P	153/160 (96%)	149 (97%)	4 (3%)	46 56
1	Q	153/160 (96%)	144 (94%)	9 (6%)	19 22
1	R	153/160 (96%)	147 (96%)	6 (4%)	32 41
1	S	153/160 (96%)	148 (97%)	5 (3%)	38 46

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	T	153/160 (96%)	148 (97%)	5 (3%)	38	46
1	U	153/160 (96%)	147 (96%)	6 (4%)	32	41
1	V	153/160 (96%)	147 (96%)	6 (4%)	32	41
1	W	153/160 (96%)	150 (98%)	3 (2%)	55	66
1	X	153/160 (96%)	149 (97%)	4 (3%)	46	56
All	All	3672/3840 (96%)	3555 (97%)	117 (3%)	39	47

All (117) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	TYR
1	A	117	LEU
1	B	32	TYR
1	B	58	GLN
1	B	91	ASP
1	B	113	SER
1	B	117	LEU
1	C	14	GLN
1	C	32	TYR
1	C	119	LYS
1	D	32	TYR
1	D	61	GLU
1	D	95	SER
1	E	32	TYR
1	E	61	GLU
1	E	86	GLN
1	E	117	LEU
1	F	14	GLN
1	F	32	TYR
1	F	68	LYS
1	F	117	LEU
1	F	146	LYS
1	F	171	ASP
1	F	177	ASP
1	G	6	SER
1	G	32	TYR
1	G	44	ASP
1	H	7	GLN
1	H	32	TYR
1	H	91	ASP

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Mol	Chain	Res	Type
1	H	117	LEU
1	H	157	LYS
1	H	171	ASP
1	H	177	ASP
1	I	32	TYR
1	I	44	ASP
1	I	64	GLU
1	I	157	LYS
1	J	4	SER
1	J	86	GLN
1	J	117	LEU
1	K	4	SER
1	K	6	SER
1	K	32	TYR
1	K	61	GLU
1	K	68	LYS
1	K	117	LEU
1	L	5	THR
1	L	32	TYR
1	L	49	LYS
1	L	53	LYS
1	L	94	GLU
1	L	117	LEU
1	L	140	GLU
1	L	171	ASP
1	M	32	TYR
1	M	123	ASP
1	M	134	GLU
1	M	171	ASP
1	N	4	SER
1	N	32	TYR
1	N	86	GLN
1	N	117	LEU
1	N	125	ASN
1	N	177	ASP
1	O	5	THR
1	O	32	TYR
1	O	87	LYS
1	O	140	GLU
1	P	32	TYR
1	P	53	LYS
1	P	61	GLU

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Mol	Chain	Res	Type
1	P	117	LEU
1	Q	5	THR
1	Q	8	VAL
1	Q	32	TYR
1	Q	49	LYS
1	Q	100	MET
1	Q	117	LEU
1	Q	123	ASP
1	Q	157	LYS
1	Q	177	ASP
1	R	32	TYR
1	R	61	GLU
1	R	64	GLU
1	R	117	LEU
1	R	157	LYS
1	R	171	ASP
1	S	6	SER
1	S	32	TYR
1	S	72	LEU
1	S	117	LEU
1	S	177	ASP
1	T	14	GLN
1	T	32	TYR
1	T	53	LYS
1	T	117	LEU
1	T	123	ASP
1	U	5	THR
1	U	32	TYR
1	U	44	ASP
1	U	117	LEU
1	U	119	LYS
1	U	175	LEU
1	V	4	SER
1	V	6	SER
1	V	32	TYR
1	V	117	LEU
1	V	123	ASP
1	V	171	ASP
1	W	32	TYR
1	W	49	LYS
1	W	91	ASP
1	X	32	TYR

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Mol	Chain	Res	Type
1	X	44	ASP
1	X	53	LYS
1	X	143	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	G	25	ASN
1	J	122	HIS
1	L	57	HIS
1	L	136	HIS
1	N	122	HIS
1	R	25	ASN
1	S	118	HIS
1	T	122	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 226 ligands modelled in this entry, 210 are monoatomic - leaving 16 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	V9Y	A	205[C]	2	9,13,13	4.98	4 (44%)	8,17,17	2.44	3 (37%)
4	V9Y	X	203[D]	2	9,13,13	4.76	4 (44%)	8,17,17	2.56	4 (50%)
4	V9Y	A	205[D]	2	9,13,13	5.01	3 (33%)	8,17,17	2.16	2 (25%)
4	V9Y	O	204[C]	2	9,13,13	5.06	3 (33%)	8,17,17	1.97	2 (25%)
4	V9Y	R	203[C]	2	9,13,13	5.04	4 (44%)	8,17,17	2.33	3 (37%)
4	V9Y	O	204[D]	2	9,13,13	4.78	3 (33%)	8,17,17	2.80	3 (37%)
4	V9Y	R	203[D]	2	9,13,13	5.06	4 (44%)	8,17,17	1.62	2 (25%)
4	V9Y	t	203[C]	2	9,13,13	5.02	4 (44%)	8,17,17	2.16	4 (50%)
4	V9Y	D	204[C]	2	9,13,13	4.86	4 (44%)	8,17,17	2.60	5 (62%)
4	V9Y	D	204[D]	2	9,13,13	4.88	4 (44%)	8,17,17	2.30	3 (37%)
4	V9Y	t	203[D]	2	9,13,13	4.82	3 (33%)	8,17,17	1.84	2 (25%)
4	V9Y	X	203[C]	2	9,13,13	5.17	3 (33%)	8,17,17	1.60	2 (25%)
4	V9Y	L	203[C]	2	9,13,13	4.82	5 (55%)	8,17,17	2.76	5 (62%)
4	V9Y	I	203[C]	2	9,13,13	4.89	3 (33%)	8,17,17	2.10	3 (37%)
4	V9Y	L	203[D]	2	9,13,13	5.08	5 (55%)	8,17,17	2.49	3 (37%)
4	V9Y	I	203[D]	2	9,13,13	5.15	4 (44%)	8,17,17	2.16	3 (37%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	V9Y	A	205[C]	2	-	1/6/12/12	0/1/1/1
4	V9Y	X	203[D]	2	-	2/6/12/12	0/1/1/1
4	V9Y	A	205[D]	2	-	1/6/12/12	0/1/1/1
4	V9Y	O	204[C]	2	-	2/6/12/12	0/1/1/1
4	V9Y	R	203[C]	2	-	0/6/12/12	0/1/1/1
4	V9Y	O	204[D]	2	-	2/6/12/12	0/1/1/1
4	V9Y	R	203[D]	2	-	2/6/12/12	0/1/1/1
4	V9Y	t	203[C]	2	-	0/6/12/12	0/1/1/1
4	V9Y	D	204[C]	2	-	2/6/12/12	0/1/1/1
4	V9Y	D	204[D]	2	-	1/6/12/12	0/1/1/1
4	V9Y	t	203[D]	2	-	2/6/12/12	0/1/1/1
4	V9Y	X	203[C]	2	-	2/6/12/12	0/1/1/1
4	V9Y	L	203[C]	2	-	1/6/12/12	0/1/1/1
4	V9Y	I	203[C]	2	-	2/6/12/12	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	V9Y	L	203[D]	2	-	1/6/12/12	0/1/1/1
4	V9Y	I	203[D]	2	-	0/6/12/12	0/1/1/1

All (60) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	I	203[C]	V9Y	C02-N03	10.74	1.46	1.32
4	X	203[C]	V9Y	C02-N03	10.66	1.46	1.32
4	L	203[D]	V9Y	C02-N03	10.62	1.46	1.32
4	A	205[D]	V9Y	C02-N03	10.60	1.46	1.32
4	X	203[D]	V9Y	C02-N03	10.59	1.46	1.32
4	R	203[C]	V9Y	C02-N03	10.59	1.46	1.32
4	O	204[C]	V9Y	C02-N03	10.58	1.46	1.32
4	A	205[C]	V9Y	C02-N03	10.58	1.46	1.32
4	I	203[D]	V9Y	C02-N03	10.48	1.46	1.32
4	O	204[D]	V9Y	C02-N03	10.48	1.46	1.32
4	t	203[C]	V9Y	C02-N03	10.42	1.46	1.32
4	R	203[D]	V9Y	C02-N03	10.41	1.46	1.32
4	I	203[D]	V9Y	C09-N10	10.36	1.46	1.32
4	X	203[C]	V9Y	C09-N10	10.35	1.46	1.32
4	t	203[D]	V9Y	C02-N03	10.33	1.46	1.32
4	R	203[D]	V9Y	C09-N10	10.25	1.46	1.32
4	D	204[C]	V9Y	C02-N03	10.06	1.45	1.32
4	O	204[C]	V9Y	C09-N10	10.02	1.45	1.32
4	L	203[D]	V9Y	C09-N10	9.94	1.45	1.32
4	t	203[C]	V9Y	C09-N10	9.94	1.45	1.32
4	R	203[C]	V9Y	C09-N10	9.91	1.45	1.32
4	D	204[D]	V9Y	C02-N03	9.88	1.45	1.32
4	L	203[C]	V9Y	C02-N03	9.88	1.45	1.32
4	D	204[D]	V9Y	C09-N10	9.75	1.45	1.32
4	A	205[D]	V9Y	C09-N10	9.64	1.45	1.32
4	D	204[C]	V9Y	C09-N10	9.45	1.45	1.32
4	A	205[C]	V9Y	C09-N10	9.41	1.45	1.32
4	I	203[C]	V9Y	C09-N10	9.32	1.44	1.32
4	L	203[C]	V9Y	C09-N10	9.32	1.44	1.32
4	t	203[D]	V9Y	C09-N10	9.30	1.44	1.32
4	O	204[D]	V9Y	C09-N10	8.88	1.44	1.32
4	X	203[D]	V9Y	C09-N10	8.53	1.43	1.32
4	L	203[C]	V9Y	C05-C02	3.26	1.54	1.49
4	A	205[D]	V9Y	C08-C09	3.21	1.54	1.49
4	I	203[D]	V9Y	C08-C09	3.20	1.54	1.49
4	X	203[C]	V9Y	C08-C09	3.14	1.54	1.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	205[C]	V9Y	C08-C09	3.12	1.54	1.49
4	O	204[C]	V9Y	C08-C09	3.03	1.54	1.49
4	D	204[D]	V9Y	C08-C09	2.98	1.54	1.49
4	D	204[C]	V9Y	C05-C02	2.90	1.54	1.49
4	X	203[D]	V9Y	C05-C02	2.86	1.54	1.49
4	L	203[D]	V9Y	C08-C09	2.81	1.54	1.49
4	t	203[C]	V9Y	C08-C09	2.78	1.54	1.49
4	O	204[D]	V9Y	C08-C09	2.72	1.53	1.49
4	A	205[C]	V9Y	C05-C02	2.59	1.53	1.49
4	D	204[C]	V9Y	C08-C09	2.56	1.53	1.49
4	X	203[D]	V9Y	C08-C09	2.56	1.53	1.49
4	R	203[D]	V9Y	C08-C09	2.53	1.53	1.49
4	L	203[C]	V9Y	C08-C09	2.49	1.53	1.49
4	t	203[C]	V9Y	C05-C02	2.42	1.53	1.49
4	R	203[C]	V9Y	C08-C09	2.41	1.53	1.49
4	t	203[D]	V9Y	C08-C09	2.24	1.53	1.49
4	D	204[D]	V9Y	C05-C02	2.19	1.53	1.49
4	I	203[C]	V9Y	C08-C09	2.14	1.53	1.49
4	L	203[C]	V9Y	O12-C09	-2.07	1.19	1.23
4	R	203[D]	V9Y	C05-C02	2.04	1.52	1.49
4	R	203[C]	V9Y	C05-C02	2.03	1.52	1.49
4	L	203[D]	V9Y	C05-C02	2.02	1.52	1.49
4	L	203[D]	V9Y	O12-C09	-2.02	1.19	1.23
4	I	203[D]	V9Y	C05-C02	2.00	1.52	1.49

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	O	204[D]	V9Y	C08-C09-N10	5.10	123.65	115.48
4	L	203[D]	V9Y	C08-C09-N10	4.98	123.45	115.48
4	A	205[C]	V9Y	O12-C09-N10	-4.34	114.84	122.94
4	O	204[D]	V9Y	O12-C09-N10	-4.30	114.91	122.94
4	A	205[D]	V9Y	C08-C09-N10	4.20	122.22	115.48
4	X	203[D]	V9Y	O12-C09-N10	-4.12	115.24	122.94
4	I	203[D]	V9Y	C08-C09-N10	4.11	122.07	115.48
4	L	203[C]	V9Y	C08-C09-N10	4.06	121.98	115.48
4	L	203[C]	V9Y	O01-C02-N03	-4.04	115.40	122.94
4	D	204[C]	V9Y	C08-C09-N10	3.97	121.84	115.48
4	A	205[C]	V9Y	C08-C09-N10	3.95	121.81	115.48
4	L	203[D]	V9Y	O12-C09-N10	-3.94	115.59	122.94
4	R	203[C]	V9Y	C05-C02-N03	3.83	121.62	115.48
4	A	205[D]	V9Y	O12-C09-N10	-3.83	115.80	122.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	204[D]	V9Y	C08-C09-N10	3.82	121.60	115.48
4	R	203[C]	V9Y	C08-C09-N10	3.75	121.48	115.48
4	O	204[C]	V9Y	C08-C09-N10	3.70	121.41	115.48
4	X	203[D]	V9Y	O11-N10-C09	-3.62	110.70	119.64
4	D	204[D]	V9Y	O12-C09-N10	-3.56	116.29	122.94
4	D	204[C]	V9Y	O12-C09-N10	-3.51	116.39	122.94
4	O	204[D]	V9Y	O11-N10-C09	-3.49	111.01	119.64
4	L	203[C]	V9Y	O12-C09-N10	-3.47	116.47	122.94
4	X	203[D]	V9Y	C08-C09-N10	3.44	120.99	115.48
4	I	203[D]	V9Y	O12-C09-N10	-3.40	116.59	122.94
4	t	203[C]	V9Y	C05-C02-N03	3.38	120.89	115.48
4	I	203[C]	V9Y	O12-C09-N10	-3.37	116.66	122.94
4	R	203[D]	V9Y	C08-C09-N10	3.28	120.74	115.48
4	O	204[C]	V9Y	O12-C09-N10	-3.24	116.89	122.94
4	I	203[C]	V9Y	C08-C09-N10	3.21	120.62	115.48
4	t	203[C]	V9Y	C08-C09-N10	3.21	120.62	115.48
4	D	204[D]	V9Y	C05-C02-N03	3.20	120.61	115.48
4	t	203[C]	V9Y	O12-C09-N10	-3.15	117.07	122.94
4	D	204[C]	V9Y	O01-C02-N03	-3.11	117.13	122.94
4	I	203[C]	V9Y	O11-N10-C09	-3.09	112.02	119.64
4	R	203[C]	V9Y	O12-C09-N10	-3.00	117.35	122.94
4	D	204[C]	V9Y	C05-C02-N03	2.94	120.19	115.48
4	t	203[D]	V9Y	C08-C09-N10	2.91	120.15	115.48
4	t	203[D]	V9Y	O12-C09-N10	-2.82	117.68	122.94
4	X	203[C]	V9Y	O12-C09-N10	-2.80	117.72	122.94
4	X	203[C]	V9Y	C08-C09-N10	2.73	119.86	115.48
4	L	203[C]	V9Y	C05-C02-N03	2.69	119.78	115.48
4	D	204[C]	V9Y	O11-N10-C09	-2.54	113.36	119.64
4	L	203[C]	V9Y	O11-N10-C09	-2.40	113.70	119.64
4	I	203[D]	V9Y	C05-C02-N03	2.34	119.23	115.48
4	A	205[C]	V9Y	O11-N10-C09	-2.28	114.01	119.64
4	R	203[D]	V9Y	O12-C09-N10	-2.17	118.89	122.94
4	X	203[D]	V9Y	O01-C02-N03	-2.11	119.00	122.94
4	t	203[C]	V9Y	O01-C02-N03	-2.04	119.13	122.94
4	L	203[D]	V9Y	O11-N10-C09	-2.02	114.64	119.64

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	205[C]	V9Y	C07-C08-C09-O12
4	A	205[D]	V9Y	O01-C02-C05-C06

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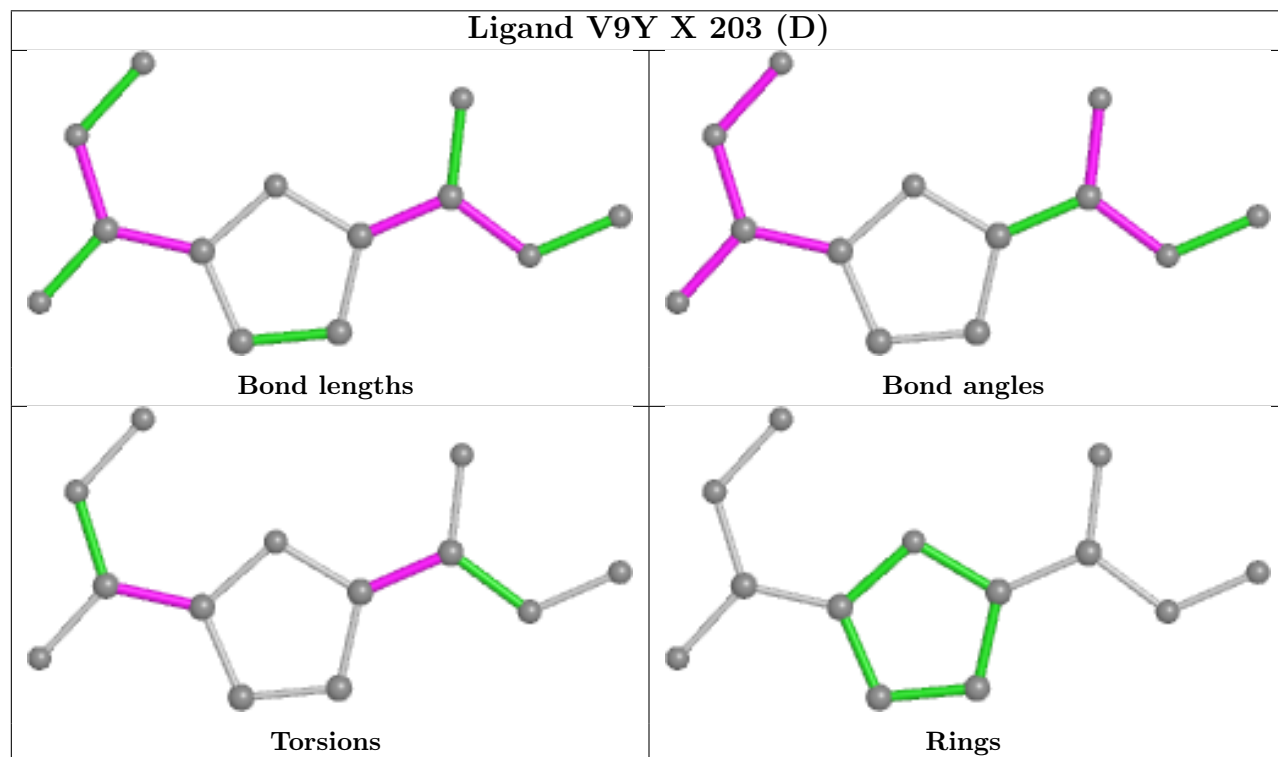
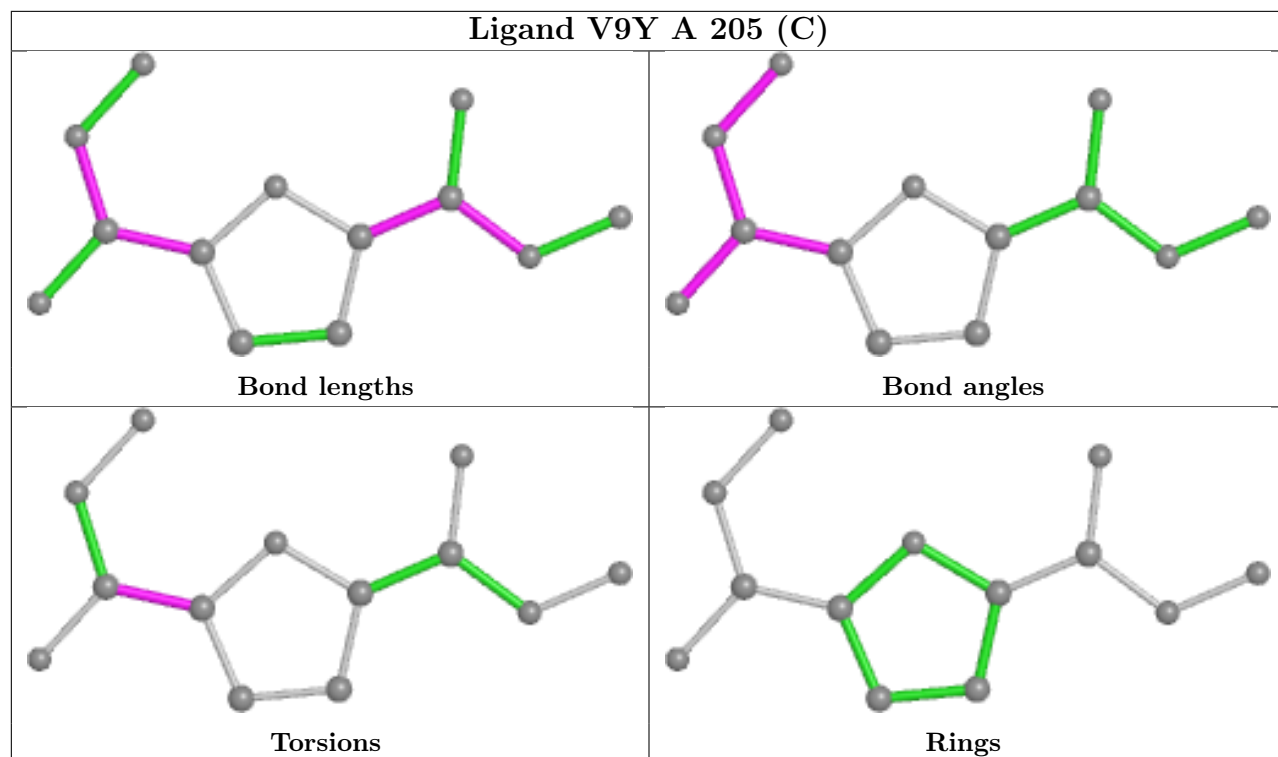
Mol	Chain	Res	Type	Atoms
4	D	204[C]	V9Y	O01-C02-C05-C06
4	D	204[C]	V9Y	C07-C08-C09-O12
4	D	204[D]	V9Y	C07-C08-C09-O12
4	I	203[C]	V9Y	O01-C02-C05-C06
4	I	203[C]	V9Y	C07-C08-C09-O12
4	L	203[C]	V9Y	C07-C08-C09-O12
4	L	203[D]	V9Y	O01-C02-C05-C06
4	O	204[C]	V9Y	O01-C02-C05-C06
4	O	204[C]	V9Y	C07-C08-C09-O12
4	O	204[D]	V9Y	O01-C02-C05-C06
4	O	204[D]	V9Y	C07-C08-C09-O12
4	R	203[D]	V9Y	O01-C02-C05-C06
4	R	203[D]	V9Y	C07-C08-C09-O12
4	X	203[C]	V9Y	O01-C02-C05-C06
4	X	203[C]	V9Y	C07-C08-C09-O12
4	X	203[D]	V9Y	O01-C02-C05-C06
4	X	203[D]	V9Y	C07-C08-C09-O12
4	t	203[D]	V9Y	O01-C02-C05-C06
4	t	203[D]	V9Y	C07-C08-C09-O12

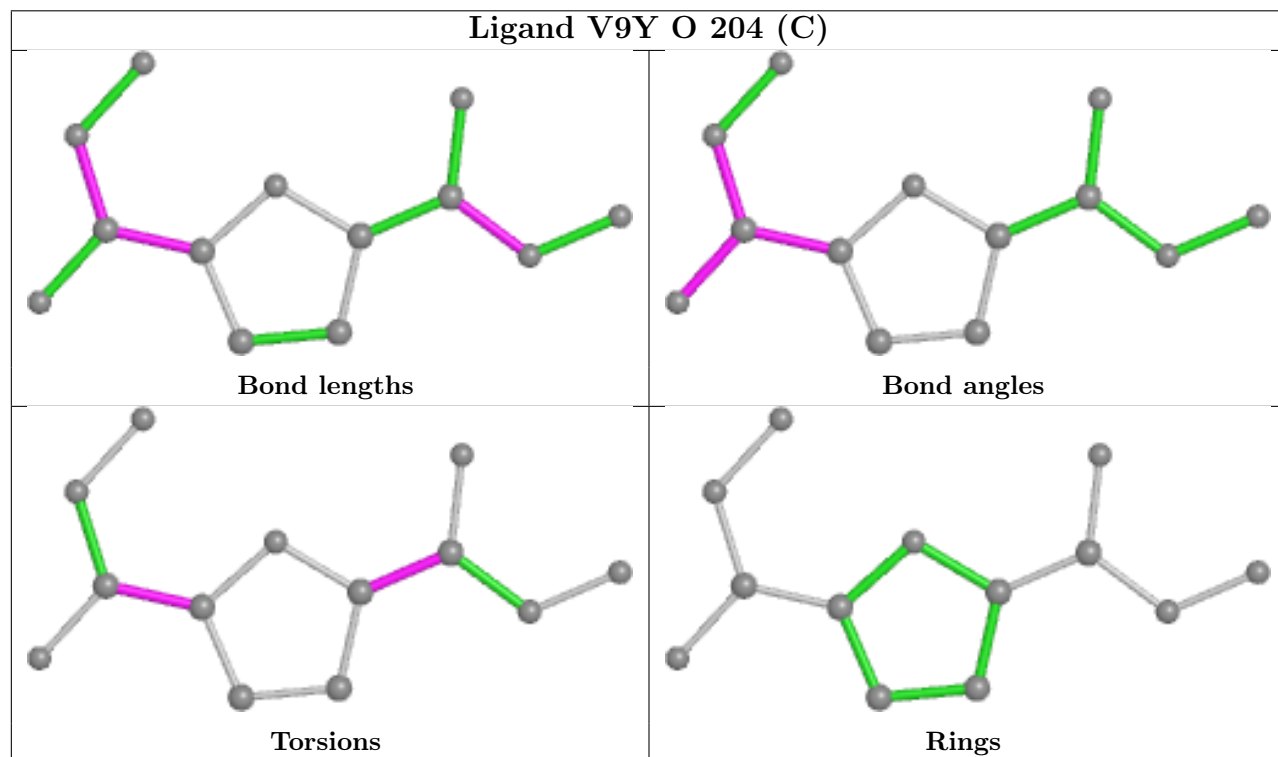
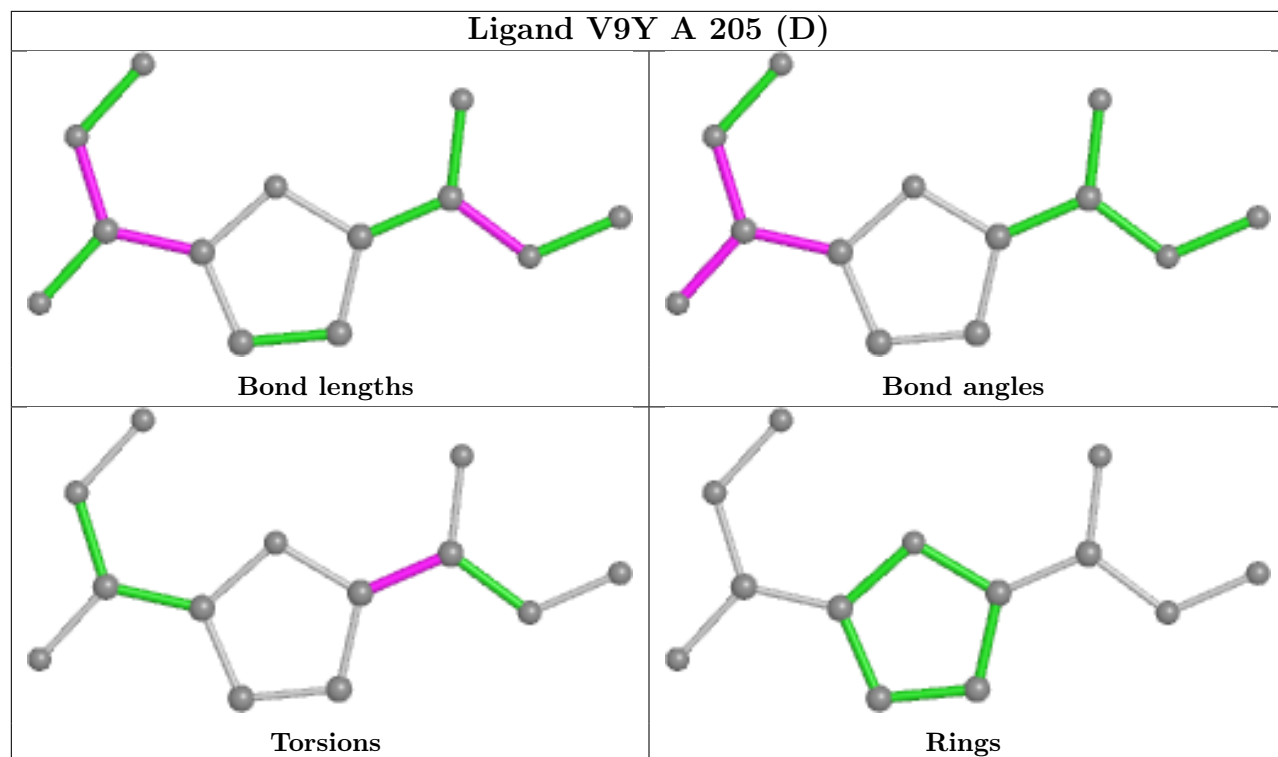
There are no ring outliers.

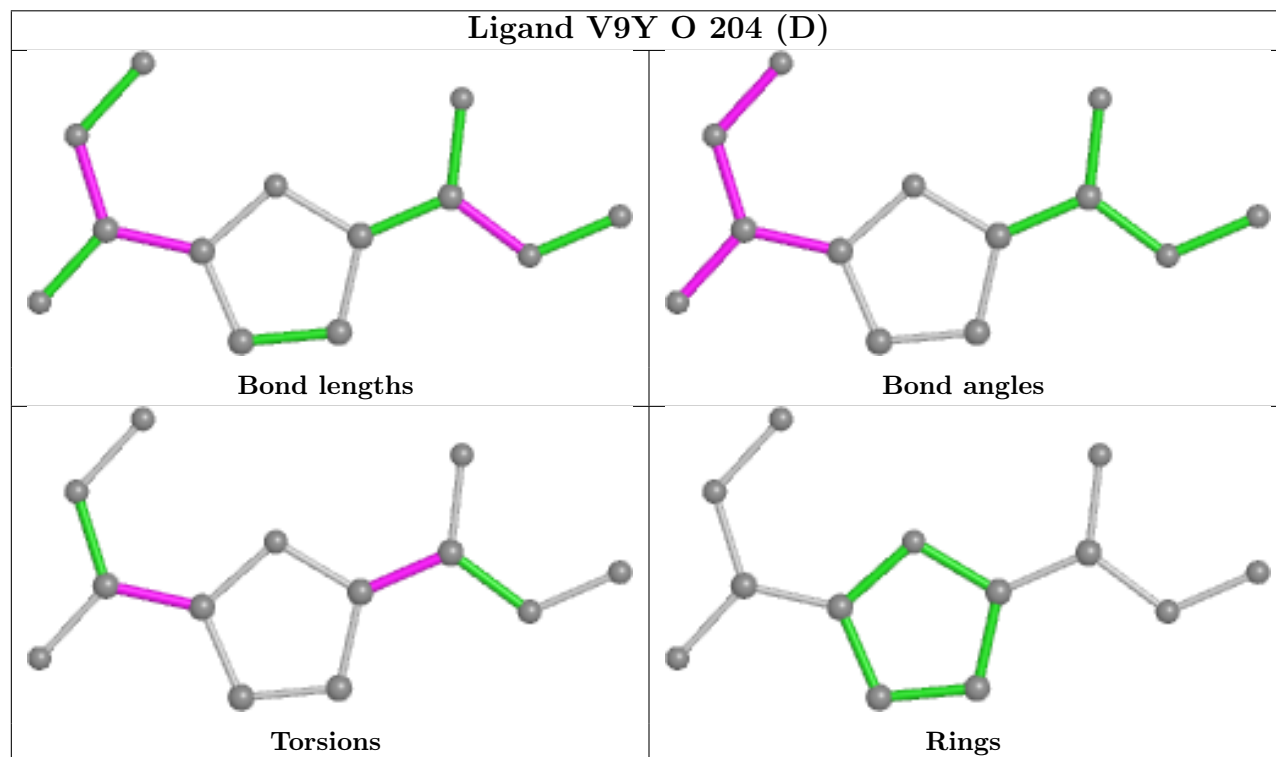
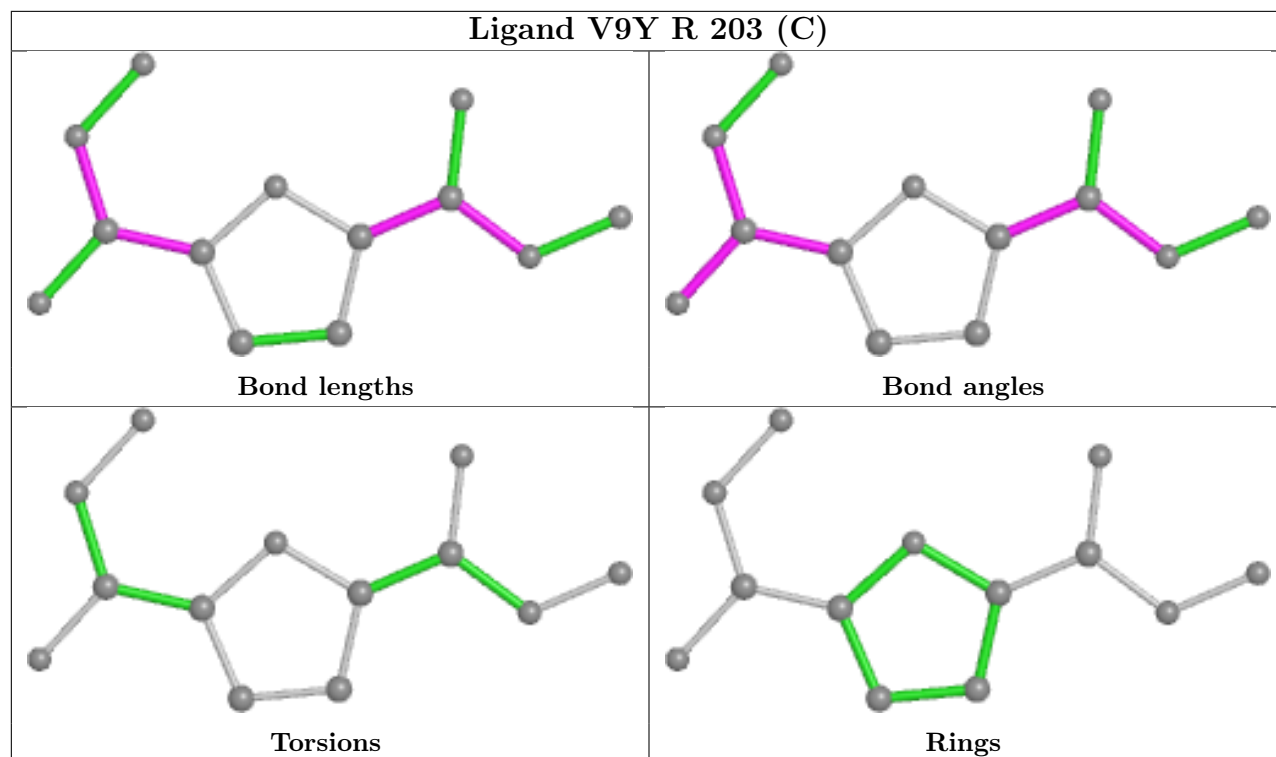
3 monomers are involved in 5 short contacts:

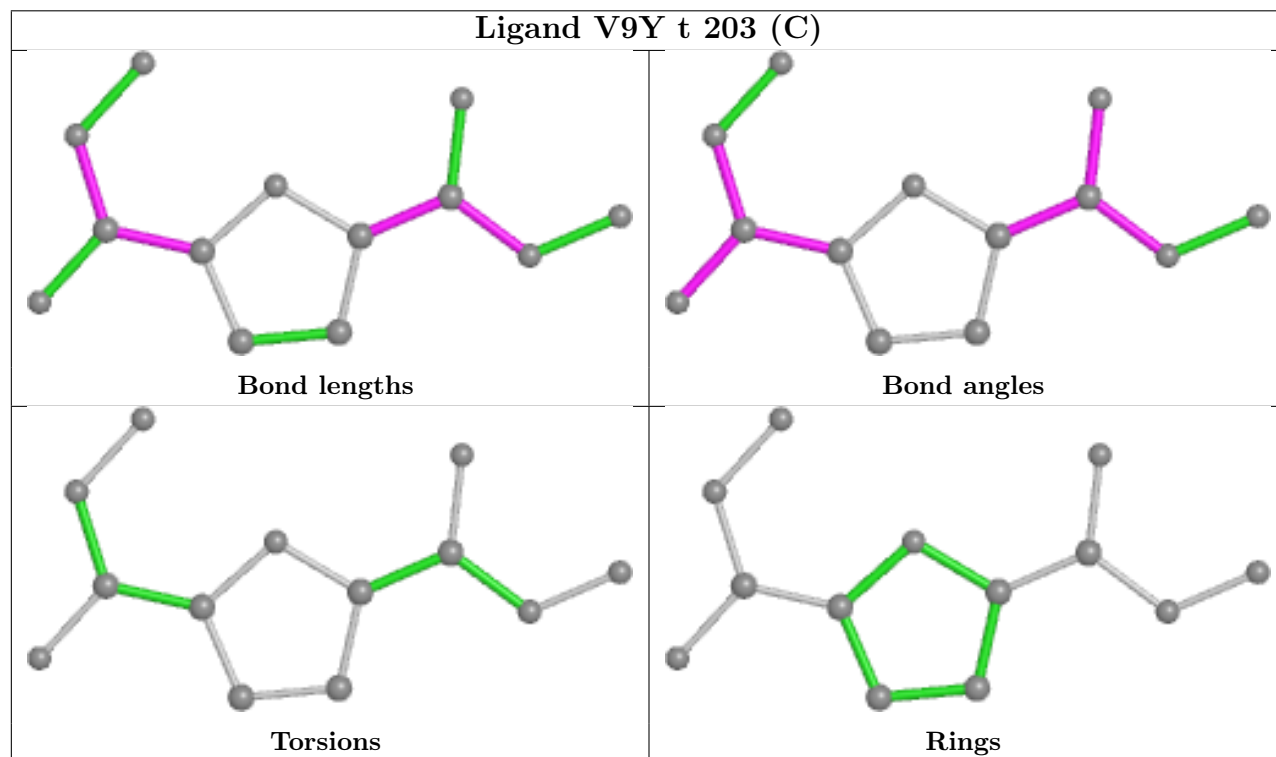
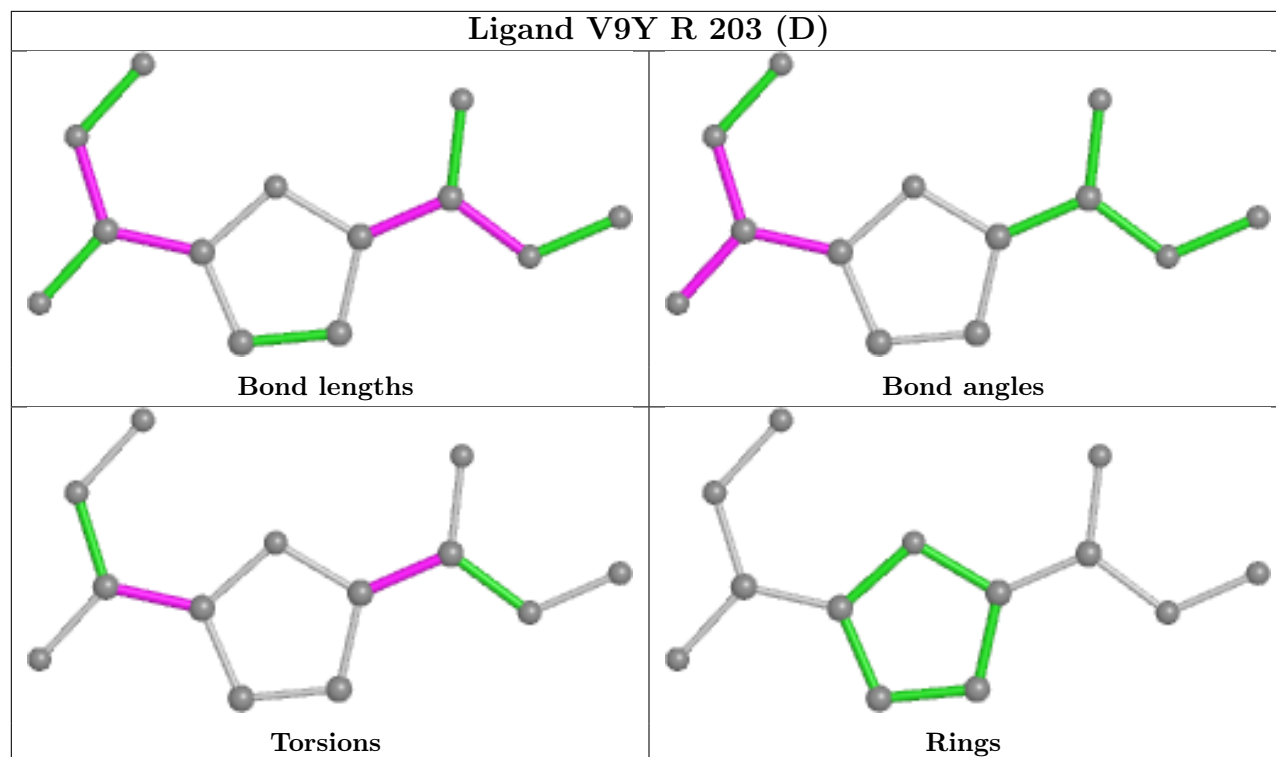
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	R	203[D]	V9Y	2	0
4	D	204[C]	V9Y	1	0
4	I	203[C]	V9Y	2	0

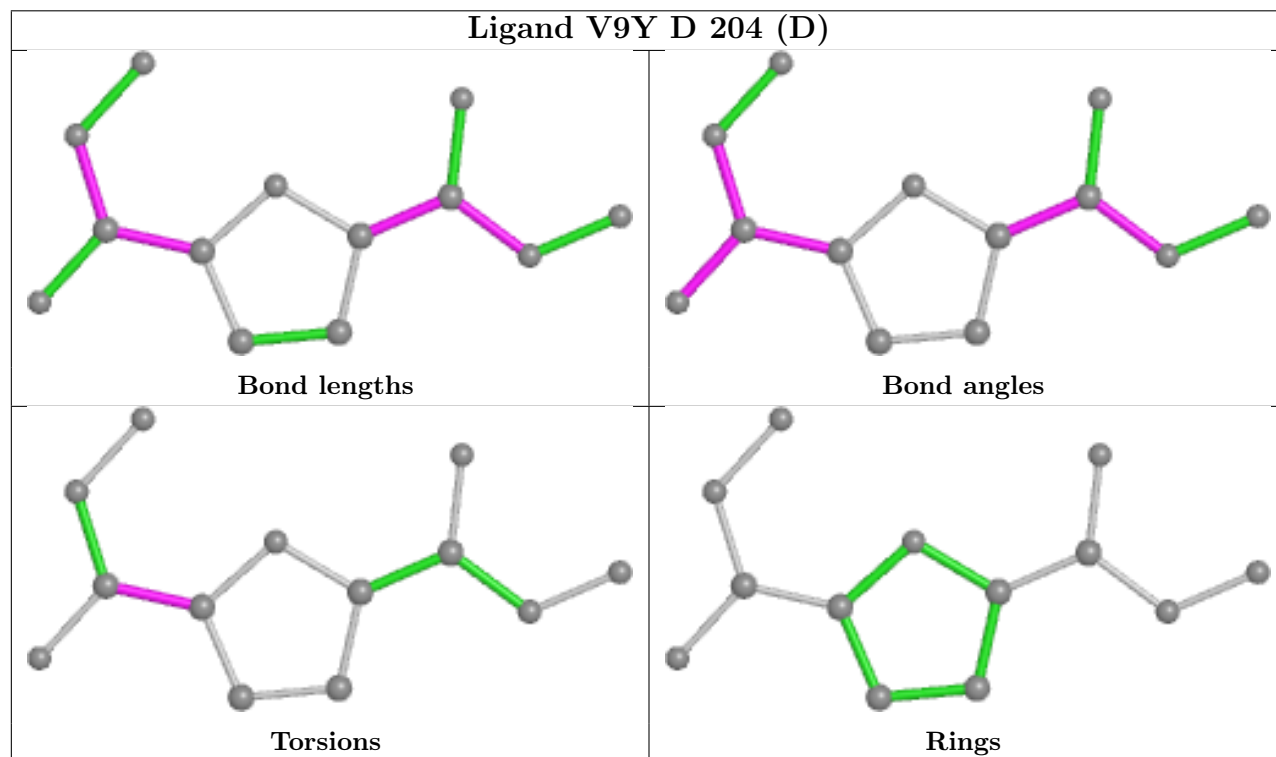
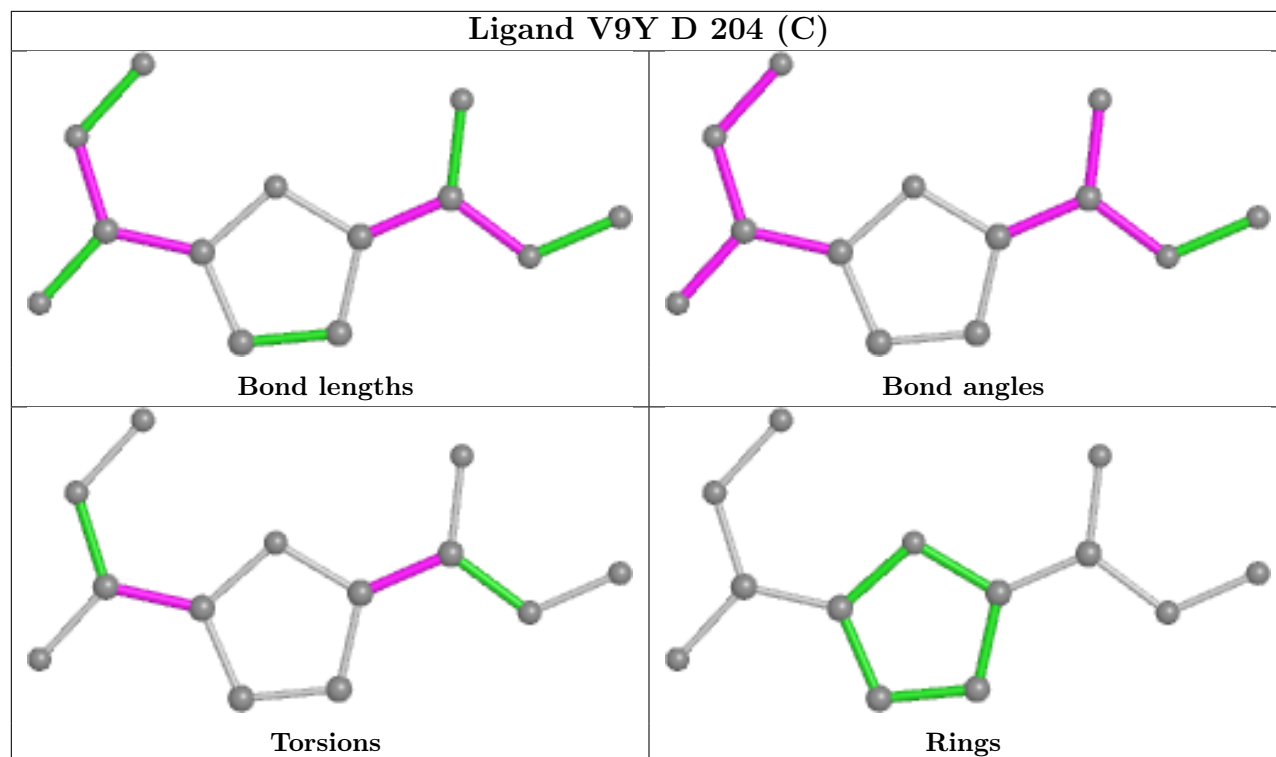
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

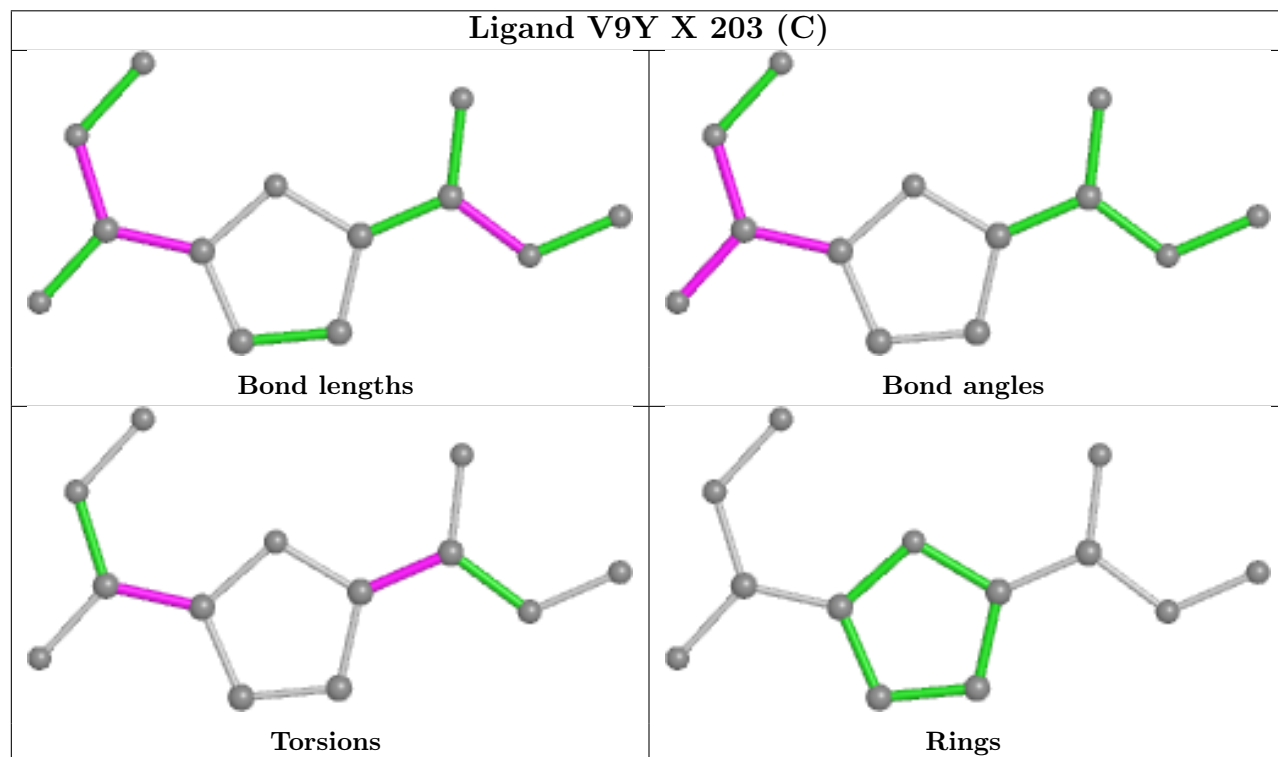
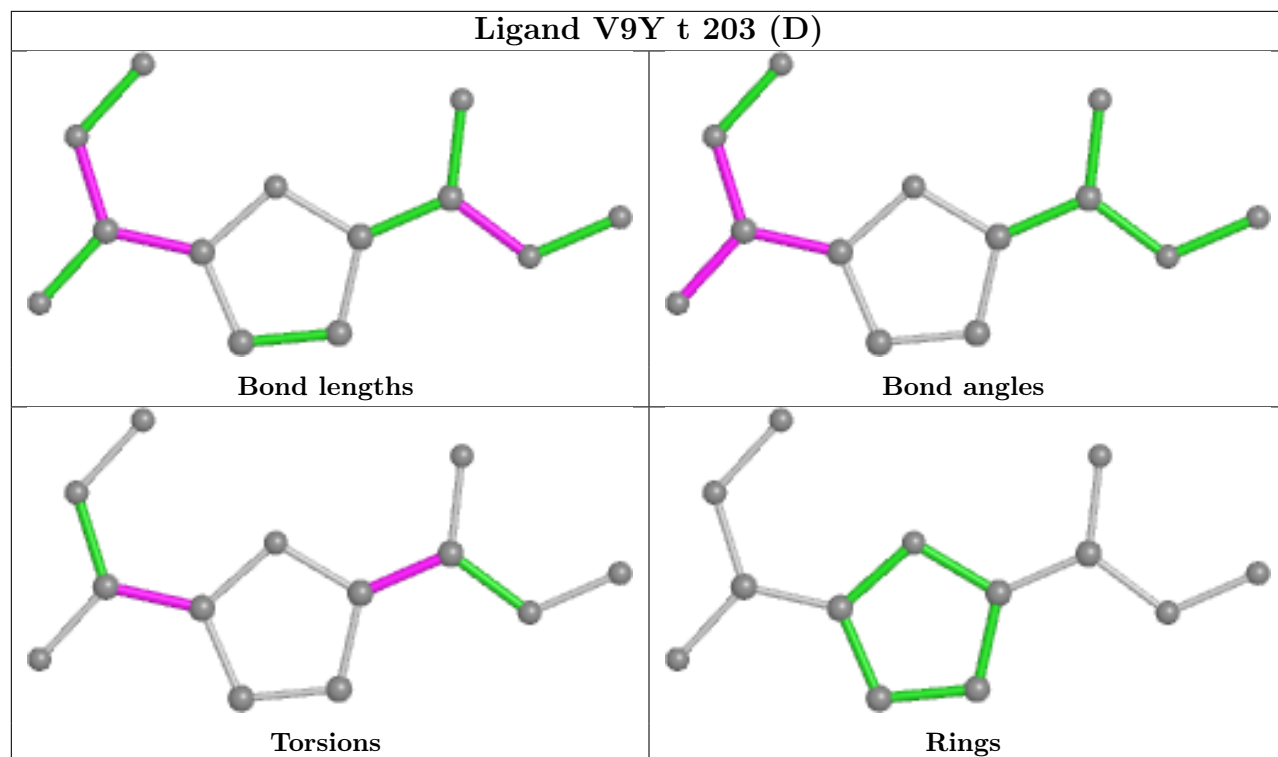


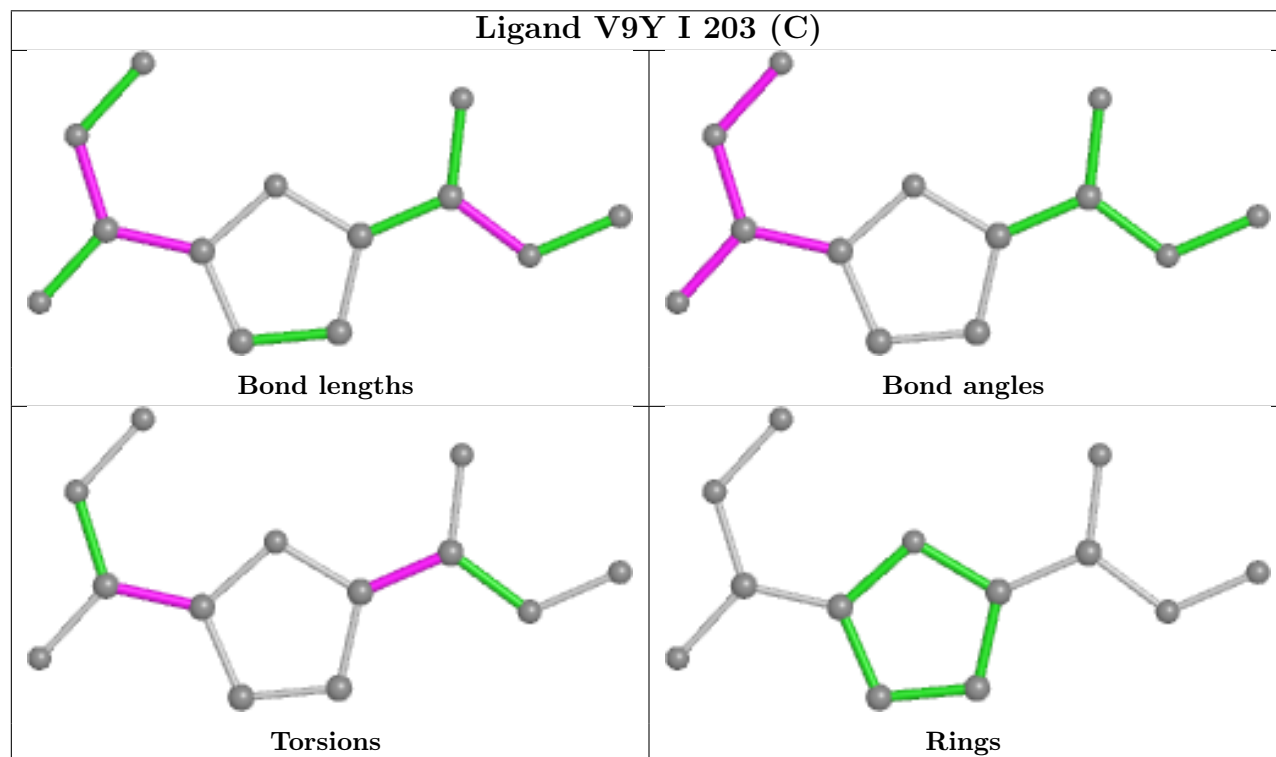
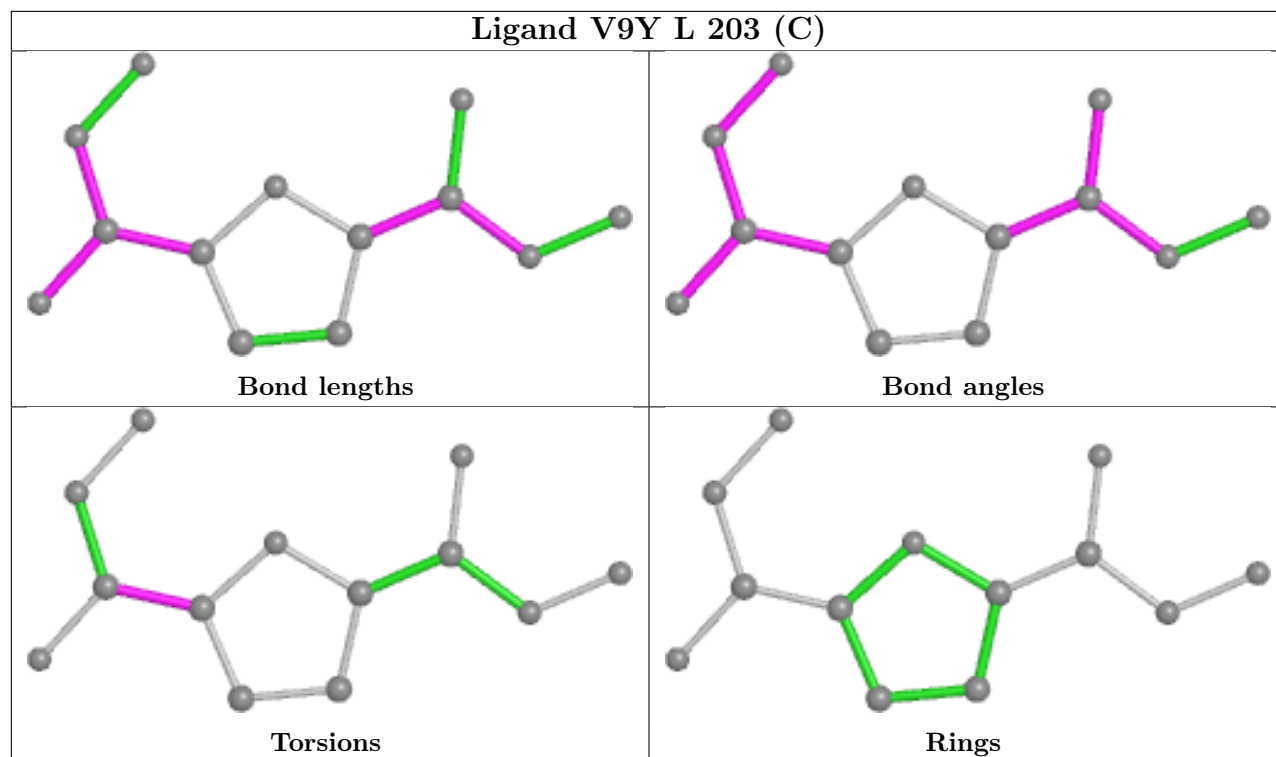


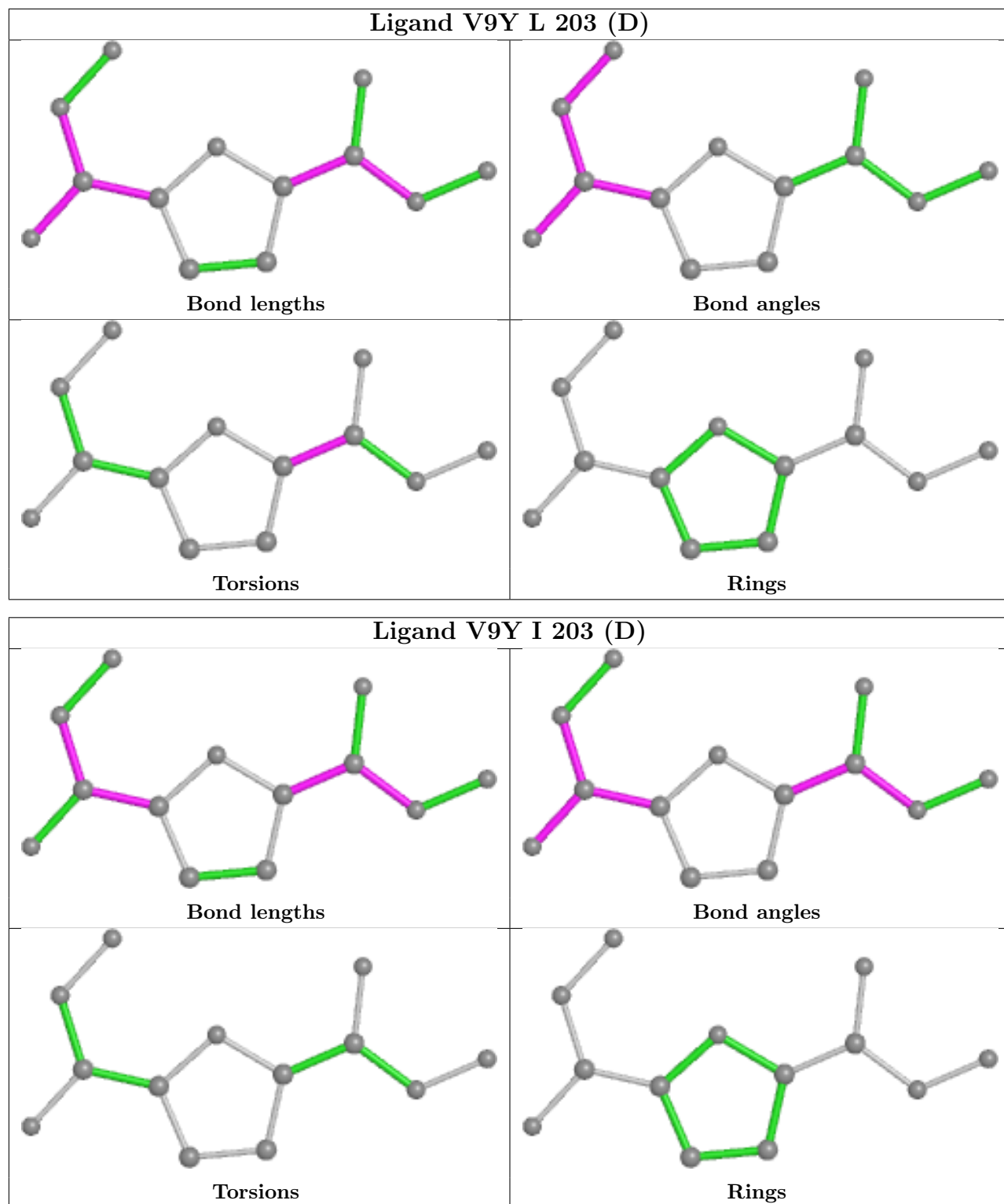












5.7 Other polymers [\(i\)](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data i

6.1 Protein, DNA and RNA chains i

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	174/182 (95%)	-0.18	1 (0%) 89 93	22, 31, 48, 75	0
1	B	174/182 (95%)	-0.14	1 (0%) 89 93	23, 32, 50, 77	0
1	C	174/182 (95%)	-0.17	1 (0%) 89 93	23, 32, 46, 80	0
1	D	174/182 (95%)	-0.10	0 100 100	26, 36, 53, 85	0
1	E	174/182 (95%)	-0.01	1 (0%) 89 93	28, 39, 56, 82	0
1	F	174/182 (95%)	0.05	3 (1%) 70 78	30, 42, 58, 85	0
1	G	174/182 (95%)	-0.31	0 100 100	24, 37, 54, 75	0
1	H	174/182 (95%)	-0.09	2 (1%) 80 86	27, 40, 53, 84	0
1	I	174/182 (95%)	-0.04	2 (1%) 80 86	29, 41, 58, 80	0
1	J	174/182 (95%)	-0.24	0 100 100	23, 30, 46, 66	0
1	K	174/182 (95%)	-0.34	1 (0%) 89 93	25, 33, 53, 75	0
1	L	174/182 (95%)	-0.25	0 100 100	24, 31, 48, 78	0
1	M	174/182 (95%)	-0.12	1 (0%) 89 93	23, 31, 46, 77	0
1	N	174/182 (95%)	-0.32	1 (0%) 89 93	26, 34, 53, 73	0
1	O	174/182 (95%)	-0.28	0 100 100	23, 32, 49, 76	0
1	P	174/182 (95%)	-0.24	0 100 100	26, 38, 55, 75	0
1	Q	174/182 (95%)	-0.06	2 (1%) 80 86	28, 41, 57, 78	0
1	R	174/182 (95%)	0.09	4 (2%) 60 69	31, 42, 61, 88	0
1	S	174/182 (95%)	-0.01	1 (0%) 89 93	23, 36, 51, 77	0
1	T	174/182 (95%)	0.13	7 (4%) 38 49	29, 40, 56, 82	0
1	U	174/182 (95%)	0.05	2 (1%) 80 86	32, 42, 60, 80	0
1	V	174/182 (95%)	-0.17	0 100 100	22, 31, 46, 73	0
1	W	174/182 (95%)	-0.03	1 (0%) 89 93	23, 32, 51, 80	0
1	X	174/182 (95%)	-0.19	1 (0%) 89 93	21, 31, 46, 83	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9	
1	a	174/182 (95%)	1.29	41 (23%)	0	1	29, 33, 42, 56	0
1	b	174/182 (95%)	1.37	44 (25%)	0	1	26, 31, 39, 51	0
1	c	174/182 (95%)	1.27	41 (23%)	0	1	28, 33, 42, 59	0
1	d	174/182 (95%)	1.43	42 (24%)	0	1	28, 33, 42, 55	0
1	e	174/182 (95%)	1.59	51 (29%)	0	0	28, 34, 42, 55	0
1	f	174/182 (95%)	1.29	39 (22%)	0	1	26, 32, 42, 62	0
1	g	174/182 (95%)	1.35	46 (26%)	0	0	27, 32, 40, 52	0
1	h	174/182 (95%)	1.30	38 (21%)	0	1	28, 33, 41, 53	0
1	i	174/182 (95%)	1.20	37 (21%)	0	1	27, 32, 44, 58	0
1	j	174/182 (95%)	1.36	41 (23%)	0	1	28, 33, 41, 55	0
1	k	174/182 (95%)	1.35	40 (22%)	0	1	27, 32, 40, 51	0
1	l	174/182 (95%)	1.38	42 (24%)	0	1	28, 34, 45, 63	0
1	m	174/182 (95%)	1.28	38 (21%)	0	1	30, 35, 45, 58	0
1	n	174/182 (95%)	1.32	38 (21%)	0	1	30, 34, 42, 52	0
1	o	174/182 (95%)	1.59	53 (30%)	0	0	28, 34, 42, 51	0
1	p	174/182 (95%)	1.50	50 (28%)	0	0	29, 35, 47, 70	0
1	q	174/182 (95%)	1.41	44 (25%)	0	1	28, 33, 41, 54	0
1	r	174/182 (95%)	1.45	45 (25%)	0	1	30, 34, 43, 55	0
1	s	174/182 (95%)	1.30	37 (21%)	0	1	27, 32, 39, 54	0
1	t	174/182 (95%)	1.50	50 (28%)	0	0	28, 32, 41, 52	0
1	u	174/182 (95%)	1.49	53 (30%)	0	0	28, 35, 43, 55	0
1	v	174/182 (95%)	1.39	43 (24%)	0	1	29, 34, 43, 55	0
1	w	174/182 (95%)	1.37	44 (25%)	0	1	28, 32, 40, 53	0
1	x	174/182 (95%)	1.28	29 (16%)	1	2	26, 30, 39, 54	0
All	All	8352/8736 (95%)	0.63	1058 (12%)	3	6	21, 34, 51, 88	0

All (1058) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	d	4[C]	SER	9.5
1	d	164[C]	GLY	7.9
1	o	129[C]	LEU	7.3
1	p	24[C]	ILE	7.3
1	o	110[C]	VAL	7.2

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Mol	Chain	Res	Type	RSRZ
1	n	127[C]	PRO	6.7
1	q	4[C]	SER	6.6
1	j	133[C]	ILE	6.6
1	p	4[C]	SER	6.5
1	u	129[C]	LEU	6.4
1	n	114[C]	LEU	6.3
1	f	20[C]	ILE	5.9
1	a	166[C]	ALA	5.9
1	m	104[C]	LEU	5.7
1	w	133[C]	ILE	5.6
1	u	24[C]	ILE	5.6
1	a	176[C]	GLY	5.5
1	Q	4	SER	5.4
1	j	165[C]	LEU	5.4
1	q	8[C]	VAL	5.3
1	t	8[C]	VAL	5.3
1	k	24[C]	ILE	5.3
1	o	24[C]	ILE	5.3
1	u	114[C]	LEU	5.3
1	p	104[C]	LEU	5.3
1	r	129[C]	LEU	5.2
1	i	130[C]	ALA	5.2
1	b	110[C]	VAL	5.1
1	j	164[C]	GLY	5.1
1	h	129[C]	LEU	5.1
1	d	177[C]	ASP	5.1
1	j	166[C]	ALA	5.1
1	w	165[C]	LEU	5.0
1	o	133[C]	ILE	5.0
1	q	164[C]	GLY	5.0
1	n	165[C]	LEU	4.9
1	o	130[C]	ALA	4.9
1	x	133[C]	ILE	4.9
1	i	133[C]	ILE	4.9
1	g	4[C]	SER	4.8
1	g	165[C]	LEU	4.8
1	x	24[C]	ILE	4.8
1	l	119[C]	LYS	4.8
1	n	8[C]	VAL	4.8
1	v	28[C]	LEU	4.8
1	h	24[C]	ILE	4.7
1	r	33[C]	VAL	4.7

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Mol	Chain	Res	Type	RSRZ
1	p	20[C]	ILE	4.7
1	r	161[C]	PRO	4.7
1	c	130[C]	ALA	4.7
1	k	104[C]	LEU	4.7
1	x	129[C]	LEU	4.7
1	g	164[C]	GLY	4.6
1	f	24[C]	ILE	4.6
1	g	114[C]	LEU	4.6
1	c	20[C]	ILE	4.6
1	l	24[C]	ILE	4.6
1	p	127[C]	PRO	4.6
1	u	133[C]	ILE	4.6
1	t	129[C]	LEU	4.6
1	W	4	SER	4.6
1	k	148[C]	LEU	4.6
1	b	20[C]	ILE	4.5
1	g	166[C]	ALA	4.5
1	t	133[C]	ILE	4.5
1	f	127[C]	PRO	4.5
1	n	169[C]	LEU	4.5
1	q	133[C]	ILE	4.5
1	t	114[C]	LEU	4.5
1	p	41[C]	PHE	4.4
1	p	114[C]	LEU	4.4
1	a	165[C]	LEU	4.4
1	b	123[C]	ASP	4.4
1	c	5[C]	THR	4.3
1	k	20[C]	ILE	4.3
1	b	33[C]	VAL	4.3
1	e	104[C]	LEU	4.3
1	r	105[C]	HIS	4.3
1	l	28[C]	LEU	4.3
1	t	104[C]	LEU	4.3
1	w	8[C]	VAL	4.3
1	e	158[C]	MET	4.2
1	j	129[C]	LEU	4.2
1	r	162[C]	GLU	4.2
1	w	153[C]	THR	4.2
1	g	8[C]	VAL	4.2
1	l	20[C]	ILE	4.2
1	w	158[C]	MET	4.2
1	f	129[C]	LEU	4.2

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Mol	Chain	Res	Type	RSRZ
1	r	127[C]	PRO	4.2
1	o	114[C]	LEU	4.2
1	e	33[C]	VAL	4.2
1	r	144[C]	ALA	4.2
1	s	73[C]	GLN	4.1
1	s	129[C]	LEU	4.1
1	h	33[C]	VAL	4.1
1	t	105[C]	HIS	4.1
1	t	153[C]	THR	4.1
1	d	129[C]	LEU	4.1
1	k	176[C]	GLY	4.1
1	t	176[C]	GLY	4.1
1	t	4[C]	SER	4.1
1	t	76[C]	ARG	4.1
1	g	24[C]	ILE	4.1
1	l	152[C]	VAL	4.1
1	l	32[C]	TYR	4.1
1	p	97[C]	LEU	4.1
1	s	105[C]	HIS	4.0
1	v	97[C]	LEU	4.0
1	o	66[C]	ALA	4.0
1	w	166[C]	ALA	4.0
1	c	127[C]	PRO	4.0
1	x	110[C]	VAL	4.0
1	T	4	SER	4.0
1	q	6[C]	SER	4.0
1	d	72[C]	LEU	4.0
1	d	82[C]	LEU	4.0
1	o	28[C]	LEU	4.0
1	t	115[C]	LEU	4.0
1	v	129[C]	LEU	4.0
1	e	130[C]	ALA	4.0
1	f	130[C]	ALA	4.0
1	j	28[C]	LEU	3.9
1	m	97[C]	LEU	3.9
1	w	76[C]	ARG	3.9
1	e	24[C]	ILE	3.9
1	n	158[C]	MET	3.9
1	r	69[C]	LEU	3.9
1	w	129[C]	LEU	3.9
1	x	130[C]	ALA	3.9
1	e	112[C]	GLN	3.9

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Mol	Chain	Res	Type	RSRZ
1	m	177[C]	ASP	3.9
1	a	34[C]	TYR	3.9
1	p	55[C]	PHE	3.9
1	q	127[C]	PRO	3.9
1	n	111[C]	ASN	3.8
1	d	24[C]	ILE	3.8
1	o	97[C]	LEU	3.8
1	a	4[C]	SER	3.8
1	e	20[C]	ILE	3.8
1	v	18[C]	ALA	3.8
1	e	129[C]	LEU	3.8
1	q	111[C]	ASN	3.8
1	a	164[C]	GLY	3.8
1	t	164[C]	GLY	3.8
1	x	105[C]	HIS	3.8
1	E	119	LYS	3.8
1	q	40[C]	TYR	3.8
1	k	103[C]	ALA	3.7
1	s	5[C]	THR	3.7
1	s	24[C]	ILE	3.7
1	e	69[C]	LEU	3.7
1	h	20[C]	ILE	3.7
1	r	104[C]	LEU	3.7
1	s	20[C]	ILE	3.7
1	e	19[C]	ALA	3.7
1	f	5[C]	THR	3.7
1	e	145[C]	ILE	3.7
1	m	73[C]	GLN	3.7
1	d	176[C]	GLY	3.7
1	h	148[C]	LEU	3.7
1	r	130[C]	ALA	3.6
1	u	130[C]	ALA	3.6
1	l	133[C]	ILE	3.6
1	d	25[C]	ASN	3.6
1	d	165[C]	LEU	3.6
1	t	165[C]	LEU	3.6
1	x	97[C]	LEU	3.6
1	h	159[C]	GLY	3.6
1	t	174[C]	THR	3.6
1	j	40[C]	TYR	3.6
1	j	158[C]	MET	3.6
1	t	28[C]	LEU	3.6

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Mol	Chain	Res	Type	RSRZ
1	c	24[C]	ILE	3.6
1	k	28[C]	LEU	3.6
1	d	166[C]	ALA	3.6
1	t	111[C]	ASN	3.6
1	p	73[C]	GLN	3.6
1	e	148[C]	LEU	3.6
1	o	4[C]	SER	3.6
1	i	152[C]	VAL	3.6
1	s	4[C]	SER	3.6
1	w	28[C]	LEU	3.6
1	x	69[C]	LEU	3.6
1	s	28[C]	LEU	3.5
1	s	152[C]	VAL	3.5
1	t	169[C]	LEU	3.5
1	w	40[C]	TYR	3.5
1	t	103[C]	ALA	3.5
1	u	144[C]	ALA	3.5
1	o	117[C]	LEU	3.5
1	t	40[C]	TYR	3.5
1	g	103[C]	ALA	3.5
1	t	19[C]	ALA	3.5
1	n	76[C]	ARG	3.5
1	o	105[C]	HIS	3.5
1	R	4	SER	3.5
1	j	25[C]	ASN	3.5
1	c	100[C]	MET	3.5
1	v	175[C]	LEU	3.5
1	n	84[C]	ASP	3.5
1	o	81[C]	PHE	3.5
1	d	169[C]	LEU	3.5
1	u	161[C]	PRO	3.5
1	s	112[C]	GLN	3.5
1	j	109[C]	ASN	3.5
1	R	159	GLY	3.5
1	q	176[C]	GLY	3.5
1	u	36[C]	SER	3.5
1	e	44[C]	ASP	3.5
1	x	114[C]	LEU	3.5
1	o	33[C]	VAL	3.5
1	f	32[C]	TYR	3.5
1	r	36[C]	SER	3.5
1	a	112[C]	GLN	3.5

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Mol	Chain	Res	Type	RSRZ
1	g	177[C]	ASP	3.4
1	v	51[C]	PHE	3.4
1	a	169[C]	LEU	3.4
1	b	148[C]	LEU	3.4
1	k	69[C]	LEU	3.4
1	k	117[C]	LEU	3.4
1	t	26[C]	LEU	3.4
1	x	5[C]	THR	3.4
1	a	129[C]	LEU	3.4
1	k	129[C]	LEU	3.4
1	n	104[C]	LEU	3.4
1	o	126[C]	ASP	3.4
1	e	162[C]	GLU	3.4
1	p	29[C]	TYR	3.4
1	h	114[C]	LEU	3.4
1	n	51[C]	PHE	3.4
1	p	28[C]	LEU	3.4
1	s	8[C]	VAL	3.4
1	d	29[C]	TYR	3.4
1	e	5[C]	THR	3.4
1	p	18[C]	ALA	3.4
1	j	76[C]	ARG	3.4
1	k	33[C]	VAL	3.4
1	q	177[C]	ASP	3.4
1	u	60[C]	HIS	3.4
1	g	39[C]	TYR	3.4
1	s	6[C]	SER	3.4
1	n	153[C]	THR	3.4
1	e	103[C]	ALA	3.4
1	e	114[C]	LEU	3.4
1	g	72[C]	LEU	3.4
1	j	24[C]	ILE	3.4
1	k	97[C]	LEU	3.4
1	o	80[C]	ILE	3.4
1	w	36[C]	SER	3.3
1	h	97[C]	LEU	3.3
1	r	24[C]	ILE	3.3
1	r	81[C]	PHE	3.3
1	v	24[C]	ILE	3.3
1	w	164[C]	GLY	3.3
1	f	119[C]	LYS	3.3
1	u	33[C]	VAL	3.3

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Mol	Chain	Res	Type	RSRZ
1	g	40[C]	TYR	3.3
1	k	99[C]	ALA	3.3
1	r	66[C]	ALA	3.3
1	t	177[C]	ASP	3.3
1	e	72[C]	LEU	3.3
1	r	148[C]	LEU	3.3
1	f	78[C]	GLY	3.3
1	o	144[C]	ALA	3.3
1	l	175[C]	LEU	3.3
1	p	169[C]	LEU	3.3
1	t	72[C]	LEU	3.3
1	u	69[C]	LEU	3.3
1	s	133[C]	ILE	3.3
1	i	119[C]	LYS	3.3
1	b	103[C]	ALA	3.3
1	l	102[C]	ALA	3.3
1	r	19[C]	ALA	3.3
1	w	66[C]	ALA	3.3
1	g	82[C]	LEU	3.3
1	m	129[C]	LEU	3.3
1	w	115[C]	LEU	3.3
1	I	4	SER	3.3
1	a	145[C]	ILE	3.3
1	w	4[C]	SER	3.3
1	h	130[C]	ALA	3.2
1	g	36[C]	SER	3.2
1	v	29[C]	TYR	3.2
1	x	104[C]	LEU	3.2
1	u	162[C]	GLU	3.2
1	p	135[C]	THR	3.2
1	p	101[C]	GLU	3.2
1	v	64[C]	GLU	3.2
1	w	6[C]	SER	3.2
1	l	72[C]	LEU	3.2
1	u	138[C]	LEU	3.2
1	k	110[C]	VAL	3.2
1	H	13	HIS	3.2
1	i	129[C]	LEU	3.2
1	o	35[C]	LEU	3.2
1	c	176[C]	GLY	3.2
1	p	105[C]	HIS	3.2
1	o	69[C]	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
1	q	169[C]	LEU	3.2
1	p	93[C]	TRP	3.2
1	v	93[C]	TRP	3.2
1	c	73[C]	GLN	3.2
1	l	66[C]	ALA	3.2
1	p	129[C]	LEU	3.2
1	d	133[C]	ILE	3.2
1	r	173[C]	HIS	3.2
1	j	112[C]	GLN	3.2
1	s	109[C]	ASN	3.2
1	o	106[C]	LEU	3.2
1	l	39[C]	TYR	3.1
1	c	88[C]	PRO	3.1
1	p	78[C]	GLY	3.1
1	c	135[C]	THR	3.1
1	h	110[C]	VAL	3.1
1	v	39[C]	TYR	3.1
1	a	28[C]	LEU	3.1
1	c	129[C]	LEU	3.1
1	q	114[C]	LEU	3.1
1	t	170[C]	PHE	3.1
1	o	125[C]	ASN	3.1
1	g	29[C]	TYR	3.1
1	n	30[C]	ALA	3.1
1	I	164	GLY	3.1
1	f	93[C]	TRP	3.1
1	m	114[C]	LEU	3.1
1	n	72[C]	LEU	3.1
1	q	28[C]	LEU	3.1
1	o	140[C]	GLU	3.1
1	f	133[C]	ILE	3.1
1	e	127[C]	PRO	3.1
1	k	8[C]	VAL	3.1
1	q	102[C]	ALA	3.1
1	v	7[C]	GLN	3.1
1	e	161[C]	PRO	3.1
1	m	20[C]	ILE	3.1
1	g	81[C]	PHE	3.1
1	i	53[C]	LYS	3.1
1	j	49[C]	LYS	3.1
1	c	133[C]	ILE	3.1
1	i	20[C]	ILE	3.1

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Mol	Chain	Res	Type	RSRZ
1	e	84[C]	ASP	3.0
1	m	29[C]	TYR	3.0
1	g	25[C]	ASN	3.0
1	l	4[C]	SER	3.0
1	d	28[C]	LEU	3.0
1	n	164[C]	GLY	3.0
1	t	127[C]	PRO	3.0
1	w	127[C]	PRO	3.0
1	v	4[C]	SER	3.0
1	j	114[C]	LEU	3.0
1	r	97[C]	LEU	3.0
1	x	138[C]	LEU	3.0
1	d	47[C]	ALA	3.0
1	n	85[C]	ILE	3.0
1	b	161[C]	PRO	3.0
1	o	103[C]	ALA	3.0
1	s	130[C]	ALA	3.0
1	g	34[C]	TYR	3.0
1	c	114[C]	LEU	3.0
1	x	28[C]	LEU	3.0
1	m	60[C]	HIS	3.0
1	g	30[C]	ALA	3.0
1	h	66[C]	ALA	3.0
1	q	24[C]	ILE	3.0
1	u	4[C]	SER	3.0
1	j	55[C]	PHE	3.0
1	e	23[C]	GLN	3.0
1	e	142[C]	VAL	3.0
1	i	32[C]	TYR	3.0
1	p	40[C]	TYR	3.0
1	d	151[C]	HIS	3.0
1	t	158[C]	MET	3.0
1	q	124[C]	LYS	3.0
1	i	73[C]	GLN	3.0
1	b	170[C]	PHE	3.0
1	m	93[C]	TRP	3.0
1	p	142[C]	VAL	3.0
1	p	156[C]	ARG	3.0
1	f	114[C]	LEU	2.9
1	n	28[C]	LEU	2.9
1	p	39[C]	TYR	2.9
1	s	114[C]	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	N	4	SER	2.9
1	e	52[C]	ALA	2.9
1	h	133[C]	ILE	2.9
1	m	80[C]	ILE	2.9
1	u	145[C]	ILE	2.9
1	i	51[C]	PHE	2.9
1	v	127[C]	PRO	2.9
1	h	105[C]	HIS	2.9
1	l	36[C]	SER	2.9
1	i	66[C]	ALA	2.9
1	a	133[C]	ILE	2.9
1	l	53[C]	LYS	2.9
1	a	13[C]	HIS	2.9
1	h	128[C]	HIS	2.9
1	s	51[C]	PHE	2.9
1	m	110[C]	VAL	2.9
1	n	46[C]	VAL	2.9
1	a	120[C]	LEU	2.9
1	d	114[C]	LEU	2.9
1	m	4[C]	SER	2.9
1	q	158[C]	MET	2.9
1	a	76[C]	ARG	2.9
1	o	145[C]	ILE	2.9
1	d	105[C]	HIS	2.9
1	h	112[C]	GLN	2.9
1	c	41[C]	PHE	2.9
1	e	36[C]	SER	2.9
1	k	76[C]	ARG	2.9
1	c	66[C]	ALA	2.9
1	j	34[C]	TYR	2.9
1	a	24[C]	ILE	2.9
1	n	133[C]	ILE	2.9
1	q	119[C]	LYS	2.9
1	u	81[C]	PHE	2.9
1	o	31[C]	SER	2.9
1	b	114[C]	LEU	2.9
1	g	129[C]	LEU	2.9
1	j	72[C]	LEU	2.9
1	j	169[C]	LEU	2.9
1	v	72[C]	LEU	2.9
1	w	47[C]	ALA	2.9
1	o	32[C]	TYR	2.9

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Mol	Chain	Res	Type	RSRZ
1	h	145[C]	ILE	2.9
1	m	24[C]	ILE	2.9
1	j	8[C]	VAL	2.9
1	u	28[C]	LEU	2.9
1	v	119[C]	LYS	2.9
1	b	130[C]	ALA	2.9
1	b	144[C]	ALA	2.9
1	g	76[C]	ARG	2.9
1	d	113[C]	SER	2.8
1	o	59[C]	SER	2.8
1	g	111[C]	ASN	2.8
1	x	176[C]	GLY	2.8
1	b	8[C]	VAL	2.8
1	u	152[C]	VAL	2.8
1	M	177	ASP	2.8
1	l	56[C]	LEU	2.8
1	q	138[C]	LEU	2.8
1	g	161[C]	PRO	2.8
1	k	12[C]	TYR	2.8
1	u	32[C]	TYR	2.8
1	p	51[C]	PHE	2.8
1	b	76[C]	ARG	2.8
1	c	106[C]	LEU	2.8
1	f	97[C]	LEU	2.8
1	i	69[C]	LEU	2.8
1	l	148[C]	LEU	2.8
1	m	127[C]	PRO	2.8
1	j	4[C]	SER	2.8
1	m	113[C]	SER	2.8
1	o	12[C]	TYR	2.8
1	e	126[C]	ASP	2.8
1	v	71[C]	LYS	2.8
1	x	33[C]	VAL	2.8
1	e	105[C]	HIS	2.8
1	t	15[C]	ASP	2.8
1	b	168[C]	TYR	2.8
1	u	97[C]	LEU	2.8
1	U	159	GLY	2.8
1	j	124[C]	LYS	2.8
1	n	101[C]	GLU	2.8
1	t	101[C]	GLU	2.8
1	f	41[C]	PHE	2.8

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Mol	Chain	Res	Type	RSRZ
1	j	102[C]	ALA	2.8
1	j	110[C]	VAL	2.8
1	t	30[C]	ALA	2.8
1	b	129[C]	LEU	2.8
1	g	123[C]	ASP	2.8
1	l	114[C]	LEU	2.8
1	p	100[C]	MET	2.8
1	u	5[C]	THR	2.8
1	c	40[C]	TYR	2.8
1	p	118[C]	HIS	2.8
1	r	29[C]	TYR	2.8
1	a	49[C]	LYS	2.8
1	h	124[C]	LYS	2.8
1	m	41[C]	PHE	2.7
1	b	28[C]	LEU	2.7
1	w	82[C]	LEU	2.7
1	e	173[C]	HIS	2.7
1	g	133[C]	ILE	2.7
1	h	161[C]	PRO	2.7
1	f	135[C]	THR	2.7
1	q	5[C]	THR	2.7
1	u	105[C]	HIS	2.7
1	i	127[C]	PRO	2.7
1	s	127[C]	PRO	2.7
1	c	33[C]	VAL	2.7
1	u	8[C]	VAL	2.7
1	b	117[C]	LEU	2.7
1	e	117[C]	LEU	2.7
1	j	138[C]	LEU	2.7
1	l	129[C]	LEU	2.7
1	t	35[C]	LEU	2.7
1	r	76[C]	ARG	2.7
1	e	125[C]	ASN	2.7
1	w	151[C]	HIS	2.7
1	b	145[C]	ILE	2.7
1	n	112[C]	GLN	2.7
1	e	29[C]	TYR	2.7
1	b	98[C]	ASN	2.7
1	h	48[C]	LEU	2.7
1	r	133[C]	ILE	2.7
1	e	66[C]	ALA	2.7
1	l	130[C]	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	s	40[C]	TYR	2.7
1	a	105[C]	HIS	2.7
1	g	28[C]	LEU	2.7
1	i	72[C]	LEU	2.7
1	v	152[C]	VAL	2.7
1	p	5[C]	THR	2.7
1	l	31[C]	SER	2.7
1	k	130[C]	ALA	2.7
1	o	121[C]	ALA	2.7
1	n	40[C]	TYR	2.7
1	k	170[C]	PHE	2.7
1	x	170[C]	PHE	2.7
1	i	156[C]	ARG	2.6
1	o	63[C]	ARG	2.6
1	a	127[C]	PRO	2.6
1	d	64[C]	GLU	2.6
1	e	15[C]	ASP	2.6
1	r	126[C]	ASP	2.6
1	h	103[C]	ALA	2.6
1	t	102[C]	ALA	2.6
1	b	128[C]	HIS	2.6
1	o	128[C]	HIS	2.6
1	d	34[C]	TYR	2.6
1	C	164	GLY	2.6
1	d	8[C]	VAL	2.6
1	r	114[C]	LEU	2.6
1	r	165[C]	LEU	2.6
1	w	51[C]	PHE	2.6
1	s	100[C]	MET	2.6
1	H	4	SER	2.6
1	b	27[C]	GLU	2.6
1	b	31[C]	SER	2.6
1	q	53[C]	LYS	2.6
1	w	101[C]	GLU	2.6
1	h	126[C]	ASP	2.6
1	w	111[C]	ASN	2.6
1	p	173[C]	HIS	2.6
1	i	148[C]	LEU	2.6
1	k	114[C]	LEU	2.6
1	n	82[C]	LEU	2.6
1	n	115[C]	LEU	2.6
1	q	82[C]	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	q	132[C]	PHE	2.6
1	r	132[C]	PHE	2.6
1	t	82[C]	LEU	2.6
1	g	31[C]	SER	2.6
1	q	103[C]	ALA	2.6
1	g	176[C]	GLY	2.6
1	s	53[C]	LYS	2.6
1	u	126[C]	ASP	2.6
1	u	177[C]	ASP	2.6
1	x	126[C]	ASP	2.6
1	e	4[C]	SER	2.6
1	f	51[C]	PHE	2.6
1	m	112[C]	GLN	2.6
1	T	161	PRO	2.6
1	a	119[C]	LYS	2.6
1	k	108[C]	LYS	2.6
1	o	18[C]	ALA	2.6
1	x	173[C]	HIS	2.6
1	m	176[C]	GLY	2.6
1	o	14[C]	GLN	2.6
1	d	120[C]	LEU	2.6
1	w	97[C]	LEU	2.6
1	p	37[C]	MET	2.6
1	g	119[C]	LYS	2.6
1	g	151[C]	HIS	2.6
1	r	128[C]	HIS	2.6
1	a	30[C]	ALA	2.6
1	r	166[C]	ALA	2.6
1	q	36[C]	SER	2.6
1	R	165	LEU	2.6
1	a	72[C]	LEU	2.6
1	c	104[C]	LEU	2.6
1	l	138[C]	LEU	2.6
1	o	127[C]	PRO	2.6
1	d	55[C]	PHE	2.6
1	l	51[C]	PHE	2.6
1	o	132[C]	PHE	2.6
1	x	132[C]	PHE	2.6
1	d	168[C]	TYR	2.6
1	o	60[C]	HIS	2.6
1	v	118[C]	HIS	2.6
1	d	123[C]	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	t	92[C]	ASP	2.6
1	f	103[C]	ALA	2.6
1	l	144[C]	ALA	2.6
1	w	149[C]	GLY	2.6
1	t	64[C]	GLU	2.6
1	l	127[C]	PRO	2.5
1	f	175[C]	LEU	2.5
1	e	55[C]	PHE	2.5
1	r	110[C]	VAL	2.5
1	j	29[C]	TYR	2.5
1	s	39[C]	TYR	2.5
1	o	124[C]	LYS	2.5
1	v	53[C]	LYS	2.5
1	i	109[C]	ASN	2.5
1	n	166[C]	ALA	2.5
1	u	66[C]	ALA	2.5
1	a	85[C]	ILE	2.5
1	f	174[C]	THR	2.5
1	k	5[C]	THR	2.5
1	l	135[C]	THR	2.5
1	l	145[C]	ILE	2.5
1	a	162[C]	GLU	2.5
1	c	138[C]	LEU	2.5
1	m	169[C]	LEU	2.5
1	o	176[C]	GLY	2.5
1	p	33[C]	VAL	2.5
1	f	156[C]	ARG	2.5
1	k	121[C]	ALA	2.5
1	o	160[C]	ALA	2.5
1	u	103[C]	ALA	2.5
1	e	133[C]	ILE	2.5
1	u	153[C]	THR	2.5
1	w	112[C]	GLN	2.5
1	s	119[C]	LYS	2.5
1	g	13[C]	HIS	2.5
1	a	8[C]	VAL	2.5
1	p	8[C]	VAL	2.5
1	v	33[C]	VAL	2.5
1	v	142[C]	VAL	2.5
1	R	16	SER	2.5
1	j	32[C]	TYR	2.5
1	m	64[C]	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	w	162[C]	GLU	2.5
1	u	174[C]	THR	2.5
1	x	80[C]	ILE	2.5
1	d	13[C]	HIS	2.5
1	j	173[C]	HIS	2.5
1	b	97[C]	LEU	2.5
1	e	165[C]	LEU	2.5
1	i	155[C]	LEU	2.5
1	l	106[C]	LEU	2.5
1	g	55[C]	PHE	2.5
1	v	170[C]	PHE	2.5
1	e	30[C]	ALA	2.5
1	m	103[C]	ALA	2.5
1	t	166[C]	ALA	2.5
1	s	32[C]	TYR	2.5
1	r	145[C]	ILE	2.5
1	k	128[C]	HIS	2.5
1	e	48[C]	LEU	2.5
1	i	28[C]	LEU	2.5
1	w	49[C]	LYS	2.5
1	h	44[C]	ASP	2.5
1	f	76[C]	ARG	2.5
1	w	32[C]	TYR	2.5
1	n	151[C]	HIS	2.5
1	m	172[C]	LYS	2.5
1	t	85[C]	ILE	2.5
1	t	119[C]	LYS	2.5
1	h	14[C]	GLN	2.5
1	t	112[C]	GLN	2.5
1	a	177[C]	ASP	2.5
1	g	69[C]	LEU	2.5
1	v	106[C]	LEU	2.5
1	w	169[C]	LEU	2.5
1	f	4[C]	SER	2.4
1	a	55[C]	PHE	2.4
1	c	46[C]	VAL	2.4
1	c	51[C]	PHE	2.4
1	e	110[C]	VAL	2.4
1	e	144[C]	ALA	2.4
1	f	102[C]	ALA	2.4
1	l	55[C]	PHE	2.4
1	q	170[C]	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
1	c	71[C]	LYS	2.4
1	g	49[C]	LYS	2.4
1	l	105[C]	HIS	2.4
1	p	77[C]	GLY	2.4
1	r	32[C]	TYR	2.4
1	u	173[C]	HIS	2.4
1	k	112[C]	GLN	2.4
1	e	76[C]	ARG	2.4
1	s	138[C]	LEU	2.4
1	t	31[C]	SER	2.4
1	v	76[C]	ARG	2.4
1	h	117[C]	LEU	2.4
1	q	129[C]	LEU	2.4
1	g	102[C]	ALA	2.4
1	m	111[C]	ASN	2.4
1	i	39[C]	TYR	2.4
1	k	39[C]	TYR	2.4
1	u	44[C]	ASP	2.4
1	q	80[C]	ILE	2.4
1	f	169[C]	LEU	2.4
1	k	72[C]	LEU	2.4
1	u	37[C]	MET	2.4
1	n	103[C]	ALA	2.4
1	b	176[C]	GLY	2.4
1	l	118[C]	HIS	2.4
1	r	151[C]	HIS	2.4
1	s	57[C]	HIS	2.4
1	d	76[C]	ARG	2.4
1	s	29[C]	TYR	2.4
1	m	133[C]	ILE	2.4
1	q	165[C]	LEU	2.4
1	w	104[C]	LEU	2.4
1	e	159[C]	GLY	2.4
1	j	161[C]	PRO	2.4
1	k	15[C]	ASP	2.4
1	t	32[C]	TYR	2.4
1	x	4[C]	SER	2.4
1	i	112[C]	GLN	2.4
1	r	11[C]	ASN	2.4
1	h	162[C]	GLU	2.4
1	c	148[C]	LEU	2.4
1	g	115[C]	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	k	115[C]	LEU	2.4
1	n	129[C]	LEU	2.4
1	p	69[C]	LEU	2.4
1	F	161	PRO	2.4
1	f	100[C]	MET	2.4
1	w	102[C]	ALA	2.4
1	w	24[C]	ILE	2.4
1	x	20[C]	ILE	2.4
1	k	164[C]	GLY	2.4
1	a	114[C]	LEU	2.4
1	g	117[C]	LEU	2.4
1	o	104[C]	LEU	2.4
1	p	120[C]	LEU	2.4
1	s	104[C]	LEU	2.4
1	w	117[C]	LEU	2.4
1	b	99[C]	ALA	2.4
1	e	47[C]	ALA	2.4
1	w	100[C]	MET	2.4
1	h	5[C]	THR	2.4
1	h	8[C]	VAL	2.4
1	j	33[C]	VAL	2.4
1	l	110[C]	VAL	2.4
1	a	31[C]	SER	2.4
1	h	4[C]	SER	2.4
1	p	145[C]	ILE	2.3
1	k	105[C]	HIS	2.3
1	o	57[C]	HIS	2.3
1	w	173[C]	HIS	2.3
1	a	167[C]	GLU	2.3
1	f	106[C]	LEU	2.3
1	i	144[C]	ALA	2.3
1	e	14[C]	GLN	2.3
1	l	7[C]	GLN	2.3
1	g	125[C]	ASN	2.3
1	h	125[C]	ASN	2.3
1	k	81[C]	PHE	2.3
1	c	105[C]	HIS	2.3
1	n	162[C]	GLU	2.3
1	w	64[C]	GLU	2.3
1	c	32[C]	TYR	2.3
1	T	119	LYS	2.3
1	o	92[C]	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	t	38[C]	SER	2.3
1	r	142[C]	VAL	2.3
1	s	33[C]	VAL	2.3
1	c	112[C]	GLN	2.3
1	f	88[C]	PRO	2.3
1	e	39[C]	TYR	2.3
1	u	175[C]	LEU	2.3
1	v	104[C]	LEU	2.3
1	q	153[C]	THR	2.3
1	e	38[C]	SER	2.3
1	t	93[C]	TRP	2.3
1	w	140[C]	GLU	2.3
1	d	49[C]	LYS	2.3
1	o	170[C]	PHE	2.3
1	q	57[C]	HIS	2.3
1	x	112[C]	GLN	2.3
1	x	128[C]	HIS	2.3
1	v	80[C]	ILE	2.3
1	v	145[C]	ILE	2.3
1	g	64[C]	GLU	2.3
1	j	47[C]	ALA	2.3
1	k	19[C]	ALA	2.3
1	m	61[C]	GLU	2.3
1	o	123[C]	ASP	2.3
1	i	138[C]	LEU	2.3
1	r	169[C]	LEU	2.3
1	i	31[C]	SER	2.3
1	T	49	LYS	2.3
1	m	119[C]	LYS	2.3
1	j	14[C]	GLN	2.3
1	m	118[C]	HIS	2.3
1	v	110[C]	VAL	2.3
1	m	154[C]	ASN	2.3
1	t	18[C]	ALA	2.3
1	a	5[C]	THR	2.3
1	u	14[C]	GLN	2.3
1	w	34[C]	TYR	2.3
1	j	151[C]	HIS	2.3
1	c	110[C]	VAL	2.3
1	v	55[C]	PHE	2.3
1	i	4[C]	SER	2.3
1	p	30[C]	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	u	38[C]	SER	2.3
1	v	112[C]	GLN	2.3
1	U	26	LEU	2.2
1	c	97[C]	LEU	2.2
1	f	138[C]	LEU	2.2
1	i	114[C]	LEU	2.2
1	o	56[C]	LEU	2.2
1	q	97[C]	LEU	2.2
1	v	5[C]	THR	2.2
1	p	76[C]	ARG	2.2
1	a	168[C]	TYR	2.2
1	b	100[C]	MET	2.2
1	x	29[C]	TYR	2.2
1	a	125[C]	ASN	2.2
1	d	109[C]	ASN	2.2
1	q	25[C]	ASN	2.2
1	u	15[C]	ASP	2.2
1	d	170[C]	PHE	2.2
1	b	127[C]	PRO	2.2
1	d	167[C]	GLU	2.2
1	k	38[C]	SER	2.2
1	b	19[C]	ALA	2.2
1	b	160[C]	ALA	2.2
1	r	52[C]	ALA	2.2
1	t	66[C]	ALA	2.2
1	d	155[C]	LEU	2.2
1	g	169[C]	LEU	2.2
1	h	104[C]	LEU	2.2
1	i	77[C]	GLY	2.2
1	v	174[C]	THR	2.2
1	t	13[C]	HIS	2.2
1	f	39[C]	TYR	2.2
1	l	109[C]	ASN	2.2
1	c	8[C]	VAL	2.2
1	d	14[C]	GLN	2.2
1	l	142[C]	VAL	2.2
1	F	4	SER	2.2
1	j	36[C]	SER	2.2
1	w	93[C]	TRP	2.2
1	K	119	LYS	2.2
1	l	103[C]	ALA	2.2
1	u	52[C]	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	b	69[C]	LEU	2.2
1	c	60[C]	HIS	2.2
1	q	105[C]	HIS	2.2
1	v	60[C]	HIS	2.2
1	h	29[C]	TYR	2.2
1	p	140[C]	GLU	2.2
1	r	168[C]	TYR	2.2
1	x	127[C]	PRO	2.2
1	Q	119	LYS	2.2
1	k	124[C]	LYS	2.2
1	x	142[C]	VAL	2.2
1	i	78[C]	GLY	2.2
1	u	176[C]	GLY	2.2
1	a	174[C]	THR	2.2
1	b	15[C]	ASP	2.2
1	n	102[C]	ALA	2.2
1	e	128[C]	HIS	2.2
1	h	98[C]	ASN	2.2
1	i	151[C]	HIS	2.2
1	m	145[C]	ILE	2.2
1	o	20[C]	ILE	2.2
1	j	82[C]	LEU	2.2
1	s	76[C]	ARG	2.2
1	u	100[C]	MET	2.2
1	u	127[C]	PRO	2.2
1	v	32[C]	TYR	2.2
1	a	46[C]	VAL	2.2
1	o	142[C]	VAL	2.2
1	f	55[C]	PHE	2.2
1	l	81[C]	PHE	2.2
1	s	55[C]	PHE	2.2
1	t	51[C]	PHE	2.2
1	b	11[C]	ASN	2.2
1	g	57[C]	HIS	2.2
1	n	19[C]	ALA	2.2
1	o	141[C]	GLN	2.2
1	r	73[C]	GLN	2.2
1	u	57[C]	HIS	2.2
1	u	73[C]	GLN	2.2
1	q	93[C]	TRP	2.2
1	c	69[C]	LEU	2.2
1	v	48[C]	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	b	126[C]	ASP	2.2
1	c	168[C]	TYR	2.2
1	e	54[C]	TYR	2.2
1	j	177[C]	ASP	2.2
1	n	29[C]	TYR	2.2
1	q	92[C]	ASP	2.2
1	u	67[C]	GLU	2.2
1	v	78[C]	GLY	2.2
1	a	111[C]	ASN	2.2
1	b	124[C]	LYS	2.2
1	i	105[C]	HIS	2.2
1	m	5[C]	THR	2.2
1	b	24[C]	ILE	2.2
1	c	145[C]	ILE	2.2
1	l	80[C]	ILE	2.2
1	f	28[C]	LEU	2.2
1	v	117[C]	LEU	2.2
1	a	36[C]	SER	2.2
1	n	6[C]	SER	2.2
1	T	14	GLN	2.2
1	b	14[C]	GLN	2.2
1	d	112[C]	GLN	2.2
1	f	73[C]	GLN	2.2
1	g	124[C]	LYS	2.1
1	n	119[C]	LYS	2.1
1	r	112[C]	GLN	2.2
1	t	124[C]	LYS	2.1
1	u	109[C]	ASN	2.1
1	h	39[C]	TYR	2.1
1	m	55[C]	PHE	2.1
1	v	41[C]	PHE	2.1
1	g	52[C]	ALA	2.1
1	b	138[C]	LEU	2.1
1	n	97[C]	LEU	2.1
1	p	117[C]	LEU	2.1
1	s	42[C]	ASP	2.1
1	A	4	SER	2.1
1	b	4[C]	SER	2.1
1	b	119[C]	LYS	2.1
1	c	4[C]	SER	2.1
1	o	36[C]	SER	2.1
1	p	112[C]	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	q	14[C]	GLN	2.1
1	q	154[C]	ASN	2.1
1	k	29[C]	TYR	2.1
1	m	46[C]	VAL	2.1
1	t	33[C]	VAL	2.1
1	r	103[C]	ALA	2.1
1	k	126[C]	ASP	2.1
1	o	119[C]	LYS	2.1
1	r	44[C]	ASP	2.1
1	u	124[C]	LYS	2.1
1	i	24[C]	ILE	2.1
1	X	4	SER	2.1
1	d	111[C]	ASN	2.1
1	h	69[C]	LEU	2.1
1	k	98[C]	ASN	2.1
1	u	117[C]	LEU	2.1
1	t	149[C]	GLY	2.1
1	i	67[C]	GLU	2.1
1	p	116[C]	GLU	2.1
1	s	118[C]	HIS	2.1
1	a	29[C]	TYR	2.1
1	a	124[C]	LYS	2.1
1	b	12[C]	TYR	2.1
1	c	142[C]	VAL	2.1
1	i	110[C]	VAL	2.1
1	j	168[C]	TYR	2.1
1	h	81[C]	PHE	2.1
1	k	66[C]	ALA	2.1
1	q	34[C]	TYR	2.1
1	r	5[C]	THR	2.1
1	c	155[C]	LEU	2.1
1	f	72[C]	LEU	2.1
1	o	120[C]	LEU	2.1
1	s	116[C]	GLU	2.1
1	o	37[C]	MET	2.1
1	r	158[C]	MET	2.1
1	f	105[C]	HIS	2.1
1	q	13[C]	HIS	2.1
1	d	40[C]	TYR	2.1
1	e	160[C]	ALA	2.1
1	f	152[C]	VAL	2.1
1	h	32[C]	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	m	32[C]	TYR	2.1
1	t	39[C]	TYR	2.1
1	n	55[C]	PHE	2.1
1	p	111[C]	ASN	2.1
1	q	64[C]	GLU	2.1
1	u	170[C]	PHE	2.1
1	v	156[C]	ARG	2.1
1	j	176[C]	GLY	2.1
1	k	127[C]	PRO	2.1
1	w	85[C]	ILE	2.1
1	B	119	LYS	2.1
1	c	28[C]	LEU	2.1
1	p	157[C]	LYS	2.1
1	r	106[C]	LEU	2.1
1	u	148[C]	LEU	2.1
1	c	13[C]	HIS	2.1
1	q	100[C]	MET	2.1
1	p	45[C]	ASP	2.1
1	u	76[C]	ARG	2.1
1	u	160[C]	ALA	2.1
1	l	33[C]	VAL	2.1
1	m	8[C]	VAL	2.1
1	i	41[C]	PHE	2.1
1	x	38[C]	SER	2.1
1	j	88[C]	PRO	2.1
1	F	120	LEU	2.1
1	p	80[C]	ILE	2.1
1	b	173[C]	HIS	2.1
1	f	118[C]	HIS	2.1
1	p	14[C]	GLN	2.1
1	s	82[C]	LEU	2.1
1	w	105[C]	HIS	2.1
1	e	147[C]	GLU	2.1
1	l	63[C]	ARG	2.1
1	p	109[C]	ASN	2.1
1	c	53[C]	LYS	2.1
1	d	119[C]	LYS	2.1
1	f	53[C]	LYS	2.1
1	h	47[C]	ALA	2.1
1	q	166[C]	ALA	2.1
1	s	103[C]	ALA	2.1
1	p	149[C]	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	d	81[C]	PHE	2.1
1	u	29[C]	TYR	2.1
1	n	167[C]	GLU	2.0
1	u	107[C]	GLU	2.0
1	T	106	LEU	2.0
1	d	117[C]	LEU	2.0
1	f	104[C]	LEU	2.0
1	g	138[C]	LEU	2.0
1	r	138[C]	LEU	2.0
1	b	158[C]	MET	2.0
1	r	109[C]	ASN	2.0
1	d	103[C]	ALA	2.0
1	i	99[C]	ALA	2.0
1	v	103[C]	ALA	2.0
1	l	95[C]	SER	2.0
1	m	33[C]	VAL	2.0
1	a	40[C]	TYR	2.0
1	b	13[C]	HIS	2.0
1	b	105[C]	HIS	2.0
1	l	151[C]	HIS	2.0
1	m	122[C]	HIS	2.0
1	T	53	LYS	2.0
1	o	109[C]	ASN	2.0
1	x	37[C]	MET	2.0
1	S	112	GLN	2.0
1	j	159[C]	GLY	2.0
1	j	127[C]	PRO	2.0
1	w	92[C]	ASP	2.0
1	f	136[C]	HIS	2.0
1	h	152[C]	VAL	2.0
1	p	46[C]	VAL	2.0
1	q	151[C]	HIS	2.0
1	w	33[C]	VAL	2.0
1	g	170[C]	PHE	2.0
1	u	132[C]	PHE	2.0
1	c	39[C]	TYR	2.0
1	s	54[C]	TYR	2.0
1	v	137[C]	TYR	2.0
1	v	133[C]	ILE	2.0
1	a	86[C]	GLN	2.0
1	g	14[C]	GLN	2.0
1	n	86[C]	GLN	2.0

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Mol	Chain	Res	Type	RSRZ
1	i	95[C]	SER	2.0
1	i	135[C]	THR	2.0
1	m	30[C]	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ZN	v	203[C]	1/1	0.61	0.15	33,33,33,33	1
2	ZN	v	203[D]	1/1	0.61	0.15	35,35,35,35	1
2	ZN	d	202[C]	1/1	0.74	0.11	38,38,38,38	1
2	ZN	d	202[D]	1/1	0.74	0.11	50,50,50,50	1
2	ZN	d	204[C]	1/1	0.83	0.11	37,37,37,37	1
2	ZN	d	204[D]	1/1	0.83	0.11	36,36,36,36	1
3	NA	G	204	1/1	0.83	0.20	46,46,46,46	0
2	ZN	n	202[D]	1/1	0.85	0.11	48,48,48,48	1
2	ZN	l	203[C]	1/1	0.85	0.12	32,32,32,32	1
2	ZN	l	203[D]	1/1	0.85	0.12	30,30,30,30	1
2	ZN	n	202[C]	1/1	0.85	0.11	46,46,46,46	1
3	NA	g	204[C]	1/1	0.85	0.18	28,28,28,28	1
3	NA	g	204[D]	1/1	0.85	0.18	32,32,32,32	1
2	ZN	g	201[D]	1/1	0.86	0.15	36,36,36,36	1
2	ZN	d	201[C]	1/1	0.86	0.13	34,34,34,34	1
2	ZN	d	201[D]	1/1	0.86	0.13	40,40,40,40	1
2	ZN	H	202	1/1	0.86	0.17	70,70,70,70	1
2	ZN	g	201[C]	1/1	0.86	0.15	27,27,27,27	1
3	NA	p	204[C]	1/1	0.86	0.15	36,36,36,36	1
3	NA	p	204[D]	1/1	0.86	0.15	33,33,33,33	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ZN	d	203[D]	1/1	0.87	0.19	52,52,52,52	1
2	ZN	d	203[C]	1/1	0.87	0.19	56,56,56,56	1
2	ZN	S	202	1/1	0.88	0.13	35,35,35,35	1
2	ZN	e	201[C]	1/1	0.88	0.14	29,29,29,29	1
2	ZN	i	201[C]	1/1	0.88	0.12	29,29,29,29	1
2	ZN	i	201[D]	1/1	0.88	0.12	34,34,34,34	1
2	ZN	e	201[D]	1/1	0.88	0.14	34,34,34,34	1
4	V9Y	O	204[C]	13/13	0.88	0.22	45,50,63,63	13
4	V9Y	O	204[D]	13/13	0.88	0.22	43,53,63,63	13
2	ZN	g	202[C]	1/1	0.89	0.08	42,42,42,42	1
2	ZN	g	202[D]	1/1	0.89	0.08	48,48,48,48	1
2	ZN	h	201[C]	1/1	0.89	0.09	34,34,34,34	1
2	ZN	h	201[D]	1/1	0.89	0.09	36,36,36,36	1
2	ZN	T	202	1/1	0.89	0.08	54,54,54,54	0
2	ZN	r	202[C]	1/1	0.89	0.09	34,34,34,34	1
2	ZN	r	202[D]	1/1	0.89	0.09	40,40,40,40	1
4	V9Y	A	205[C]	13/13	0.89	0.20	41,46,60,61	13
4	V9Y	A	205[D]	13/13	0.89	0.20	43,49,60,61	13
2	ZN	s	201[C]	1/1	0.89	0.10	28,28,28,28	1
2	ZN	s	201[D]	1/1	0.89	0.10	33,33,33,33	1
2	ZN	w	201[D]	1/1	0.90	0.14	31,31,31,31	1
2	ZN	a	201[C]	1/1	0.90	0.10	30,30,30,30	1
3	NA	a	204[C]	1/1	0.90	0.20	34,34,34,34	1
3	NA	a	204[D]	1/1	0.90	0.20	31,31,31,31	1
2	ZN	a	201[D]	1/1	0.90	0.10	34,34,34,34	1
2	ZN	g	203[C]	1/1	0.90	0.14	32,32,32,32	1
2	ZN	g	203[D]	1/1	0.90	0.14	40,40,40,40	1
2	ZN	v	201[C]	1/1	0.90	0.09	38,38,38,38	1
2	ZN	v	201[D]	1/1	0.90	0.09	33,33,33,33	1
2	ZN	a	203[C]	1/1	0.90	0.10	33,33,33,33	1
2	ZN	a	203[D]	1/1	0.90	0.10	34,34,34,34	1
2	ZN	w	201[C]	1/1	0.90	0.14	35,35,35,35	1
2	ZN	D	201	1/1	0.91	0.10	34,34,34,34	0
2	ZN	a	202[C]	1/1	0.91	0.09	42,42,42,42	1
2	ZN	b	201[C]	1/1	0.91	0.11	30,30,30,30	1
2	ZN	o	202[C]	1/1	0.91	0.10	38,38,38,38	1
3	NA	A	204	1/1	0.91	0.22	33,33,33,33	0
2	ZN	o	202[D]	1/1	0.91	0.10	34,34,34,34	1
2	ZN	r	201[C]	1/1	0.91	0.10	31,31,31,31	1
2	ZN	r	201[D]	1/1	0.91	0.10	34,34,34,34	1
2	ZN	b	201[D]	1/1	0.91	0.11	31,31,31,31	1
2	ZN	k	201[C]	1/1	0.91	0.09	31,31,31,31	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ZN	k	201[D]	1/1	0.91	0.09	37,37,37,37	1
2	ZN	l	201[C]	1/1	0.91	0.11	26,26,26,26	1
2	ZN	s	202[C]	1/1	0.91	0.08	40,40,40,40	1
2	ZN	s	202[D]	1/1	0.91	0.08	44,44,44,44	1
4	V9Y	D	204[C]	13/13	0.91	0.21	49,57,65,67	13
4	V9Y	D	204[D]	13/13	0.91	0.21	45,49,64,66	13
4	V9Y	I	203[C]	13/13	0.91	0.18	48,53,66,66	13
4	V9Y	I	203[D]	13/13	0.91	0.18	44,51,64,66	13
2	ZN	l	201[D]	1/1	0.91	0.11	34,34,34,34	1
2	ZN	a	202[D]	1/1	0.91	0.09	50,50,50,50	1
4	V9Y	X	203[C]	13/13	0.91	0.20	40,47,57,58	13
4	V9Y	X	203[D]	13/13	0.91	0.20	34,45,57,58	13
3	NA	n	204[D]	1/1	0.92	0.14	36,36,36,36	1
2	ZN	f	201[C]	1/1	0.92	0.12	26,26,26,26	1
2	ZN	f	201[D]	1/1	0.92	0.12	38,38,38,38	1
2	ZN	t	201[C]	1/1	0.92	0.08	35,35,35,35	1
2	ZN	t	201[D]	1/1	0.92	0.08	34,34,34,34	1
3	NA	D	203	1/1	0.92	0.21	48,48,48,48	0
2	ZN	t	202[C]	1/1	0.92	0.11	35,35,35,35	1
3	NA	P	204	1/1	0.92	0.20	45,45,45,45	0
2	ZN	t	202[D]	1/1	0.92	0.11	31,31,31,31	1
2	ZN	f	203[C]	1/1	0.92	0.12	26,26,26,26	1
2	ZN	f	203[D]	1/1	0.92	0.12	42,42,42,42	1
2	ZN	C	202	1/1	0.92	0.05	71,71,71,71	1
3	NA	n	204[C]	1/1	0.92	0.14	36,36,36,36	1
4	V9Y	t	203[C]	13/13	0.92	0.18	43,46,65,67	13
4	V9Y	t	203[D]	13/13	0.92	0.18	51,56,67,68	13
3	NA	x	204[C]	1/1	0.93	0.18	31,31,31,31	1
3	NA	x	204[D]	1/1	0.93	0.18	33,33,33,33	1
2	ZN	i	202[C]	1/1	0.93	0.10	29,29,29,29	1
3	NA	S	203	1/1	0.93	0.27	45,45,45,45	0
2	ZN	i	202[D]	1/1	0.93	0.10	36,36,36,36	1
2	ZN	s	203[C]	1/1	0.93	0.08	31,31,31,31	1
2	ZN	s	203[D]	1/1	0.93	0.08	32,32,32,32	1
2	ZN	w	204[C]	1/1	0.93	0.12	32,32,32,32	1
3	NA	j	203[C]	1/1	0.93	0.14	28,28,28,28	1
3	NA	j	203[D]	1/1	0.93	0.14	33,33,33,33	1
4	V9Y	R	203[C]	13/13	0.93	0.20	45,56,70,70	13
4	V9Y	R	203[D]	13/13	0.93	0.20	47,58,70,71	13
2	ZN	w	204[D]	1/1	0.93	0.12	37,37,37,37	1
2	ZN	P	202	1/1	0.93	0.09	55,55,55,55	0
2	ZN	v	202[C]	1/1	0.93	0.18	44,44,44,44	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	ZN	v	202[D]	1/1	0.93	0.18	44,44,44,44	1
2	ZN	I	202	1/1	0.94	0.13	44,44,44,44	1
2	ZN	q	201[C]	1/1	0.94	0.08	36,36,36,36	1
2	ZN	q	201[D]	1/1	0.94	0.08	32,32,32,32	1
2	ZN	x	201[C]	1/1	0.94	0.11	35,35,35,35	1
2	ZN	x	201[D]	1/1	0.94	0.11	30,30,30,30	1
2	ZN	q	202[C]	1/1	0.94	0.08	35,35,35,35	1
2	ZN	q	202[D]	1/1	0.94	0.08	37,37,37,37	1
2	ZN	G	202	1/1	0.94	0.10	49,49,49,49	0
2	ZN	u	202[C]	1/1	0.94	0.08	40,40,40,40	1
2	ZN	u	202[D]	1/1	0.94	0.08	35,35,35,35	1
2	ZN	f	202[C]	1/1	0.94	0.11	38,38,38,38	1
4	V9Y	L	203[C]	13/13	0.94	0.19	42,46,67,68	13
4	V9Y	L	203[D]	13/13	0.94	0.19	43,50,67,68	13
2	ZN	j	201[C]	1/1	0.94	0.12	31,31,31,31	1
3	NA	d	205[C]	1/1	0.94	0.16	31,31,31,31	1
3	NA	d	205[D]	1/1	0.94	0.16	36,36,36,36	1
2	ZN	j	201[D]	1/1	0.94	0.12	39,39,39,39	1
2	ZN	f	202[D]	1/1	0.94	0.11	45,45,45,45	1
2	ZN	o	201[C]	1/1	0.94	0.09	32,32,32,32	1
2	ZN	o	201[D]	1/1	0.94	0.09	33,33,33,33	1
2	ZN	U	201	1/1	0.94	0.12	43,43,43,43	0
2	ZN	e	203[C]	1/1	0.95	0.08	38,38,38,38	1
2	ZN	e	203[D]	1/1	0.95	0.08	36,36,36,36	1
2	ZN	x	203[C]	1/1	0.95	0.11	39,39,39,39	1
2	ZN	x	203[D]	1/1	0.95	0.11	30,30,30,30	1
2	ZN	n	203[C]	1/1	0.95	0.07	33,33,33,33	1
2	ZN	n	203[D]	1/1	0.95	0.07	31,31,31,31	1
2	ZN	l	202[C]	1/1	0.95	0.09	36,36,36,36	1
3	NA	J	203	1/1	0.95	0.23	32,32,32,32	0
3	NA	M	203	1/1	0.95	0.17	28,28,28,28	0
2	ZN	l	202[D]	1/1	0.95	0.09	47,47,47,47	1
2	ZN	I	201	1/1	0.95	0.14	43,43,43,43	0
3	NA	V	203	1/1	0.95	0.22	32,32,32,32	0
2	ZN	F	204	1/1	0.95	0.12	45,45,45,45	1
2	ZN	p	202[C]	1/1	0.95	0.07	51,51,51,51	1
2	ZN	p	202[D]	1/1	0.95	0.07	39,39,39,39	1
2	ZN	p	203[C]	1/1	0.95	0.06	32,32,32,32	1
2	ZN	p	203[D]	1/1	0.95	0.06	31,31,31,31	1
2	ZN	w	203[C]	1/1	0.95	0.10	44,44,44,44	1
2	ZN	w	203[D]	1/1	0.95	0.10	50,50,50,50	1
2	ZN	m	201[C]	1/1	0.95	0.08	40,40,40,40	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	ZN	m	201[D]	1/1	0.95	0.08	33,33,33,33	1
2	ZN	b	202[C]	1/1	0.96	0.11	28,28,28,28	1
2	ZN	b	202[D]	1/1	0.96	0.11	35,35,35,35	1
2	ZN	w	202[C]	1/1	0.96	0.11	42,42,42,42	1
2	ZN	w	202[D]	1/1	0.96	0.11	38,38,38,38	1
2	ZN	c	201[C]	1/1	0.96	0.08	30,30,30,30	1
2	ZN	c	201[D]	1/1	0.96	0.08	38,38,38,38	1
2	ZN	k	202[C]	1/1	0.96	0.08	34,34,34,34	1
2	ZN	k	202[D]	1/1	0.96	0.08	33,33,33,33	1
2	ZN	c	202[C]	1/1	0.96	0.08	30,30,30,30	1
2	ZN	c	202[D]	1/1	0.96	0.08	27,27,27,27	1
2	ZN	x	202[C]	1/1	0.96	0.23	52,52,52,52	1
2	ZN	x	202[D]	1/1	0.96	0.23	51,51,51,51	1
2	ZN	p	201[C]	1/1	0.96	0.06	32,32,32,32	1
2	ZN	p	201[D]	1/1	0.96	0.06	32,32,32,32	1
2	ZN	R	202	1/1	0.96	0.08	45,45,45,45	1
2	ZN	H	203	1/1	0.96	0.12	37,37,37,37	1
2	ZN	u	201[C]	1/1	0.96	0.09	37,37,37,37	1
2	ZN	u	201[D]	1/1	0.96	0.09	34,34,34,34	1
2	ZN	E	201	1/1	0.96	0.12	44,44,44,44	0
2	ZN	F	202	1/1	0.96	0.10	45,45,45,45	0
2	ZN	F	203	1/1	0.96	0.10	76,76,76,76	1
2	ZN	Q	201	1/1	0.96	0.10	42,42,42,42	0
2	ZN	m	202[C]	1/1	0.96	0.08	30,30,30,30	1
2	ZN	m	202[D]	1/1	0.96	0.08	30,30,30,30	1
2	ZN	n	201[C]	1/1	0.96	0.10	37,37,37,37	1
2	ZN	n	201[D]	1/1	0.96	0.10	32,32,32,32	1
2	ZN	j	202[C]	1/1	0.97	0.11	33,33,33,33	1
2	ZN	j	202[D]	1/1	0.97	0.11	34,34,34,34	1
2	ZN	R	201	1/1	0.97	0.11	43,43,43,43	0
2	ZN	B	203	1/1	0.97	0.13	34,34,34,34	1
2	ZN	S	201	1/1	0.97	0.08	33,33,33,33	0
2	ZN	E	202	1/1	0.97	0.10	79,79,79,79	1
2	ZN	T	201	1/1	0.97	0.09	42,42,42,42	0
2	ZN	C	201	1/1	0.97	0.10	36,36,36,36	0
2	ZN	B	201	1/1	0.97	0.13	33,33,33,33	0
2	ZN	V	201	1/1	0.97	0.14	33,33,33,33	0
2	ZN	W	201	1/1	0.97	0.14	35,35,35,35	0
2	ZN	X	202	1/1	0.97	0.09	36,36,36,36	1
2	ZN	N	203	1/1	0.97	0.09	37,37,37,37	1
2	ZN	B	202	1/1	0.97	0.06	44,44,44,44	0
2	ZN	D	202	1/1	0.97	0.12	35,35,35,35	1

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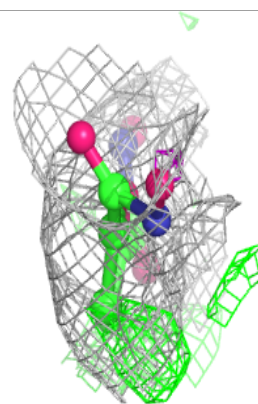
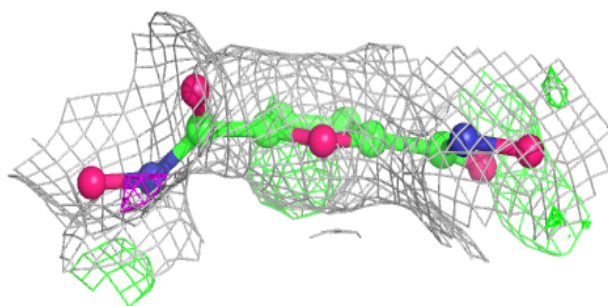
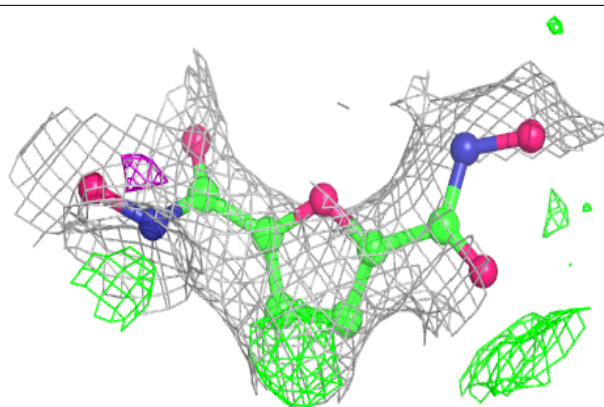
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	ZN	Q	202	1/1	0.97	0.09	38,38,38,38	1
2	ZN	W	203	1/1	0.98	0.10	38,38,38,38	1
3	NA	s	204[C]	1/1	0.98	0.10	33,33,33,33	1
3	NA	s	204[D]	1/1	0.98	0.10	27,27,27,27	1
2	ZN	X	201	1/1	0.98	0.12	34,34,34,34	0
2	ZN	A	202	1/1	0.98	0.04	48,48,48,48	1
2	ZN	J	201	1/1	0.98	0.17	32,32,32,32	0
2	ZN	K	202	1/1	0.98	0.08	42,42,42,42	0
2	ZN	K	203	1/1	0.98	0.05	36,36,36,36	1
2	ZN	M	201	1/1	0.98	0.13	42,42,42,42	0
2	ZN	M	202	1/1	0.98	0.12	35,35,35,35	1
2	ZN	C	203	1/1	0.98	0.12	30,30,30,30	1
2	ZN	O	202	1/1	0.98	0.08	71,71,71,71	1
2	ZN	T	203	1/1	0.98	0.09	41,41,41,41	1
2	ZN	P	201	1/1	0.98	0.07	38,38,38,38	0
2	ZN	e	202[C]	1/1	0.98	0.10	43,43,43,43	1
2	ZN	h	202[C]	1/1	0.98	0.06	37,37,37,37	1
2	ZN	h	202[D]	1/1	0.98	0.06	43,43,43,43	1
2	ZN	e	202[D]	1/1	0.98	0.10	50,50,50,50	1
2	ZN	H	201	1/1	0.98	0.09	43,43,43,43	0
2	ZN	V	202	1/1	0.98	0.17	31,31,31,31	1
2	ZN	P	203	1/1	0.98	0.10	39,39,39,39	1
2	ZN	L	201	1/1	0.99	0.09	31,31,31,31	0
2	ZN	L	202	1/1	0.99	0.10	30,30,30,30	1
2	ZN	F	201	1/1	0.99	0.06	45,45,45,45	0
2	ZN	G	203	1/1	0.99	0.09	39,39,39,39	1
2	ZN	N	201	1/1	0.99	0.05	34,34,34,34	0
2	ZN	N	202	1/1	0.99	0.08	41,41,41,41	0
2	ZN	A	201	1/1	0.99	0.14	35,35,35,35	0
2	ZN	O	201	1/1	0.99	0.11	30,30,30,30	0
2	ZN	J	202	1/1	0.99	0.16	34,34,34,34	1
2	ZN	O	203	1/1	0.99	0.13	33,33,33,33	1
2	ZN	A	203	1/1	0.99	0.10	33,33,33,33	1
2	ZN	U	202	1/1	0.99	0.12	44,44,44,44	1
2	ZN	E	203	1/1	0.99	0.10	40,40,40,40	1
2	ZN	W	202	1/1	1.00	0.14	40,40,40,40	0
2	ZN	K	201	1/1	1.00	0.12	33,33,33,33	0
2	ZN	G	201	1/1	1.00	0.11	34,34,34,34	0

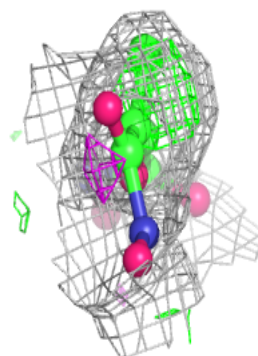
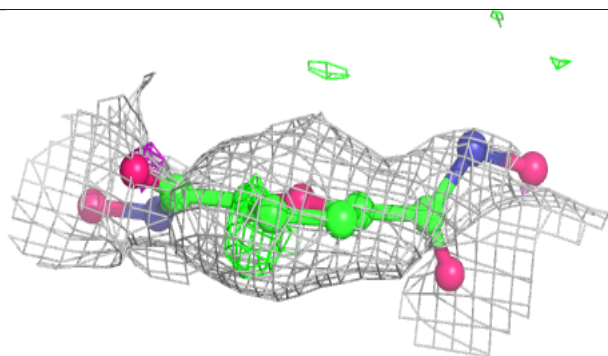
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around V9Y O 204 (C):

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

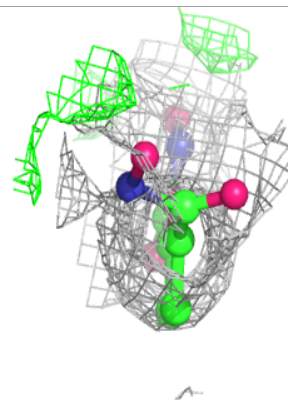
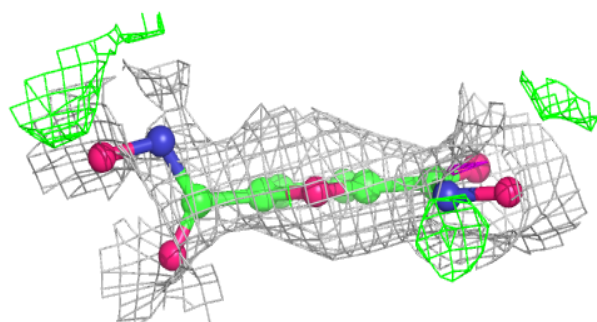
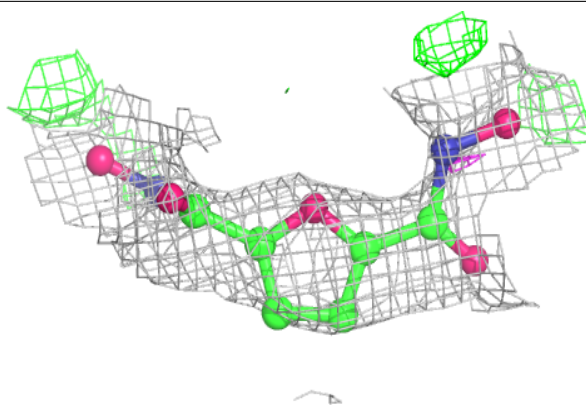
**Electron density around V9Y O 204 (D):**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

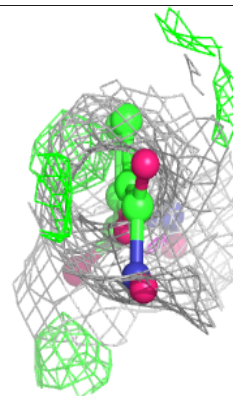
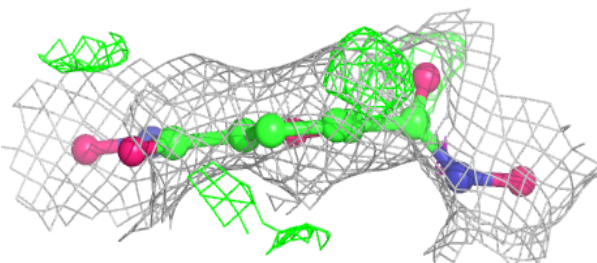
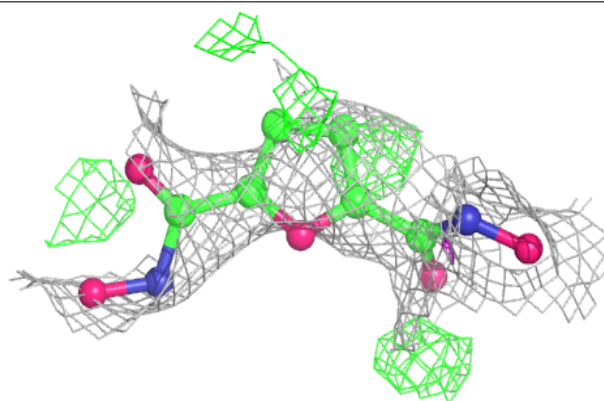


Electron density around V9Y A 205 (C):

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

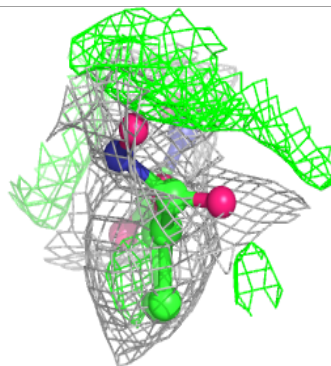
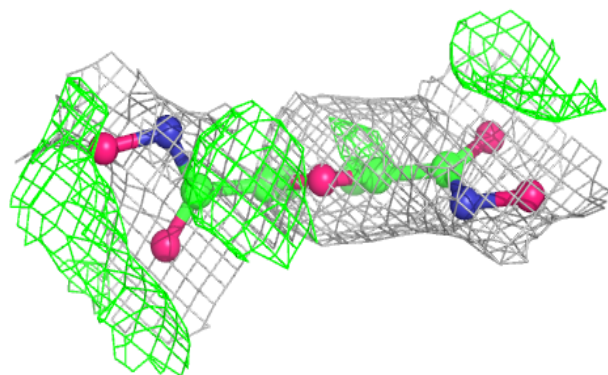
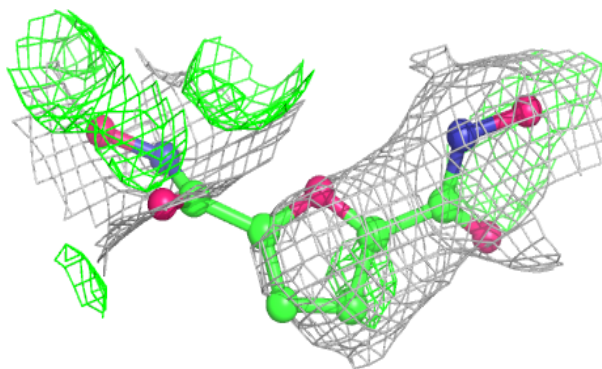
**Electron density around V9Y A 205 (D):**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

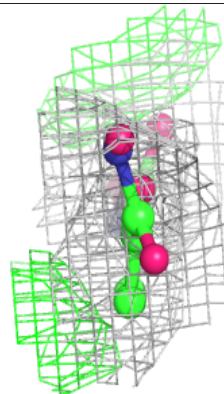
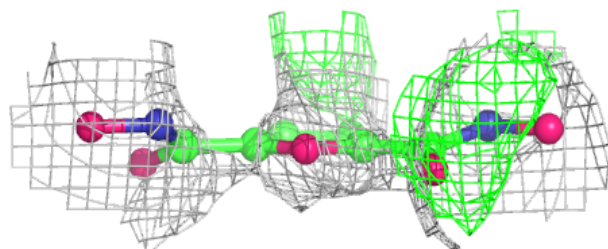
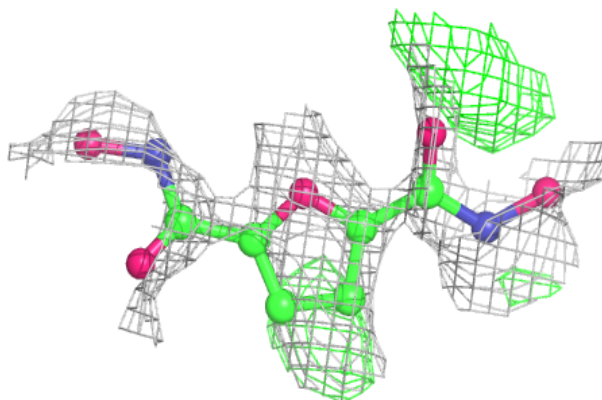


Electron density around V9Y D 204 (C):

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

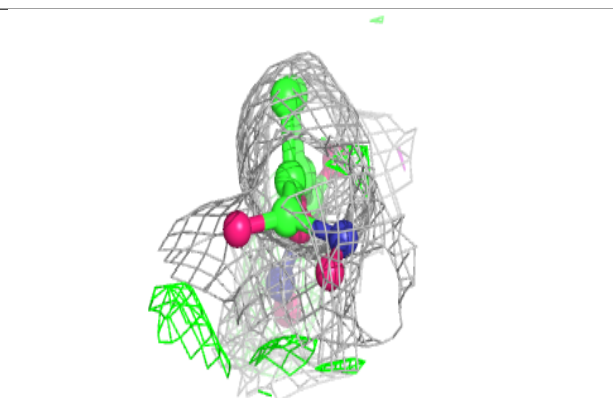
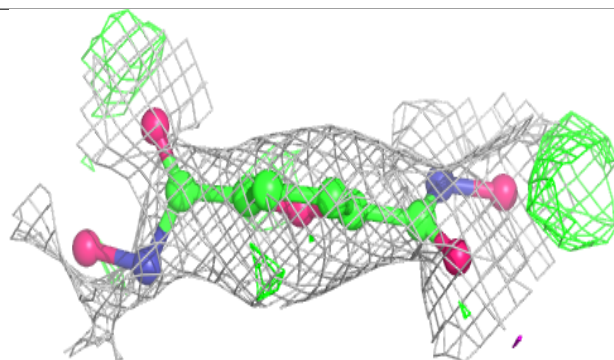
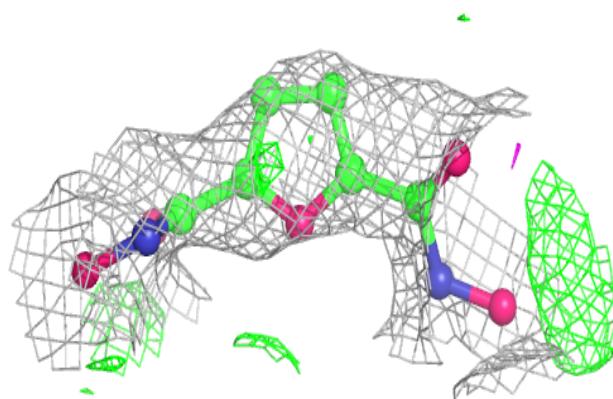
**Electron density around V9Y D 204 (D):**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

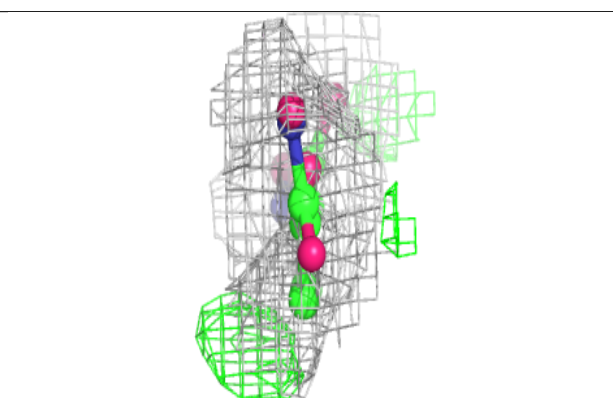
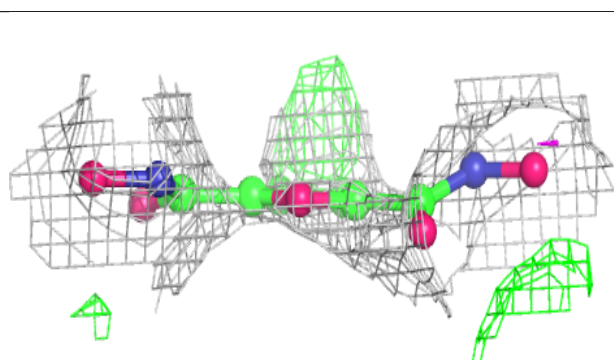
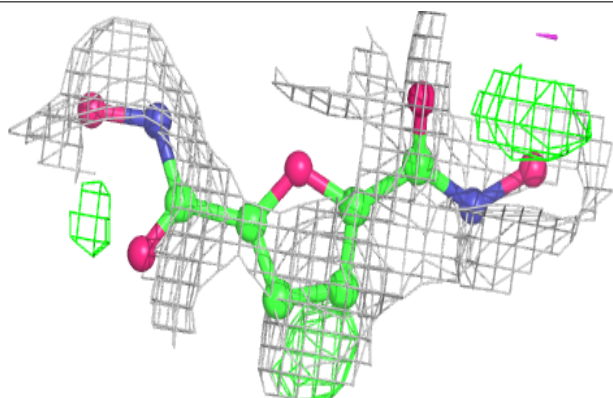


Electron density around V9Y I 203 (C):

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

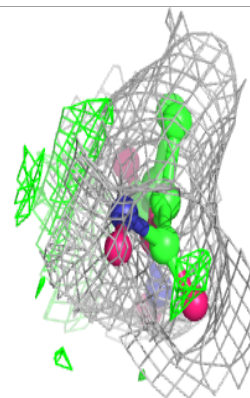
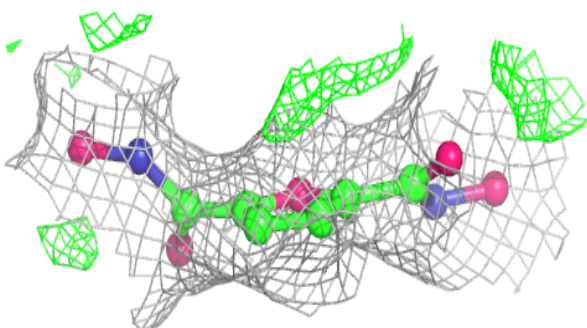
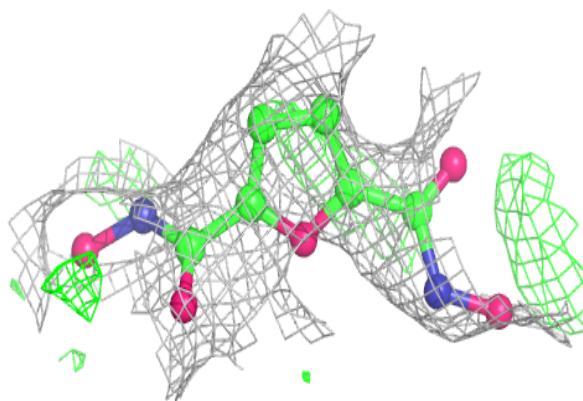
**Electron density around V9Y I 203 (D):**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

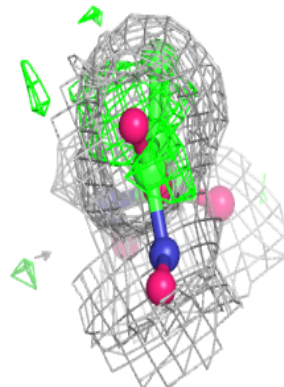
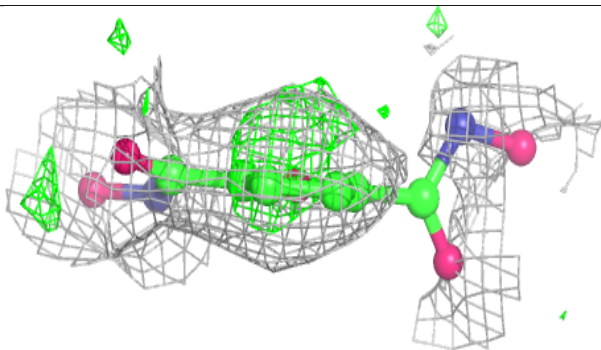
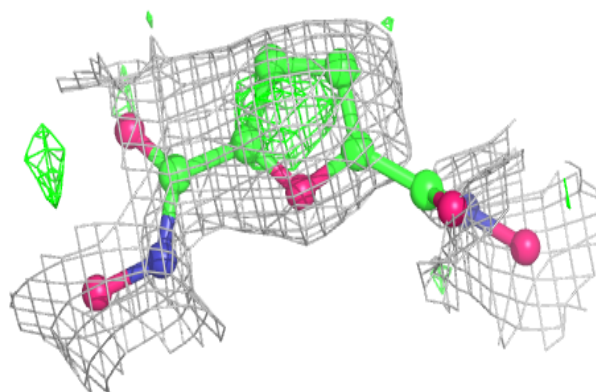


Electron density around V9Y X 203 (C):

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

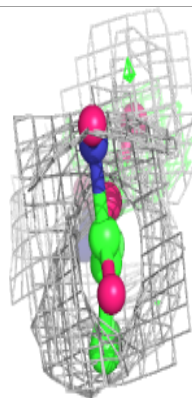
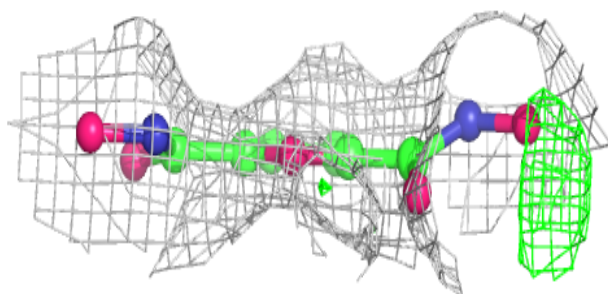
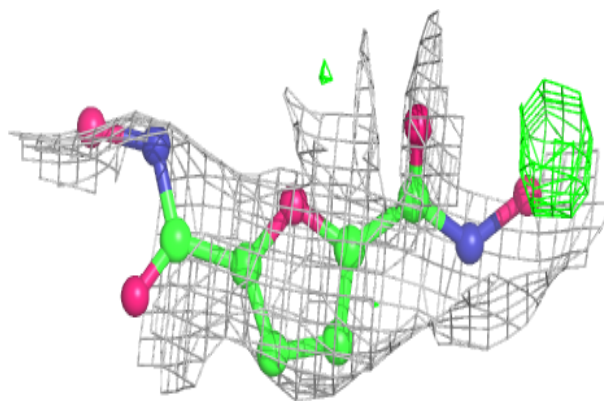
**Electron density around V9Y X 203 (D):**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

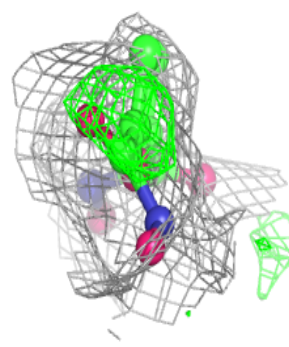
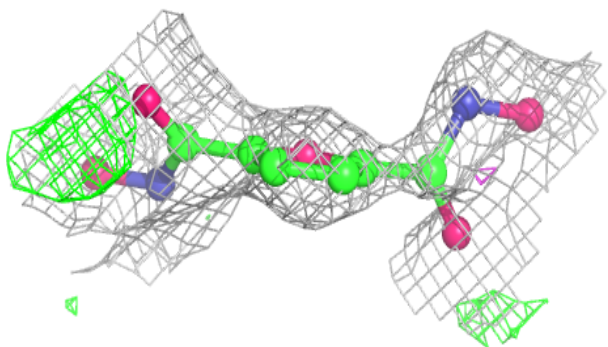
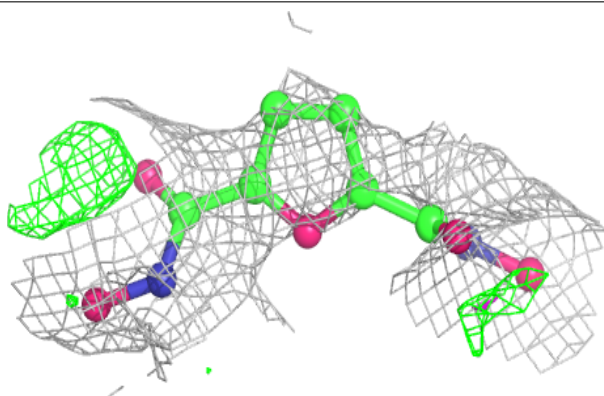


Electron density around V9Y t 203 (C):

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

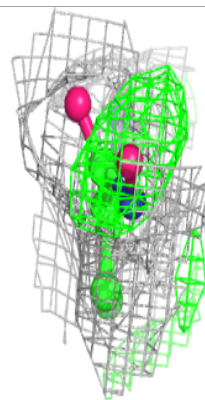
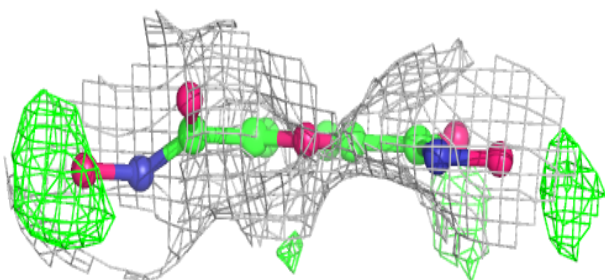
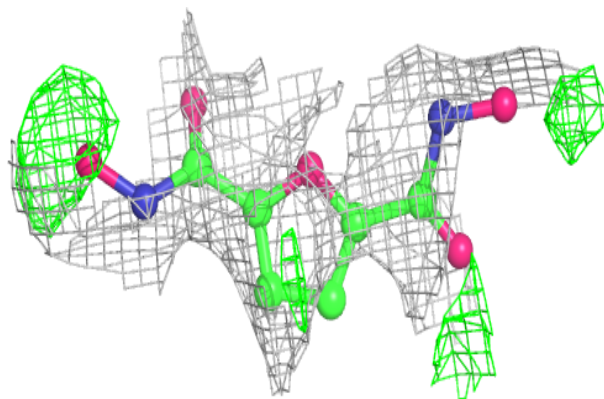
**Electron density around V9Y t 203 (D):**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

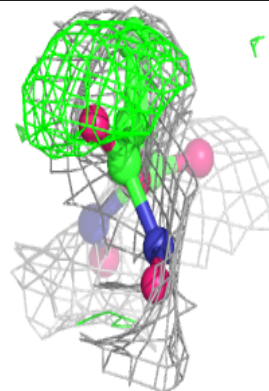
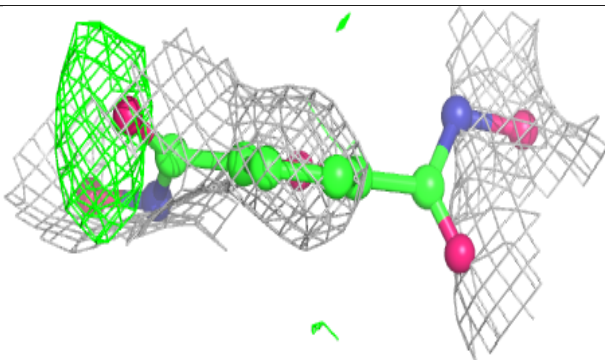
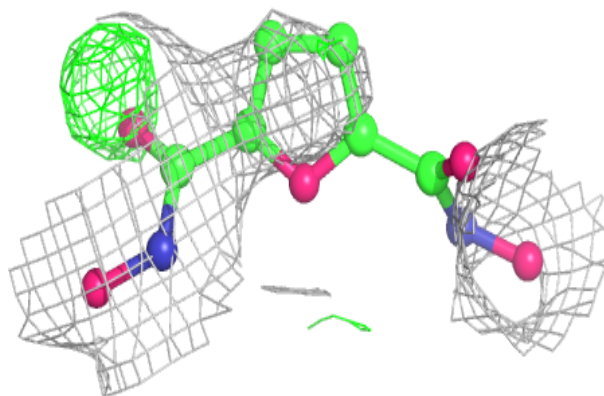


Electron density around V9Y R 203 (C):

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

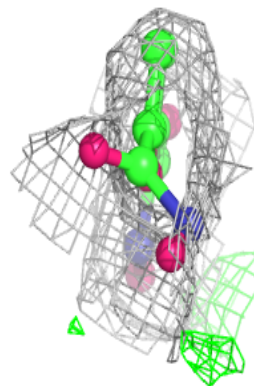
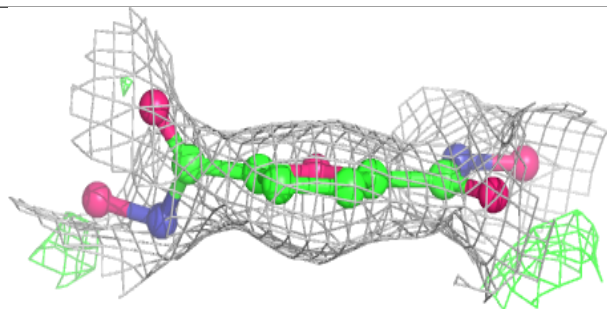
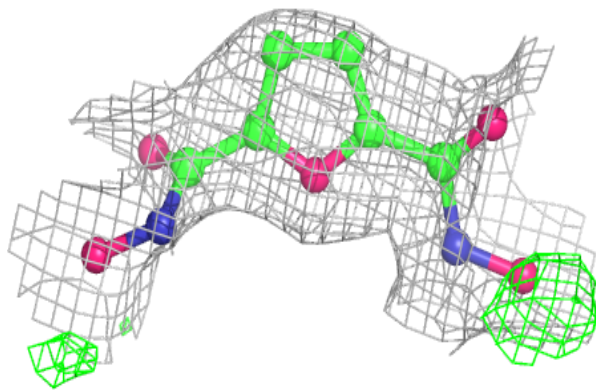
**Electron density around V9Y R 203 (D):**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

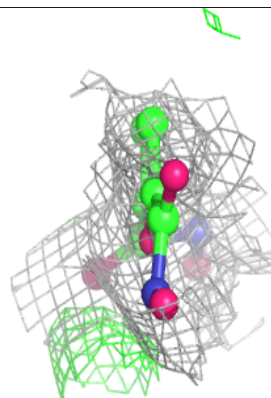
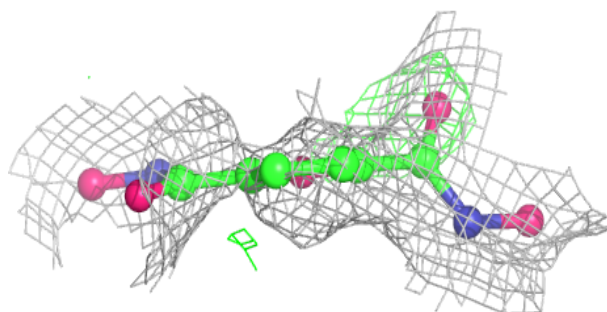
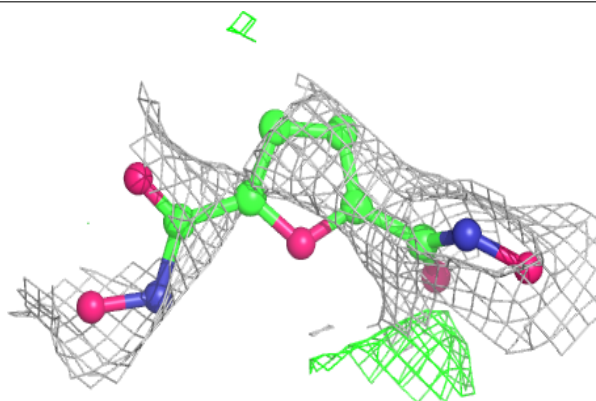


Electron density around V9Y L 203 (C):

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around V9Y L 203 (D):**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.