Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 08:37 PM GMT

PDB ID	:	4TNW
Title	:	C. elegans glutamate-gated chloride channel (GluCl) in complex with Fab and POPC in a lipid-modulated conformation
Authors	:	Althoff, T.; Hibbs, R.E.; Banerjee, S.; Gouaux, E.
Deposited on	:	2014-06-05
Resolution	:	3.20 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

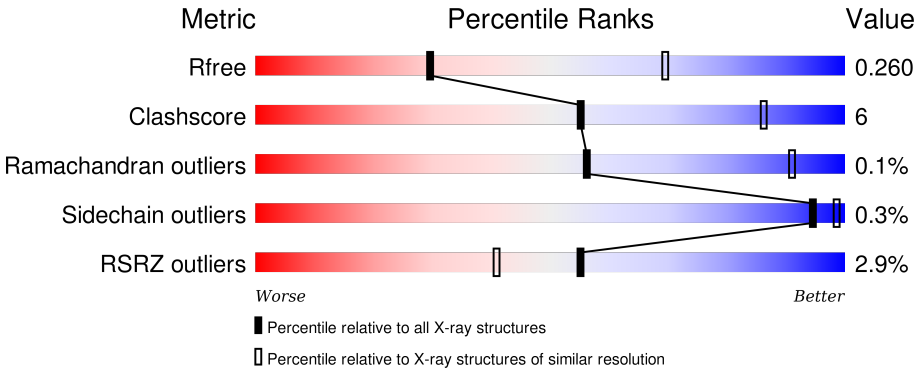
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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}	91344	1124 (3.24-3.16)
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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 on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. 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 <=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	347	81% 18%
1	B	347	% 79% 19%
1	C	347	% 77% 21%
1	D	347	80% 18%
1	E	347	% 80% 18%

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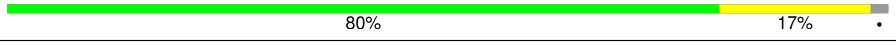
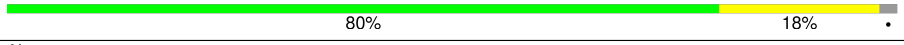
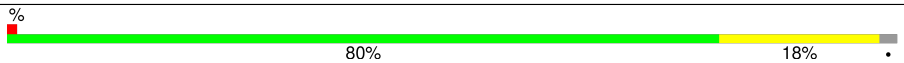
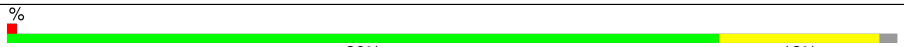
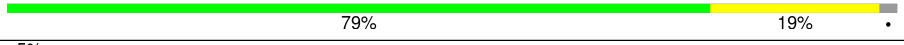
Mol	Chain	Length	Quality of chain
1	P	347	
1	Q	347	
1	R	347	
1	S	347	
1	T	347	
2	F	224	
2	G	224	
2	H	224	
2	I	224	
2	J	224	
2	U	224	
2	V	224	
2	W	224	
2	X	224	
2	Y	224	
3	K	215	
3	L	215	
3	M	215	
3	N	215	
3	O	215	
3	Z	215	
3	f	215	
3	g	215	
3	h	215	
3	i	215	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	POV	A	402	-	-	-	X
5	POV	D	401	-	-	-	X
5	POV	P	403	-	-	-	X
5	POV	Q	401	-	-	-	X
5	POV	R	401	-	-	-	X
5	POV	R	402	-	-	-	X
5	POV	T	401	-	-	-	X
8	LMT	B	403	-	-	-	X
8	LMT	D	403	-	-	-	X
8	LMT	P	402	-	-	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 60818 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 Avermectin-sensitive glutamate-gated chloride channel GluCl alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	342	Total	C	N	O	S	0	0	0
			2744	1786	449	494	15			
1	B	341	Total	C	N	O	S	0	0	0
			2734	1780	446	493	15			
1	C	340	Total	C	N	O	S	0	0	0
			2724	1774	443	492	15			
1	D	340	Total	C	N	O	S	0	0	0
			2724	1774	443	492	15			
1	E	340	Total	C	N	O	S	0	0	0
			2724	1774	443	492	15			
1	P	340	Total	C	N	O	S	0	0	0
			2724	1774	443	492	15			
1	Q	340	Total	C	N	O	S	0	0	0
			2724	1774	443	492	15			
1	R	340	Total	C	N	O	S	0	0	0
			2724	1774	443	492	15			
1	S	340	Total	C	N	O	S	0	0	0
			2724	1774	443	492	15			
1	T	340	Total	C	N	O	S	0	0	0
			2724	1774	443	492	15			

There are 110 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	303	ALA	-	linker	UNP G5EBR3
A	304	GLY	-	linker	UNP G5EBR3
A	305	THR	-	linker	UNP G5EBR3
A	340	HIS	-	expression tag	UNP G5EBR3
A	341	HIS	-	expression tag	UNP G5EBR3
A	342	HIS	-	expression tag	UNP G5EBR3
A	343	HIS	-	expression tag	UNP G5EBR3
A	344	HIS	-	expression tag	UNP G5EBR3

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Chain	Residue	Modelled	Actual	Comment	Reference
A	345	HIS	-	expression tag	UNP G5EBR3
A	346	HIS	-	expression tag	UNP G5EBR3
A	347	HIS	-	expression tag	UNP G5EBR3
B	303	ALA	-	linker	UNP G5EBR3
B	304	GLY	-	linker	UNP G5EBR3
B	305	THR	-	linker	UNP G5EBR3
B	340	HIS	-	expression tag	UNP G5EBR3
B	341	HIS	-	expression tag	UNP G5EBR3
B	342	HIS	-	expression tag	UNP G5EBR3
B	343	HIS	-	expression tag	UNP G5EBR3
B	344	HIS	-	expression tag	UNP G5EBR3
B	345	HIS	-	expression tag	UNP G5EBR3
B	346	HIS	-	expression tag	UNP G5EBR3
B	347	HIS	-	expression tag	UNP G5EBR3
C	303	ALA	-	linker	UNP G5EBR3
C	304	GLY	-	linker	UNP G5EBR3
C	305	THR	-	linker	UNP G5EBR3
C	340	HIS	-	expression tag	UNP G5EBR3
C	341	HIS	-	expression tag	UNP G5EBR3
C	342	HIS	-	expression tag	UNP G5EBR3
C	343	HIS	-	expression tag	UNP G5EBR3
C	344	HIS	-	expression tag	UNP G5EBR3
C	345	HIS	-	expression tag	UNP G5EBR3
C	346	HIS	-	expression tag	UNP G5EBR3
C	347	HIS	-	expression tag	UNP G5EBR3
D	303	ALA	-	linker	UNP G5EBR3
D	304	GLY	-	linker	UNP G5EBR3
D	305	THR	-	linker	UNP G5EBR3
D	340	HIS	-	expression tag	UNP G5EBR3
D	341	HIS	-	expression tag	UNP G5EBR3
D	342	HIS	-	expression tag	UNP G5EBR3
D	343	HIS	-	expression tag	UNP G5EBR3
D	344	HIS	-	expression tag	UNP G5EBR3
D	345	HIS	-	expression tag	UNP G5EBR3
D	346	HIS	-	expression tag	UNP G5EBR3
D	347	HIS	-	expression tag	UNP G5EBR3
E	303	ALA	-	linker	UNP G5EBR3
E	304	GLY	-	linker	UNP G5EBR3
E	305	THR	-	linker	UNP G5EBR3
E	340	HIS	-	expression tag	UNP G5EBR3
E	341	HIS	-	expression tag	UNP G5EBR3
E	342	HIS	-	expression tag	UNP G5EBR3

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Chain	Residue	Modelled	Actual	Comment	Reference
E	343	HIS	-	expression tag	UNP G5EBR3
E	344	HIS	-	expression tag	UNP G5EBR3
E	345	HIS	-	expression tag	UNP G5EBR3
E	346	HIS	-	expression tag	UNP G5EBR3
E	347	HIS	-	expression tag	UNP G5EBR3
P	303	ALA	-	linker	UNP G5EBR3
P	304	GLY	-	linker	UNP G5EBR3
P	305	THR	-	linker	UNP G5EBR3
P	340	HIS	-	expression tag	UNP G5EBR3
P	341	HIS	-	expression tag	UNP G5EBR3
P	342	HIS	-	expression tag	UNP G5EBR3
P	343	HIS	-	expression tag	UNP G5EBR3
P	344	HIS	-	expression tag	UNP G5EBR3
P	345	HIS	-	expression tag	UNP G5EBR3
P	346	HIS	-	expression tag	UNP G5EBR3
P	347	HIS	-	expression tag	UNP G5EBR3
Q	303	ALA	-	linker	UNP G5EBR3
Q	304	GLY	-	linker	UNP G5EBR3
Q	305	THR	-	linker	UNP G5EBR3
Q	340	HIS	-	expression tag	UNP G5EBR3
Q	341	HIS	-	expression tag	UNP G5EBR3
Q	342	HIS	-	expression tag	UNP G5EBR3
Q	343	HIS	-	expression tag	UNP G5EBR3
Q	344	HIS	-	expression tag	UNP G5EBR3
Q	345	HIS	-	expression tag	UNP G5EBR3
Q	346	HIS	-	expression tag	UNP G5EBR3
Q	347	HIS	-	expression tag	UNP G5EBR3
R	303	ALA	-	linker	UNP G5EBR3
R	304	GLY	-	linker	UNP G5EBR3
R	305	THR	-	linker	UNP G5EBR3
R	340	HIS	-	expression tag	UNP G5EBR3
R	341	HIS	-	expression tag	UNP G5EBR3
R	342	HIS	-	expression tag	UNP G5EBR3
R	343	HIS	-	expression tag	UNP G5EBR3
R	344	HIS	-	expression tag	UNP G5EBR3
R	345	HIS	-	expression tag	UNP G5EBR3
R	346	HIS	-	expression tag	UNP G5EBR3
R	347	HIS	-	expression tag	UNP G5EBR3
S	303	ALA	-	linker	UNP G5EBR3
S	304	GLY	-	linker	UNP G5EBR3
S	305	THR	-	linker	UNP G5EBR3
S	340	HIS	-	expression tag	UNP G5EBR3

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Chain	Residue	Modelled	Actual	Comment	Reference
S	341	HIS	-	expression tag	UNP G5EBR3
S	342	HIS	-	expression tag	UNP G5EBR3
S	343	HIS	-	expression tag	UNP G5EBR3
S	344	HIS	-	expression tag	UNP G5EBR3
S	345	HIS	-	expression tag	UNP G5EBR3
S	346	HIS	-	expression tag	UNP G5EBR3
S	347	HIS	-	expression tag	UNP G5EBR3
T	303	ALA	-	linker	UNP G5EBR3
T	304	GLY	-	linker	UNP G5EBR3
T	305	THR	-	linker	UNP G5EBR3
T	340	HIS	-	expression tag	UNP G5EBR3
T	341	HIS	-	expression tag	UNP G5EBR3
T	342	HIS	-	expression tag	UNP G5EBR3
T	343	HIS	-	expression tag	UNP G5EBR3
T	344	HIS	-	expression tag	UNP G5EBR3
T	345	HIS	-	expression tag	UNP G5EBR3
T	346	HIS	-	expression tag	UNP G5EBR3
T	347	HIS	-	expression tag	UNP G5EBR3

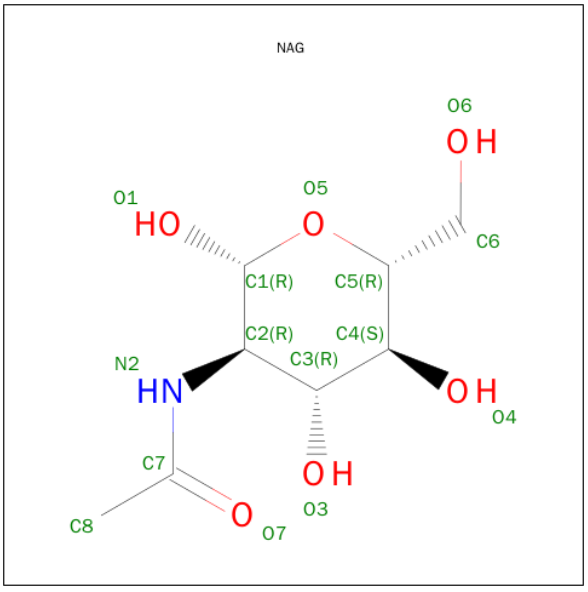
- Molecule 2 is a protein called Mouse monoclonal Fab fragment, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	221	Total	C	N	O	S	0	0	0
			1682	1067	273	334	8			
2	G	222	Total	C	N	O	S	0	0	0
			1693	1073	277	335	8			
2	H	223	Total	C	N	O	S	0	0	0
			1701	1077	278	338	8			
2	I	222	Total	C	N	O	S	0	0	0
			1693	1073	277	335	8			
2	U	221	Total	C	N	O	S	0	0	0
			1682	1067	273	334	8			
2	V	224	Total	C	N	O	S	0	0	0
			1707	1080	279	339	9			
2	W	224	Total	C	N	O	S	0	0	0
			1707	1080	279	339	9			
2	X	224	Total	C	N	O	S	0	0	0
			1707	1080	279	339	9			
2	Y	221	Total	C	N	O	S	0	0	0
			1682	1067	273	334	8			
2	J	222	Total	C	N	O	S	0	0	0
			1693	1073	277	335	8			

- Molecule 3 is a protein called Mouse monoclonal Fab fragment, light chain.

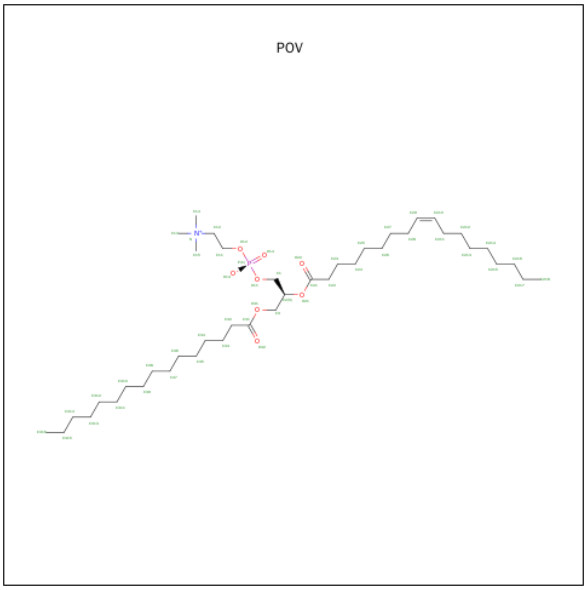
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	K	210	Total	C	N	O	S	0	0	0
			1590	999	266	319	6			
3	L	211	Total	C	N	O	S	0	0	0
			1601	1005	270	320	6			
3	N	211	Total	C	N	O	S	0	0	0
			1601	1005	270	320	6			
3	O	211	Total	C	N	O	S	0	0	0
			1601	1005	270	320	6			
3	Z	215	Total	C	N	O	S	0	0	0
			1626	1018	274	327	7			
3	f	215	Total	C	N	O	S	0	0	0
			1626	1018	274	327	7			
3	g	211	Total	C	N	O	S	0	0	0
			1601	1005	270	320	6			
3	h	211	Total	C	N	O	S	0	0	0
			1601	1005	270	320	6			
3	i	214	Total	C	N	O	S	0	0	0
			1620	1015	273	325	7			
3	M	210	Total	C	N	O	S	0	0	0
			1590	999	266	319	6			

- Molecule 4 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N O 14 8 1 5	0	0
4	B	1	Total C N O 14 8 1 5	0	0
4	C	1	Total C N O 14 8 1 5	0	0
4	D	1	Total C N O 14 8 1 5	0	0
4	E	1	Total C N O 14 8 1 5	0	0
4	P	1	Total C N O 14 8 1 5	0	0
4	Q	1	Total C N O 14 8 1 5	0	0
4	S	1	Total C N O 14 8 1 5	0	0
4	T	1	Total C N O 14 8 1 5	0	0

- Molecule 5 is (2S)-3-(hexadecanoyloxy)-2-[(9Z)-octadec-9-enoyloxy]propyl 2-(trimethylammonio)ethyl phosphate (three-letter code: POV) (formula: $C_{42}H_{82}NO_8P$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C N O P 23 14 1 7 1	0	0
5	D	1	Total C N O P 26 16 1 8 1	0	0

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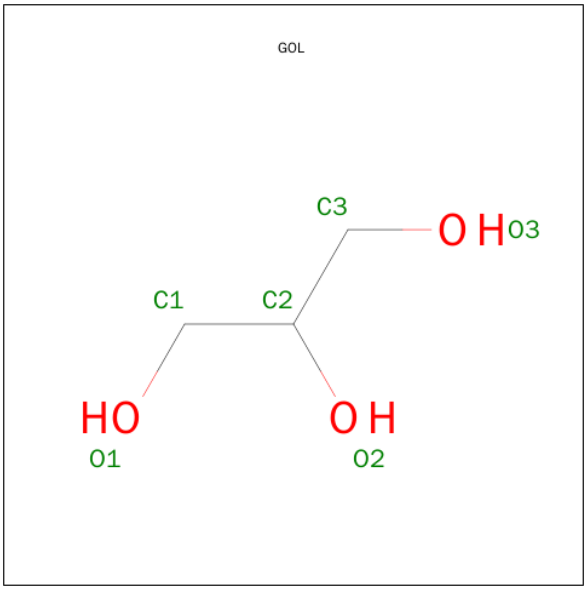
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	P	1	Total C N O P 42 32 1 8 1	0	0
5	Q	1	Total C N O P 46 36 1 8 1	0	0
5	R	1	Total C N O P 34 24 1 8 1	0	0
5	R	1	Total C O P 35 26 8 1	0	0
5	T	1	Total C N O P 43 33 1 8 1	0	0

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	P	1	Total Cl 1 1	0	0
6	A	1	Total Cl 1 1	0	0

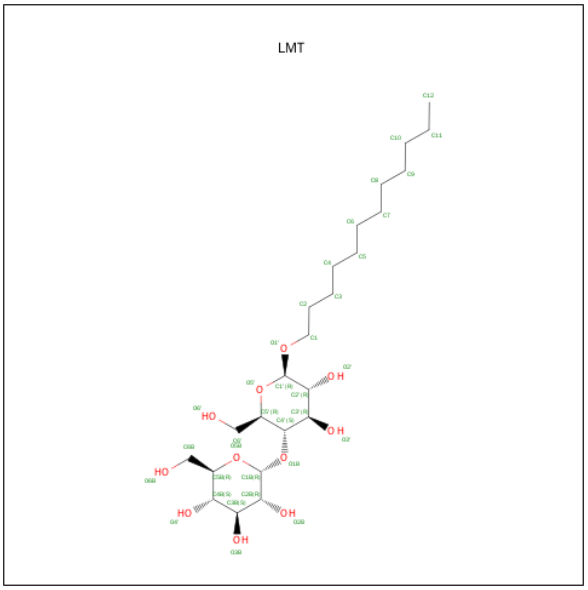
- Molecule 7 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	B	1	Total C O 6 3 3	0	0

- Molecule 8 is DODECYL-BETA-D-MALTOSIDE (three-letter code: LMT) (formula:

$C_{24}H_{46}O_{11}$).



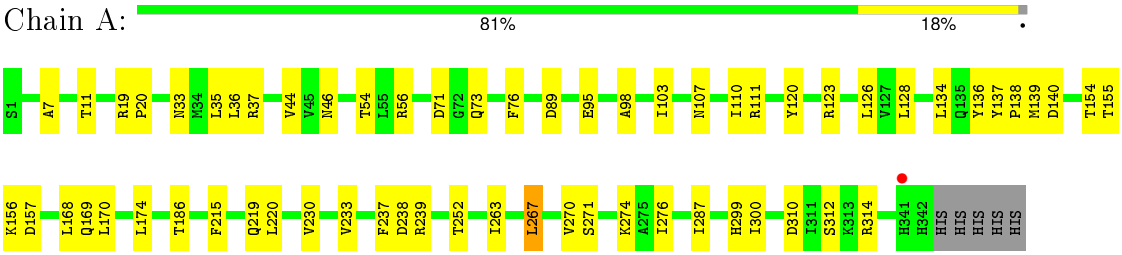
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	B	1	Total C O 23 12 11	0	0
8	C	1	Total C O 23 12 11	0	0
8	D	1	Total C O 23 12 11	0	0
8	E	1	Total C O 23 12 11	0	0
8	P	1	Total C O 23 12 11	0	0
8	S	1	Total C O 23 12 11	0	0
8	T	1	Total C O 23 12 11	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

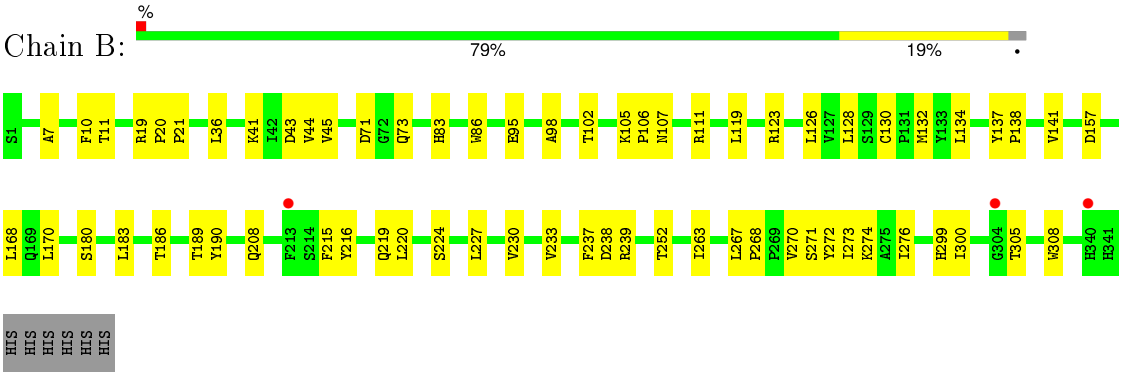
- Molecule 1: Avermectin-sensitive glutamate-gated chloride channel GluCl alpha

Chain A:



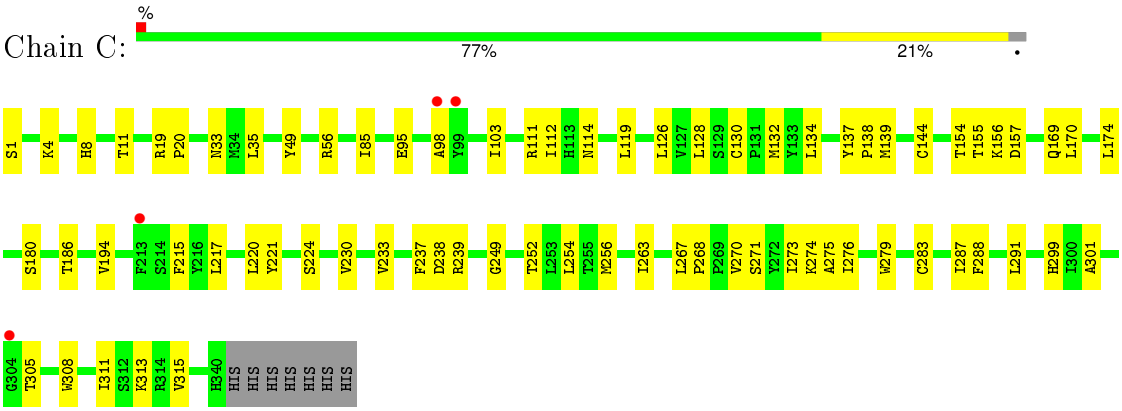
- Molecule 1: Avermectin-sensitive glutamate-gated chloride channel GluCl alpha

Chain B:



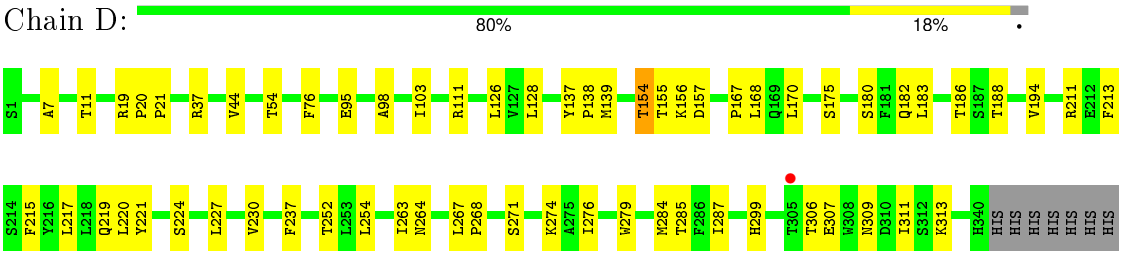
- Molecule 1: Avermectin-sensitive glutamate-gated chloride channel GluCl alpha

Chain C:



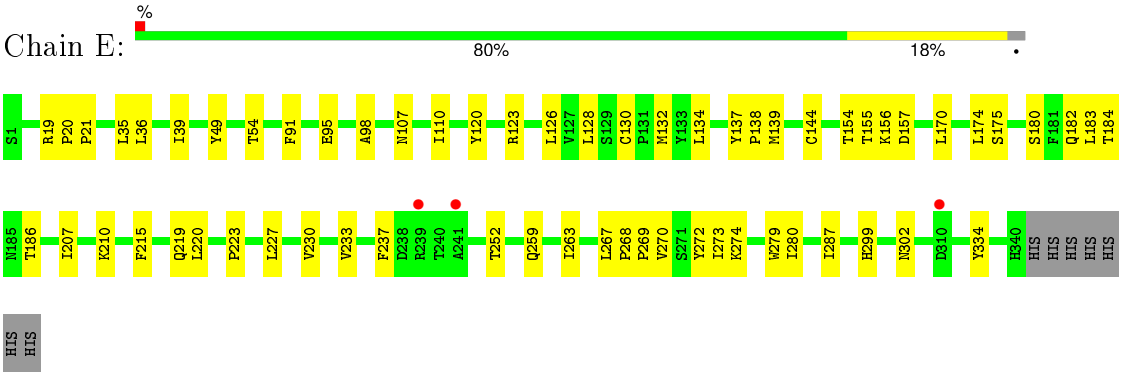
- Molecule 1: Avermectin-sensitive glutamate-gated chloride channel GluCl alpha

Chain D:



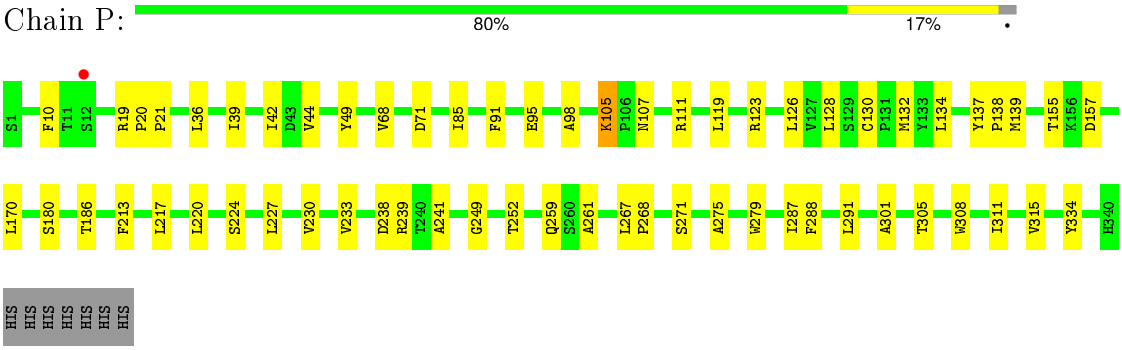
• Molecule 1: Avermectin-sensitive glutamate-gated chloride channel GluCl alpha

Chain E:



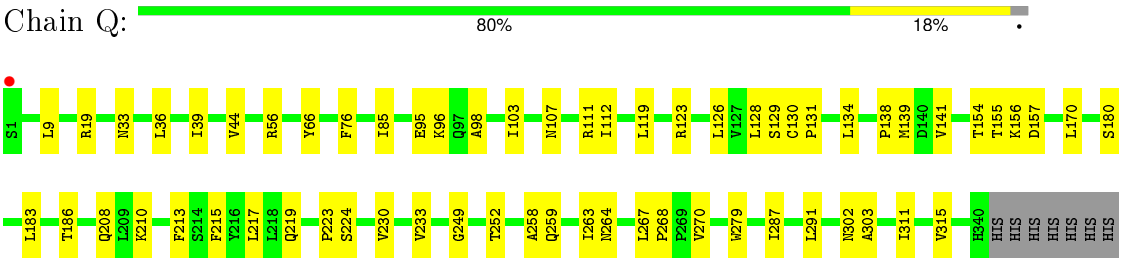
• Molecule 1: Avermectin-sensitive glutamate-gated chloride channel GluCl alpha

Chain P:



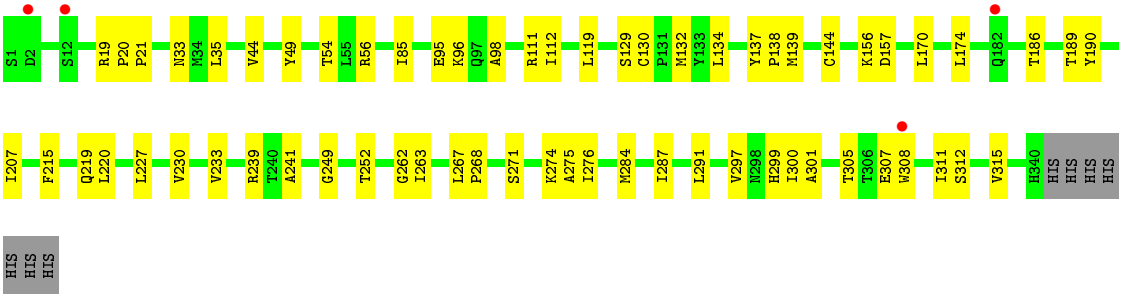
• Molecule 1: Avermectin-sensitive glutamate-gated chloride channel GluCl alpha

Chain Q:



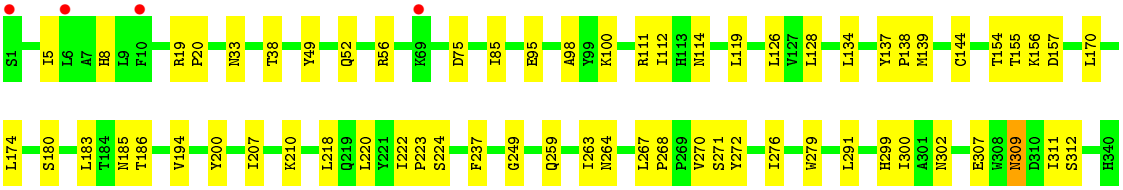
• Molecule 1: Avermectin-sensitive glutamate-gated chloride channel GluCl alpha

Chain R:



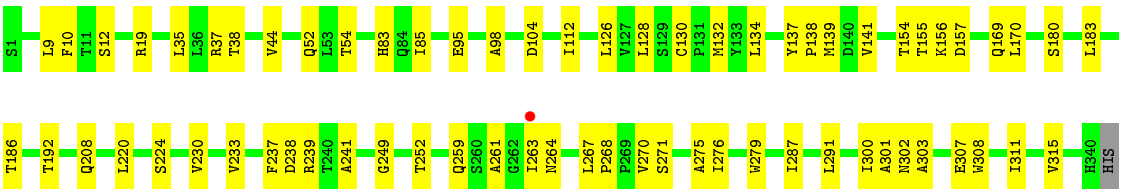
• Molecule 1: Avermectin-sensitive glutamate-gated chloride channel GluCl alpha

Chain S:



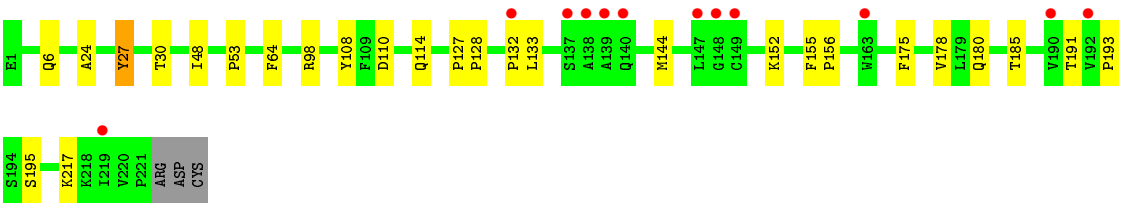
• Molecule 1: Avermectin-sensitive glutamate-gated chloride channel GluCl alpha

Chain T:



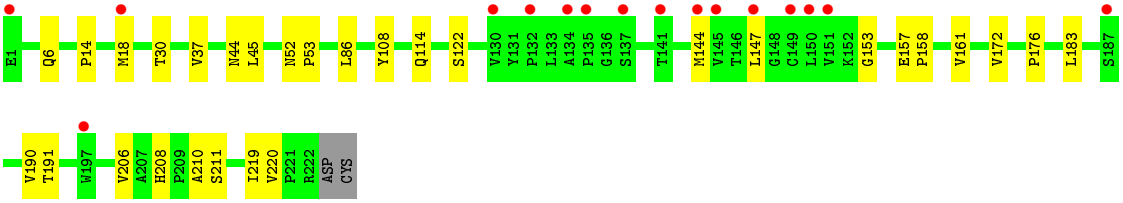
• Molecule 2: Mouse monoclonal Fab fragment, heavy chain

Chain F:

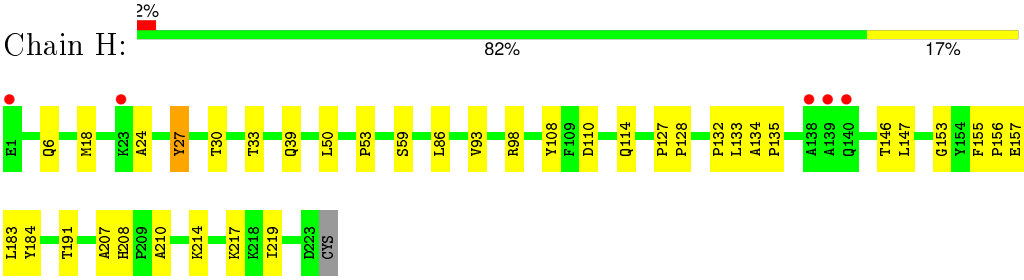


• Molecule 2: Mouse monoclonal Fab fragment, heavy chain

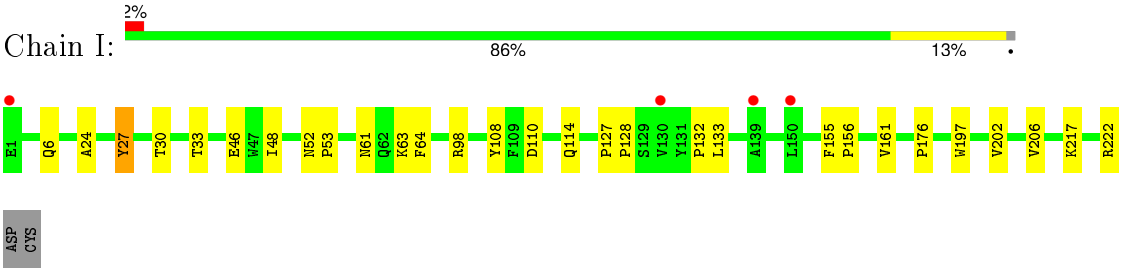
Chain G:



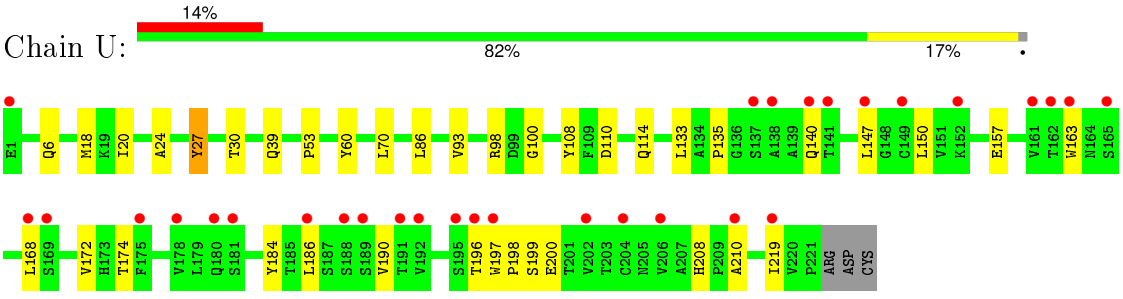
• Molecule 2: Mouse monoclonal Fab fragment, heavy chain



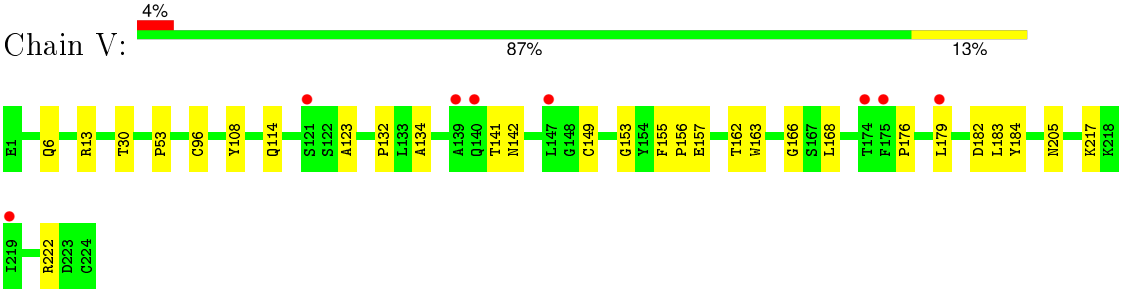
• Molecule 2: Mouse monoclonal Fab fragment, heavy chain



• Molecule 2: Mouse monoclonal Fab fragment, heavy chain

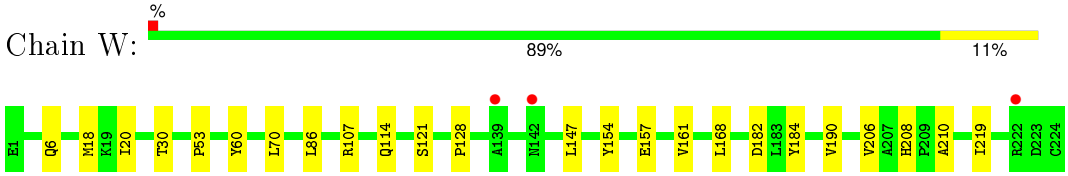


• Molecule 2: Mouse monoclonal Fab fragment, heavy chain



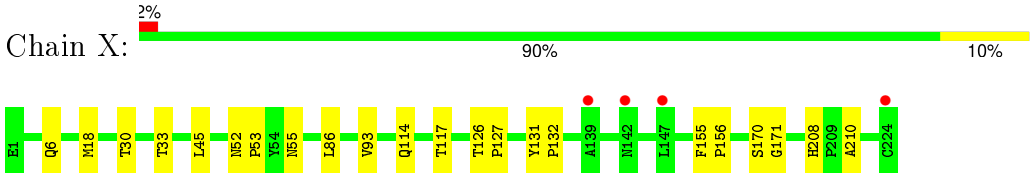
• Molecule 2: Mouse monoclonal Fab fragment, heavy chain

Chain W:



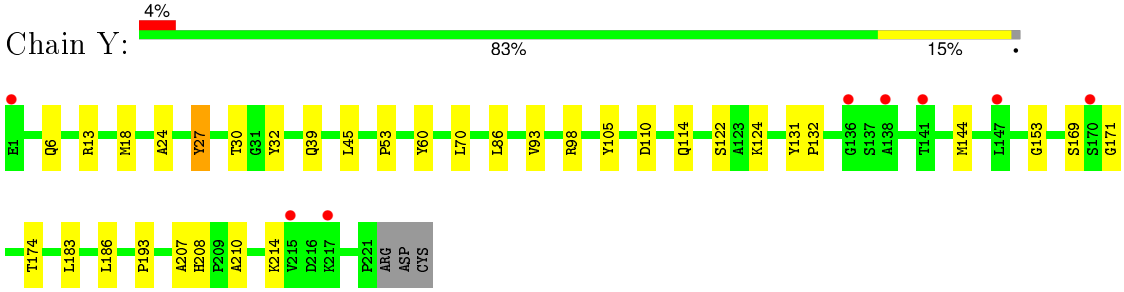
• Molecule 2: Mouse monoclonal Fab fragment, heavy chain

Chain X:



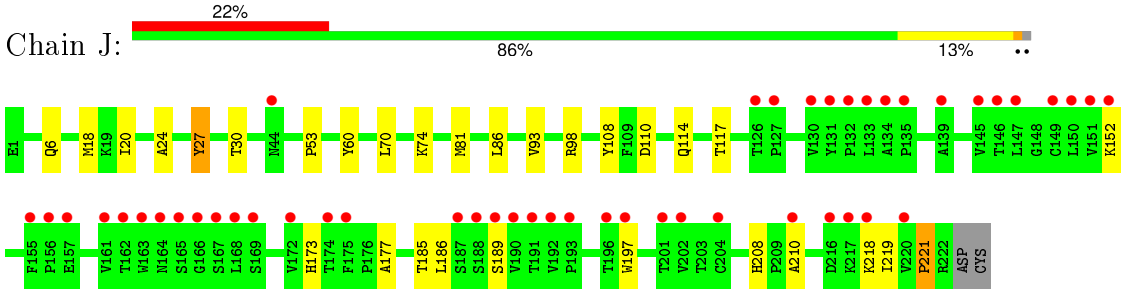
• Molecule 2: Mouse monoclonal Fab fragment, heavy chain

Chain Y:



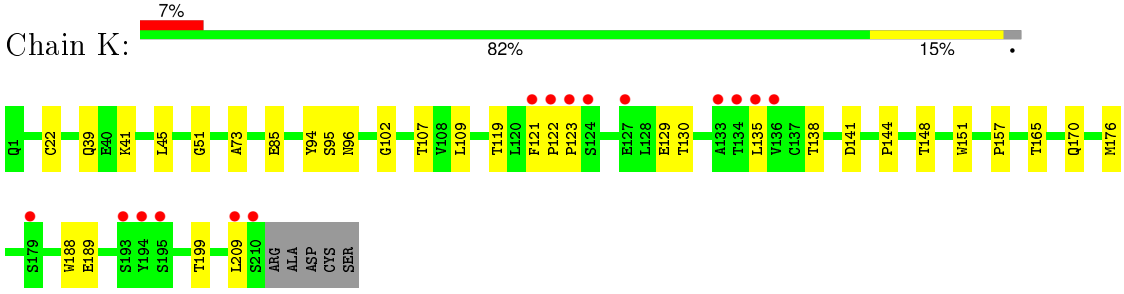
• Molecule 2: Mouse monoclonal Fab fragment, heavy chain

Chain J:



• Molecule 3: Mouse monoclonal Fab fragment, light chain

Chain K:



• Molecule 3: Mouse monoclonal Fab fragment, light chain

Chain L:

• Molecule 3: Mouse monoclonal Fab fragment, light chain

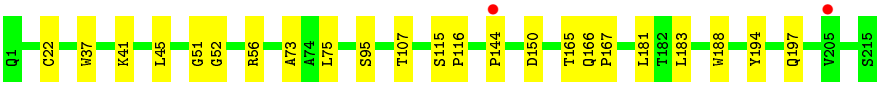
Chain N:

• Molecule 3: Mouse monoclonal Fab fragment, light chain

Chain O:

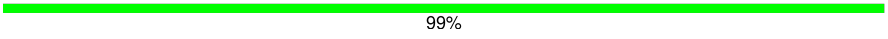
• Molecule 3: Mouse monoclonal Fab fragment, light chain

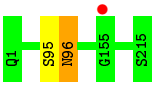
Chain Z:



• Molecule 3: Mouse monoclonal Fab fragment, light chain

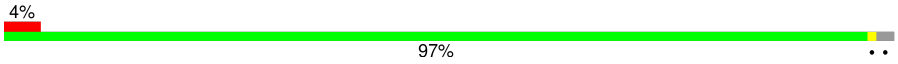
Chain f:

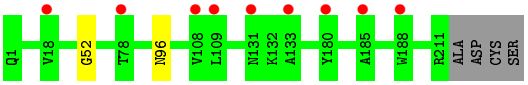




• Molecule 3: Mouse monoclonal Fab fragment, light chain

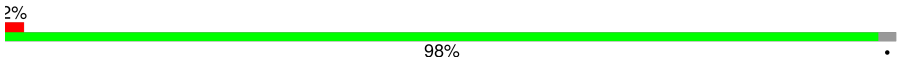
Chain g:





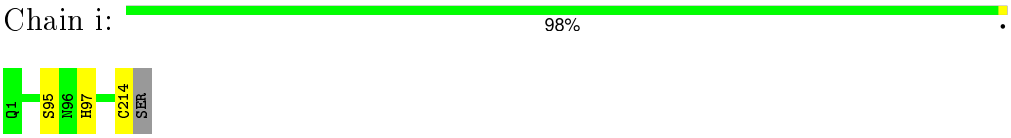
• Molecule 3: Mouse monoclonal Fab fragment, light chain

Chain h:



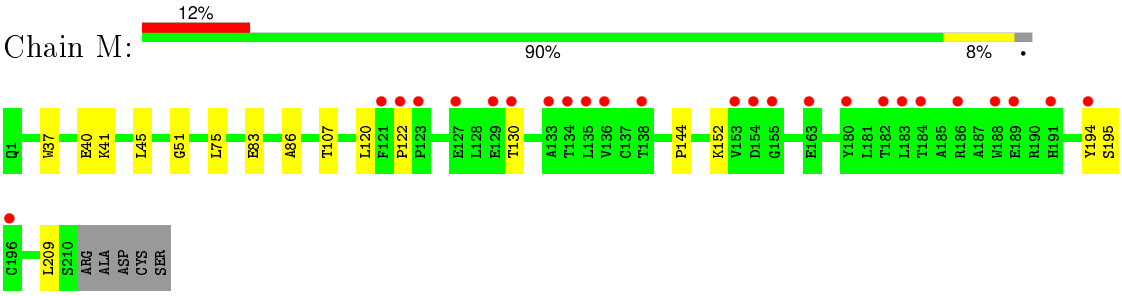
• Molecule 3: Mouse monoclonal Fab fragment, light chain

Chain i:



• Molecule 3: Mouse monoclonal Fab fragment, light chain

Chain M:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	453.75Å 192.87Å 196.14Å 90.00° 92.27° 90.00°	Depositor
Resolution (Å)	58.70 – 3.20 59.51 – 3.00	Depositor EDS
% Data completeness (in resolution range)	77.6 (58.70-3.20) 88.8 (59.51-3.00)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.25 (at 3.01Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.227 , 0.251 0.240 , 0.260	Depositor DCC
R_{free} test set	12812 reflections (5.26%)	DCC
Wilson B-factor (Å^2)	84.0	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.27 , 48.9	EDS
Estimated twinning fraction	0.021 for -h,-k,l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.43$, $< L^2 > = 0.26$	Xtriage
Outliers	0 of 297382 reflections	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	60818	wwPDB-VP
Average B, all atoms (Å^2)	115.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.31% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: POV, GOL, LMT, NAG, CL

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	0.24	0/2819	0.44	0/3847
1	B	0.23	0/2808	0.43	0/3832
1	C	0.23	0/2797	0.43	0/3817
1	D	0.25	0/2797	0.48	0/3817
1	E	0.24	0/2797	0.44	0/3817
1	P	0.24	0/2797	0.45	0/3817
1	Q	0.24	0/2797	0.45	0/3817
1	R	0.24	0/2797	0.43	0/3817
1	S	0.23	0/2797	0.43	0/3817
1	T	0.24	0/2797	0.45	0/3817
2	F	0.21	0/1728	0.40	0/2360
2	G	0.22	0/1739	0.41	0/2374
2	H	0.22	0/1747	0.41	0/2385
2	I	0.21	0/1739	0.40	0/2374
2	J	0.23	0/1739	0.42	0/2374
2	U	0.21	0/1728	0.40	0/2360
2	V	0.21	0/1753	0.40	0/2393
2	W	0.21	0/1753	0.40	0/2393
2	X	0.22	0/1753	0.42	0/2393
2	Y	0.21	0/1728	0.41	0/2360
3	K	0.22	0/1628	0.42	0/2226
3	L	0.22	0/1639	0.43	0/2240
3	M	0.22	0/1628	0.43	0/2226
3	N	0.23	0/1639	0.43	0/2240
3	O	0.22	0/1639	0.42	0/2240
3	Z	0.23	0/1664	0.43	0/2274
3	f	0.22	0/1664	0.42	0/2274
3	g	0.22	0/1639	0.42	0/2240
3	h	0.22	0/1639	0.41	0/2240
3	i	0.22	0/1658	0.43	0/2266
All	All	0.23	0/61847	0.43	0/84447

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	2744	0	2744	45	0
1	B	2734	0	2737	48	0
1	C	2724	0	2729	51	0
1	D	2724	0	2729	45	0
1	E	2724	0	2730	46	0
1	P	2724	0	2730	41	0
1	Q	2724	0	2730	41	0
1	R	2724	0	2731	43	0
1	S	2724	0	2730	50	0
1	T	2724	0	2730	49	0
2	F	1682	0	1632	24	0
2	G	1693	0	1645	20	0
2	H	1701	0	1649	26	0
2	I	1693	0	1645	23	0
2	J	1693	0	1645	23	0
2	U	1682	0	1632	22	0
2	V	1707	0	1653	20	0
2	W	1707	0	1653	14	0
2	X	1707	0	1653	13	0
2	Y	1682	0	1632	21	0
3	K	1590	0	1542	23	0
3	L	1601	0	1555	19	0
3	M	1590	0	1542	11	0
3	N	1601	0	1555	21	0
3	O	1601	0	1555	16	0
3	Z	1626	0	1573	13	0
3	f	1626	0	1573	0	0
3	g	1601	0	1555	0	0
3	h	1601	0	1555	0	0
3	i	1620	0	1568	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	14	0	13	0	0
4	B	14	0	13	0	0
4	C	14	0	13	0	0
4	D	14	0	13	0	0
4	E	14	0	13	0	0
4	P	14	0	13	0	0
4	Q	14	0	13	0	0
4	S	14	0	13	1	0
4	T	14	0	13	0	0
5	A	23	0	26	1	0
5	D	26	0	26	5	0
5	P	42	0	61	3	0
5	Q	46	0	67	5	0
5	R	69	0	83	5	0
5	T	43	0	63	3	0
6	A	1	0	0	0	0
6	P	1	0	0	0	0
7	B	6	0	8	0	0
8	B	23	0	21	0	0
8	C	23	0	21	0	0
8	D	23	0	21	1	0
8	E	23	0	21	0	0
8	P	23	0	21	0	0
8	S	23	0	21	1	0
8	T	23	0	21	1	0
All	All	60818	0	59930	703	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 6.

All (703) 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:D:306:THR:HA	1:D:309:ASN:HB2	1.53	0.90
1:Q:302:ASN:HD22	1:R:241:ALA:HB2	1.36	0.89
1:T:37:ARG:NH2	1:T:54:THR:OG1	2.10	0.84
1:C:305:THR:HG22	1:C:308:TRP:HD1	1.43	0.83
1:R:156:LYS:HA	3:Z:95:SER:HB2	1.61	0.83
1:D:156:LYS:HA	3:O:95:SER:HB2	1.61	0.81
2:Y:13:ARG:HH22	2:Y:124:LYS:HD2	1.46	0.81
1:Q:302:ASN:ND2	1:R:241:ALA:HB2	1.96	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:37:ARG:NH1	1:D:54:THR:OG1	2.17	0.78
1:Q:141:VAL:HG22	1:Q:210:LYS:HG2	1.67	0.76
1:P:241:ALA:HB2	1:T:302:ASN:HD22	1.52	0.75
1:P:91:PHE:HA	1:Q:103:ILE:HD11	1.69	0.75
1:B:220:LEU:HD22	1:B:263:ILE:HG13	1.68	0.74
3:M:194:TYR:HB2	3:M:209:LEU:HB3	1.69	0.74
1:B:44:VAL:HG11	1:B:134:LEU:HD11	1.69	0.74
1:T:10:PHE:HE2	1:T:83:HIS:HB3	1.52	0.74
1:Q:44:VAL:HG11	1:Q:134:LEU:HD11	1.69	0.73
3:M:107:THR:HG21	3:M:144:PRO:HB3	1.73	0.71
1:Q:267:LEU:HD12	1:Q:268:PRO:HD2	1.72	0.70
1:B:71:ASP:OD2	1:B:73:GLN:NE2	2.23	0.70
1:C:305:THR:HG22	1:C:308:TRP:CD1	2.25	0.70
1:P:267:LEU:HD12	1:P:268:PRO:HD2	1.73	0.69
2:F:152:LYS:HB2	2:F:195:SER:HB2	30.38	0.69
2:V:153:GLY:HA2	2:V:183:LEU:HB3	1.73	0.69
2:I:6:GLN:H	2:I:114:GLN:HE22	1.41	0.69
2:V:6:GLN:H	2:V:114:GLN:HE22	1.40	0.69
1:T:267:LEU:HD12	1:T:268:PRO:HD2	1.74	0.68
1:R:267:LEU:HD12	1:R:268:PRO:HD2	1.74	0.68
1:S:111:ARG:HD2	1:S:119:LEU:HD23	1.76	0.68
2:V:108:TYR:HB2	3:Z:51:GLY:HA3	1.76	0.68
2:I:108:TYR:HB2	3:O:51:GLY:HA3	1.74	0.68
1:T:10:PHE:CE2	1:T:83:HIS:HB3	2.28	0.68
1:A:238:ASP:OD1	1:A:239:ARG:N	2.27	0.68
2:J:219:ILE:HG22	2:J:221:PRO:HD3	1.76	0.67
2:W:6:GLN:H	2:W:114:GLN:HE22	1.42	0.67
1:S:210:LYS:NZ	8:S:402:LMT:O4'	2.27	0.67
2:J:108:TYR:HB2	3:M:51:GLY:HA3	1.76	0.67
2:G:108:TYR:HB2	3:K:51:GLY:HA3	1.77	0.67
3:K:107:THR:HG21	3:K:144:PRO:HB3	1.76	0.67
1:C:267:LEU:HD21	1:C:274:LYS:HE3	1.77	0.67
2:H:108:TYR:HB2	3:L:51:GLY:HA3	1.78	0.66
1:E:186:THR:HG22	1:E:207:ILE:HG22	1.78	0.66
1:D:267:LEU:HD11	1:D:274:LYS:HE3	1.78	0.66
5:D:401:POV:O13	1:E:259:GLN:NE2	2.28	0.66
1:A:37:ARG:NH1	1:A:54:THR:OG1	2.29	0.66
1:Q:249:GLY:HA3	1:Q:291:LEU:HD13	1.78	0.66
2:U:6:GLN:H	2:U:114:GLN:HE22	1.43	0.66
1:E:134:LEU:HD12	1:E:270:VAL:HG21	1.78	0.66
3:O:135:LEU:HD12	3:O:181:LEU:HD23	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:103:ILE:HG21	1:E:98:ALA:HB3	1.78	0.66
1:T:238:ASP:OD1	1:T:239:ARG:N	2.28	0.65
3:N:107:THR:HG21	3:N:144:PRO:HB3	1.78	0.65
2:J:173:HIS:HB2	2:J:189:SER:HB2	1.78	0.65
2:V:134:ALA:HB3	2:V:222:ARG:HG3	1.78	0.65
2:W:30:THR:HA	2:W:53:PRO:HB2	1.78	0.65
1:Q:36:LEU:HD21	1:Q:39:ILE:HD11	1.78	0.65
3:K:41:LYS:HB2	3:K:45:LEU:HB3	1.77	0.65
1:B:267:LEU:HD21	1:B:274:LYS:HE3	1.78	0.65
1:R:44:VAL:HG11	1:R:134:LEU:HD11	1.78	0.65
1:D:307:GLU:O	1:D:311:ILE:HG13	1.96	0.64
1:T:249:GLY:HA3	1:T:291:LEU:HD13	1.79	0.64
1:C:220:LEU:HD22	1:C:263:ILE:HG13	1.78	0.64
1:A:238:ASP:OD2	1:E:302:ASN:ND2	2.31	0.64
2:G:6:GLN:H	2:G:114:GLN:HE22	1.44	0.64
1:T:300:ILE:HG23	1:T:308:TRP:HE3	1.62	0.64
2:J:6:GLN:H	2:J:114:GLN:HE22	1.45	0.64
2:H:6:GLN:H	2:H:114:GLN:HE22	1.46	0.64
2:Y:6:GLN:H	2:Y:114:GLN:HE22	1.44	0.63
1:P:238:ASP:OD1	1:P:239:ARG:N	2.31	0.63
1:B:10:PHE:CE2	1:B:83:HIS:HB3	2.33	0.63
1:P:44:VAL:HG11	1:P:134:LEU:HD11	1.80	0.63
5:D:401:POV:H3A	1:E:223:PRO:HB3	1.79	0.62
1:B:238:ASP:OD1	1:B:239:ARG:N	2.32	0.62
2:I:222:ARG:NH2	3:O:123:PRO:O	2.32	0.62
2:F:144:MET:HB2	2:F:191:THR:HG22	1.81	0.62
3:O:96:ASN:O	3:O:96:ASN:ND2	2.28	0.62
2:F:6:GLN:H	2:F:114:GLN:HE22	1.45	0.62
3:L:107:THR:HG21	3:L:144:PRO:HB3	1.82	0.61
1:T:19:ARG:HH11	1:T:157:ASP:HA	1.65	0.61
1:Q:141:VAL:HG11	1:Q:208:GLN:HE21	1.66	0.61
1:P:180:SER:HA	1:T:271:SER:HB2	1.81	0.61
2:I:30:THR:HA	2:I:53:PRO:HB2	1.82	0.61
1:S:154:THR:OG1	1:S:155:THR:N	2.34	0.61
1:A:36:LEU:HD13	1:A:168:LEU:HD11	1.82	0.61
1:S:267:LEU:HD12	1:S:268:PRO:HD2	1.83	0.61
2:Y:24:ALA:HB1	2:Y:27:TYR:HE1	1.66	0.61
3:K:22:CYS:HB3	3:K:73:ALA:HB3	1.82	0.61
1:D:44:VAL:HG21	1:D:268:PRO:HG3	1.82	0.61
1:C:170:LEU:HD23	1:C:174:LEU:HD23	1.82	0.60
2:V:30:THR:HA	2:V:53:PRO:HB2	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:98:ARG:NH2	2:I:110:ASP:OD2	2.27	0.60
2:H:33:THR:HG21	2:H:50:LEU:HD12	1.84	0.60
1:D:213:PHE:CE2	1:D:217:LEU:HD12	2.37	0.60
1:D:220:LEU:HD22	1:D:263:ILE:HG13	1.83	0.60
1:A:44:VAL:HG11	1:A:134:LEU:HD11	1.83	0.60
2:F:98:ARG:NH2	2:F:110:ASP:OD2	2.29	0.60
1:E:267:LEU:HD21	1:E:274:LYS:HE3	1.84	0.59
1:T:35:LEU:HB3	1:T:54:THR:HB	1.84	0.59
2:Y:24:ALA:HB1	2:Y:27:TYR:CE1	2.38	0.59
1:A:134:LEU:HD12	1:A:270:VAL:HG21	1.83	0.59
2:I:24:ALA:HB1	2:I:27:TYR:HE1	1.67	0.59
1:R:170:LEU:HD23	1:R:174:LEU:HD23	1.85	0.58
1:P:155:THR:OG1	2:W:107:ARG:NH2	2.37	0.58
1:S:218:LEU:HD23	1:S:222:ILE:HG13	1.83	0.58
3:N:94:TYR:O	3:N:96:ASN:N	2.36	0.58
2:H:147:LEU:HD13	2:H:219:ILE:HG13	1.84	0.58
2:Y:30:THR:HA	2:Y:53:PRO:HB2	1.86	0.58
1:P:249:GLY:HA3	1:P:291:LEU:HD13	1.85	0.58
1:S:134:LEU:HD12	1:S:270:VAL:HG21	1.86	0.58
1:D:7:ALA:O	1:D:11:THR:HG23	2.04	0.58
1:C:98:ALA:HB3	1:D:103:ILE:HG13	1.85	0.57
2:H:98:ARG:NH2	2:H:110:ASP:OD2	2.33	0.57
1:T:9:LEU:O	1:T:12:SER:OG	2.21	0.57
3:K:39:GLN:NE2	3:K:41:LYS:HE3	2.18	0.57
1:R:239:ARG:NH1	1:R:301:ALA:HB1	2.19	0.57
1:D:19:ARG:HH11	1:D:157:ASP:HA	1.69	0.57
3:L:135:LEU:HB2	3:L:181:LEU:HB3	1.86	0.57
1:A:7:ALA:O	1:A:11:THR:HG23	2.03	0.57
1:B:36:LEU:HD13	1:B:168:LEU:HD11	1.85	0.57
1:C:238:ASP:OD1	1:C:239:ARG:N	2.37	0.57
2:G:30:THR:HA	2:G:53:PRO:HB2	1.85	0.57
1:P:107:ASN:OD1	1:P:123:ARG:NH1	2.38	0.57
3:Z:41:LYS:HB2	3:Z:45:LEU:HB2	1.86	0.57
1:P:227:LEU:HD13	5:P:403:POV:H22A	1.87	0.57
2:X:93:VAL:HG22	2:X:117:THR:HG22	1.86	0.57
3:N:135:LEU:HB2	3:N:181:LEU:HB3	1.87	0.57
2:X:30:THR:HA	2:X:53:PRO:HB2	1.86	0.56
2:W:128:PRO:HB3	2:W:154:TYR:HB3	1.86	0.56
2:F:27:TYR:CE2	2:F:98:ARG:HD2	2.41	0.56
2:H:24:ALA:HB1	2:H:27:TYR:HE1	1.68	0.56
3:M:41:LYS:HB2	3:M:45:LEU:HB2	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:9:LEU:HD11	1:Q:66:TYR:HB3	1.86	0.56
3:N:152:LYS:HE3	3:N:155:GLY:HA2	1.86	0.56
1:C:154:THR:OG1	1:C:155:THR:N	2.38	0.56
1:A:239:ARG:HD3	1:A:312:SER:HB3	1.87	0.56
1:T:154:THR:OG1	1:T:155:THR:N	2.38	0.56
1:B:98:ALA:HB3	1:C:103:ILE:HG13	1.88	0.56
1:A:156:LYS:HA	3:L:95:SER:HB2	1.87	0.56
2:V:13:ARG:NH2	2:V:123:ALA:O	2.39	0.56
1:Q:258:ALA:HB1	5:Q:401:POV:H13B	1.87	0.56
1:T:220:LEU:HD22	1:T:263:ILE:HG13	1.87	0.55
2:U:30:THR:HA	2:U:53:PRO:HB2	1.89	0.55
1:Q:154:THR:OG1	1:Q:155:THR:N	2.38	0.55
2:F:24:ALA:HB1	2:F:27:TYR:HE1	1.69	0.55
2:J:93:VAL:HG22	2:J:117:THR:HG22	1.87	0.55
1:P:36:LEU:HD21	1:P:39:ILE:HD11	1.87	0.55
2:H:24:ALA:HB1	2:H:27:TYR:CE1	2.42	0.55
1:T:302:ASN:OD1	1:T:303:ALA:N	2.40	0.55
2:X:6:GLN:H	2:X:114:GLN:HE22	1.53	0.55
1:B:271:SER:HB2	1:C:180:SER:HA	1.89	0.55
2:J:98:ARG:NH2	2:J:110:ASP:OD2	2.32	0.55
1:A:126:LEU:HD13	1:A:128:LEU:HD21	1.87	0.55
1:A:219:GLN:NE2	5:A:402:POV:H12	2.21	0.55
2:U:60:TYR:HE1	2:U:70:LEU:HG	1.72	0.55
2:I:24:ALA:HB1	2:I:27:TYR:CE1	2.41	0.55
1:T:44:VAL:HG11	1:T:134:LEU:HD11	1.88	0.55
2:U:172:VAL:HA	2:U:190:VAL:HG12	1.87	0.55
1:P:42:ILE:HG12	1:P:49:TYR:HB2	1.87	0.55
1:R:137:TYR:HB3	1:R:274:LYS:HB3	1.88	0.55
1:Q:111:ARG:HD2	1:Q:119:LEU:HD23	1.89	0.55
1:B:107:ASN:OD1	1:B:123:ARG:NH1	2.39	0.55
1:E:184:THR:HG21	1:E:210:LYS:HD2	1.88	0.55
3:O:94:TYR:HB2	3:O:97:HIS:HB3	1.89	0.54
1:R:111:ARG:HD2	1:R:119:LEU:HD23	1.89	0.54
2:Y:39:GLN:HB3	2:Y:93:VAL:HG13	1.89	0.54
1:P:95:GLU:HG3	1:P:98:ALA:HB2	1.88	0.54
1:E:154:THR:OG1	1:E:155:THR:N	2.40	0.54
1:B:141:VAL:HG11	1:B:208:GLN:HE21	1.73	0.54
2:V:108:TYR:CB	3:Z:51:GLY:HA3	2.37	0.54
1:D:220:LEU:HD23	1:D:276:ILE:HD11	1.90	0.54
1:P:261:ALA:HA	5:Q:401:POV:H14B	1.90	0.54
1:E:126:LEU:HD13	1:E:128:LEU:HD21	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:172:VAL:HA	2:G:190:VAL:HG12	1.90	0.54
1:P:19:ARG:HH11	1:P:157:ASP:HA	1.73	0.54
1:C:252:THR:HG22	1:C:287:ILE:HD13	1.88	0.54
1:E:252:THR:HG22	1:E:287:ILE:HD13	1.90	0.54
1:D:284:MET:HG2	5:D:401:POV:H1	1.88	0.54
1:A:103:ILE:HD11	1:E:91:PHE:HA	1.90	0.54
1:E:95:GLU:HG3	1:E:98:ALA:HB2	1.89	0.54
1:S:249:GLY:HA3	1:S:291:LEU:HD13	1.90	0.53
1:D:139:MET:HG2	8:D:403:LMT:H6D	1.90	0.53
1:Q:302:ASN:OD1	1:Q:303:ALA:N	2.42	0.53
1:T:85:ILE:HD11	1:T:112:ILE:HD11	1.90	0.53
1:D:95:GLU:HG3	1:D:98:ALA:HB2	1.89	0.53
3:K:170:GLN:HE21	3:K:176:MET:HB3	1.73	0.53
2:U:24:ALA:HB1	2:U:27:TYR:HE1	1.73	0.53
3:L:152:LYS:HB2	3:L:195:SER:HB3	1.90	0.53
1:D:167:PRO:HB2	1:D:188:THR:HG21	1.90	0.53
2:G:108:TYR:CB	3:K:51:GLY:HA3	2.38	0.53
3:O:135:LEU:HB2	3:O:181:LEU:HB3	1.90	0.53
1:C:85:ILE:HD11	1:C:112:ILE:HD11	1.90	0.53
1:A:95:GLU:HG3	1:A:98:ALA:HB2	1.90	0.53
1:E:107:ASN:OD1	1:E:123:ARG:NH1	2.41	0.53
3:K:119:THR:OG1	3:K:138:THR:OG1	2.26	0.53
2:Y:98:ARG:NH2	2:Y:110:ASP:OD2	2.40	0.53
1:C:126:LEU:HD13	1:C:128:LEU:HD21	1.91	0.53
3:K:123:PRO:HD3	3:K:135:LEU:HG	1.91	0.53
1:D:267:LEU:HD12	1:D:268:PRO:HD2	1.90	0.53
1:B:134:LEU:HD12	1:B:270:VAL:HG21	1.91	0.53
5:D:401:POV:P	1:E:259:GLN:HE22	2.32	0.53
1:A:154:THR:OG1	1:A:155:THR:N	2.38	0.52
3:O:85:GLU:HG3	3:O:107:THR:HA	1.89	0.52
1:B:126:LEU:HD13	1:B:128:LEU:HD21	1.90	0.52
1:A:71:ASP:OD2	1:A:73:GLN:NE2	2.29	0.52
2:J:108:TYR:CB	3:M:51:GLY:HA3	2.39	0.52
2:V:132:PRO:HG3	2:V:217:LYS:HB3	1.92	0.52
2:F:108:TYR:HB2	3:N:51:GLY:HA2	1.91	0.52
1:A:46:ASN:ND2	1:B:41:LYS:HE2	2.25	0.52
2:H:30:THR:HA	2:H:53:PRO:HB2	1.91	0.52
1:S:126:LEU:HD13	1:S:128:LEU:HD21	1.92	0.52
1:T:134:LEU:HD12	1:T:270:VAL:HG21	1.92	0.52
3:N:131:ASN:HA	3:N:185:ALA:HB2	1.92	0.52
2:H:39:GLN:HB3	2:H:93:VAL:HG13	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:207:ALA:HB2	2:H:214:LYS:HD3	1.92	0.51
1:Q:252:THR:HG22	1:Q:287:ILE:HD13	1.92	0.51
1:A:220:LEU:HD22	1:A:263:ILE:HG13	1.92	0.51
1:A:267:LEU:HD11	1:A:274:LYS:HE3	1.91	0.51
3:Z:37:TRP:CE2	3:Z:75:LEU:HB2	2.46	0.51
2:V:176:PRO:HG2	3:Z:165:THR:HB	1.92	0.51
2:I:6:GLN:N	2:I:114:GLN:HE22	2.08	0.51
1:S:220:LEU:HD22	1:S:263:ILE:HG13	1.92	0.51
1:R:239:ARG:HH11	1:R:301:ALA:HB1	1.76	0.51
1:T:263:ILE:HG21	1:T:276:ILE:HG12	1.93	0.51
2:G:147:LEU:HD13	2:G:219:ILE:HD13	1.93	0.51
1:E:138:PRO:HD2	1:E:139:MET:SD	2.50	0.51
2:U:39:GLN:HB3	2:U:93:VAL:HG13	1.92	0.51
1:C:249:GLY:HA3	1:C:291:LEU:HD13	1.92	0.51
2:F:24:ALA:HB1	2:F:27:TYR:CE1	2.45	0.51
1:P:213:PHE:CE2	1:P:217:LEU:HD12	2.45	0.51
1:Q:213:PHE:CE2	1:Q:217:LEU:HD12	2.45	0.51
3:L:37:TRP:CE2	3:L:75:LEU:HB2	2.45	0.51
1:T:252:THR:HG22	1:T:287:ILE:HD13	1.93	0.51
1:P:105:LYS:N	1:P:105:LYS:HD2	2.26	0.51
2:J:30:THR:HA	2:J:53:PRO:HB2	1.93	0.51
1:Q:170:LEU:HD11	1:Q:186:THR:HG21	1.93	0.51
1:C:95:GLU:HG3	1:C:98:ALA:HB2	1.92	0.51
3:L:22:CYS:HB3	3:L:73:ALA:HB3	1.92	0.51
1:A:89:ASP:HB2	1:B:105:LYS:HD3	1.93	0.51
1:S:154:THR:HG23	1:S:156:LYS:HG2	1.93	0.51
1:P:239:ARG:NH1	1:P:301:ALA:HB1	2.26	0.51
2:U:24:ALA:HB1	2:U:27:TYR:CE1	2.45	0.51
1:S:95:GLU:HG3	1:S:98:ALA:HB2	1.92	0.51
2:H:108:TYR:CB	3:L:51:GLY:HA3	2.41	0.51
1:B:7:ALA:O	1:B:11:THR:HG23	2.10	0.51
2:W:161:VAL:HG22	2:W:206:VAL:HG22	1.92	0.51
1:R:85:ILE:HD11	1:R:112:ILE:HD11	1.92	0.51
1:T:154:THR:HG23	1:T:156:LYS:HG2	1.93	0.51
3:L:93:TRP:NE1	3:L:95:SER:HA	2.26	0.51
2:J:24:ALA:HB1	2:J:27:TYR:CE1	2.46	0.51
1:Q:19:ARG:HH11	1:Q:157:ASP:HA	1.76	0.51
1:Q:85:ILE:HD11	1:Q:112:ILE:HD11	1.93	0.51
1:D:183:LEU:HD21	1:D:186:THR:HG23	1.93	0.51
1:T:37:ARG:HH11	1:T:52:GLN:HE21	1.58	0.50
2:I:27:TYR:CE2	2:I:98:ARG:HD2	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:T:95:GLU:HG3	1:T:98:ALA:HB2	1.92	0.50
1:R:262:GLY:HA3	5:R:401:POV:H12	1.92	0.50
1:B:95:GLU:HG3	1:B:98:ALA:HB2	1.93	0.50
2:U:98:ARG:NH2	2:U:110:ASP:OD2	2.36	0.50
1:S:38:THR:HG22	1:S:52:GLN:HB3	1.92	0.50
1:C:33:ASN:HB3	1:C:56:ARG:HB2	1.93	0.50
1:T:38:THR:HG22	1:T:52:GLN:HB3	1.92	0.50
3:N:170:GLN:NE2	3:N:172:ASN:OD1	2.45	0.50
2:Y:27:TYR:CE2	2:Y:98:ARG:HD2	2.46	0.50
2:I:132:PRO:HD3	2:I:217:LYS:HE2	1.93	0.50
1:C:256:MET:HE3	1:C:283:CYS:HB3	1.94	0.50
2:U:133:LEU:HD23	2:U:150:LEU:HB2	1.92	0.50
1:T:126:LEU:HD13	1:T:128:LEU:HD21	1.92	0.50
1:C:154:THR:HG23	1:C:156:LYS:HG2	1.92	0.50
1:E:36:LEU:HD21	1:E:39:ILE:HD11	1.94	0.50
1:E:215:PHE:CE1	1:E:219:GLN:HG3	2.46	0.50
1:R:95:GLU:HG3	1:R:98:ALA:HB2	1.94	0.50
1:D:170:LEU:HD11	1:D:186:THR:HG21	1.94	0.50
1:R:137:TYR:O	1:R:275:ALA:HB3	2.12	0.50
1:S:19:ARG:HH11	1:S:157:ASP:HA	1.75	0.50
1:P:10:PHE:HZ	1:P:85:ILE:HG22	1.77	0.50
1:S:307:GLU:HG2	1:S:309:ASN:H	1.76	0.50
1:D:154:THR:HG22	1:D:155:THR:H	1.77	0.49
1:S:194:VAL:O	2:X:55:ASN:ND2	2.43	0.49
1:B:183:LEU:HD21	1:B:186:THR:HG23	1.93	0.49
2:V:162:THR:HG23	2:V:205:ASN:HB2	1.94	0.49
1:Q:126:LEU:HD13	1:Q:128:LEU:HD21	1.95	0.49
2:V:123:ALA:HB2	2:V:182:ASP:HB3	1.94	0.49
1:E:215:PHE:CZ	1:E:219:GLN:HG3	2.47	0.49
2:U:147:LEU:HD13	2:U:219:ILE:HD12	1.93	0.49
1:E:220:LEU:HD13	1:E:263:ILE:HG13	1.93	0.49
1:Q:138:PRO:HD2	1:Q:139:MET:SD	2.52	0.49
1:Q:95:GLU:HG3	1:Q:98:ALA:HB2	1.93	0.49
1:B:137:TYR:HD2	1:B:276:ILE:HB	1.77	0.49
1:R:35:LEU:HB3	1:R:54:THR:HB	1.95	0.49
3:N:37:TRP:CE2	3:N:75:LEU:HB2	2.46	0.49
2:Y:60:TYR:HE1	2:Y:70:LEU:HG	1.78	0.49
1:R:284:MET:HG2	5:R:402:POV:H1	1.95	0.49
1:D:217:LEU:O	1:D:221:TYR:HB2	2.12	0.49
1:S:134:LEU:N	1:S:272:TYR:OH	2.42	0.49
1:E:220:LEU:HD11	1:E:280:ILE:HD11	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:154:THR:HG23	1:Q:156:LYS:HG2	1.93	0.49
1:P:271:SER:HB2	1:Q:180:SER:HA	1.95	0.49
1:A:154:THR:HG23	1:A:156:LYS:HG2	1.95	0.49
2:I:197:TRP:CZ3	2:I:202:VAL:HG22	2.48	0.49
1:T:141:VAL:HG11	1:T:208:GLN:HE21	1.78	0.49
3:M:37:TRP:CE2	3:M:75:LEU:HB2	2.48	0.49
1:P:241:ALA:HB2	1:T:302:ASN:ND2	2.24	0.48
1:P:98:ALA:HB3	1:Q:103:ILE:HG21	1.95	0.48
1:C:170:LEU:HD11	1:C:186:THR:HG21	1.94	0.48
2:J:27:TYR:CE2	2:J:98:ARG:HD2	2.47	0.48
1:Q:215:PHE:CE1	1:Q:219:GLN:HG3	2.48	0.48
1:P:170:LEU:HD11	1:P:186:THR:HG21	1.95	0.48
1:T:220:LEU:HB2	1:T:263:ILE:HD11	1.95	0.48
1:E:19:ARG:HH11	1:E:157:ASP:HA	1.78	0.48
1:D:126:LEU:HD13	1:D:128:LEU:HD21	1.95	0.48
1:S:111:ARG:HB2	1:S:119:LEU:HB3	1.94	0.48
2:G:14:PRO:HG2	2:G:122:SER:HB3	1.95	0.48
2:Y:207:ALA:HB2	2:Y:214:LYS:HD3	1.95	0.48
1:C:224:SER:OG	1:C:283:CYS:SG	2.72	0.48
1:S:264:ASN:HD21	5:T:401:POV:H14B	1.78	0.48
3:Z:107:THR:HG21	3:Z:144:PRO:HB3	1.94	0.48
2:I:176:PRO:HG2	3:O:165:THR:HB	1.94	0.48
1:S:264:ASN:ND2	5:T:401:POV:H14B	2.28	0.48
2:H:157:GLU:HG3	2:H:184:TYR:CD2	2.49	0.48
1:C:1:SER:HA	1:C:4:LYS:HD2	1.94	0.48
1:R:138:PRO:HD2	1:R:139:MET:SD	2.53	0.48
3:N:170:GLN:NE2	3:N:174:LYS:HB2	2.29	0.48
3:L:148:THR:HB	3:L:199:THR:HB	1.94	0.48
2:U:6:GLN:N	2:U:114:GLN:HE22	2.12	0.48
1:C:111:ARG:HD2	1:C:119:LEU:HD23	1.94	0.48
1:B:134:LEU:N	1:B:272:TYR:OH	2.44	0.48
1:A:110:ILE:HG12	1:A:120:TYR:HD1	1.79	0.48
1:R:271:SER:HB2	1:S:180:SER:HA	1.96	0.48
2:J:152:LYS:HA	2:J:185:THR:HG23	1.96	0.48
3:L:85:GLU:HG3	3:L:107:THR:HA	1.96	0.47
1:T:311:ILE:O	1:T:315:VAL:HG23	2.14	0.47
1:P:220:LEU:O	1:P:224:SER:OG	2.23	0.47
1:C:288:PHE:HE1	1:D:230:VAL:HG11	1.78	0.47
1:C:271:SER:HB2	1:D:180:SER:HA	1.96	0.47
1:D:271:SER:HB2	1:E:180:SER:HA	1.95	0.47
2:W:168:LEU:HD13	2:W:190:VAL:HG21	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:24:ALA:HB1	2:J:27:TYR:HE1	1.77	0.47
1:B:300:ILE:O	1:B:308:TRP:HB3	2.13	0.47
1:D:285:THR:OG1	5:D:401:POV:O22	2.28	0.47
1:S:267:LEU:HG	1:S:268:PRO:O	2.15	0.47
1:P:126:LEU:HD13	1:P:128:LEU:HD21	1.97	0.47
1:A:170:LEU:HD22	1:A:174:LEU:HD23	1.96	0.47
2:G:157:GLU:HB2	2:G:158:PRO:HA	1.96	0.47
1:Q:259:GLN:O	1:Q:263:ILE:HG12	2.15	0.47
1:R:299:HIS:CE1	1:S:237:PHE:HA	2.49	0.47
1:B:220:LEU:O	1:B:224:SER:OG	2.21	0.47
2:H:6:GLN:N	2:H:114:GLN:HE22	2.12	0.47
1:S:137:TYR:HD2	1:S:276:ILE:HB	1.80	0.47
1:Q:223:PRO:HB3	5:Q:401:POV:H3	1.97	0.47
1:Q:215:PHE:CZ	1:Q:219:GLN:HG3	2.49	0.47
1:S:85:ILE:HD11	1:S:112:ILE:HD11	1.97	0.47
1:C:230:VAL:O	1:C:233:VAL:HG22	2.14	0.47
1:E:137:TYR:CE1	1:E:267:LEU:HD13	2.50	0.47
1:E:268:PRO:HA	1:E:269:PRO:HD3	1.77	0.47
2:G:161:VAL:HG22	2:G:206:VAL:HG22	1.96	0.47
2:G:44:ASN:ND2	3:K:102:GLY:HA2	2.30	0.47
2:X:155:PHE:HA	2:X:156:PRO:HA	1.76	0.47
3:M:120:LEU:HD13	3:M:209:LEU:HD22	1.97	0.47
1:T:259:GLN:NE2	5:T:401:POV:O13	2.48	0.47
1:R:230:VAL:O	1:R:233:VAL:HG22	2.14	0.47
1:R:220:LEU:HD22	1:R:263:ILE:HG13	1.96	0.47
1:T:170:LEU:HD11	1:T:186:THR:HG21	1.96	0.47
1:E:184:THR:CG2	1:E:210:LYS:HD2	2.45	0.47
2:F:175:PHE:CZ	3:N:138:THR:HG23	2.50	0.47
1:S:185:ASN:HD22	4:S:401:NAG:H61	1.80	0.47
1:Q:230:VAL:O	1:Q:233:VAL:HG22	2.14	0.47
2:H:133:LEU:HD13	3:L:121:PHE:CG	2.50	0.47
2:G:6:GLN:N	2:G:114:GLN:HE22	2.10	0.46
2:U:163:TRP:HB3	2:U:168:LEU:HD12	1.98	0.46
1:R:215:PHE:CZ	1:R:219:GLN:HG3	2.50	0.46
2:U:157:GLU:HG3	2:U:184:TYR:CD2	2.50	0.46
1:E:154:THR:HG23	1:E:156:LYS:HG2	1.95	0.46
1:R:227:LEU:HD13	5:R:401:POV:H34A	1.96	0.46
1:B:215:PHE:O	1:B:219:GLN:N	2.46	0.46
2:G:208:HIS:CE1	2:G:211:SER:HG	2.30	0.46
1:T:230:VAL:O	1:T:233:VAL:HG22	2.15	0.46
5:Q:401:POV:H13B	5:Q:401:POV:H11	1.64	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:197:TRP:CZ2	2:J:221:PRO:HG3	2.50	0.46
2:H:27:TYR:CE2	2:H:98:ARG:HD2	2.50	0.46
1:E:175:SER:HB2	1:E:182:GLN:OE1	2.16	0.46
2:G:208:HIS:NE2	2:G:210:ALA:HB3	2.31	0.46
1:Q:33:ASN:HB3	1:Q:56:ARG:HB2	1.98	0.46
1:R:297:VAL:HG13	1:R:312:SER:HB2	1.97	0.46
2:F:175:PHE:HZ	3:N:138:THR:HG23	1.81	0.46
2:Y:18:MET:HG3	2:Y:86:LEU:HD11	1.98	0.46
3:K:122:PRO:HG3	3:K:209:LEU:HD11	1.98	0.46
2:Y:208:HIS:CE1	2:Y:210:ALA:HB3	2.51	0.46
1:A:170:LEU:HD11	1:A:186:THR:HG21	1.98	0.46
2:U:18:MET:HG3	2:U:86:LEU:HD11	1.97	0.46
1:D:299:HIS:CE1	1:E:237:PHE:HA	2.51	0.46
3:K:151:TRP:O	3:K:157:PRO:HA	2.15	0.46
2:V:6:GLN:HE22	2:V:96:CYS:H	1.64	0.46
2:Y:174:THR:HG23	2:Y:186:LEU:HD21	1.97	0.46
1:B:299:HIS:CE1	1:C:237:PHE:HA	2.51	0.46
2:I:155:PHE:HA	2:I:156:PRO:HA	1.75	0.46
1:T:137:TYR:HD2	1:T:276:ILE:HB	1.81	0.46
1:C:35:LEU:HD12	1:C:169:GLN:O	2.16	0.46
1:A:35:LEU:HB3	1:A:54:THR:HB	1.98	0.45
2:G:37:VAL:HG11	2:G:45:LEU:HD23	1.98	0.45
1:B:305:THR:CG2	1:B:308:TRP:HD1	2.28	0.45
2:I:108:TYR:CB	3:O:51:GLY:HA3	2.44	0.45
1:R:111:ARG:HB2	1:R:119:LEU:HB3	1.99	0.45
2:H:132:PRO:HG3	2:H:217:LYS:HE2	1.97	0.45
1:R:186:THR:HG22	1:R:207:ILE:HG22	1.98	0.45
1:S:271:SER:HB2	1:T:180:SER:HA	1.98	0.45
1:C:19:ARG:HG2	1:C:20:PRO:HD2	1.98	0.45
2:F:132:PRO:HD3	2:F:217:LYS:HE2	1.97	0.45
3:Z:22:CYS:HB3	3:Z:73:ALA:HB3	1.99	0.45
2:I:6:GLN:H	2:I:114:GLN:NE2	2.12	0.45
2:V:6:GLN:N	2:V:114:GLN:HE22	2.11	0.45
3:K:141:ASP:H	3:K:170:GLN:HE22	1.64	0.45
3:O:107:THR:HG21	3:O:144:PRO:HB3	1.99	0.45
1:S:170:LEU:HD22	1:S:174:LEU:HD23	1.99	0.45
1:D:227:LEU:HG	1:D:252:THR:HG23	1.99	0.45
2:H:127:PRO:HA	2:H:128:PRO:HD3	1.83	0.45
2:F:6:GLN:N	2:F:114:GLN:HE22	2.14	0.45
2:H:50:LEU:HD21	2:H:59:SER:HB3	1.97	0.45
1:B:98:ALA:CB	1:C:103:ILE:HG13	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:130:CYS:O	1:B:132:MET:HG3	2.17	0.45
3:Z:150:ASP:O	3:Z:197:GLN:N	2.46	0.45
1:A:252:THR:HG22	1:A:287:ILE:HD13	1.98	0.45
1:E:183:LEU:HD12	1:E:183:LEU:O	2.17	0.45
1:T:239:ARG:HD2	1:T:301:ALA:HB2	1.98	0.45
3:N:85:GLU:HG3	3:N:107:THR:HA	1.99	0.45
2:J:6:GLN:N	2:J:114:GLN:HE22	2.12	0.45
1:C:137:TYR:HA	1:C:138:PRO:HA	1.74	0.45
1:S:138:PRO:HD2	1:S:139:MET:SD	2.57	0.45
1:B:138:PRO:HG3	1:B:216:TYR:CG	2.51	0.45
1:S:154:THR:HA	1:S:200:TYR:CD1	2.52	0.45
2:V:6:GLN:NE2	2:V:96:CYS:H	2.15	0.45
1:D:137:TYR:HA	1:D:138:PRO:HA	1.80	0.45
2:V:141:THR:OG1	2:V:142:ASN:N	2.49	0.45
2:U:174:THR:HG23	2:U:186:LEU:HD21	1.98	0.45
1:C:239:ARG:HD2	1:C:301:ALA:HB1	1.98	0.45
1:R:215:PHE:CE1	1:R:219:GLN:HG3	2.52	0.45
1:D:20:PRO:HA	1:D:21:PRO:HD3	1.88	0.45
2:F:48:ILE:HA	2:F:64:PHE:HD2	1.81	0.45
1:A:215:PHE:HA	1:E:273:ILE:HD11	1.99	0.45
1:C:254:LEU:HD11	1:D:254:LEU:HD23	1.99	0.45
1:E:259:GLN:O	1:E:263:ILE:HG12	2.17	0.45
1:C:138:PRO:HD2	1:C:139:MET:SD	2.57	0.45
1:A:230:VAL:O	1:A:233:VAL:HG22	2.17	0.45
1:B:137:TYR:HA	1:B:138:PRO:HA	1.69	0.45
1:B:230:VAL:O	1:B:233:VAL:HG22	2.16	0.45
2:G:176:PRO:HG2	3:K:165:THR:HB	1.98	0.45
3:K:122:PRO:HB3	3:K:209:LEU:HD21	1.98	0.44
2:I:46:GLU:OE1	2:I:63:LYS:HE2	2.16	0.44
1:C:313:LYS:HD3	1:C:313:LYS:HA	1.80	0.44
3:N:22:CYS:HB3	3:N:73:ALA:HB3	1.99	0.44
2:F:30:THR:HA	2:F:53:PRO:HB2	1.99	0.44
2:X:131:TYR:HA	2:X:132:PRO:HD3	1.83	0.44
2:J:18:MET:HE2	2:J:18:MET:HB2	1.78	0.44
3:O:37:TRP:CE2	3:O:75:LEU:HB2	2.52	0.44
1:P:239:ARG:HH11	1:P:301:ALA:HB1	1.81	0.44
1:E:267:LEU:HD12	1:E:268:PRO:HD2	1.99	0.44
1:S:259:GLN:O	1:S:263:ILE:HG12	2.18	0.44
2:U:199:SER:OG	2:U:200:GLU:OE1	2.28	0.44
2:V:163:TRP:HB2	2:V:168:LEU:HD11	1.99	0.44
1:S:299:HIS:CE1	1:T:237:PHE:HA	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:20:ILE:HD13	2:J:81:MET:HE3	1.99	0.44
1:P:137:TYR:HA	1:P:138:PRO:HA	1.78	0.44
2:H:146:THR:HG22	2:H:191:THR:HG22	2.00	0.44
1:R:267:LEU:HG	1:R:268:PRO:O	2.17	0.44
3:K:109:LEU:HD21	3:K:144:PRO:HG3	2.00	0.44
1:Q:224:SER:HB2	1:Q:279:TRP:CZ3	2.53	0.44
1:A:107:ASN:OD1	1:A:123:ARG:NH1	2.50	0.44
2:Y:6:GLN:N	2:Y:114:GLN:HE22	2.12	0.44
3:N:172:ASN:OD1	3:N:172:ASN:N	2.48	0.44
2:I:127:PRO:HA	2:I:128:PRO:HD3	1.83	0.44
2:J:60:TYR:HE1	2:J:70:LEU:HG	1.83	0.44
1:R:252:THR:HG22	1:R:287:ILE:HD13	2.00	0.44
1:A:33:ASN:HB3	1:A:56:ARG:HB2	1.99	0.44
1:T:192:THR:O	2:Y:105:TYR:N	2.49	0.44
1:R:305:THR:HG22	1:R:307:GLU:OE1	2.18	0.44
5:P:403:POV:H13B	5:P:403:POV:H11A	1.77	0.44
1:E:49:TYR:CE1	1:E:144:CYS:HB3	2.52	0.44
2:J:208:HIS:CE1	2:J:210:ALA:HB3	2.53	0.44
1:S:33:ASN:HB3	1:S:56:ARG:HB2	2.00	0.44
1:Q:311:ILE:O	1:Q:315:VAL:HG23	2.18	0.44
2:X:208:HIS:CE1	2:X:210:ALA:HB3	2.53	0.44
2:G:6:GLN:H	2:G:114:GLN:NE2	2.13	0.44
1:D:194:VAL:HG13	2:I:52:ASN:HD22	1.82	0.44
1:C:194:VAL:HG13	2:G:52:ASN:HD22	1.82	0.44
1:D:175:SER:HB2	1:D:182:GLN:OE1	2.18	0.44
1:A:310:ASP:O	1:A:314:ARG:HG3	2.18	0.44
2:G:153:GLY:HA2	2:G:183:LEU:HD22	2.00	0.44
1:Q:111:ARG:HB2	1:Q:119:LEU:HB3	1.99	0.44
3:L:37:TRP:CZ3	3:L:90:CYS:HB3	2.53	0.44
2:W:157:GLU:HG3	2:W:184:TYR:CD2	2.53	0.44
1:P:68:VAL:N	1:P:71:ASP:OD2	2.49	0.44
2:W:60:TYR:CE1	2:W:70:LEU:HG	2.53	0.44
1:R:96:LYS:HD2	1:R:129:SER:HB3	2.00	0.44
2:Y:144:MET:HG2	2:Y:193:PRO:HA	2.00	0.44
1:S:309:ASN:OD1	1:S:311:ILE:HG12	2.18	0.43
1:C:19:ARG:HH11	1:C:157:ASP:HA	1.82	0.43
3:O:37:TRP:CZ3	3:O:90:CYS:HB3	2.53	0.43
3:M:152:LYS:HB2	3:M:195:SER:HB3	1.99	0.43
1:A:299:HIS:CE1	1:B:237:PHE:HA	2.53	0.43
1:R:311:ILE:O	1:R:315:VAL:HG23	2.18	0.43
2:J:177:ALA:HB2	2:J:186:LEU:HD23	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:W:6:GLN:N	2:W:114:GLN:HE22	2.11	0.43
1:T:137:TYR:O	1:T:275:ALA:HB3	2.18	0.43
2:W:208:HIS:CE1	2:W:210:ALA:HB3	2.53	0.43
1:E:230:VAL:O	1:E:233:VAL:HG22	2.18	0.43
3:M:41:LYS:HE2	3:M:83:GLU:O	2.18	0.43
2:H:153:GLY:HA2	2:H:183:LEU:HB3	1.99	0.43
1:S:100:LYS:HE2	1:T:104:ASP:H	1.83	0.43
1:A:300:ILE:HG13	1:A:312:SER:OG	2.18	0.43
1:R:220:LEU:HD23	1:R:276:ILE:HD11	2.01	0.43
1:D:252:THR:HG22	1:D:287:ILE:HD13	2.00	0.43
1:P:230:VAL:O	1:P:233:VAL:HG22	2.18	0.43
1:S:224:SER:HB2	1:S:279:TRP:CZ3	2.53	0.43
1:P:259:GLN:HE22	5:P:403:POV:P	2.42	0.43
2:F:108:TYR:HB3	3:N:52:GLY:H	1.84	0.43
1:S:170:LEU:HD11	1:S:186:THR:HG21	2.01	0.43
2:F:48:ILE:HG12	2:F:64:PHE:CE2	2.54	0.43
1:P:305:THR:HB	1:P:308:TRP:HD1	1.83	0.43
1:S:183:LEU:HD11	1:S:207:ILE:HB	2.01	0.43
1:R:249:GLY:HA3	1:R:291:LEU:HD13	1.99	0.43
3:N:199:THR:HA	3:N:204:THR:HA	2.01	0.43
8:T:403:LMT:H5B	8:T:403:LMT:H6D	2.01	0.43
3:K:188:TRP:CG	3:K:189:GLU:N	2.87	0.43
2:I:161:VAL:HG22	2:I:206:VAL:HG22	1.99	0.43
1:C:311:ILE:O	1:C:315:VAL:HG23	2.19	0.43
3:L:41:LYS:HB2	3:L:45:LEU:HB2	1.99	0.43
2:U:27:TYR:CE2	2:U:98:ARG:HD2	2.53	0.43
1:P:137:TYR:O	1:P:275:ALA:HB3	2.18	0.43
1:A:19:ARG:HH11	1:A:157:ASP:HA	1.83	0.43
2:H:164:ASN:HB3	2:H:167:SER:HB3	1.99	0.43
1:Q:267:LEU:HA	1:Q:268:PRO:HD2	1.87	0.43
1:S:222:ILE:HB	1:S:223:PRO:HD3	2.00	0.43
2:F:108:TYR:CB	3:N:51:GLY:HA2	2.49	0.43
1:A:220:LEU:HD23	1:A:276:ILE:HD11	2.01	0.43
1:Q:183:LEU:O	1:Q:183:LEU:HD12	2.19	0.43
3:K:85:GLU:HG3	3:K:107:THR:HA	1.99	0.43
1:S:220:LEU:C	1:S:223:PRO:HD2	2.38	0.43
1:R:137:TYR:HA	1:R:138:PRO:HA	1.79	0.43
1:P:138:PRO:HD2	1:P:139:MET:SD	2.59	0.43
1:S:49:TYR:CE1	1:S:144:CYS:HB3	2.54	0.43
1:B:267:LEU:HD12	1:B:268:PRO:HD2	2.00	0.43
1:R:20:PRO:HA	1:R:21:PRO:HD3	1.86	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:U:135:PRO:HB2	2:U:140:GLN:HE22	1.84	0.43
2:X:18:MET:HG3	2:X:86:LEU:HD11	2.00	0.43
3:L:139:ILE:HD12	3:L:198:VAL:HG21	2.00	0.43
2:F:133:LEU:HD13	3:N:121:PHE:CG	2.53	0.43
1:B:111:ARG:HD2	1:B:119:LEU:HD23	2.00	0.43
1:Q:134:LEU:HD12	1:Q:270:VAL:HG21	2.00	0.42
1:D:267:LEU:HA	1:D:268:PRO:HD2	1.92	0.42
1:D:138:PRO:HD2	1:D:139:MET:SD	2.59	0.42
1:B:215:PHE:CZ	1:B:219:GLN:HG2	2.54	0.42
2:G:144:MET:HB2	2:G:191:THR:HG22	2.01	0.42
2:U:208:HIS:CE1	2:U:210:ALA:HB3	2.54	0.42
1:B:20:PRO:HA	1:B:21:PRO:HD3	1.86	0.42
1:R:300:ILE:HB	1:R:308:TRP:HE3	1.84	0.42
3:O:45:LEU:HD13	3:O:45:LEU:HA	1.87	0.42
3:M:122:PRO:HG3	3:M:209:LEU:HD11	2.00	0.42
1:B:267:LEU:HA	1:B:268:PRO:HD2	1.90	0.42
1:B:10:PHE:HE2	1:B:83:HIS:HB3	1.80	0.42
1:R:33:ASN:HB3	1:R:56:ARG:HB2	2.00	0.42
2:F:127:PRO:HA	2:F:128:PRO:HD3	1.85	0.42
1:A:136:TYR:O	1:A:140:ASP:HB3	2.19	0.42
1:A:138:PRO:HD2	1:A:139:MET:SD	2.59	0.42
1:R:19:ARG:HH11	1:R:157:ASP:HA	1.84	0.42
1:C:299:HIS:CE1	1:D:237:PHE:HA	2.54	0.42
1:B:170:LEU:HD11	1:B:186:THR:HG21	2.01	0.42
1:P:311:ILE:O	1:P:315:VAL:HG23	2.19	0.42
1:D:313:LYS:HA	1:D:313:LYS:HD3	1.88	0.42
1:P:20:PRO:HA	1:P:21:PRO:HD3	1.88	0.42
5:R:402:POV:H28A	5:R:402:POV:H211	1.73	0.42
1:B:227:LEU:HG	1:B:252:THR:HG23	2.01	0.42
1:R:189:THR:OG1	1:R:190:TYR:N	2.52	0.42
1:D:224:SER:HB2	1:D:279:TRP:CZ3	2.55	0.42
1:C:114:ASN:N	1:C:114:ASN:OD1	2.53	0.42
1:C:130:CYS:O	1:C:132:MET:HG3	2.20	0.42
1:B:220:LEU:HD23	1:B:276:ILE:HD11	2.01	0.42
1:C:267:LEU:HD12	1:C:268:PRO:HD2	2.02	0.42
2:X:33:THR:HA	2:X:53:PRO:HD3	2.00	0.42
3:K:141:ASP:H	3:K:170:GLN:NE2	2.18	0.42
1:C:217:LEU:HA	1:C:221:TYR:HB2	2.01	0.42
1:T:224:SER:HB2	1:T:279:TRP:CZ3	2.55	0.42
1:S:300:ILE:HD11	1:S:312:SER:HA	2.02	0.42
3:N:118:VAL:O	3:N:207:LYS:HE3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:168:LEU:HD23	1:D:186:THR:HB	2.01	0.42
1:D:227:LEU:HD12	1:D:227:LEU:HA	1.88	0.42
1:A:76:PHE:CD2	1:A:111:ARG:HD3	2.55	0.42
1:T:138:PRO:HD2	1:T:139:MET:SD	2.59	0.42
2:I:133:LEU:HD13	3:O:121:PHE:CG	2.55	0.42
1:B:273:ILE:HD11	1:C:215:PHE:HA	2.02	0.42
1:D:76:PHE:CD2	1:D:111:ARG:HD3	2.55	0.42
1:T:261:ALA:O	1:T:264:ASN:HB2	2.20	0.42
1:C:8:HIS:O	1:C:11:THR:HG22	2.19	0.42
1:P:252:THR:HG22	1:P:287:ILE:HD13	2.01	0.42
1:S:114:ASN:N	1:S:114:ASN:OD1	2.53	0.42
1:E:267:LEU:HA	1:E:268:PRO:HD2	1.81	0.42
2:J:18:MET:HG3	2:J:86:LEU:HD11	2.01	0.42
1:D:215:PHE:CE1	1:D:219:GLN:HG3	2.55	0.42
2:I:61:ASN:HD21	3:O:97:HIS:HE1	1.66	0.42
1:S:194:VAL:HG13	2:X:52:ASN:HD22	1.85	0.42
2:F:48:ILE:HA	2:F:64:PHE:CD2	2.55	0.42
1:C:221:TYR:HA	1:C:279:TRP:CZ3	2.55	0.42
1:S:75:ASP:O	1:S:114:ASN:N	2.53	0.42
2:X:170:SER:OG	2:X:171:GLY:N	2.52	0.42
1:P:111:ARG:HD2	1:P:119:LEU:HD23	2.01	0.42
2:F:178:VAL:HG12	2:F:180:GLN:H	1.85	0.42
3:K:129:GLU:O	3:K:130:THR:OG1	2.34	0.42
2:V:155:PHE:HA	2:V:156:PRO:HA	1.76	0.42
2:W:18:MET:HE1	2:W:20:ILE:HG12	2.01	0.42
2:F:144:MET:HG2	2:F:193:PRO:HA	2.02	0.41
3:L:94:TYR:O	3:L:95:SER:OG	2.29	0.41
2:J:53:PRO:O	2:J:74:LYS:HE2	2.20	0.41
1:C:137:TYR:O	1:C:275:ALA:HB3	2.20	0.41
2:W:18:MET:HG3	2:W:86:LEU:HD11	2.01	0.41
1:B:102:THR:HA	1:B:106:PRO:HA	2.02	0.41
1:C:134:LEU:HD12	1:C:270:VAL:HG21	2.01	0.41
3:N:152:LYS:HD3	3:N:197:GLN:OE1	2.20	0.41
1:D:138:PRO:HB2	1:D:211:ARG:HD2	2.02	0.41
2:V:157:GLU:HG3	2:V:184:TYR:CD2	2.56	0.41
1:Q:130:CYS:HA	1:Q:131:PRO:HD2	1.95	0.41
1:B:189:THR:OG1	1:B:190:TYR:N	2.53	0.41
1:R:130:CYS:O	1:R:132:MET:HG3	2.19	0.41
2:U:100:GLY:HA3	2:U:108:TYR:CZ	2.55	0.41
1:E:19:ARG:HG2	1:E:20:PRO:HD2	2.02	0.41
1:C:111:ARG:HB2	1:C:119:LEU:HB3	2.00	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:T:307:GLU:O	1:T:311:ILE:HG12	2.19	0.41
2:F:155:PHE:HA	2:F:156:PRO:HA	1.76	0.41
3:K:94:TYR:O	3:K:96:ASN:N	2.53	0.41
2:H:208:HIS:NE2	2:H:210:ALA:HB3	2.34	0.41
1:S:154:THR:HA	1:S:200:TYR:HD1	1.85	0.41
1:Q:76:PHE:CD2	1:Q:111:ARG:HD3	2.56	0.41
1:C:273:ILE:HG12	1:D:215:PHE:HD1	1.85	0.41
3:Z:181:LEU:HG	3:Z:183:LEU:HD13	2.01	0.41
1:P:279:TRP:HB2	1:P:334:TYR:CE1	2.56	0.41
1:P:288:PHE:CG	5:Q:401:POV:H34	2.56	0.41
2:G:18:MET:HG3	2:G:86:LEU:HD11	2.01	0.41
1:A:237:PHE:HA	1:E:299:HIS:CE1	2.55	0.41
2:U:197:TRP:HA	2:U:198:PRO:HA	1.86	0.41
1:R:262:GLY:HA2	5:R:401:POV:H14B	2.03	0.41
1:B:19:ARG:HG2	1:B:20:PRO:HD2	2.02	0.41
1:A:137:TYR:CE2	1:A:138:PRO:HB3	2.56	0.41
3:Z:166:GLN:HA	3:Z:167:PRO:HD2	1.91	0.41
3:K:121:PHE:HA	3:K:122:PRO:HD3	1.80	0.41
2:V:166:GLY:C	2:V:168:LEU:H	2.24	0.41
3:Z:115:SER:HA	3:Z:116:PRO:HD3	1.92	0.41
3:L:187:ALA:O	3:L:191:HIS:ND1	2.51	0.41
1:A:271:SER:HB2	1:B:180:SER:HA	2.03	0.41
2:Y:13:ARG:HD2	2:Y:122:SER:HA	2.03	0.41
2:H:6:GLN:H	2:H:114:GLN:NE2	2.17	0.41
2:Y:27:TYR:HD2	2:Y:32:TYR:CD1	2.39	0.41
1:E:20:PRO:HA	1:E:21:PRO:HD3	1.85	0.41
2:J:177:ALA:HA	2:J:186:LEU:HB3	2.02	0.41
1:B:20:PRO:HD3	1:B:86:TRP:CD2	2.56	0.41
1:B:19:ARG:HH11	1:B:157:ASP:HA	1.86	0.41
2:Y:169:SER:C	2:Y:171:GLY:H	2.24	0.41
2:Y:153:GLY:HA2	2:Y:183:LEU:HB3	2.03	0.41
1:P:130:CYS:O	1:P:132:MET:HG3	2.21	0.41
1:E:130:CYS:O	1:E:132:MET:HG3	2.20	0.41
1:E:170:LEU:HD22	1:E:174:LEU:HD23	2.01	0.41
2:W:121:SER:OG	2:W:182:ASP:OD1	2.39	0.41
1:T:130:CYS:O	1:T:132:MET:HG3	2.21	0.41
1:E:279:TRP:HB2	1:E:334:TYR:CZ	2.56	0.41
3:L:128:LEU:HD21	3:L:188:TRP:CZ3	2.55	0.41
1:E:35:LEU:HB3	1:E:54:THR:HB	2.03	0.41
3:Z:188:TRP:HA	3:Z:194:TYR:OH	2.20	0.41
3:M:40:GLU:O	3:M:86:ALA:HB1	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:227:LEU:HD12	1:E:227:LEU:HA	1.89	0.41
1:R:49:TYR:CE1	1:R:144:CYS:HB3	2.56	0.41
2:J:218:LYS:HG2	2:J:219:ILE:H	1.86	0.41
1:T:220:LEU:HD23	1:T:276:ILE:HD11	2.02	0.41
2:U:18:MET:HE1	2:U:20:ILE:HG12	2.02	0.41
2:Y:131:TYR:HA	2:Y:132:PRO:HD3	1.87	0.41
1:A:35:LEU:HD12	1:A:169:GLN:O	2.21	0.40
1:A:137:TYR:HA	1:A:138:PRO:HA	1.79	0.40
1:Q:96:LYS:HD2	1:Q:129:SER:HB3	2.03	0.40
1:B:43:ASP:OD1	1:B:45:VAL:N	2.51	0.40
2:I:33:THR:HA	2:I:53:PRO:HD3	2.03	0.40
1:Q:107:ASN:OD1	1:Q:123:ARG:NH1	2.54	0.40
1:E:110:ILE:HG12	1:E:120:TYR:HD1	1.85	0.40
1:T:35:LEU:HD12	1:T:169:GLN:O	2.22	0.40
2:X:6:GLN:N	2:X:114:GLN:HE22	2.20	0.40
2:H:155:PHE:HA	2:H:156:PRO:HA	1.80	0.40
1:P:19:ARG:HA	1:P:20:PRO:HD3	1.97	0.40
2:V:179:LEU:HD12	2:V:184:TYR:HE1	1.87	0.40
3:L:188:TRP:HA	3:L:194:TYR:OH	2.21	0.40
1:C:49:TYR:CE1	1:C:144:CYS:HB3	2.56	0.40
2:W:147:LEU:HD22	2:W:219:ILE:HG21	2.03	0.40
1:S:5:ILE:O	1:S:8:HIS:HB3	2.20	0.40
3:N:53:ILE:CG2	3:N:54:ASN:N	2.84	0.40
1:S:302:ASN:ND2	1:T:241:ALA:HA	2.36	0.40
1:T:183:LEU:O	1:T:183:LEU:HD12	2.22	0.40
1:E:134:LEU:N	1:E:272:TYR:OH	2.41	0.40
1:C:220:LEU:HD23	1:C:276:ILE:HD11	2.02	0.40
1:T:300:ILE:HG23	1:T:308:TRP:CE3	2.49	0.40
1:S:19:ARG:HG2	1:S:20:PRO:HD2	2.04	0.40
1:A:19:ARG:HG2	1:A:20:PRO:HD2	2.04	0.40
2:H:18:MET:HG3	2:H:86:LEU:HD11	2.04	0.40
2:X:126:THR:HA	2:X:127:PRO:HD3	1.96	0.40
2:F:152:LYS:HA	2:F:185:THR:HG23	2.04	0.40
1:S:137:TYR:OH	1:S:267:LEU:HB2	2.21	0.40
1:A:267:LEU:HD22	1:A:276:ILE:HG21	2.04	0.40
1:S:19:ARG:NH1	1:S:157:ASP:HA	2.37	0.40
2:H:134:ALA:HA	2:H:135:PRO:HD3	1.97	0.40
3:K:148:THR:HB	3:K:199:THR:HB	2.03	0.40
2:I:48:ILE:HA	2:I:64:PHE:CD2	2.57	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	340/347 (98%)	327 (96%)	13 (4%)	0	100	100
1	B	339/347 (98%)	325 (96%)	14 (4%)	0	100	100
1	C	338/347 (97%)	324 (96%)	14 (4%)	0	100	100
1	D	338/347 (97%)	328 (97%)	10 (3%)	0	100	100
1	E	338/347 (97%)	325 (96%)	13 (4%)	0	100	100
1	P	338/347 (97%)	326 (96%)	12 (4%)	0	100	100
1	Q	338/347 (97%)	326 (96%)	12 (4%)	0	100	100
1	R	338/347 (97%)	326 (96%)	12 (4%)	0	100	100
1	S	338/347 (97%)	325 (96%)	12 (4%)	1 (0%)	46	85
1	T	338/347 (97%)	326 (96%)	12 (4%)	0	100	100
2	F	219/224 (98%)	206 (94%)	13 (6%)	0	100	100
2	G	220/224 (98%)	203 (92%)	17 (8%)	0	100	100
2	H	221/224 (99%)	206 (93%)	15 (7%)	0	100	100
2	I	220/224 (98%)	206 (94%)	14 (6%)	0	100	100
2	J	220/224 (98%)	203 (92%)	16 (7%)	1 (0%)	34	78
2	U	219/224 (98%)	205 (94%)	14 (6%)	0	100	100
2	V	222/224 (99%)	203 (91%)	19 (9%)	0	100	100
2	W	222/224 (99%)	209 (94%)	13 (6%)	0	100	100
2	X	222/224 (99%)	206 (93%)	16 (7%)	0	100	100
2	Y	219/224 (98%)	202 (92%)	17 (8%)	0	100	100
3	K	208/215 (97%)	193 (93%)	14 (7%)	1 (0%)	34	78
3	L	209/215 (97%)	198 (95%)	11 (5%)	0	100	100
3	M	208/215 (97%)	193 (93%)	15 (7%)	0	100	100
3	N	209/215 (97%)	194 (93%)	14 (7%)	1 (0%)	34	78
3	O	209/215 (97%)	194 (93%)	15 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
3	Z	213/215 (99%)	193 (91%)	19 (9%)	1 (0%)	34 78
3	f	213/215 (99%)	195 (92%)	16 (8%)	2 (1%)	21 67
3	g	209/215 (97%)	196 (94%)	11 (5%)	2 (1%)	19 65
3	h	209/215 (97%)	194 (93%)	15 (7%)	0	100 100
3	i	212/215 (99%)	199 (94%)	12 (6%)	1 (0%)	34 78
All	All	7686/7860 (98%)	7256 (94%)	420 (6%)	10 (0%)	56 91

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	K	95	SER
3	N	95	SER
3	i	95	SER
3	f	95	SER
1	S	309	ASN
3	f	96	ASN
2	J	221	PRO
3	g	96	ASN
3	g	52	GLY
3	Z	52	GLY

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	311/316 (98%)	310 (100%)	1 (0%)	94 98
1	B	310/316 (98%)	310 (100%)	0	100 100
1	C	309/316 (98%)	309 (100%)	0	100 100
1	D	309/316 (98%)	307 (99%)	2 (1%)	90 97
1	E	309/316 (98%)	309 (100%)	0	100 100
1	P	309/316 (98%)	308 (100%)	1 (0%)	94 98
1	Q	309/316 (98%)	308 (100%)	1 (0%)	94 98

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	R	309/316 (98%)	309 (100%)	0	100 100
1	S	309/316 (98%)	309 (100%)	0	100 100
1	T	309/316 (98%)	309 (100%)	0	100 100
2	F	190/193 (98%)	189 (100%)	1 (0%)	92 97
2	G	191/193 (99%)	190 (100%)	1 (0%)	92 97
2	H	192/193 (100%)	191 (100%)	1 (0%)	92 97
2	I	191/193 (99%)	190 (100%)	1 (0%)	92 97
2	J	191/193 (99%)	190 (100%)	1 (0%)	92 97
2	U	190/193 (98%)	188 (99%)	2 (1%)	80 94
2	V	193/193 (100%)	192 (100%)	1 (0%)	92 97
2	W	193/193 (100%)	193 (100%)	0	100 100
2	X	193/193 (100%)	192 (100%)	1 (0%)	92 97
2	Y	190/193 (98%)	188 (99%)	2 (1%)	80 94
3	K	178/182 (98%)	178 (100%)	0	100 100
3	L	179/182 (98%)	179 (100%)	0	100 100
3	M	178/182 (98%)	177 (99%)	1 (1%)	90 97
3	N	179/182 (98%)	179 (100%)	0	100 100
3	O	179/182 (98%)	177 (99%)	2 (1%)	80 94
3	Z	182/182 (100%)	181 (100%)	1 (0%)	92 97
3	f	182/182 (100%)	181 (100%)	1 (0%)	92 97
3	g	179/182 (98%)	179 (100%)	0	100 100
3	h	179/182 (98%)	179 (100%)	0	100 100
3	i	181/182 (100%)	179 (99%)	2 (1%)	80 94
All	All	6803/6910 (98%)	6780 (100%)	23 (0%)	94 98

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	267	LEU
1	D	154	THR
1	D	264	ASN
2	F	27	TYR
2	G	220	VAL
2	H	27	TYR

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Mol	Chain	Res	Type
2	I	27	TYR
3	O	96	ASN
3	O	97	HIS
1	P	105	LYS
1	Q	264	ASN
2	U	27	TYR
2	U	196	THR
2	V	149	CYS
2	X	45	LEU
2	Y	27	TYR
2	Y	45	LEU
3	Z	56	ARG
3	f	96	ASN
3	i	97	HIS
3	i	214	CYS
2	J	27	TYR
3	M	130	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	ASN
1	A	219	GLN
1	E	259	GLN
2	H	140	GLN
3	K	170	GLN
3	L	200	HIS
3	N	170	GLN
3	O	97	HIS
1	P	73	GLN
1	Q	299	HIS
1	S	264	ASN
3	Z	97	HIS
3	f	39	GLN
3	h	39	GLN
3	h	170	GLN

5.3.3 RNA

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 26 ligands modelled in this entry, 2 are monoatomic - leaving 24 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	NAG	A	401	1	14,14,15	0.27	0	15,19,21	1.14	1 (6%)
5	POV	A	402	-	22,22,51	1.06	1 (4%)	25,29,59	1.21	2 (8%)
7	GOL	B	401	-	5,5,5	0.34	0	5,5,5	0.22	0
4	NAG	B	402	1	14,14,15	0.24	0	15,19,21	0.69	0
8	LMT	B	403	-	24,24,36	1.08	2 (8%)	35,35,47	1.51	6 (17%)
4	NAG	C	401	1	14,14,15	0.38	0	15,19,21	0.81	1 (6%)
8	LMT	C	402	-	24,24,36	1.17	2 (8%)	35,35,47	1.14	2 (5%)
5	POV	D	401	-	25,25,51	1.31	2 (8%)	29,33,59	1.35	3 (10%)
4	NAG	D	402	1	14,14,15	0.33	0	15,19,21	1.47	1 (6%)
8	LMT	D	403	-	24,24,36	1.14	2 (8%)	35,35,47	1.21	4 (11%)
4	NAG	E	401	1	14,14,15	0.25	0	15,19,21	0.65	0
8	LMT	E	402	-	24,24,36	1.11	2 (8%)	35,35,47	1.15	3 (8%)
4	NAG	P	401	1	14,14,15	0.32	0	15,19,21	0.63	0
8	LMT	P	402	-	24,24,36	1.10	2 (8%)	35,35,47	1.23	4 (11%)
5	POV	P	403	-	41,41,51	1.03	2 (4%)	45,49,59	1.13	4 (8%)
5	POV	Q	401	-	45,45,51	0.99	2 (4%)	49,53,59	1.06	3 (6%)
4	NAG	Q	402	1	14,14,15	0.25	0	15,19,21	0.62	0
5	POV	R	401	-	33,33,51	1.13	2 (6%)	37,41,59	1.15	3 (8%)
5	POV	R	402	-	34,34,51	1.09	2 (5%)	38,39,59	1.24	3 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
4	NAG	S	401	1	14,14,15	0.33	0	15,19,21	1.40	1 (6%)
8	LMT	S	402	-	24,24,36	1.12	2 (8%)	35,35,47	1.24	5 (14%)
5	POV	T	401	-	42,42,51	1.01	2 (4%)	46,50,59	1.15	4 (8%)
4	NAG	T	402	1	14,14,15	0.35	0	15,19,21	0.84	1 (6%)
8	LMT	T	403	-	24,24,36	1.11	2 (8%)	35,35,47	1.29	6 (17%)

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. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	401	1	-	0/6/23/26	0/1/1/1
5	POV	A	402	-	-	0/25/25/55	0/0/0/0
7	GOL	B	401	-	-	0/4/4/4	0/0/0/0
4	NAG	B	402	1	-	0/6/23/26	0/1/1/1
8	LMT	B	403	-	-	0/8/48/61	0/2/2/2
4	NAG	C	401	1	-	0/6/23/26	0/1/1/1
8	LMT	C	402	-	-	0/8/48/61	0/2/2/2
5	POV	D	401	-	-	0/29/29/55	0/0/0/0
4	NAG	D	402	1	-	0/6/23/26	0/1/1/1
8	LMT	D	403	-	-	0/8/48/61	0/2/2/2
4	NAG	E	401	1	-	0/6/23/26	0/1/1/1
8	LMT	E	402	-	-	0/8/48/61	0/2/2/2
4	NAG	P	401	1	-	0/6/23/26	0/1/1/1
8	LMT	P	402	-	-	0/8/48/61	0/2/2/2
5	POV	P	403	-	-	0/45/45/55	0/0/0/0
5	POV	Q	401	-	-	0/49/49/55	0/0/0/0
4	NAG	Q	402	1	-	0/6/23/26	0/1/1/1
5	POV	R	401	-	-	1/37/37/55	0/0/0/0
5	POV	R	402	-	-	0/36/36/55	0/0/0/0
4	NAG	S	401	1	-	0/6/23/26	0/1/1/1
8	LMT	S	402	-	-	0/8/48/61	0/2/2/2
5	POV	T	401	-	-	0/46/46/55	0/0/0/0
4	NAG	T	402	1	-	1/6/23/26	0/1/1/1
8	LMT	T	403	-	-	0/8/48/61	0/2/2/2

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	403	LMT	O5'-C1'	2.55	1.47	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	E	402	LMT	O5B-C1B	2.57	1.48	1.41
8	T	403	LMT	O5B-C1B	2.65	1.48	1.41
8	E	402	LMT	O5'-C1'	2.69	1.48	1.43
8	S	402	LMT	O5B-C1B	2.70	1.48	1.41
8	D	403	LMT	O5'-C1'	2.75	1.48	1.43
8	P	402	LMT	O5'-C1'	2.82	1.48	1.43
8	P	402	LMT	O5B-C1B	2.83	1.49	1.41
8	B	403	LMT	O5B-C1B	2.86	1.49	1.41
8	C	402	LMT	O5'-C1'	2.93	1.48	1.43
8	T	403	LMT	O5'-C1'	2.93	1.48	1.43
8	D	403	LMT	O5B-C1B	2.96	1.49	1.41
8	S	402	LMT	O5'-C1'	2.96	1.48	1.43
8	C	402	LMT	O5B-C1B	3.00	1.49	1.41
5	D	401	POV	O21-C21	3.72	1.45	1.34
5	R	401	POV	O21-C21	3.80	1.45	1.34
5	P	403	POV	O21-C21	3.83	1.45	1.34
5	T	401	POV	O21-C21	3.83	1.45	1.34
5	Q	401	POV	O21-C21	3.84	1.45	1.34
5	A	402	POV	O21-C21	3.89	1.45	1.34
5	R	402	POV	O21-C21	3.91	1.46	1.34
5	R	402	POV	O31-C31	4.07	1.45	1.33
5	T	401	POV	O31-C31	4.10	1.45	1.33
5	R	401	POV	O31-C31	4.13	1.45	1.33
5	Q	401	POV	O31-C31	4.20	1.46	1.33
5	P	403	POV	O31-C31	4.21	1.46	1.33
5	D	401	POV	O31-C31	4.34	1.46	1.33

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	D	402	NAG	C1-O5-C5	-4.69	106.29	112.25
4	A	401	NAG	C1-O5-C5	-3.82	107.41	112.25
5	D	401	POV	C2-O21-C21	-3.40	109.72	117.89
5	R	401	POV	C2-O21-C21	-3.37	109.79	117.89
8	S	402	LMT	C1B-O1B-C4'	-3.28	109.45	118.01
8	P	402	LMT	C1B-O1B-C4'	-3.25	109.52	118.01
5	P	403	POV	C2-O21-C21	-2.97	110.75	117.89
5	T	401	POV	C2-O21-C21	-2.88	110.99	117.89
8	T	403	LMT	C1B-O1B-C4'	-2.79	110.71	118.01
5	R	402	POV	C2-O21-C21	-2.61	111.62	117.89
5	A	402	POV	C2-O21-C21	-2.56	111.75	117.89
8	D	403	LMT	C1B-O1B-C4'	-2.55	111.35	118.01

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	Q	401	POV	C2-O21-C21	-2.36	112.24	117.89
5	T	401	POV	C11-C12-N	-2.34	108.50	116.03
5	P	403	POV	C11-C12-N	-2.33	108.53	116.03
8	B	403	LMT	O3'-C3'-C2'	-2.18	105.42	110.34
8	C	402	LMT	C1B-O1B-C4'	-2.17	112.33	118.01
8	E	402	LMT	C1B-O1B-C4'	-2.06	112.62	118.01
8	S	402	LMT	C3B-C4B-C5B	2.01	113.70	110.20
8	E	402	LMT	C3'-C4'-C5'	2.06	115.50	110.84
8	D	403	LMT	C1B-C2B-C3B	2.10	114.11	109.97
8	S	402	LMT	C1B-C2B-C3B	2.13	114.17	109.97
8	S	402	LMT	O5'-C1'-C2'	2.14	113.21	109.80
8	B	403	LMT	C1B-C2B-C3B	2.15	114.21	109.97
4	T	402	NAG	C1-O5-C5	2.20	115.04	112.25
8	S	402	LMT	C4B-C3B-C2B	2.20	114.90	110.79
8	B	403	LMT	C3B-C4B-C5B	2.23	114.09	110.20
8	T	403	LMT	O5B-C5B-C4B	2.25	113.90	109.68
8	C	402	LMT	C1B-C2B-C3B	2.29	114.49	109.97
8	D	403	LMT	O5'-C1'-C2'	2.31	113.48	109.80
4	C	401	NAG	C1-O5-C5	2.32	115.19	112.25
8	T	403	LMT	O5'-C5'-C4'	2.36	114.72	109.75
8	T	403	LMT	C2'-C3'-C4'	2.39	114.84	109.60
8	T	403	LMT	C1'-C2'-C3'	2.46	114.09	110.43
5	T	401	POV	O31-C31-C32	2.48	119.45	111.90
8	P	402	LMT	O5'-C1'-C2'	2.49	113.77	109.80
8	T	403	LMT	O5'-C1'-C2'	2.55	113.86	109.80
5	P	403	POV	O31-C31-C32	2.62	119.89	111.90
8	P	402	LMT	C2'-C3'-C4'	2.63	115.37	109.60
8	B	403	LMT	O5B-C5B-C4B	2.65	114.65	109.68
8	P	402	LMT	C1'-C2'-C3'	2.65	114.37	110.43
8	D	403	LMT	C1'-C2'-C3'	2.73	114.48	110.43
5	D	401	POV	O31-C31-C32	2.73	120.21	111.90
5	R	402	POV	O31-C31-C32	2.74	120.25	111.90
5	R	401	POV	O31-C31-C32	2.77	120.35	111.90
8	B	403	LMT	C2'-C3'-C4'	2.90	115.97	109.60
5	Q	401	POV	O31-C31-C32	3.04	121.15	111.90
5	Q	401	POV	O21-C21-C22	3.22	118.53	111.53
5	R	401	POV	O21-C21-C22	3.23	118.55	111.53
8	E	402	LMT	O5'-C5'-C4'	3.59	117.32	109.75
8	B	403	LMT	C1'-C2'-C3'	3.70	115.93	110.43
5	P	403	POV	O21-C21-C22	3.79	119.76	111.53
5	D	401	POV	O21-C21-C22	4.00	120.23	111.53
5	A	402	POV	O21-C21-C22	4.03	120.28	111.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	T	401	POV	O21-C21-C22	4.14	120.53	111.53
5	R	402	POV	O21-C21-C22	4.26	120.79	111.53
4	S	401	NAG	C1-O5-C5	4.78	118.32	112.25

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	T	402	NAG	O7-C7-N2-C2
5	R	401	POV	C2-O21-C21-C22

There are no ring outliers.

11 monomers are involved in 26 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	402	POV	1	0
5	D	401	POV	5	0
8	D	403	LMT	1	0
5	P	403	POV	3	0
5	Q	401	POV	5	0
5	R	401	POV	3	0
5	R	402	POV	2	0
4	S	401	NAG	1	0
8	S	402	LMT	1	0
5	T	401	POV	3	0
8	T	403	LMT	1	0

5.7 Other polymers

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

6.1 Protein, DNA and RNA chains

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, 95^{th} 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	342/347 (98%)	0.08	1 (0%) 94 93	56, 102, 165, 226	0
1	B	341/347 (98%)	0.05	3 (0%) 85 78	58, 109, 175, 270	0
1	C	340/347 (97%)	0.05	4 (1%) 81 69	58, 106, 185, 251	0
1	D	340/347 (97%)	0.01	1 (0%) 94 93	54, 106, 170, 263	0
1	E	340/347 (97%)	0.04	3 (0%) 85 78	57, 105, 170, 230	0
1	P	340/347 (97%)	-0.01	1 (0%) 94 93	51, 93, 156, 239	0
1	Q	340/347 (97%)	-0.03	1 (0%) 94 93	58, 99, 167, 242	0
1	R	340/347 (97%)	-0.05	4 (1%) 81 69	58, 97, 158, 274	0
1	S	340/347 (97%)	0.10	4 (1%) 81 69	53, 89, 144, 247	0
1	T	340/347 (97%)	0.02	1 (0%) 94 93	49, 88, 151, 210	0
2	F	221/224 (98%)	0.15	12 (5%) 29 17	68, 113, 202, 265	0
2	G	222/224 (99%)	0.43	16 (7%) 18 10	75, 142, 230, 309	0
2	H	223/224 (99%)	0.02	5 (2%) 65 50	52, 100, 158, 275	0
2	I	222/224 (99%)	0.01	4 (1%) 71 58	63, 105, 192, 239	0
2	J	222/224 (99%)	0.86	49 (22%) 1 1	64, 146, 281, 313	0
2	U	221/224 (98%)	0.68	31 (14%) 4 2	75, 131, 220, 317	0
2	V	224/224 (100%)	0.15	8 (3%) 46 31	70, 117, 169, 269	0
2	W	224/224 (100%)	-0.04	3 (1%) 79 67	68, 102, 155, 257	0
2	X	224/224 (100%)	0.03	4 (1%) 71 58	57, 98, 169, 249	0
2	Y	221/224 (98%)	0.25	8 (3%) 46 31	64, 127, 195, 277	0
3	K	210/215 (97%)	0.18	15 (7%) 19 10	59, 114, 222, 295	0
3	L	211/215 (98%)	-0.07	2 (0%) 85 78	49, 91, 170, 224	0
3	M	210/215 (97%)	0.49	25 (11%) 6 3	69, 130, 248, 270	0
3	N	211/215 (98%)	-0.05	0 100 100	66, 116, 158, 200	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
3	O	211/215 (98%)	-0.03	0 100 100	65, 108, 143, 157	0
3	Z	215/215 (100%)	0.22	2 (0%) 85 78	71, 120, 160, 199	0
3	f	215/215 (100%)	-0.10	1 (0%) 91 87	69, 104, 153, 231	0
3	g	211/215 (98%)	0.26	9 (4%) 39 25	71, 127, 173, 218	0
3	h	211/215 (98%)	0.28	4 (1%) 70 55	75, 137, 177, 249	0
3	i	214/215 (99%)	-0.04	0 100 100	60, 99, 135, 165	0
All	All	7746/7860 (98%)	0.11	221 (2%) 55 41	49, 107, 191, 317	0

All (221) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	J	165	SER	9.0
2	V	140	GLN	8.8
2	X	142	ASN	7.0
2	J	147	LEU	6.9
2	J	218	LYS	5.9
3	M	122	PRO	5.7
2	J	166	GLY	5.7
2	J	151	VAL	5.7
2	G	149	CYS	5.5
3	M	123	PRO	5.3
2	J	134	ALA	5.3
2	J	149	CYS	5.1
2	J	168	LEU	5.1
2	U	138	ALA	5.0
2	J	150	LEU	5.0
2	U	191	THR	4.9
3	K	123	PRO	4.7
1	B	304	GLY	4.6
2	Y	136	GLY	4.6
2	J	161	VAL	4.6
3	M	130	THR	4.5
3	M	188	TRP	4.5
2	J	197	TRP	4.5
3	M	136	VAL	4.5
2	U	192	VAL	4.4
2	V	139	ALA	4.3
2	J	187	SER	4.2
2	J	188	SER	4.2
3	M	194	TYR	4.2

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Mol	Chain	Res	Type	RSRZ
3	M	183	LEU	4.2
2	J	174	THR	4.2
2	G	150	LEU	4.1
2	F	192	VAL	4.1
2	U	149	CYS	4.1
2	U	140	GLN	4.1
2	U	161	VAL	4.0
2	J	190	VAL	4.0
3	M	134	THR	4.0
2	J	133	LEU	3.9
1	R	12	SER	3.9
2	U	147	LEU	3.9
2	J	172	VAL	3.9
3	M	189	GLU	3.9
3	K	194	TYR	3.8
2	J	201	THR	3.8
2	U	137	SER	3.8
2	J	163	TRP	3.8
2	F	147	LEU	3.8
2	J	220	VAL	3.8
2	J	192	VAL	3.8
2	X	224	CYS	3.8
2	G	137	SER	3.7
2	J	217	LYS	3.7
2	F	149	CYS	3.6
3	M	153	VAL	3.6
2	U	1	GLU	3.6
2	U	175	PHE	3.6
2	J	132	PRO	3.6
2	J	189	SER	3.6
3	Z	205	VAL	3.5
2	G	135	PRO	3.5
2	U	165	SER	3.4
2	J	167	SER	3.3
3	h	144	PRO	3.3
2	V	121	SER	3.3
2	F	139	ALA	3.3
3	K	122	PRO	3.3
2	J	157	GLU	3.3
2	F	140	GLN	3.3
3	M	135	LEU	3.1
2	J	204	CYS	3.1

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Mol	Chain	Res	Type	RSRZ
2	G	132	PRO	3.1
3	g	188	TRP	3.1
2	F	137	SER	3.1
3	K	195	SER	3.1
1	S	69	LYS	3.0
3	M	186	ARG	3.0
1	B	213	PHE	3.0
2	Y	147	LEU	3.0
2	G	134	ALA	3.0
2	H	1	GLU	2.9
2	F	163	TRP	2.9
2	H	138	ALA	2.9
2	V	219	ILE	2.9
2	U	206	VAL	2.9
3	h	165	THR	2.9
3	Z	144	PRO	2.9
3	M	127	GLU	2.9
2	U	141	THR	2.8
2	U	196	THR	2.8
2	U	188	SER	2.8
3	M	121	PHE	2.8
3	K	193	SER	2.8
2	J	162	THR	2.8
1	B	340	HIS	2.8
3	K	121	PHE	2.8
2	J	193	PRO	2.8
3	K	135	LEU	2.8
2	G	141	THR	2.8
2	U	162	THR	2.8
2	J	146	THR	2.8
2	U	195	SER	2.7
2	W	139	ALA	2.7
1	S	10	PHE	2.7
3	K	124	SER	2.7
2	Y	138	ALA	2.7
3	M	182	THR	2.7
1	C	213	PHE	2.6
2	J	175	PHE	2.6
3	g	133	ALA	2.6
2	X	147	LEU	2.6
2	U	181	SER	2.6
2	G	147	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
2	I	1	GLU	2.6
2	V	175	PHE	2.6
3	M	154	ASP	2.6
2	G	18	MET	2.6
2	U	163	TRP	2.6
2	G	130	VAL	2.6
2	G	187	SER	2.6
2	I	139	ALA	2.6
2	J	196	THR	2.6
2	J	126	THR	2.5
2	J	135	PRO	2.5
2	F	148	GLY	2.5
2	W	142	ASN	2.5
2	J	216	ASP	2.5
2	J	145	VAL	2.5
1	S	6	LEU	2.5
2	G	151	VAL	2.5
2	U	210	ALA	2.5
1	A	341	HIS	2.5
2	U	152	LYS	2.5
3	f	155	GLY	2.5
1	T	263	ILE	2.5
2	U	202	VAL	2.5
2	H	23	LYS	2.5
2	J	139	ALA	2.4
2	V	174	THR	2.4
2	J	131	TYR	2.4
3	L	1	GLN	2.4
2	J	152	LYS	2.4
2	F	132	PRO	2.4
2	U	169	SER	2.4
1	D	305	THR	2.4
2	U	189	SER	2.4
2	U	204	CYS	2.4
3	M	191	HIS	2.4
2	U	186	LEU	2.4
3	K	136	VAL	2.4
3	K	133	ALA	2.4
1	E	310	ASP	2.3
3	K	179	SER	2.3
2	G	1	GLU	2.3
1	E	241	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
2	J	156	PRO	2.3
2	J	210	ALA	2.3
2	U	197	TRP	2.3
2	Y	170	SER	2.3
3	h	178	SER	2.3
3	g	18	VAL	2.3
2	F	219	ILE	2.3
3	K	127	GLU	2.3
3	g	78	THR	2.3
2	J	164	ASN	2.3
2	J	155	PHE	2.3
1	R	308	TRP	2.3
2	X	139	ALA	2.3
2	W	222	ARG	2.3
1	P	12	SER	2.3
2	V	147	LEU	2.3
2	Y	217	LYS	2.3
3	M	138	THR	2.2
1	R	2	ASP	2.2
2	Y	1	GLU	2.2
3	M	129	GLU	2.2
2	U	178	VAL	2.2
2	G	197	TRP	2.2
1	Q	1	SER	2.2
2	H	140	GLN	2.2
2	U	180	GLN	2.2
3	K	209	LEU	2.2
2	J	169	SER	2.2
3	M	196	CYS	2.2
2	G	145	VAL	2.2
2	Y	141	THR	2.2
3	g	180	TYR	2.2
2	J	44	ASN	2.2
3	h	152	LYS	2.2
3	M	155	GLY	2.2
3	g	109	LEU	2.2
3	g	108	VAL	2.2
2	J	127	PRO	2.1
3	M	163	GLU	2.1
1	E	239	ARG	2.1
2	J	191	THR	2.1
1	C	99	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
3	M	184	THR	2.1
2	Y	215	VAL	2.1
2	J	202	VAL	2.1
3	K	134	THR	2.1
2	V	179	LEU	2.1
2	G	144	MET	2.1
3	M	180	TYR	2.1
2	H	139	ALA	2.1
1	R	182	GLN	2.1
2	U	219	ILE	2.1
3	g	131	ASN	2.1
3	L	186	ARG	2.1
3	g	185	ALA	2.1
2	U	168	LEU	2.1
2	J	130	VAL	2.1
3	K	210	SER	2.1
2	F	138	ALA	2.0
2	F	190	VAL	2.0
1	S	1	SER	2.0
2	I	150	LEU	2.0
1	C	304	GLY	2.0
1	C	98	ALA	2.0
2	I	130	VAL	2.0
3	M	133	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95^{th} 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	LLDF	B-factors(Å^2)	Q<0.9
5	POV	Q	401	46/52	0.81	0.67	11.09	108,149,198,237	0
5	POV	R	402	35/52	0.81	0.50	5.92	96,123,186,208	0
5	POV	A	402	23/52	0.73	0.62	5.72	137,177,208,241	0
5	POV	D	401	26/52	0.76	0.63	5.68	125,176,201,334	0
5	POV	T	401	43/52	0.77	0.47	4.77	77,129,177,205	0
8	LMT	D	403	23/35	0.79	0.58	4.02	196,218,239,247	0
5	POV	P	403	42/52	0.78	0.43	3.95	82,125,179,206	0
5	POV	R	401	34/52	0.73	0.43	3.78	99,146,200,535	0
8	LMT	P	402	23/35	0.56	0.43	1.99	221,221,221,221	0
8	LMT	B	403	23/35	0.70	0.55	1.45	176,217,235,242	0
7	GOL	B	401	6/6	0.58	0.35	1.29	115,145,151,167	0
8	LMT	C	402	23/35	0.79	0.32	-0.10	179,179,179,179	0
4	NAG	S	401	14/15	0.80	0.22	-	153,184,189,195	0
6	CL	P	404	1/1	0.89	0.20	-	90,90,90,90	0
4	NAG	A	401	14/15	0.86	0.23	-	137,181,192,196	0
4	NAG	T	402	14/15	0.69	0.34	-	153,200,210,211	0
8	LMT	S	402	23/35	0.61	0.51	-	169,214,224,235	0
8	LMT	E	402	23/35	0.63	0.63	-	218,239,254,257	0
6	CL	A	403	1/1	0.54	0.61	-	119,119,119,119	0
4	NAG	E	401	14/15	0.90	0.17	-	157,190,206,212	0
4	NAG	B	402	14/15	0.87	0.25	-	139,168,176,177	0
4	NAG	C	401	14/15	0.84	0.26	-	191,210,218,223	0
4	NAG	D	402	14/15	0.80	0.20	-	153,197,214,221	0
8	LMT	T	403	23/35	0.58	0.38	-	156,202,217,222	0
4	NAG	Q	402	14/15	0.82	0.23	-	146,184,202,207	0
4	NAG	P	401	14/15	0.85	0.24	-	133,182,192,201	0

6.5 Other polymers (i)

There are no such residues in this entry.