



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 19, 2026 – 02:46 AM UTC

PDB ID : 6ELU / pdb_00006elu
Title : Structure of Serum Resistance Associated protein from T. b. rhodesiense
Authors : Zoll, S.; Higgins, M.K.
Deposited on : 2017-09-29
Resolution : 2.30 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	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

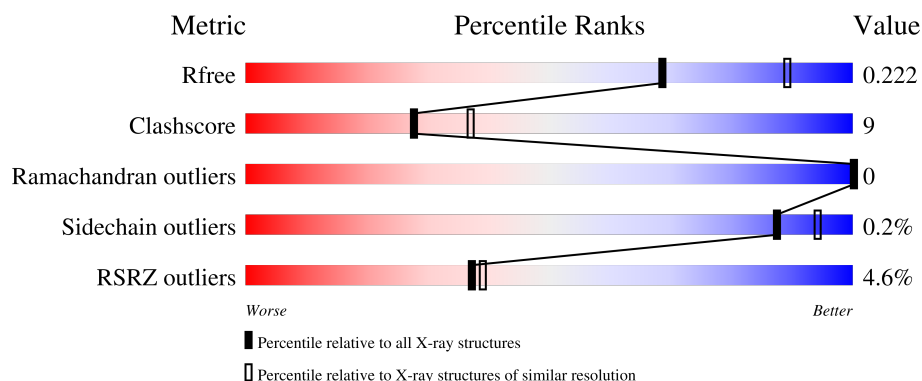
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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.



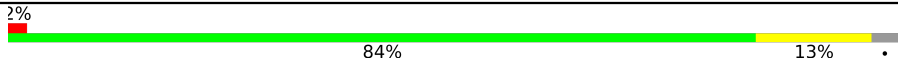
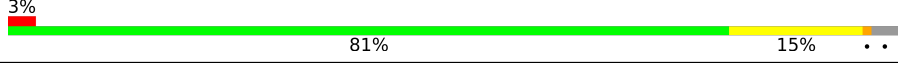
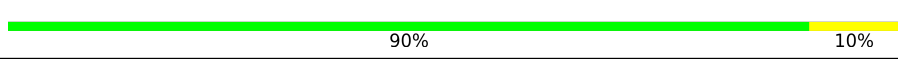
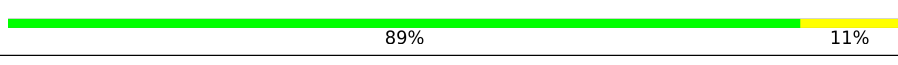
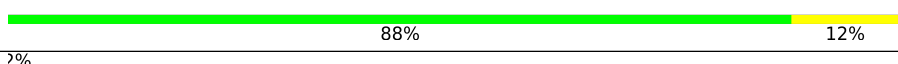
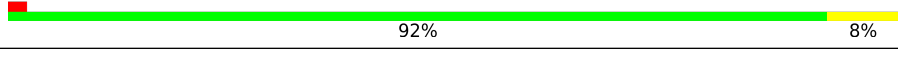
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	233	9% 61% 22% 17%
1	D	233	11% 48% 21% 31%
1	G	233	6% 52% 20% 28%
1	J	233	11% 54% 19% 26%
2	B	229	2% 83% 13% •

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Mol	Chain	Length	Quality of chain
2	E	229	
2	H	229	
2	K	229	
3	C	217	
3	F	217	
3	I	217	
3	L	217	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 19153 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 Serum resistance associated; VSG protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	193	Total	C	N	O	S	0	0	0
			1413	883	247	280	3			
1	D	161	Total	C	N	O	S	0	0	0
			1188	745	208	232	3			
1	G	168	Total	C	N	O	S	0	0	0
			1236	775	216	242	3			
1	J	172	Total	C	N	O	S	0	0	0
			1264	793	221	247	3			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	29	GLY	-	expression tag	UNP Q8T309
A	30	SER	-	expression tag	UNP Q8T309
A	31	HIS	-	expression tag	UNP Q8T309
A	32	MET	-	expression tag	UNP Q8T309
D	29	GLY	-	expression tag	UNP Q8T309
D	30	SER	-	expression tag	UNP Q8T309
D	31	HIS	-	expression tag	UNP Q8T309
D	32	MET	-	expression tag	UNP Q8T309
G	29	GLY	-	expression tag	UNP Q8T309
G	30	SER	-	expression tag	UNP Q8T309
G	31	HIS	-	expression tag	UNP Q8T309
G	32	MET	-	expression tag	UNP Q8T309
J	29	GLY	-	expression tag	UNP Q8T309
J	30	SER	-	expression tag	UNP Q8T309
J	31	HIS	-	expression tag	UNP Q8T309
J	32	MET	-	expression tag	UNP Q8T309

- Molecule 2 is a protein called G10_3 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	222	Total 1686	C 1064	N 280	O 334	S 8	0	0	0
2	E	222	Total 1686	C 1064	N 280	O 334	S 8	0	0	0
2	H	222	Total 1686	C 1064	N 280	O 334	S 8	0	0	0
2	K	222	Total 1686	C 1064	N 280	O 334	S 8	0	0	0

- Molecule 3 is a protein called G10_3 Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	217	Total 1682	C 1046	N 282	O 347	S 7	0	0	0
3	F	217	Total 1682	C 1046	N 282	O 347	S 7	0	0	0
3	I	217	Total 1682	C 1046	N 282	O 347	S 7	0	0	0
3	L	217	Total 1682	C 1046	N 282	O 347	S 7	0	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	O 1	0	0
4	B	49	Total 49	O 49	0	0
4	C	83	Total 83	O 83	0	0
4	E	50	Total 50	O 50	0	0
4	F	77	Total 77	O 77	0	0
4	G	4	Total 4	O 4	0	0
4	H	86	Total 86	O 86	0	0
4	I	91	Total 91	O 91	0	0
4	J	3	Total 3	O 3	0	0
4	K	74	Total 74	O 74	0	0

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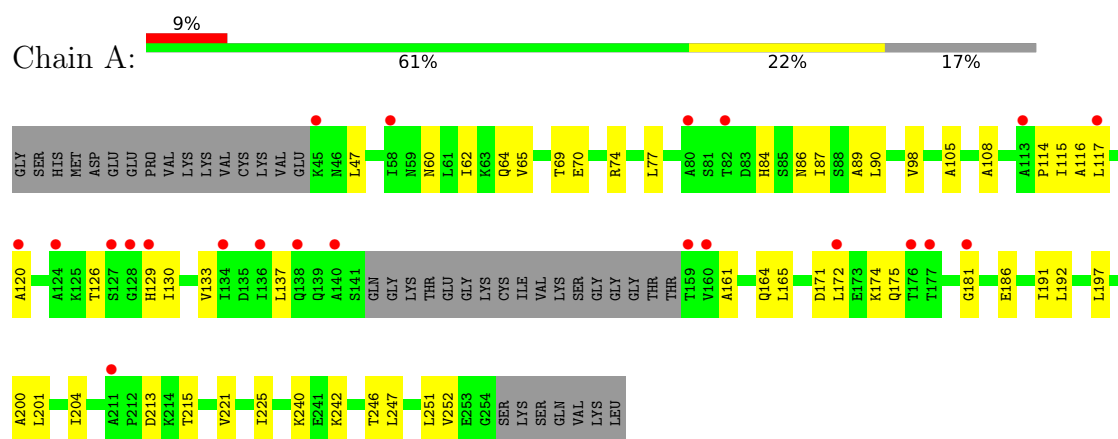
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	62	Total	O	0	0
			62	62		

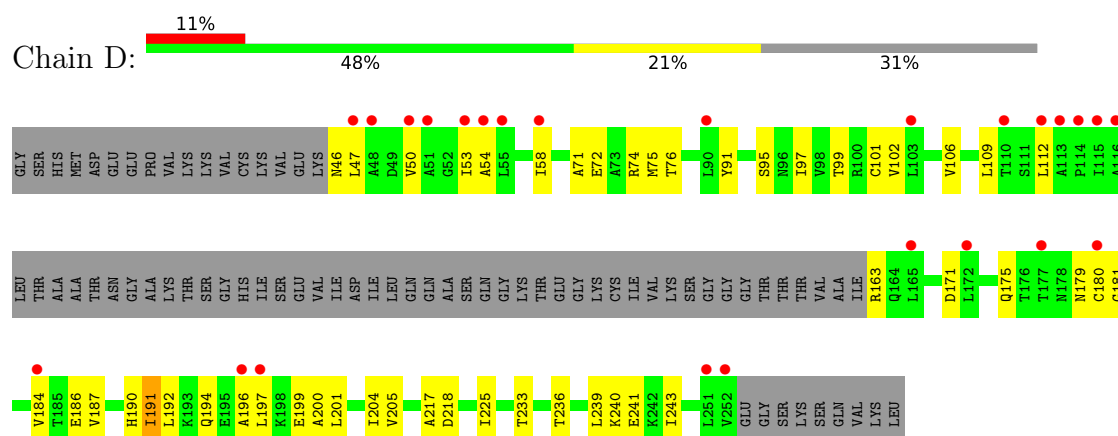
3 Residue-property plots [i](#)

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.

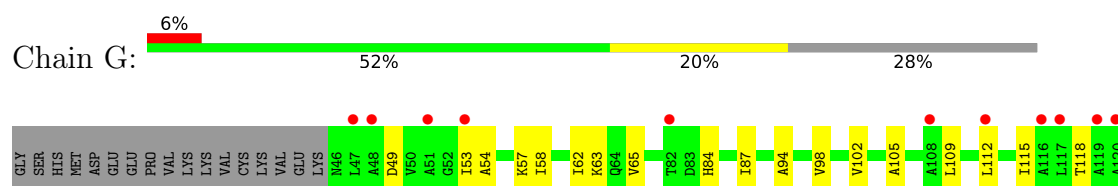
- Molecule 1: Serum resistance associated; VSG protein

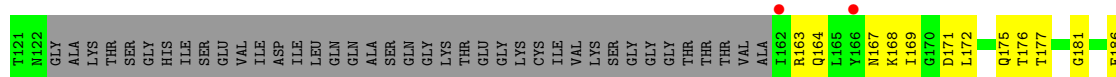


- Molecule 1: Serum resistance associated; VSG protein



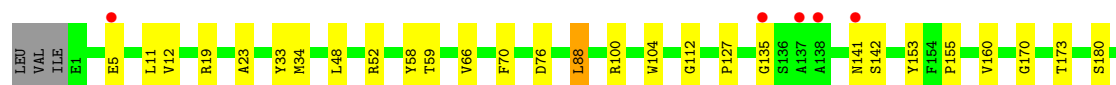
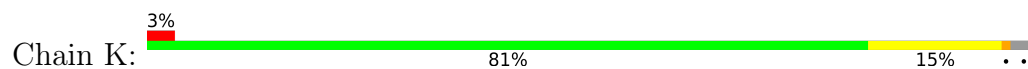
- Molecule 1: Serum resistance associated; VSG protein







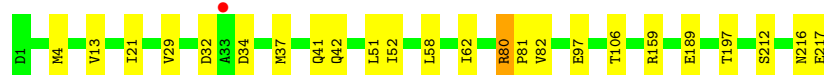
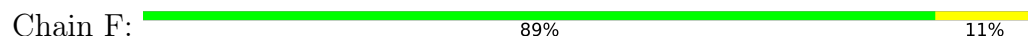
- Molecule 2: G10_3 heavy chain



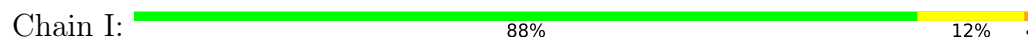
- Molecule 3: G10_3 Light chain



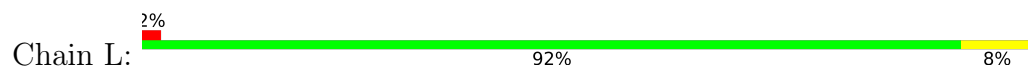
- Molecule 3: G10_3 Light chain



- Molecule 3: G10_3 Light chain



- Molecule 3: G10_3 Light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	127.60Å 103.55Å 150.64Å 90.00° 114.58° 90.00°	Depositor
Resolution (Å)	48.43 – 2.30 48.43 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.43-2.30) 99.9 (48.43-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.74 (at 2.29Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.191 , 0.220 0.192 , 0.222	Depositor DCC
R_{free} test set	7914 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	46.9	Xtriage
Anisotropy	0.209	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 63.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.026 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	19153	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% 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

5.1 Standard geometry

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	0.33	0/1422	0.64	0/1932
1	D	0.25	0/1196	0.49	1/1625 (0.1%)
1	G	0.31	0/1244	0.64	0/1692
1	J	0.34	0/1272	0.71	3/1730 (0.2%)
2	B	0.33	1/1731 (0.1%)	0.60	3/2364 (0.1%)
2	E	0.28	1/1731 (0.1%)	0.55	1/2364 (0.0%)
2	H	0.22	0/1731	0.49	0/2364
2	K	0.24	0/1731	0.55	1/2364 (0.0%)
3	C	0.23	0/1722	0.49	1/2342 (0.0%)
3	F	0.31	1/1722 (0.1%)	0.58	3/2342 (0.1%)
3	I	0.22	0/1722	0.48	1/2342 (0.0%)
3	L	0.19	0/1722	0.44	0/2342
All	All	0.27	3/18946 (0.0%)	0.55	14/25803 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	40	PRO	CA-C	-6.34	1.48	1.51
2	E	40	PRO	CA-C	5.45	1.54	1.51
3	F	80	ARG	CG-CD	-5.35	1.36	1.52

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	80	ARG	CA-CB-CG	-9.17	95.75	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	80	ARG	CG-CD-NE	-8.22	93.91	112.00
1	J	175	GLN	N-CA-C	-7.78	102.40	111.03
2	E	113	GLN	CB-CA-C	-7.08	97.08	110.67
3	I	160	GLN	CA-CB-CG	-6.98	100.14	114.10
2	B	3	LYS	CG-CD-CE	-6.96	95.28	111.30
3	F	97	GLU	N-CA-CB	6.78	122.63	110.83
3	F	97	GLU	CB-CA-C	-6.00	97.92	109.37
1	D	191	ILE	CG1-CB-CG2	-5.87	93.08	110.70
2	B	123	LYS	CB-CG-CD	-5.49	98.67	111.30
1	J	172	LEU	CA-C-N	-5.23	112.36	120.31
1	J	172	LEU	C-N-CA	-5.23	112.36	120.31
2	K	199	GLU	N-CA-CB	-5.16	102.25	111.08
2	B	54	LYS	CB-CG-CD	5.04	122.90	111.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	140	THR	Peptide

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	1413	0	1493	51	0
1	D	1188	0	1263	53	0
1	G	1236	0	1313	38	0
1	J	1264	0	1347	43	0
2	B	1686	0	1629	20	0
2	E	1686	0	1629	24	0
2	H	1686	0	1629	25	0
2	K	1686	0	1629	27	0
3	C	1682	0	1593	13	0
3	F	1682	0	1593	19	0
3	I	1682	0	1593	17	0
3	L	1682	0	1593	10	0
4	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	49	0	0	0	0
4	C	83	0	0	0	0
4	E	50	0	0	0	0
4	F	77	0	0	0	0
4	G	4	0	0	1	0
4	H	86	0	0	1	0
4	I	91	0	0	0	0
4	J	3	0	0	1	0
4	K	74	0	0	2	0
4	L	62	0	0	0	0
All	All	19153	0	18304	330	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 (330) 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:A:161:ALA:O	1:A:164:GLN:CG	2.05	1.04
1:A:161:ALA:HA	1:A:164:GLN:HG2	1.42	1.00
3:L:193:HIS:O	3:L:215:ARG:NH2	1.94	1.00
1:A:161:ALA:O	1:A:164:GLN:HG2	1.64	0.97
1:J:192:LEU:HB3	1:J:243:ILE:HD12	1.49	0.94
1:G:57:LYS:HE2	1:G:172:LEU:HD21	1.50	0.94
1:A:161:ALA:CA	1:A:164:GLN:HG2	2.02	0.90
2:K:135:GLY:HA2	2:K:221:ARG:HD3	1.54	0.89
1:A:161:ALA:C	1:A:164:GLN:HG2	1.98	0.88
3:F:80:ARG:HG3	3:F:81:PRO:HA	1.55	0.85
1:A:77:LEU:O	1:A:77:LEU:HD23	1.76	0.85
1:A:161:ALA:O	1:A:164:GLN:HG3	1.73	0.85
1:J:173:GLU:HG3	1:J:174:LYS:N	1.91	0.84
1:A:181:GLY:HA3	1:A:186:GLU:HB3	1.60	0.83
1:G:167:ASN:O	1:G:171:ASP:N	2.11	0.82
1:D:58:ILE:HD11	1:D:112:LEU:HD11	1.62	0.81
2:K:5:GLU:OE1	2:K:23:ALA:HB3	1.81	0.80
1:A:115:ILE:O	1:A:164:GLN:NE2	2.14	0.80
1:A:164:GLN:HG3	1:A:165:LEU:N	1.99	0.78
1:D:46:ASN:HB2	1:D:163:ARG:NH2	2.00	0.77
1:A:129:HIS:O	1:A:129:HIS:ND1	2.17	0.77
1:A:221:VAL:O	1:A:225:ILE:HG13	1.84	0.76
2:H:37:VAL:HG21	2:H:108:MET:HE1	1.68	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:141:ASN:OD1	2:B:142:SER:N	2.19	0.76
3:I:216:ASN:O	3:I:216:ASN:ND2	2.19	0.76
1:A:98:VAL:HG23	1:A:191:ILE:HD11	1.68	0.75
1:A:171:ASP:HA	1:A:174:LYS:HD2	1.67	0.74
1:G:65:VAL:HG13	1:G:102:VAL:HG13	1.70	0.74
1:D:171:ASP:O	1:D:175:GLN:NE2	2.22	0.73
1:A:77:LEU:HD23	1:A:77:LEU:C	2.13	0.73
3:F:80:ARG:HG3	3:F:81:PRO:CA	2.19	0.72
1:D:233:THR:HG23	1:D:236:THR:H	1.54	0.71
1:D:46:ASN:HB2	1:D:163:ARG:HH22	1.56	0.71
1:J:167:ASN:O	1:J:171:ASP:N	2.23	0.71
1:G:57:LYS:CE	1:G:172:LEU:HD21	2.21	0.71
1:G:112:LEU:HD11	1:G:169:ILE:HG22	1.71	0.70
1:A:108:ALA:HB1	1:A:172:LEU:HD21	1.71	0.70
2:H:52:ARG:HE	2:H:59:THR:CG2	2.05	0.69
1:A:114:PRO:HB3	1:A:126:THR:HG23	1.75	0.69
2:K:195:THR:O	2:K:199:GLU:HB2	1.93	0.69
1:D:197:LEU:CA	1:D:199:GLU:OE2	2.40	0.68
2:B:23:ALA:HA	2:B:80:ILE:HD13	1.73	0.68
2:H:33:TYR:CE2	2:H:52:ARG:HG2	2.28	0.68
1:D:72:GLU:OE2	1:D:76:THR:HG23	1.94	0.68
2:K:141:ASN:OD1	2:K:142:SER:N	2.26	0.67
1:D:181:GLY:HA3	1:D:186:GLU:HB3	1.76	0.67
1:D:192:LEU:HA	1:D:197:LEU:HD11	1.76	0.67
2:H:52:ARG:HE	2:H:59:THR:HG23	1.58	0.67
3:F:159:ARG:NH2	3:F:189:GLU:OE2	2.21	0.67
2:E:24:THR:HG21	2:E:29:PHE:CD1	2.30	0.67
3:C:216:ASN:O	3:C:216:ASN:ND2	2.28	0.67
1:J:171:ASP:O	1:J:175:GLN:HG2	1.95	0.66
1:D:47:LEU:HD13	1:D:47:LEU:O	1.95	0.66
2:E:24:THR:HG22	2:E:79:SER:O	1.94	0.66
1:J:233:THR:HG23	1:J:236:THR:H	1.59	0.66
2:E:113:GLN:HG2	2:E:114:GLY:N	2.11	0.65
1:D:54:ALA:O	1:D:58:ILE:HD13	1.95	0.65
1:D:197:LEU:HA	1:D:199:GLU:OE2	1.97	0.65
1:J:64:GLN:NE2	4:J:301:HOH:O	2.25	0.65
3:I:40:TYR:HE1	3:I:93:GLN:HG2	1.61	0.65
2:B:4:LEU:HD23	2:B:98:CYS:SG	2.36	0.65
1:A:191:ILE:HG22	1:A:197:LEU:HD22	1.79	0.65
2:B:48:LEU:HD22	2:B:66:VAL:HG21	1.77	0.64
1:G:84:HIS:HB2	1:G:87:ILE:HG22	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:191:ILE:HG22	1:G:197:LEU:HD12	1.80	0.64
1:J:115:ILE:HD12	1:J:165:LEU:HA	1.79	0.63
2:H:33:TYR:CZ	2:H:52:ARG:HG2	2.33	0.62
3:L:159:ARG:NH2	3:L:189:GLU:OE1	2.33	0.61
1:A:225:ILE:HD12	1:A:240:LYS:HB2	1.83	0.61
1:J:218:ASP:O	1:J:222:THR:HG23	2.01	0.61
3:I:158:GLU:OE1	3:I:160:GLN:NE2	2.33	0.61
1:A:171:ASP:O	1:A:174:LYS:HB2	2.01	0.60
1:J:57:LYS:HE2	1:J:172:LEU:HD12	1.82	0.60
2:E:39:GLN:HE22	3:F:42:GLN:HE22	1.49	0.60
2:E:48:LEU:HD22	2:E:66:VAL:HG21	1.82	0.60
1:G:241:GLU:OE2	2:H:52:ARG:HD3	2.00	0.60
2:H:69:ARG:NH2	2:H:92:ASP:OD2	2.36	0.59
2:H:85:MET:HE2	2:H:88:LEU:HD21	1.86	0.58
1:A:117:LEU:HD13	1:A:130:ILE:HB	1.86	0.58
1:D:187:VAL:O	1:D:191:ILE:HD12	2.04	0.57
2:H:52:ARG:NH2	2:H:61:GLU:OE2	2.35	0.57
1:A:47:LEU:HG	1:A:137:LEU:HD11	1.86	0.57
3:C:41:GLN:HB2	3:C:51:LEU:HD11	1.86	0.57
3:F:41:GLN:HB2	3:F:51:LEU:HD11	1.86	0.57
1:G:175:GLN:NE2	1:G:177:THR:HG23	2.20	0.57
1:J:47:LEU:HD21	1:J:162:ILE:HD11	1.86	0.57
1:A:47:LEU:HD11	1:A:120:ALA:HA	1.88	0.56
1:A:86:ASN:N	1:A:86:ASN:OD1	2.39	0.56
2:B:54:LYS:HG3	2:B:58:TYR:CZ	2.40	0.56
3:L:41:GLN:HB2	3:L:51:LEU:HD11	1.86	0.56
2:E:33:TYR:CZ	2:E:52:ARG:HG2	2.40	0.56
1:G:163:ARG:H	1:G:163:ARG:HD3	1.69	0.56
2:K:180:SER:O	4:K:301:HOH:O	2.17	0.55
3:C:16:GLY:HA2	3:C:81:PRO:HB2	1.87	0.55
3:C:165:ASN:HB3	3:C:179:MET:HE1	1.88	0.55
1:D:191:ILE:O	1:D:197:LEU:HD22	2.07	0.55
1:D:199:GLU:OE2	1:D:199:GLU:N	2.40	0.55
3:C:20:THR:HG22	3:C:78:ASN:OD1	2.06	0.54
1:A:70:GLU:O	1:A:74:ARG:HG3	2.08	0.54
1:D:241:GLU:HG2	2:E:104:TRP:HB3	1.89	0.54
1:G:207:LYS:HG3	1:G:210:GLY:H	1.72	0.54
1:A:164:GLN:HG3	1:A:165:LEU:H	1.73	0.54
3:F:52:ILE:HD13	3:F:58:LEU:HA	1.90	0.54
1:J:70:GLU:OE1	1:J:73:ALA:HB3	2.08	0.54
1:G:164:GLN:HA	1:G:167:ASN:OD1	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:160:VAL:HG22	2:K:205:VAL:HG22	1.90	0.53
3:F:13:VAL:HG23	3:F:82:VAL:HG21	1.90	0.53
1:D:191:ILE:HD12	1:D:191:ILE:H	1.73	0.53
2:K:12:VAL:HG11	2:K:88:LEU:HD23	1.91	0.53
2:E:52:ARG:NH2	2:E:61:GLU:OE2	2.41	0.53
1:G:115:ILE:O	1:G:118:THR:HG22	2.09	0.52
1:J:173:GLU:HG3	1:J:174:LYS:H	1.71	0.52
2:E:207:HIS:HB3	2:E:212:THR:OG1	2.08	0.52
1:G:112:LEU:HD11	1:G:169:ILE:CG2	2.39	0.52
1:G:193:LYS:O	4:G:301:HOH:O	2.19	0.52
2:B:75:ASP:CG	2:B:78:GLN:HG3	2.35	0.52
2:H:85:MET:HB3	2:H:88:LEU:HD21	1.90	0.52
1:J:65:VAL:HG12	1:J:105:ALA:HB3	1.91	0.52
3:L:6:GLN:NE2	3:L:106:THR:HG23	2.23	0.52
3:C:112:ARG:HG3	3:C:113:ALA:O	2.10	0.52
1:A:84:HIS:HB3	1:A:86:ASN:OD1	2.10	0.51
1:D:101:CYS:CB	1:D:180:CYS:HG	2.22	0.51
1:D:71:ALA:O	1:D:75:MET:HG3	2.10	0.51
1:G:188:LEU:HA	1:G:191:ILE:HD12	1.93	0.51
2:K:48:LEU:HD22	2:K:66:VAL:HG21	1.93	0.51
2:B:1:GLU:HG3	2:B:2:VAL:N	2.25	0.51
1:D:192:LEU:C	1:D:197:LEU:HD21	2.35	0.51
2:E:5:GLU:HA	2:E:113:GLN:OE1	2.11	0.51
3:F:29:VAL:HG21	3:F:37:MET:HE2	1.93	0.51
1:G:192:LEU:HA	1:G:197:LEU:HD13	1.93	0.51
1:D:191:ILE:HD12	1:D:191:ILE:N	2.26	0.51
2:E:200:THR:HG22	2:E:217:LYS:HE3	1.93	0.50
1:J:115:ILE:HD11	1:J:168:LYS:HB3	1.94	0.50
1:J:168:LYS:HA	1:J:171:ASP:HB2	1.92	0.50
3:F:13:VAL:HG21	3:F:82:VAL:HG11	1.94	0.50
1:J:233:THR:O	1:J:236:THR:HG22	2.12	0.50
1:A:192:LEU:HA	1:A:197:LEU:HD23	1.94	0.50
2:H:108:MET:HE2	2:H:111:TRP:CZ2	2.46	0.50
2:B:70:PHE:HB3	2:B:83:LEU:HD11	1.94	0.50
1:G:200:ALA:O	1:G:204:ILE:HG13	2.12	0.50
3:C:29:VAL:HG21	3:C:37:MET:HE2	1.94	0.49
2:B:33:TYR:CZ	2:B:52:ARG:HG2	2.46	0.49
2:K:33:TYR:CE2	2:K:52:ARG:HG2	2.47	0.49
2:K:52:ARG:HD2	2:K:59:THR:OG1	2.12	0.49
3:C:112:ARG:NH1	3:C:113:ALA:O	2.45	0.49
3:I:41:GLN:HB2	3:I:51:LEU:HD11	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:GLN:OE1	1:A:175:GLN:HA	2.12	0.49
2:K:11:LEU:HB2	2:K:155:PRO:HG3	1.93	0.49
1:G:98:VAL:O	1:G:102:VAL:HG23	2.13	0.49
1:J:62:ILE:O	1:J:65:VAL:HG22	2.11	0.49
1:J:200:ALA:O	1:J:204:ILE:HG13	2.13	0.49
2:E:85:MET:HB3	2:E:88:LEU:HD21	1.94	0.49
2:K:173:THR:HG22	2:K:187:SER:OG	2.13	0.49
1:A:77:LEU:C	1:A:77:LEU:CD2	2.86	0.48
1:A:84:HIS:HB2	1:A:87:ILE:HG22	1.93	0.48
2:B:66:VAL:HG13	2:B:70:PHE:HB2	1.94	0.48
1:G:191:ILE:HG22	1:G:197:LEU:CD1	2.43	0.48
1:G:241:GLU:HG2	2:H:104:TRP:HB3	1.95	0.48
1:G:49:ASP:O	1:G:53:ILE:HD13	2.14	0.48
1:A:137:LEU:HD13	1:A:137:LEU:O	2.13	0.48
1:A:213:ASP:OD1	1:A:215:THR:HG22	2.13	0.48
1:J:50:VAL:CG1	1:J:165:LEU:HD22	2.43	0.48
1:J:192:LEU:HA	1:J:197:LEU:HD23	1.94	0.48
1:G:181:GLY:HA3	1:G:186:GLU:HB3	1.95	0.48
2:H:48:LEU:HD22	2:H:66:VAL:HG11	1.96	0.48
1:D:200:ALA:O	1:D:204:ILE:HG13	2.14	0.48
1:J:59:ASN:HB3	1:J:63:LYS:NZ	2.28	0.48
2:K:127:PRO:HB3	2:K:153:TYR:HB3	1.95	0.48
1:D:101:CYS:CB	1:D:180:CYS:SG	3.02	0.48
1:D:194:GLN:O	1:D:197:LEU:HD23	2.14	0.48
1:D:233:THR:HG23	1:D:236:THR:N	2.27	0.48
1:J:191:ILE:HG22	1:J:197:LEU:HD22	1.96	0.48
1:J:178:ASN:O	1:J:190:HIS:NE2	2.46	0.47
2:K:33:TYR:CZ	2:K:52:ARG:HG2	2.49	0.47
2:K:112:GLY:O	3:L:47:PRO:HB3	2.14	0.47
3:L:29:VAL:HG21	3:L:37:MET:HE2	1.95	0.47
1:D:75:MET:HE2	1:D:91:TYR:HD1	1.80	0.47
3:I:18:ARG:NH2	3:I:80:ARG:HD3	2.28	0.47
1:G:65:VAL:HG22	1:G:105:ALA:HB3	1.97	0.47
1:J:46:ASN:O	1:J:50:VAL:HG23	2.15	0.47
2:B:127:PRO:HB3	2:B:153:TYR:HB3	1.95	0.47
2:H:99:ALA:CB	2:H:108:MET:HE3	2.45	0.47
1:A:130:ILE:O	1:A:133:VAL:HG12	2.14	0.47
1:D:197:LEU:N	1:D:199:GLU:OE2	2.49	0.46
2:H:135:GLY:HA2	2:H:221:ARG:HD3	1.96	0.46
2:K:205:VAL:HB	2:K:214:VAL:HG13	1.97	0.46
1:J:50:VAL:HG11	1:J:165:LEU:HD22	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:197:THR:HG22	3:F:212:SER:CB	2.46	0.46
3:F:197:THR:HG22	3:F:212:SER:HB3	1.96	0.46
1:A:161:ALA:HA	1:A:164:GLN:CG	2.29	0.46
3:L:191:GLU:C	3:L:215:ARG:NH2	2.74	0.46
1:D:180:CYS:HB3	1:D:190:HIS:CG	2.51	0.46
1:D:201:LEU:O	1:D:205:VAL:HG22	2.16	0.46
1:D:218:ASP:CG	1:D:240:LYS:HZ1	2.23	0.46
3:I:167:TRP:CZ2	3:I:179:MET:HE2	2.51	0.46
2:K:34:MET:HE1	2:K:100:ARG:HG3	1.97	0.46
2:K:66:VAL:HG13	2:K:70:PHE:HB2	1.97	0.46
2:B:54:LYS:HG3	2:B:58:TYR:OH	2.16	0.46
1:D:102:VAL:O	1:D:106:VAL:HG12	2.16	0.46
3:I:217:GLU:N	3:I:217:GLU:OE1	2.49	0.46
2:K:19:ARG:NH1	4:K:309:HOH:O	2.39	0.46
1:D:225:ILE:HG21	1:D:236:THR:HG23	1.98	0.45
3:C:199:GLU:HG2	3:C:210:VAL:HG22	1.97	0.45
1:D:95:SER:O	1:D:99:THR:HG23	2.16	0.45
1:G:201:LEU:O	1:G:205:VAL:HG22	2.16	0.45
1:J:233:THR:CG2	1:J:236:THR:H	2.29	0.45
2:E:39:GLN:HE22	3:F:42:GLN:NE2	2.13	0.45
3:I:4:MET:HE1	3:I:93:GLN:CA	2.46	0.45
3:I:29:VAL:HG21	3:I:37:MET:HE2	1.98	0.45
3:I:159:ARG:NH2	3:I:189:GLU:OE2	2.46	0.45
2:K:141:ASN:CG	2:K:142:SER:N	2.75	0.45
3:C:197:THR:HG22	3:C:212:SER:HB3	1.99	0.45
1:A:65:VAL:O	1:A:69:THR:HG23	2.17	0.45
2:E:33:TYR:CE2	2:E:52:ARG:HG2	2.51	0.45
2:K:197:PRO:HB3	2:K:220:PRO:HG3	1.99	0.45
1:A:89:ALA:HB3	1:A:204:ILE:HD11	1.99	0.44
1:G:57:LYS:HD2	1:G:57:LYS:HA	1.55	0.44
1:J:95:SER:HA	1:J:98:VAL:HG12	1.99	0.44
1:A:108:ALA:HB1	1:A:172:LEU:CD2	2.43	0.44
1:A:116:ALA:HB3	1:A:130:ILE:HD11	1.99	0.44
1:D:97:ILE:HD11	1:D:196:ALA:HB3	1.99	0.44
2:B:23:ALA:CA	2:B:80:ILE:HD13	2.46	0.44
1:D:192:LEU:O	1:D:197:LEU:HD21	2.17	0.44
3:F:21:ILE:HG12	3:F:106:THR:HG21	1.98	0.44
2:B:33:TYR:CE2	2:B:52:ARG:HG2	2.53	0.43
1:G:94:ALA:O	1:G:98:VAL:HG13	2.18	0.43
1:J:233:THR:H	1:J:236:THR:HG22	1.82	0.43
2:E:66:VAL:HG13	2:E:70:PHE:CG	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:58:ILE:O	1:G:62:ILE:HD13	2.18	0.43
1:J:71:ALA:HB3	1:J:98:VAL:HG21	2.00	0.43
1:A:116:ALA:HA	1:A:165:LEU:HD22	2.00	0.43
1:G:57:LYS:HE2	1:G:172:LEU:CD2	2.35	0.43
2:E:162:TRP:CZ3	2:E:203:CYS:HB3	2.54	0.43
1:G:216:ALA:O	1:G:220:LEU:HG	2.18	0.43
3:I:144:TYR:CG	3:I:145:PRO:HA	2.53	0.43
1:A:62:ILE:O	1:A:65:VAL:HG22	2.18	0.43
3:F:58:LEU:HD21	3:F:62:ILE:O	2.18	0.43
2:H:51:ILE:HD13	2:H:74:ARG:HG2	2.01	0.43
1:J:46:ASN:OD1	1:J:46:ASN:N	2.50	0.43
2:K:33:TYR:O	2:K:34:MET:HE2	2.18	0.43
2:B:52:ARG:NH1	2:B:59:THR:OG1	2.49	0.43
1:D:205:VAL:HG11	1:D:217:ALA:HB2	2.00	0.43
1:D:239:LEU:O	1:D:243:ILE:HD12	2.19	0.43
2:H:127:PRO:HB3	2:H:153:TYR:HB3	2.01	0.43
3:I:13:VAL:HG12	3:I:108:LEU:HD11	1.99	0.43
1:A:242:LYS:O	1:A:246:THR:HG23	2.18	0.43
2:H:167:LEU:HD23	2:H:167:LEU:HA	1.90	0.43
1:J:84:HIS:HB2	1:J:87:ILE:HG22	2.00	0.43
3:L:95:SER:HA	3:L:100:TRP:CE3	2.53	0.43
1:D:192:LEU:HD23	1:D:197:LEU:HD11	2.01	0.43
1:D:197:LEU:C	1:D:199:GLU:OE2	2.61	0.43
1:J:165:LEU:C	1:J:165:LEU:HD23	2.43	0.43
3:F:32:ASP:C	3:F:34:ASP:H	2.26	0.43
2:B:204:ASN:ND2	2:B:215:ASP:OD1	2.45	0.43
3:F:4:MET:HE1	3:F:37:MET:HE1	2.01	0.43
1:A:225:ILE:CD1	1:A:240:LYS:HB2	2.47	0.42
1:D:240:LYS:HE3	2:E:56:ASN:OD1	2.19	0.42
2:H:196:TRP:O	4:H:301:HOH:O	2.21	0.42
1:D:187:VAL:O	1:D:191:ILE:CD1	2.66	0.42
1:A:47:LEU:HG	1:A:137:LEU:CD1	2.48	0.42
1:D:191:ILE:HG22	1:D:197:LEU:HD13	2.01	0.42
3:C:126:SER:O	3:C:130:THR:HG23	2.19	0.42
3:C:201:THR:HG22	3:C:208:PRO:HB3	2.01	0.42
1:J:59:ASN:HB3	1:J:63:LYS:HZ2	1.84	0.42
2:K:66:VAL:HG13	2:K:70:PHE:CG	2.54	0.42
1:A:200:ALA:O	1:A:204:ILE:HG13	2.19	0.42
1:A:247:LEU:O	1:A:251:LEU:HD12	2.19	0.42
1:D:50:VAL:HG13	1:D:163:ARG:NH1	2.34	0.42
1:G:164:GLN:O	1:G:168:LYS:N	2.51	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:70:GLU:OE1	1:J:74:ARG:HG2	2.20	0.42
1:D:47:LEU:HD13	1:D:47:LEU:C	2.43	0.42
1:D:101:CYS:HB3	1:D:180:CYS:SG	2.60	0.42
1:D:225:ILE:HD13	1:D:240:LYS:HB2	2.01	0.42
2:E:1:GLU:O	2:E:26:GLY:HA3	2.19	0.42
2:H:131:PRO:HD3	2:H:216:LYS:HG2	2.02	0.42
3:I:159:ARG:C	3:I:160:GLN:HG2	2.44	0.42
3:I:122:PHE:HA	3:I:123:PRO:HD3	1.94	0.42
3:C:112:ARG:HG3	3:C:112:ARG:HH11	1.84	0.42
1:G:233:THR:H	1:G:236:THR:HB	1.85	0.42
2:B:39:GLN:HB2	2:B:45:LEU:HD23	2.02	0.42
1:G:62:ILE:HD12	1:G:109:LEU:HD11	2.02	0.42
1:J:173:GLU:O	1:J:177:THR:HG23	2.20	0.42
1:J:235:GLN:OE1	1:J:235:GLN:N	2.47	0.42
1:A:65:VAL:HG12	1:A:105:ALA:HB3	2.01	0.41
1:D:109:LEU:HD23	1:D:109:LEU:HA	1.86	0.41
2:E:39:GLN:NE2	3:F:42:GLN:HE22	2.16	0.41
3:F:216:ASN:O	3:F:217:GLU:HG3	2.20	0.41
2:K:170:GLY:O	2:K:189:VAL:HA	2.20	0.41
3:L:179:MET:HG2	3:L:180:SER:N	2.34	0.41
2:H:216:LYS:HE2	3:I:127:GLU:OE1	2.19	0.41
1:A:252:VAL:HG12	1:A:252:VAL:O	2.19	0.41
1:D:187:VAL:CG1	1:D:191:ILE:HD11	2.51	0.41
1:G:63:LYS:HD3	1:G:63:LYS:HA	1.78	0.41
2:H:85:MET:HE2	2:H:88:LEU:CD2	2.50	0.41
1:J:70:GLU:O	1:J:71:ALA:C	2.63	0.41
3:L:199:GLU:HG2	3:L:210:VAL:HG22	2.01	0.41
2:B:162:TRP:CZ3	2:B:203:CYS:HB3	2.55	0.41
1:D:74:ARG:HH21	1:D:184:VAL:HG21	1.85	0.41
3:I:199:GLU:HG2	3:I:210:VAL:HG12	2.02	0.41
1:D:50:VAL:HG13	1:D:163:ARG:HH12	1.86	0.41
1:D:187:VAL:HG12	1:D:191:ILE:CD1	2.51	0.41
1:G:115:ILE:HA	1:G:118:THR:HG22	2.03	0.41
1:G:168:LYS:HA	1:G:171:ASP:HB2	2.03	0.41
2:K:58:TYR:OH	2:K:76:ASP:OD2	2.25	0.41
1:A:90:LEU:HD21	1:A:201:LEU:HB2	2.03	0.41
2:E:24:THR:CG2	2:E:79:SER:HB3	2.49	0.41
1:J:70:GLU:OE1	1:J:70:GLU:O	2.39	0.41
1:A:60:ASN:O	1:A:64:GLN:HG3	2.21	0.41
1:A:137:LEU:HD13	1:A:137:LEU:C	2.46	0.41
3:I:18:ARG:NH2	3:I:80:ARG:HH11	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:5:GLU:O	2:K:5:GLU:HG2	2.19	0.41
2:B:66:VAL:HG13	2:B:70:PHE:CG	2.56	0.41
2:E:113:GLN:HG2	2:E:114:GLY:H	1.86	0.41
2:E:127:PRO:HB3	2:E:153:TYR:HB3	2.03	0.41
1:J:241:GLU:HG2	2:K:104:TRP:HB3	2.03	0.41
2:B:66:VAL:CG1	2:B:70:PHE:HB2	2.52	0.40
1:D:197:LEU:HA	1:D:199:GLU:CD	2.46	0.40
3:F:80:ARG:CG	3:F:81:PRO:N	2.81	0.40
1:J:176:THR:O	1:J:179:ASN:HB3	2.21	0.40
1:D:53:ILE:HD12	1:D:54:ALA:N	2.35	0.40
2:H:5:GLU:O	2:H:22:CYS:HA	2.21	0.40
1:J:178:ASN:OD1	1:J:178:ASN:N	2.54	0.40
1:G:176:THR:O	1:G:176:THR:HG22	2.20	0.40
1:J:233:THR:H	1:J:236:THR:CG2	2.33	0.40
2:E:47:TRP:CZ2	2:E:49:GLY:HA2	2.57	0.40
2:E:66:VAL:HG13	2:E:70:PHE:HB2	2.03	0.40
2:H:204:ASN:ND2	2:H:215:ASP:OD1	2.43	0.40
1:G:54:ALA:O	1:G:57:LYS:HB3	2.22	0.40
2:H:99:ALA:HB3	2:H:108:MET:HE3	2.02	0.40
1:J:201:LEU:O	1:J:205:VAL:HG22	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	189/233 (81%)	185 (98%)	4 (2%)	0	100	100
1	D	157/233 (67%)	153 (98%)	4 (2%)	0	100	100
1	G	164/233 (70%)	160 (98%)	4 (2%)	0	100	100
1	J	168/233 (72%)	163 (97%)	5 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	220/229 (96%)	215 (98%)	5 (2%)	0	100	100
2	E	220/229 (96%)	215 (98%)	5 (2%)	0	100	100
2	H	220/229 (96%)	216 (98%)	4 (2%)	0	100	100
2	K	220/229 (96%)	215 (98%)	5 (2%)	0	100	100
3	C	215/217 (99%)	209 (97%)	6 (3%)	0	100	100
3	F	215/217 (99%)	208 (97%)	7 (3%)	0	100	100
3	I	215/217 (99%)	209 (97%)	6 (3%)	0	100	100
3	L	215/217 (99%)	209 (97%)	6 (3%)	0	100	100
All	All	2418/2716 (89%)	2357 (98%)	61 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	154/188 (82%)	154 (100%)	0	100	100
1	D	130/188 (69%)	129 (99%)	1 (1%)	73	86
1	G	135/188 (72%)	135 (100%)	0	100	100
1	J	138/188 (73%)	138 (100%)	0	100	100
2	B	188/194 (97%)	188 (100%)	0	100	100
2	E	188/194 (97%)	188 (100%)	0	100	100
2	H	188/194 (97%)	188 (100%)	0	100	100
2	K	188/194 (97%)	187 (100%)	1 (0%)	81	90
3	C	190/190 (100%)	189 (100%)	1 (0%)	81	90
3	F	190/190 (100%)	190 (100%)	0	100	100
3	I	190/190 (100%)	188 (99%)	2 (1%)	65	81
3	L	190/190 (100%)	190 (100%)	0	100	100
All	All	2069/2288 (90%)	2064 (100%)	5 (0%)	87	94

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	179	MET
1	D	179	ASN
3	I	73	THR
3	I	93	GLN
2	K	88	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	GLN
1	A	190	HIS
1	A	194	GLN
3	C	41	GLN
3	C	149	ASN
3	C	214	ASN
1	D	175	GLN
1	D	178	ASN
1	D	179	ASN
1	D	194	GLN
2	E	13	GLN
2	E	39	GLN
2	E	139	GLN
3	F	149	ASN
2	H	86	ASN
3	I	41	GLN
1	J	122	ASN
1	J	226	ASN
2	K	39	GLN
3	L	42	GLN
3	L	96	ASN
3	L	216	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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, 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	193/233 (82%)	0.93	22 (11%) 10 11	65, 98, 165, 185	0
1	D	161/233 (69%)	1.08	25 (15%) 5 6	74, 113, 188, 213	0
1	G	168/233 (72%)	0.80	15 (8%) 15 17	53, 97, 177, 196	0
1	J	172/233 (73%)	0.85	26 (15%) 5 6	49, 94, 183, 210	0
2	B	222/229 (96%)	0.03	5 (2%) 61 63	32, 58, 101, 148	0
2	E	222/229 (96%)	-0.10	4 (1%) 67 69	30, 54, 88, 141	1 (0%)
2	H	222/229 (96%)	-0.19	2 (0%) 81 82	32, 51, 90, 170	1 (0%)
2	K	222/229 (96%)	-0.11	6 (2%) 56 58	35, 52, 88, 165	0
3	C	217/217 (100%)	-0.24	0 100 100	32, 52, 84, 122	0
3	F	217/217 (100%)	-0.22	1 (0%) 87 87	31, 51, 86, 114	0
3	I	217/217 (100%)	-0.35	1 (0%) 87 87	30, 47, 79, 145	0
3	L	217/217 (100%)	-0.02	5 (2%) 61 63	36, 56, 103, 152	0
All	All	2450/2716 (90%)	0.15	112 (4%) 37 39	30, 60, 156, 213	2 (0%)

All (112) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	196	ALA	5.4
1	D	252	VAL	4.8
2	B	137	ALA	4.5
1	D	115	ILE	4.2
1	D	112	LEU	4.1
1	G	252	VAL	4.0
1	A	128	GLY	4.0
1	D	58	ILE	4.0
1	J	47	LEU	3.9
1	J	117	LEU	3.9
1	A	134	ILE	3.8

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Mol	Chain	Res	Type	RSRZ
2	E	113	GLN	3.8
1	A	160	VAL	3.7
1	G	48	ALA	3.6
1	J	120	ALA	3.6
1	G	162	ILE	3.6
1	D	48	ALA	3.6
1	G	119	ALA	3.6
2	E	5	GLU	3.5
1	J	112	LEU	3.5
1	J	176	THR	3.4
1	D	165	LEU	3.3
1	A	113	ALA	3.3
1	G	112	LEU	3.3
1	D	114	PRO	3.3
1	D	113	ALA	3.3
2	H	5	GLU	3.3
1	D	50	VAL	3.2
1	G	211	ALA	3.2
1	D	47	LEU	3.1
1	J	121	THR	3.1
1	A	140	ALA	3.1
1	A	124	ALA	3.0
2	K	135	GLY	3.0
1	A	136	ILE	3.0
1	J	48	ALA	2.9
2	E	137	ALA	2.9
3	L	165	ASN	2.9
1	J	82	THR	2.9
1	A	127	SER	2.8
3	L	217	GLU	2.8
3	L	216	ASN	2.8
1	A	45	LYS	2.8
1	J	252	VAL	2.8
1	J	51	ALA	2.7
2	K	5	GLU	2.7
2	B	138	ALA	2.7
2	K	137	ALA	2.7
3	I	4	MET	2.7
1	D	172	LEU	2.7
1	G	117	LEU	2.7
1	J	211	ALA	2.6
1	D	110	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	181	GLY	2.6
1	D	51	ALA	2.6
1	G	120	ALA	2.6
1	D	251	LEU	2.6
1	G	47	LEU	2.6
1	A	176	THR	2.6
1	D	55	LEU	2.5
2	K	138	ALA	2.5
1	D	54	ALA	2.5
1	J	116	ALA	2.5
2	B	179	GLN	2.5
1	A	129	HIS	2.5
1	G	166	TYR	2.5
1	D	53	ILE	2.4
1	A	117	LEU	2.4
1	A	58	ILE	2.4
1	D	103	LEU	2.4
1	J	71	ALA	2.3
1	D	116	ALA	2.3
1	G	108	ALA	2.3
2	E	138	ALA	2.3
2	B	136	SER	2.3
1	J	53	ILE	2.3
1	A	172	LEU	2.3
1	D	180	CYS	2.3
1	G	116	ALA	2.3
1	A	159	THR	2.3
1	A	120	ALA	2.2
1	G	51	ALA	2.2
1	D	177	THR	2.2
1	J	159	THR	2.2
2	H	140	THR	2.2
1	D	197	LEU	2.2
1	J	50	VAL	2.2
1	J	86	ASN	2.2
1	A	211	ALA	2.2
1	A	82	THR	2.2
1	J	119	ALA	2.2
1	D	184	VAL	2.2
2	K	222	ASP	2.2
1	J	180	CYS	2.2
1	J	162	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	138	GLN	2.1
1	A	80	ALA	2.1
1	G	53	ILE	2.1
1	J	58	ILE	2.1
1	J	110	THR	2.1
1	J	118	THR	2.1
3	L	4	MET	2.1
1	J	90	LEU	2.1
3	L	164	LEU	2.1
1	J	160	VAL	2.1
1	A	177	THR	2.0
2	B	98	CYS	2.0
3	F	33	ALA	2.0
1	J	166	TYR	2.0
1	G	82	THR	2.0
1	D	90	LEU	2.0
2	K	141	ASN	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.