



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:56 PM GMT

PDB ID : 1Z3G  
Title : Crystal structure of complex between Pvs25 and Fab fragment of malaria transmission blocking antibody 2A8  
Authors : Saxena, A.K.; Singh, K.; Su, H.P.; Klein, M.M.; Stowers, A.W.; Saul, A.J.; Long, C.A.; Garboczi, D.N.  
Deposited on : 2005-03-12  
Resolution : 3.30 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) 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

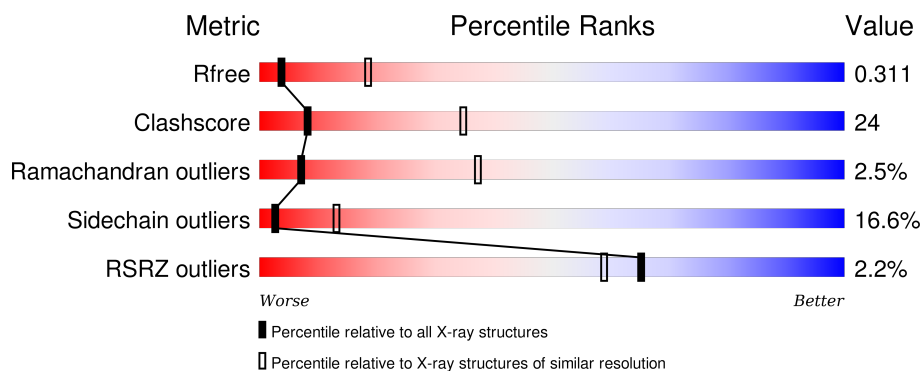
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.30 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))
$R_{free}$	91344	2060 (3.40-3.20)
Clashscore	102246	1058 (3.38-3.22)
Ramachandran outliers	100387	1038 (3.38-3.22)
Sidechain outliers	100360	1037 (3.38-3.22)
RSRZ outliers	91569	2070 (3.40-3.20)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	L	213	 70% 28% •
1	M	213	 72% 25% •
2	H	216	 47% 42% 9% ••
2	I	216	 44% 44% 11% •
3	A	186	 44% 40% 8% • 7%

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Mol	Chain	Length	Quality of chain
3	B	186	<div><div>5%</div><div><div></div><div>44%</div><div>37%</div><div>13%</div><div>7%</div></div></div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9120 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 2A8 Fab Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	213	Total	C	N	O	S	0	0	0
			1634	1011	274	339	10			
1	M	213	Total	C	N	O	S	0	0	0
			1634	1011	274	339	10			

- Molecule 2 is a protein called 2A8 Fab Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	214	Total	C	N	O	S	0	0	0
			1617	1027	267	315	8			
2	I	214	Total	C	N	O	S	0	0	0
			1617	1027	267	315	8			

- Molecule 3 is a protein called ookinete surface protein Pvs25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	173	Total	C	N	O	S	0	0	0
			1309	795	216	272	26			
3	B	173	Total	C	N	O	S	0	0	0
			1309	795	216	272	26			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLU	-	CLONING ARTIFACT	UNP O96555
A	-3	ALA	-	CLONING ARTIFACT	UNP O96555
A	-2	GLU	-	CLONING ARTIFACT	UNP O96555
A	-1	ALA	-	CLONING ARTIFACT	UNP O96555
A	0	SER	-	CLONING ARTIFACT	UNP O96555
A	174	GLY	-	CLONING ARTIFACT	UNP O96555
A	175	PRO	-	CLONING ARTIFACT	UNP O96555
A	176	HIS	-	EXPRESSION TAG	UNP O96555

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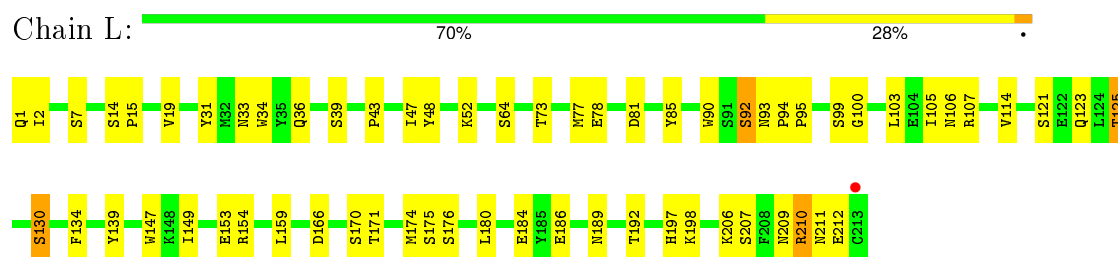
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Chain	Residue	Modelled	Actual	Comment	Reference
A	177	HIS	-	EXPRESSION TAG	UNP O96555
A	178	HIS	-	EXPRESSION TAG	UNP O96555
A	179	HIS	-	EXPRESSION TAG	UNP O96555
A	180	HIS	-	EXPRESSION TAG	UNP O96555
A	181	HIS	-	EXPRESSION TAG	UNP O96555
B	-4	GLU	-	CLONING ARTIFACT	UNP O96555
B	-3	ALA	-	CLONING ARTIFACT	UNP O96555
B	-2	GLU	-	CLONING ARTIFACT	UNP O96555
B	-1	ALA	-	CLONING ARTIFACT	UNP O96555
B	0	SER	-	CLONING ARTIFACT	UNP O96555
B	174	GLY	-	CLONING ARTIFACT	UNP O96555
B	175	PRO	-	CLONING ARTIFACT	UNP O96555
B	176	HIS	-	EXPRESSION TAG	UNP O96555
B	177	HIS	-	EXPRESSION TAG	UNP O96555
B	178	HIS	-	EXPRESSION TAG	UNP O96555
B	179	HIS	-	EXPRESSION TAG	UNP O96555
B	180	HIS	-	EXPRESSION TAG	UNP O96555
B	181	HIS	-	EXPRESSION TAG	UNP O96555

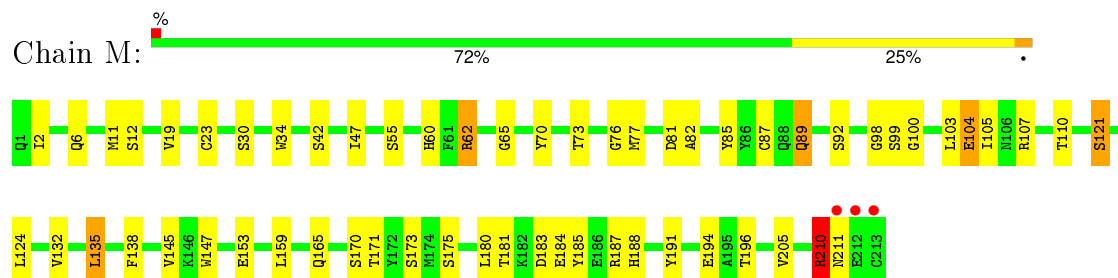
### 3 Residue-property plots

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.

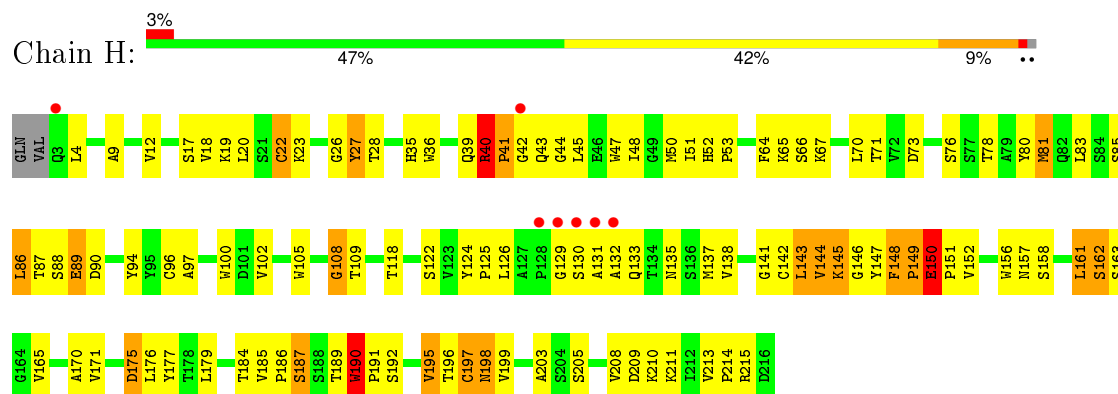
- Molecule 1: 2A8 Fab Light Chain



- Molecule 1: 2A8 Fab Light Chain

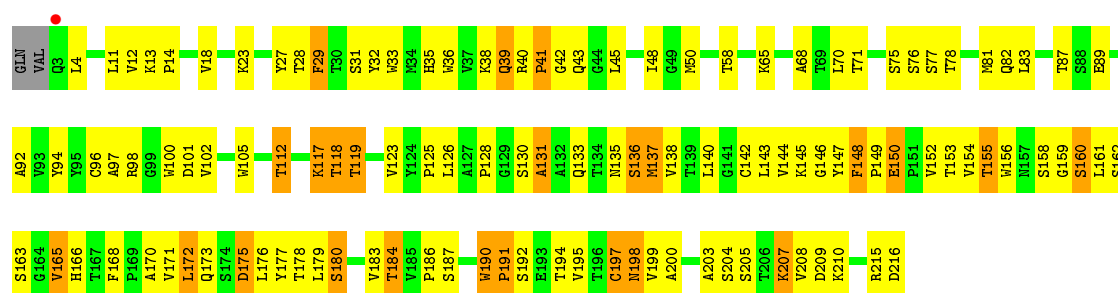


- Molecule 2: 2A8 Fab Heavy Chain

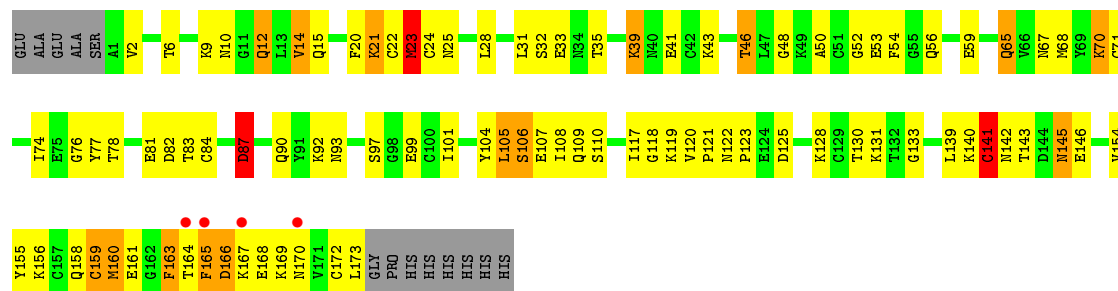


- Molecule 2: 2A8 Fab Heavy Chain

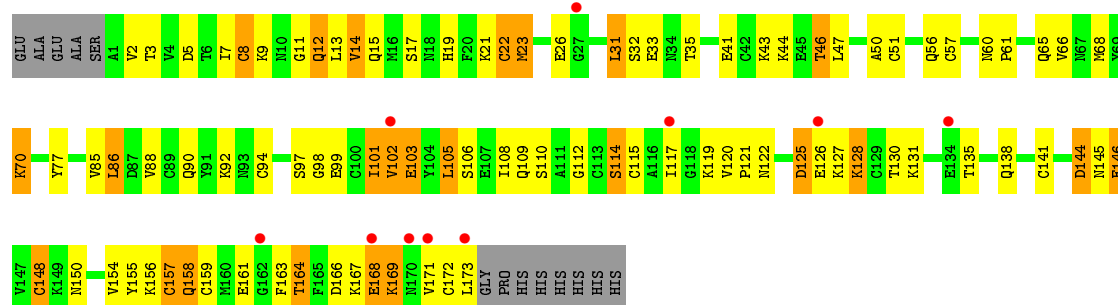
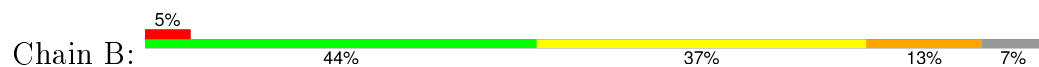




• Molecule 3: ookinete surface protein Pvs25



• Molecule 3: ookinete surface protein Pvs25



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	86.26 Å   61.65 Å   142.65 Å 90.00°   101.68°   90.00°	Depositor
Resolution (Å)	49.15 – 3.30 49.16 – 3.28	Depositor EDS
% Data completeness (in resolution range)	94.8 (49.15-3.30) 94.9 (49.16-3.28)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.28	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.38 (at 3.25 Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.272   ,   0.327 0.258   ,   0.311	Depositor DCC
$R_{free}$ test set	1095 reflections (5.38%)	DCC
Wilson B-factor (Å <sup>2</sup> )	50.7	Xtriage
Anisotropy	0.380	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.26 , 11.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.44$ , $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	1 of 21698 reflections (0.005%)	Xtriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	9120	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	2.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.16% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  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

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	L	0.52	0/1673	0.68	1/2274 (0.0%)
1	M	0.57	0/1673	0.73	1/2274 (0.0%)
2	H	0.65	1/1664 (0.1%)	0.87	3/2278 (0.1%)
2	I	0.67	0/1664	0.83	0/2278
3	A	0.91	0/1323	1.06	3/1776 (0.2%)
3	B	0.77	0/1323	0.96	2/1776 (0.1%)
All	All	0.68	1/9320 (0.0%)	0.85	10/12656 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	190	TRP	NE1-CE2	-5.15	1.30	1.37

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	215	ARG	NE-CZ-NH2	8.46	124.53	120.30
2	H	40	ARG	NE-CZ-NH2	7.33	123.96	120.30
1	L	210	ARG	NE-CZ-NH2	7.17	123.89	120.30
1	M	210	ARG	NE-CZ-NH2	6.63	123.61	120.30
3	A	87	ASP	CB-CG-OD2	6.10	123.79	118.30
3	A	160	MET	CG-SD-CE	6.10	109.95	100.20
3	B	23	MET	CG-SD-CE	5.82	109.51	100.20
3	A	23	MET	CG-SD-CE	5.51	109.02	100.20
2	H	81	MET	CG-SD-CE	5.49	108.99	100.20
3	B	144	ASP	CB-CG-OD2	5.39	123.15	118.30

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	L	1634	0	1549	50	0
1	M	1634	0	1549	34	0
2	H	1617	0	1578	116	0
2	I	1617	0	1578	105	0
3	A	1309	0	1242	79	0
3	B	1309	0	1242	72	0
All	All	9120	0	8738	433	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 24.

All (433) 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)
3:B:120:VAL:CG1	3:B:121:PRO:HD2	1.38	1.48
3:A:125:ASP:OD1	3:A:130:THR:HG21	1.35	1.21
2:I:190:TRP:HB3	2:I:191:PRO:CD	1.75	1.14
3:B:120:VAL:CG1	3:B:121:PRO:CD	2.24	1.13
2:H:148:PHE:HB3	2:H:149:PRO:HD3	1.24	1.12
2:H:190:TRP:HB3	2:H:191:PRO:HD3	1.30	1.10
3:B:120:VAL:HG12	3:B:121:PRO:HD2	1.32	1.10
2:H:40:ARG:HB3	2:H:41:PRO:HD2	1.17	1.10
3:A:142:ASN:O	3:A:146:GLU:O	1.71	1.08
2:H:138:VAL:HG23	2:H:187:SER:HA	1.32	1.08
3:B:120:VAL:HG13	3:B:121:PRO:CD	1.84	1.06
2:I:190:TRP:HB3	2:I:191:PRO:HD3	1.32	1.05
1:L:125:THR:HG21	2:I:119:THR:HG22	1.37	1.04
2:H:40:ARG:HB3	2:H:41:PRO:CD	1.87	1.03
2:H:186:PRO:O	2:H:189:THR:HG22	1.58	1.01
1:L:149:ILE:HD12	1:L:154:ARG:HG3	1.43	1.00
3:B:122:ASN:HD22	3:B:130:THR:HB	1.27	0.99
2:H:213:VAL:HG13	2:H:214:PRO:HD2	1.43	0.98
2:I:153:THR:HG23	2:I:200:ALA:HB3	1.45	0.97
2:H:138:VAL:CG2	2:H:187:SER:HA	1.94	0.96
3:B:120:VAL:HG13	3:B:121:PRO:HD2	0.97	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:148:PHE:HB3	2:H:149:PRO:CD	1.97	0.94
2:I:172:LEU:O	2:I:173:GLN:HG3	1.67	0.94
2:H:40:ARG:HD2	2:H:41:PRO:HD3	1.49	0.93
3:B:120:VAL:HG12	3:B:121:PRO:CD	1.95	0.92
1:L:192:THR:HG23	1:L:207:SER:HB3	1.53	0.91
2:I:190:TRP:O	2:I:192:SER:N	2.01	0.91
3:A:139:LEU:HB3	3:A:141:CYS:SG	2.10	0.91
2:I:190:TRP:CB	2:I:191:PRO:HD3	1.99	0.91
2:H:190:TRP:HB3	2:H:191:PRO:CD	2.00	0.91
2:I:137:MET:HA	2:I:186:PRO:HA	1.52	0.91
2:I:172:LEU:CD1	2:I:173:GLN:H	1.83	0.91
3:A:25:ASN:HB2	3:A:28:LEU:HD12	1.51	0.90
1:L:210:ARG:NH1	1:L:211:ASN:OD1	2.02	0.90
2:I:190:TRP:CB	2:I:191:PRO:CD	2.48	0.90
2:I:172:LEU:HD12	2:I:173:GLN:H	1.35	0.90
3:A:65:GLN:HB2	3:A:68:MET:HG3	1.53	0.89
3:B:94:CYS:O	3:B:128:LYS:HG2	1.72	0.89
1:M:2:ILE:HD11	1:M:92:SER:HB2	1.59	0.85
2:I:190:TRP:C	2:I:192:SER:H	1.79	0.84
3:A:43:LYS:HG2	3:A:46:THR:HG23	1.57	0.84
3:B:120:VAL:HG22	3:B:135:THR:HG22	1.60	0.84
3:B:98:GLY:HA3	3:B:114:SER:O	1.79	0.83
2:H:149:PRO:C	2:H:151:PRO:HD2	1.99	0.82
3:B:166:ASP:HB2	3:B:171:VAL:O	1.78	0.82
2:I:138:VAL:HG23	2:I:187:SER:HA	1.60	0.82
3:A:12:GLN:OE1	3:A:156:LYS:HD2	1.79	0.81
3:B:156:LYS:HG2	3:B:157:CYS:N	1.95	0.81
2:I:27:TYR:CE1	2:I:29:PHE:HA	2.16	0.81
3:A:43:LYS:HG2	3:A:46:THR:CG2	2.11	0.81
3:B:19:HIS:HB3	3:B:101:ILE:HD13	1.63	0.80
3:B:164:THR:O	3:B:173:LEU:N	2.13	0.80
3:A:14:VAL:HG13	3:A:154:VAL:HG21	1.64	0.80
2:H:125:PRO:HD3	2:H:210:LYS:HD3	1.63	0.79
1:L:123:GLN:HE22	1:L:130:SER:HB2	1.47	0.79
3:A:92:LYS:CG	3:A:93:ASN:H	1.94	0.79
3:A:41:GLU:O	3:A:46:THR:HG21	1.83	0.78
3:A:165:PHE:HE1	3:A:170:ASN:HA	1.48	0.78
2:H:198:ASN:HB3	2:H:209:ASP:OD2	1.83	0.78
2:I:172:LEU:HD12	2:I:173:GLN:N	1.99	0.77
3:A:32:SER:HB3	3:A:35:THR:HB	1.66	0.76
2:H:9:ALA:HB2	2:H:151:PRO:HG3	1.68	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:147:TYR:CE2	2:H:152:VAL:HG21	2.20	0.76
2:H:213:VAL:CG1	2:H:214:PRO:HD2	2.14	0.76
2:I:190:TRP:HB3	2:I:191:PRO:HD2	1.68	0.76
2:H:190:TRP:CB	2:H:191:PRO:HD3	2.12	0.75
2:H:142:CYS:HB2	2:H:156:TRP:CH2	2.21	0.75
1:M:132:VAL:HG21	2:I:126:LEU:HD21	1.66	0.75
3:B:126:GLU:O	3:B:127:LYS:HB2	1.87	0.74
3:A:23:MET:O	3:A:23:MET:HG3	1.86	0.74
2:H:40:ARG:CB	2:H:41:PRO:HD2	2.08	0.74
3:B:156:LYS:HG2	3:B:157:CYS:H	1.53	0.74
2:I:190:TRP:CG	2:I:191:PRO:HD3	2.22	0.73
2:H:150:GLU:HG2	2:H:177:TYR:CZ	2.23	0.72
1:L:210:ARG:HG3	1:L:210:ARG:HH11	1.55	0.72
2:H:9:ALA:HB2	2:H:151:PRO:CG	2.19	0.71
3:B:168:GLU:O	3:B:168:GLU:OE1	2.08	0.71
1:L:90:TRP:CH2	2:H:50:MET:HE1	2.25	0.71
3:B:19:HIS:CD2	3:B:101:ILE:HG21	2.25	0.71
3:A:158:GLN:HG3	3:A:159:CYS:H	1.55	0.71
2:H:9:ALA:CB	2:H:151:PRO:CG	2.69	0.71
2:I:130:SER:O	2:I:131:ALA:CB	2.38	0.70
3:B:166:ASP:OD1	3:B:169:LYS:HB2	1.92	0.70
1:M:183:ASP:O	1:M:187:ARG:HG3	1.91	0.69
3:B:103:GLU:OE2	3:B:112:GLY:HA3	1.92	0.69
3:A:163:PHE:CG	3:A:172:CYS:HB3	2.27	0.69
3:B:43:LYS:HG2	3:B:46:THR:HG22	1.73	0.69
2:H:147:TYR:CE2	2:H:152:VAL:CG2	2.75	0.69
2:H:175:ASP:O	2:H:176:LEU:HD23	1.92	0.69
3:A:120:VAL:HB	3:A:121:PRO:HD2	1.74	0.69
3:A:15:GLN:O	3:A:154:VAL:HG23	1.92	0.69
1:M:34:TRP:HB2	1:M:47:ILE:HB	1.74	0.68
2:H:9:ALA:CB	2:H:151:PRO:HG2	2.22	0.68
2:I:130:SER:O	2:I:131:ALA:HB3	1.92	0.68
1:L:125:THR:HG21	2:I:119:THR:CG2	2.22	0.68
3:A:14:VAL:HG12	3:A:21:LYS:HG3	1.76	0.68
2:H:151:PRO:O	2:H:152:VAL:HG23	1.92	0.68
3:B:98:GLY:CA	3:B:114:SER:O	2.42	0.68
3:A:139:LEU:CB	3:A:141:CYS:SG	2.82	0.68
2:I:138:VAL:CG2	2:I:187:SER:HA	2.24	0.67
2:I:12:VAL:HG21	2:I:18:VAL:HG22	1.76	0.67
2:H:190:TRP:C	2:H:190:TRP:CD1	2.65	0.67
3:B:117:ILE:N	3:B:117:ILE:HD13	2.09	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:23:LYS:HG2	2:H:78:THR:HG22	1.78	0.67
2:H:86:LEU:HD23	2:H:90:ASP:HB2	1.77	0.66
3:A:122:ASN:HD22	3:A:125:ASP:HB2	1.59	0.66
3:A:70:LYS:HG3	3:A:71:CYS:N	2.11	0.66
3:A:65:GLN:CB	3:A:68:MET:HG3	2.25	0.66
3:A:92:LYS:HG3	3:A:93:ASN:H	1.60	0.66
2:I:40:ARG:HD2	2:I:41:PRO:HD2	1.77	0.66
2:H:185:VAL:HB	2:H:186:PRO:HD2	1.78	0.66
3:B:101:ILE:HG12	3:B:114:SER:HB2	1.78	0.66
2:I:190:TRP:CD2	2:I:191:PRO:HD3	2.31	0.65
2:I:146:GLY:HA2	2:I:176:LEU:HD23	1.79	0.65
3:A:92:LYS:CG	3:A:93:ASN:N	2.59	0.65
3:A:145:ASN:HB2	3:A:172:CYS:SG	2.37	0.65
2:H:195:VAL:HG13	2:H:195:VAL:O	1.97	0.65
2:I:147:TYR:OH	2:I:170:ALA:HB2	1.96	0.65
2:H:48:ILE:HG21	2:H:81:MET:CE	2.26	0.65
3:B:108:ILE:HG13	3:B:109:GLN:N	2.12	0.65
3:A:74:ILE:O	3:A:77:TYR:HB2	1.97	0.65
2:H:36:TRP:CD1	2:H:70:LEU:CD2	2.80	0.64
2:I:172:LEU:CD1	2:I:173:GLN:N	2.57	0.64
2:H:130:SER:O	2:H:131:ALA:HB3	1.97	0.64
3:B:12:GLN:O	3:B:22:CYS:HA	1.97	0.64
3:A:122:ASN:N	3:A:128:LYS:O	2.31	0.64
1:M:210:ARG:HG2	1:M:211:ASN:ND2	2.13	0.64
1:M:60:HIS:NE2	1:M:81:ASP:OD2	2.27	0.64
2:H:40:ARG:CB	2:H:41:PRO:CD	2.62	0.64
1:M:107:ARG:HD3	1:M:171:THR:HG22	1.80	0.64
2:H:208:VAL:HG22	2:I:208:VAL:HG22	1.80	0.63
2:I:4:LEU:HD21	2:I:96:CYS:HB3	1.81	0.63
2:I:204:SER:O	2:I:205:SER:OG	2.14	0.63
2:H:124:TYR:HB2	2:H:143:LEU:HD23	1.79	0.63
2:I:39:GLN:O	2:I:92:ALA:HB1	1.98	0.63
3:A:125:ASP:CG	3:A:130:THR:HG21	2.19	0.63
2:I:190:TRP:CE3	2:I:191:PRO:HD3	2.33	0.63
2:I:147:TYR:OH	2:I:170:ALA:CB	2.46	0.63
1:L:149:ILE:CD1	1:L:154:ARG:HG3	2.25	0.63
3:B:120:VAL:HG12	3:B:121:PRO:N	2.13	0.62
2:H:175:ASP:N	2:H:175:ASP:OD1	2.32	0.62
2:H:36:TRP:CD1	2:H:70:LEU:HD21	2.35	0.62
2:H:170:ALA:HA	2:H:179:LEU:HB3	1.80	0.62
1:L:77:MET:HE1	1:L:103:LEU:HD21	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:149:PRO:HG2	2:H:203:ALA:HB3	1.82	0.61
2:H:209:ASP:HB2	2:I:207:LYS:HB2	1.81	0.61
2:H:40:ARG:HD2	2:H:41:PRO:CD	2.25	0.61
2:H:94:TYR:O	2:H:108:GLY:HA2	2.00	0.61
2:I:48:ILE:HG21	2:I:81:MET:CE	2.30	0.61
3:A:122:ASN:HB2	3:A:131:LYS:HB2	1.83	0.61
1:L:159:LEU:HD11	2:H:171:VAL:HG11	1.84	0.60
2:I:190:TRP:C	2:I:192:SER:N	2.51	0.60
1:M:110:THR:HG23	1:M:138:PHE:HA	1.84	0.60
2:I:154:VAL:O	2:I:155:THR:HG23	2.02	0.59
3:A:164:THR:O	3:A:173:LEU:N	2.35	0.59
1:L:77:MET:CE	1:L:103:LEU:HD21	2.32	0.59
3:B:23:MET:HG3	3:B:23:MET:O	2.03	0.59
2:H:149:PRO:O	2:H:151:PRO:HD2	2.03	0.59
3:A:118:GLY:O	3:A:133:GLY:O	2.21	0.59
2:I:172:LEU:HD13	2:I:173:GLN:H	1.66	0.58
1:L:93:ASN:OD1	1:L:94:PRO:HA	2.03	0.58
2:H:190:TRP:CB	2:H:191:PRO:CD	2.67	0.58
2:H:149:PRO:HG2	2:H:203:ALA:CB	2.33	0.58
2:I:27:TYR:CZ	2:I:98:ARG:HD2	2.39	0.58
1:L:36:GLN:HG3	1:L:85:TYR:CE2	2.39	0.58
2:H:48:ILE:HG23	2:H:64:PHE:CD2	2.39	0.58
1:M:180:LEU:HD22	1:M:184:GLU:HG2	1.85	0.58
1:L:43:PRO:HD2	2:H:105:TRP:CE3	2.39	0.58
3:B:120:VAL:HG22	3:B:135:THR:CG2	2.32	0.57
3:B:125:ASP:O	3:B:126:GLU:HB2	2.03	0.57
3:B:161:GLU:C	3:B:163:PHE:H	2.07	0.57
2:I:138:VAL:HG23	2:I:187:SER:CA	2.34	0.57
3:A:165:PHE:CE1	3:A:170:ASN:HA	2.36	0.57
2:H:9:ALA:CB	2:H:151:PRO:HG3	2.31	0.57
3:B:108:ILE:HG13	3:B:109:GLN:H	1.70	0.57
2:I:158:SER:H	2:I:198:ASN:HD21	1.54	0.56
2:I:92:ALA:HB3	2:I:94:TYR:CE2	2.40	0.56
1:M:194:GLU:HG2	1:M:205:VAL:HG22	1.87	0.56
2:I:125:PRO:HG3	2:I:210:LYS:HG2	1.86	0.56
1:M:104:GLU:HG2	1:M:165:GLN:OE1	2.04	0.56
3:A:106:SER:O	3:A:107:GLU:OE1	2.23	0.56
1:L:159:LEU:HD11	2:H:171:VAL:HG21	1.88	0.56
3:B:14:VAL:HG13	3:B:154:VAL:HG22	1.88	0.56
1:L:174:MET:HE2	1:L:176:SER:HB2	1.88	0.56
1:L:106:ASN:HA	1:L:139:TYR:OH	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:15:GLN:HG3	3:A:20:PHE:CZ	2.42	0.55
2:I:18:VAL:O	2:I:82:GLN:HA	2.05	0.55
3:B:126:GLU:O	3:B:127:LYS:CB	2.53	0.55
2:I:12:VAL:HG21	2:I:18:VAL:CG2	2.35	0.55
2:I:38:LYS:HB2	2:I:48:ILE:HD11	1.87	0.55
2:I:168:PHE:CD2	2:I:168:PHE:N	2.75	0.55
1:L:189:ASN:O	1:L:209:ASN:HA	2.07	0.54
2:I:136:SER:C	2:I:137:MET:HG2	2.28	0.54
3:B:14:VAL:HG13	3:B:154:VAL:CG2	2.37	0.54
2:I:149:PRO:HG2	2:I:203:ALA:CB	2.37	0.54
2:I:154:VAL:HG12	2:I:199:VAL:HG22	1.88	0.54
2:I:48:ILE:HG21	2:I:81:MET:HE1	1.89	0.54
2:H:150:GLU:HG2	2:H:177:TYR:CE2	2.42	0.54
3:B:146:GLU:HA	3:B:158:GLN:O	2.07	0.54
3:B:8:CYS:CB	3:B:22:CYS:SG	2.95	0.54
3:A:25:ASN:HB2	3:A:28:LEU:CD1	2.32	0.54
3:B:102:VAL:HG12	3:B:103:GLU:H	1.72	0.54
2:I:156:TRP:CZ3	2:I:197:CYS:HB3	2.43	0.54
2:H:148:PHE:O	2:H:177:TYR:HD2	1.92	0.53
3:A:122:ASN:ND2	3:A:125:ASP:HB2	2.23	0.53
3:A:15:GLN:HG3	3:A:20:PHE:CE2	2.42	0.53
2:H:23:LYS:HE3	2:H:76:SER:O	2.08	0.53
3:A:65:GLN:HB2	3:A:68:MET:CG	2.31	0.53
3:A:120:VAL:HB	3:A:121:PRO:CD	2.40	0.53
2:H:129:GLY:O	2:H:132:ALA:HB2	2.09	0.53
3:A:46:THR:OG1	3:A:46:THR:O	2.26	0.52
1:M:62:ARG:HG2	1:M:73:THR:HB	1.91	0.52
3:B:31:LEU:HD22	3:B:50:ALA:HB2	1.90	0.52
2:I:76:SER:OG	2:I:78:THR:HG23	2.10	0.52
2:I:136:SER:O	2:I:186:PRO:HB3	2.10	0.52
3:A:10:ASN:HB3	3:A:28:LEU:HD13	1.92	0.52
3:B:60:ASN:ND2	3:B:61:PRO:HD2	2.24	0.52
2:H:26:GLY:O	2:H:27:TYR:HB3	2.09	0.52
3:A:166:ASP:C	3:A:168:GLU:H	2.12	0.52
3:A:32:SER:OG	3:A:33:GLU:N	2.43	0.52
2:I:149:PRO:CG	2:I:203:ALA:CB	2.88	0.52
2:H:126:LEU:HB2	2:H:141:GLY:HA3	1.92	0.52
2:I:148:PHE:O	2:I:149:PRO:C	2.48	0.52
2:I:32:TYR:CZ	3:B:47:LEU:HD13	2.45	0.52
2:I:100:TRP:CE3	3:B:66:VAL:HA	2.45	0.52
3:A:139:LEU:C	3:A:141:CYS:H	2.14	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:14:VAL:HG22	3:A:154:VAL:HG22	1.92	0.51
3:B:163:PHE:CE1	3:B:172:CYS:O	2.63	0.51
3:A:31:LEU:HB2	3:A:50:ALA:HB2	1.92	0.51
2:H:129:GLY:O	2:H:130:SER:C	2.49	0.51
2:H:142:CYS:HB2	2:H:156:TRP:CZ2	2.45	0.51
2:I:168:PHE:H	2:I:168:PHE:HD2	1.53	0.51
2:I:149:PRO:CG	2:I:203:ALA:HB1	2.41	0.51
2:H:196:THR:HA	2:H:211:LYS:HA	1.93	0.51
3:B:60:ASN:HD22	3:B:70:LYS:HB2	1.74	0.51
2:H:129:GLY:O	2:H:132:ALA:N	2.44	0.51
2:H:158:SER:H	2:H:198:ASN:HD21	1.59	0.51
2:I:145:LYS:HD2	2:I:178:THR:OG1	2.10	0.51
1:M:34:TRP:CZ3	1:M:87:CYS:HB3	2.46	0.51
2:I:100:TRP:CZ3	3:B:66:VAL:HA	2.46	0.51
3:B:115:CYS:HB2	3:B:119:LYS:O	2.11	0.51
2:I:40:ARG:HB2	2:I:43:GLN:HB2	1.93	0.50
1:M:159:LEU:HD21	2:I:171:VAL:HG11	1.92	0.50
1:M:2:ILE:CD1	1:M:92:SER:HB2	2.37	0.50
1:L:123:GLN:HA	2:H:124:TYR:CE1	2.47	0.50
2:I:68:ALA:HB2	2:I:83:LEU:HD23	1.93	0.49
3:B:12:GLN:HE22	3:B:156:LYS:HE2	1.76	0.49
2:H:147:TYR:CE2	2:H:152:VAL:HG23	2.47	0.49
1:L:31:TYR:OH	1:L:90:TRP:HB3	2.13	0.49
1:M:85:TYR:O	1:M:100:GLY:HA2	2.12	0.49
2:I:36:TRP:CD1	2:I:70:LEU:CD2	2.96	0.49
3:B:125:ASP:O	3:B:126:GLU:CB	2.61	0.49
3:A:108:ILE:HD12	3:A:109:GLN:H	1.76	0.49
2:H:148:PHE:O	2:H:177:TYR:CD2	2.65	0.49
1:M:210:ARG:CZ	1:M:211:ASN:OD1	2.61	0.49
3:A:14:VAL:HG21	3:A:154:VAL:HG13	1.94	0.49
2:H:144:VAL:HG11	2:H:199:VAL:HG11	1.95	0.49
1:L:184:GLU:OE2	1:L:184:GLU:HA	2.13	0.49
1:L:186:GLU:HA	1:L:210:ARG:HE	1.77	0.49
2:H:156:TRP:CE3	2:H:197:CYS:HB2	2.48	0.49
3:A:81:GLU:O	3:A:82:ASP:HB2	2.13	0.49
2:H:137:MET:HB3	2:H:184:THR:HG22	1.95	0.49
2:I:48:ILE:HG21	2:I:81:MET:HE2	1.94	0.48
2:H:35:HIS:HB2	2:H:97:ALA:HB3	1.95	0.48
3:A:92:LYS:HG2	3:A:93:ASN:H	1.76	0.48
3:B:2:VAL:HG21	3:B:21:LYS:HA	1.95	0.48
1:M:185:TYR:HA	1:M:191:TYR:OH	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:139:LEU:O	3:A:141:CYS:SG	2.71	0.48
1:L:210:ARG:O	1:L:211:ASN:HB2	2.13	0.48
1:L:125:THR:CG2	2:I:119:THR:HG22	2.26	0.48
1:M:2:ILE:HD12	1:M:89:GLN:NE2	2.29	0.48
3:A:14:VAL:HG21	3:A:154:VAL:CG1	2.44	0.48
1:M:121:SER:HA	1:M:124:LEU:HD12	1.94	0.48
2:I:149:PRO:HG3	2:I:203:ALA:HB1	1.94	0.48
2:H:145:LYS:HG3	2:H:146:GLY:N	2.29	0.48
2:H:125:PRO:CD	2:H:210:LYS:HD3	2.39	0.48
1:M:23:CYS:HB2	1:M:34:TRP:CH2	2.49	0.47
2:H:130:SER:O	2:H:131:ALA:CB	2.60	0.47
1:L:147:TRP:O	1:L:153:GLU:HA	2.14	0.47
2:H:19:LYS:HG2	2:H:80:TYR:HB3	1.96	0.47
1:L:31:TYR:HE1	1:L:33:ASN:HD21	1.61	0.47
2:H:17:SER:HA	2:H:86:LEU:HD12	1.96	0.47
1:L:36:GLN:HG3	1:L:85:TYR:CZ	2.50	0.47
3:A:122:ASN:HB2	3:A:131:LYS:CB	2.44	0.47
1:M:6:GLN:OE1	1:M:98:GLY:HA3	2.14	0.47
2:H:17:SER:OG	2:H:83:LEU:O	2.21	0.47
2:I:117:LYS:HG3	2:I:118:THR:H	1.79	0.47
3:A:142:ASN:HB3	3:A:145:ASN:OD1	2.14	0.47
1:L:90:TRP:HH2	2:H:50:MET:HE1	1.76	0.47
1:L:78:GLU:O	1:L:81:ASP:HB2	2.15	0.47
1:M:173:SER:HG	2:I:166:HIS:CE1	2.31	0.47
2:I:179:LEU:HG	2:I:180:SER:H	1.80	0.47
3:B:77:TYR:HA	3:B:85:VAL:O	2.15	0.46
3:B:8:CYS:HB3	3:B:22:CYS:SG	2.55	0.46
1:M:210:ARG:NH1	1:M:211:ASN:HD21	2.14	0.46
3:A:125:ASP:O	3:A:128:LYS:HD2	2.14	0.46
3:B:8:CYS:HB2	3:B:11:GLY:O	2.16	0.46
2:I:29:PHE:C	2:I:31:SER:H	2.18	0.46
2:I:159:GLY:O	2:I:161:LEU:N	2.48	0.46
3:A:48:GLY:O	3:A:56:GLN:NE2	2.43	0.46
2:H:151:PRO:O	2:H:152:VAL:CG2	2.60	0.46
1:L:210:ARG:HG3	1:L:210:ARG:NH1	2.25	0.46
1:L:206:LYS:HA	1:L:206:LYS:HD3	1.77	0.46
3:B:94:CYS:O	3:B:128:LYS:CG	2.56	0.46
2:H:156:TRP:CZ3	2:H:197:CYS:HB2	2.51	0.46
1:L:34:TRP:HB2	1:L:47:ILE:HB	1.97	0.46
3:B:60:ASN:HA	3:B:61:PRO:HD3	1.84	0.46
3:A:104:TYR:HE1	3:A:109:GLN:HG2	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:11:LEU:HD23	2:I:112:THR:HB	1.98	0.46
2:I:146:GLY:CA	2:I:176:LEU:HD23	2.45	0.46
2:H:151:PRO:C	2:H:152:VAL:HG23	2.36	0.46
2:I:41:PRO:HG2	2:I:42:GLY:H	1.80	0.46
1:L:14:SER:HA	1:L:106:ASN:HB2	1.97	0.46
3:B:51:CYS:HB3	3:B:57:CYS:SG	2.55	0.46
1:M:135:LEU:HD21	1:M:145:VAL:HG21	1.96	0.46
3:A:76:GLY:C	3:A:87:ASP:OD1	2.54	0.46
3:A:142:ASN:HB2	3:A:146:GLU:HB2	1.98	0.45
2:H:48:ILE:HG21	2:H:81:MET:HE3	1.97	0.45
2:H:149:PRO:CG	2:H:203:ALA:CB	2.94	0.45
3:A:166:ASP:C	3:A:168:GLU:N	2.69	0.45
1:L:114:VAL:HA	1:L:134:PHE:O	2.16	0.45
2:H:149:PRO:C	2:H:151:PRO:CD	2.79	0.45
3:B:120:VAL:CG2	3:B:135:THR:HG22	2.40	0.45
2:I:179:LEU:HG	2:I:180:SER:N	2.30	0.45
2:H:138:VAL:HG23	2:H:187:SER:CA	2.23	0.45
1:L:90:TRP:CZ2	2:H:50:MET:HE1	2.51	0.45
2:I:33:TRP:CE3	2:I:50:MET:HG3	2.52	0.45
3:B:122:ASN:CG	3:B:125:ASP:HB2	2.37	0.45
2:H:148:PHE:O	2:H:150:GLU:N	2.49	0.45
2:I:40:ARG:HD3	2:I:92:ALA:HB2	1.99	0.45
3:A:14:VAL:HG23	3:A:156:LYS:HB2	1.98	0.45
2:I:147:TYR:OH	2:I:170:ALA:HB1	2.17	0.45
1:L:90:TRP:CZ2	2:H:50:MET:CE	3.00	0.45
1:M:19:VAL:O	1:M:73:THR:HA	2.18	0.44
2:I:35:HIS:CD2	2:I:35:HIS:N	2.83	0.44
3:B:148:CYS:HA	3:B:156:LYS:O	2.17	0.44
2:H:189:THR:CG2	2:H:190:TRP:N	2.80	0.44
2:I:27:TYR:CD1	2:I:29:PHE:HA	2.51	0.44
3:A:39:LYS:HD2	3:A:52:GLY:HA2	1.99	0.44
3:B:77:TYR:CE1	3:B:86:LEU:HG	2.52	0.44
1:M:11:MET:HG3	1:M:103:LEU:HD12	1.98	0.44
1:L:48:TYR:CE2	1:L:52:LYS:HD2	2.53	0.44
2:H:209:ASP:CB	2:I:207:LYS:HB2	2.48	0.44
2:I:175:ASP:O	2:I:176:LEU:HD12	2.17	0.44
2:I:147:TYR:CZ	2:I:177:TYR:HB3	2.53	0.44
1:L:95:PRO:HD2	2:H:47:TRP:CE3	2.53	0.44
2:H:52:HIS:HA	2:H:53:PRO:HD3	1.83	0.44
2:I:172:LEU:C	2:I:173:GLN:HG3	2.36	0.44
3:A:14:VAL:CG2	3:A:154:VAL:HG13	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:66:SER:OG	2:H:67:LYS:N	2.51	0.44
1:L:123:GLN:NE2	1:L:130:SER:HB2	2.23	0.44
3:B:122:ASN:ND2	3:B:131:LYS:HG3	2.32	0.44
2:H:86:LEU:HD23	2:H:90:ASP:CB	2.45	0.44
3:B:166:ASP:HB2	3:B:171:VAL:C	2.37	0.43
3:B:164:THR:O	3:B:172:CYS:HA	2.18	0.43
3:A:122:ASN:OD1	3:A:123:PRO:HD2	2.18	0.43
2:H:20:LEU:O	2:H:80:TYR:HA	2.17	0.43
2:H:41:PRO:HB2	2:H:42:GLY:H	1.53	0.43
2:I:29:PHE:CE2	2:I:77:SER:HA	2.54	0.43
3:A:158:GLN:CG	3:A:159:CYS:H	2.24	0.43
3:A:166:ASP:O	3:A:168:GLU:N	2.51	0.43
2:I:135:ASN:OD1	2:I:136:SER:N	2.52	0.43
3:B:19:HIS:CG	3:B:101:ILE:HG21	2.53	0.43
3:B:166:ASP:O	3:B:169:LYS:C	2.57	0.43
1:M:147:TRP:O	1:M:153:GLU:HA	2.19	0.43
3:B:161:GLU:C	3:B:163:PHE:N	2.71	0.43
2:I:156:TRP:CD1	2:I:165:VAL:HG21	2.54	0.43
2:H:4:LEU:HD21	2:H:22:CYS:SG	2.59	0.43
3:A:54:PHE:HD1	3:A:74:ILE:HD11	1.84	0.43
2:H:129:GLY:O	2:H:132:ALA:CB	2.66	0.43
2:H:185:VAL:HB	2:H:186:PRO:CD	2.48	0.42
2:H:40:ARG:HH22	2:H:89:GLU:HA	1.84	0.42
1:L:180:LEU:HD22	1:L:184:GLU:HG3	2.00	0.42
2:H:195:VAL:CG1	2:H:195:VAL:O	2.64	0.42
2:H:36:TRP:CD1	2:H:70:LEU:HD22	2.54	0.42
3:B:15:GLN:HB3	3:B:155:TYR:HB2	2.00	0.42
1:M:188:HIS:O	1:M:210:ARG:HD3	2.19	0.42
2:H:100:TRP:CD1	3:A:59:GLU:HG2	2.55	0.42
2:H:149:PRO:O	2:H:151:PRO:CD	2.67	0.42
2:I:136:SER:O	2:I:137:MET:HG2	2.19	0.42
3:A:70:LYS:HG3	3:A:71:CYS:H	1.84	0.42
1:M:77:MET:HE1	1:M:103:LEU:HD21	2.01	0.42
1:L:107:ARG:HD3	1:L:171:THR:HG22	2.01	0.42
2:I:133:GLN:HG3	2:I:133:GLN:O	2.19	0.42
1:L:166:ASP:O	1:L:170:SER:HA	2.19	0.42
2:H:213:VAL:CG1	2:H:214:PRO:CD	2.90	0.42
1:L:159:LEU:HD21	2:H:171:VAL:HG11	2.00	0.42
3:A:117:ILE:HG12	3:A:154:VAL:HA	2.02	0.42
2:H:73:ASP:CG	2:H:76:SER:HG	2.24	0.42
2:I:11:LEU:HD12	2:I:117:LYS:HE3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:22:CYS:O	3:A:24:CYS:N	2.53	0.42
2:H:157:ASN:HB2	2:H:161:LEU:HD12	2.01	0.42
1:L:139:TYR:O	1:L:197:HIS:HE1	2.02	0.41
2:I:13:LYS:HA	2:I:14:PRO:HD2	1.80	0.41
3:B:65:GLN:HB3	3:B:68:MET:HG3	2.02	0.41
2:H:189:THR:HG23	2:H:190:TRP:N	2.35	0.41
3:B:105:LEU:H	3:B:105:LEU:HG	1.67	0.41
2:H:94:TYR:O	2:H:108:GLY:CA	2.67	0.41
1:L:15:PRO:HD3	1:L:106:ASN:O	2.20	0.41
2:I:32:TYR:CD1	2:I:32:TYR:N	2.88	0.41
3:A:53:GLU:O	3:A:84:CYS:HB2	2.20	0.41
3:B:12:GLN:O	3:B:22:CYS:CA	2.67	0.41
1:M:34:TRP:CH2	1:M:87:CYS:HB3	2.55	0.41
2:I:40:ARG:HH21	2:I:89:GLU:HA	1.86	0.41
2:H:12:VAL:CG2	2:H:18:VAL:HG21	2.51	0.41
1:M:82:ALA:HB2	1:M:105:ILE:HB	2.02	0.41
3:A:14:VAL:HG22	3:A:155:TYR:O	2.21	0.41
3:B:41:GLU:O	3:B:46:THR:HG21	2.20	0.41
3:A:120:VAL:HG22	3:A:133:GLY:N	2.35	0.41
2:I:96:CYS:O	2:I:97:ALA:HB2	2.21	0.41
2:I:123:VAL:HB	2:I:210:LYS:HD3	2.03	0.41
1:M:65:GLY:HA3	1:M:70:TYR:HA	2.02	0.41
2:I:105:TRP:N	2:I:105:TRP:CD1	2.87	0.41
2:I:199:VAL:HG12	2:I:200:ALA:N	2.35	0.41
1:L:19:VAL:O	1:L:73:THR:HA	2.20	0.41
3:A:163:PHE:CB	3:A:172:CYS:HB3	2.50	0.41
1:L:123:GLN:HE22	1:L:130:SER:CB	2.26	0.41
1:L:85:TYR:O	1:L:100:GLY:HA2	2.21	0.41
2:I:161:LEU:O	2:I:162:SER:C	2.59	0.41
3:B:60:ASN:ND2	3:B:61:PRO:CD	2.84	0.41
3:A:105:LEU:HA	3:A:105:LEU:HD23	1.76	0.41
3:A:163:PHE:HA	3:A:163:PHE:HD1	1.71	0.40
2:H:22:CYS:O	2:H:78:THR:HA	2.20	0.40
2:H:162:SER:O	2:H:165:VAL:HG12	2.21	0.40
3:A:125:ASP:HB3	3:A:128:LYS:HB2	2.03	0.40
2:H:148:PHE:CB	2:H:149:PRO:CD	2.71	0.40
1:M:12:SER:HB3	1:M:104:GLU:HB3	2.04	0.40
2:I:183:VAL:HG22	2:I:184:THR:N	2.36	0.40
2:H:23:LYS:HA	2:H:78:THR:HA	2.03	0.40
2:I:149:PRO:HG2	2:I:203:ALA:HB3	2.02	0.40
1:L:2:ILE:CD1	1:L:92:SER:HB2	2.52	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:150:GLU:HG2	2:I:177:TYR:CZ	2.56	0.40
3:B:14:VAL:CG1	3:B:154:VAL:CG2	2.99	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	L	211/213 (99%)	197 (93%)	14 (7%)	0	100	100
1	M	211/213 (99%)	197 (93%)	12 (6%)	2 (1%)	21	60
2	H	212/216 (98%)	177 (84%)	22 (10%)	13 (6%)	2	14
2	I	212/216 (98%)	184 (87%)	20 (9%)	8 (4%)	4	26
3	A	171/186 (92%)	143 (84%)	24 (14%)	4 (2%)	8	39
3	B	171/186 (92%)	136 (80%)	32 (19%)	3 (2%)	11	46
All	All	1188/1230 (97%)	1034 (87%)	124 (10%)	30 (2%)	7	37

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	41	PRO
2	I	190	TRP
3	B	128	LYS
2	H	44	GLY
2	H	108	GLY
2	H	163	SER
1	M	210	ARG
2	I	128	PRO
2	I	131	ALA
2	I	160	SER

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Mol	Chain	Res	Type
3	B	167	LYS
2	H	148	PHE
2	I	65	LYS
3	A	23	MET
3	A	141	CYS
2	H	27	TYR
2	H	150	GLU
2	I	148	PHE
2	I	191	PRO
3	A	140	LYS
3	A	167	LYS
2	H	65	LYS
2	H	149	PRO
3	B	17	SER
2	H	40	ARG
2	H	190	TRP
1	M	76	GLY
2	I	41	PRO
2	H	102	VAL
2	H	195	VAL

### 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	L	188/188 (100%)	175 (93%)	13 (7%)	19	57
1	M	188/188 (100%)	175 (93%)	13 (7%)	19	57
2	H	183/185 (99%)	152 (83%)	31 (17%)	2	12
2	I	183/185 (99%)	144 (79%)	39 (21%)	1	5
3	A	153/163 (94%)	120 (78%)	33 (22%)	1	5
3	B	153/163 (94%)	108 (71%)	45 (29%)	0	1
All	All	1048/1072 (98%)	874 (83%)	174 (17%)	3	13

All (174) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	1	GLN
1	L	7	SER
1	L	39	SER
1	L	64	SER
1	L	92	SER
1	L	99	SER
1	L	105	ILE
1	L	121	SER
1	L	125	THR
1	L	130	SER
1	L	175	SER
1	L	198	LYS
1	L	212	GLU
2	H	22	CYS
2	H	28	THR
2	H	39	GLN
2	H	40	ARG
2	H	43	GLN
2	H	45	LEU
2	H	51	ILE
2	H	71	THR
2	H	85	SER
2	H	86	LEU
2	H	87	THR
2	H	88	SER
2	H	89	GLU
2	H	96	CYS
2	H	109	THR
2	H	118	THR
2	H	122	SER
2	H	133	GLN
2	H	135	ASN
2	H	143	LEU
2	H	144	VAL
2	H	145	LYS
2	H	150	GLU
2	H	161	LEU
2	H	162	SER
2	H	175	ASP
2	H	187	SER
2	H	192	SER
2	H	197	CYS
2	H	198	ASN

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Mol	Chain	Res	Type
2	H	205	SER
1	M	30	SER
1	M	42	SER
1	M	55	SER
1	M	62	ARG
1	M	89	GLN
1	M	99	SER
1	M	104	GLU
1	M	121	SER
1	M	135	LEU
1	M	170	SER
1	M	175	SER
1	M	181	THR
1	M	196	THR
2	I	23	LYS
2	I	28	THR
2	I	29	PHE
2	I	39	GLN
2	I	45	LEU
2	I	58	THR
2	I	71	THR
2	I	75	SER
2	I	87	THR
2	I	101	ASP
2	I	102	VAL
2	I	112	THR
2	I	117	LYS
2	I	118	THR
2	I	119	THR
2	I	136	SER
2	I	137	MET
2	I	140	LEU
2	I	142	CYS
2	I	143	LEU
2	I	144	VAL
2	I	150	GLU
2	I	152	VAL
2	I	155	THR
2	I	160	SER
2	I	163	SER
2	I	165	VAL
2	I	172	LEU

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Mol	Chain	Res	Type
2	I	175	ASP
2	I	180	SER
2	I	184	THR
2	I	194	THR
2	I	195	VAL
2	I	197	CYS
2	I	198	ASN
2	I	207	LYS
2	I	209	ASP
2	I	215	ARG
2	I	216	ASP
3	A	2	VAL
3	A	6	THR
3	A	9	LYS
3	A	12	GLN
3	A	14	VAL
3	A	21	LYS
3	A	23	MET
3	A	39	LYS
3	A	46	THR
3	A	65	GLN
3	A	67	ASN
3	A	70	LYS
3	A	78	THR
3	A	83	THR
3	A	87	ASP
3	A	90	GLN
3	A	97	SER
3	A	99	GLU
3	A	101	ILE
3	A	105	LEU
3	A	106	SER
3	A	110	SER
3	A	119	LYS
3	A	141	CYS
3	A	143	THR
3	A	145	ASN
3	A	159	CYS
3	A	160	MET
3	A	161	GLU
3	A	163	PHE
3	A	165	PHE

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Mol	Chain	Res	Type
3	A	166	ASP
3	A	169	LYS
3	B	3	THR
3	B	5	ASP
3	B	7	ILE
3	B	8	CYS
3	B	9	LYS
3	B	12	GLN
3	B	13	LEU
3	B	14	VAL
3	B	22	CYS
3	B	26	GLU
3	B	31	LEU
3	B	32	SER
3	B	33	GLU
3	B	35	THR
3	B	44	LYS
3	B	46	THR
3	B	56	GLN
3	B	70	LYS
3	B	86	LEU
3	B	88	VAL
3	B	90	GLN
3	B	92	LYS
3	B	97	SER
3	B	99	GLU
3	B	101	ILE
3	B	102	VAL
3	B	103	GLU
3	B	105	LEU
3	B	106	SER
3	B	110	SER
3	B	114	SER
3	B	125	ASP
3	B	138	GLN
3	B	141	CYS
3	B	144	ASP
3	B	145	ASN
3	B	146	GLU
3	B	148	CYS
3	B	150	ASN
3	B	157	CYS

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Mol	Chain	Res	Type
3	B	158	GLN
3	B	159	CYS
3	B	164	THR
3	B	168	GLU
3	B	169	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
3	B	142	ASN
3	B	145	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	L	213/213 (100%)	-0.25	1 (0%) 91 90	2, 2, 2, 2	0
1	M	213/213 (100%)	-0.08	3 (1%) 78 73	2, 2, 2, 2	0
2	H	214/216 (99%)	-0.26	7 (3%) 50 43	2, 2, 2, 2	0
2	I	214/216 (99%)	-0.25	1 (0%) 91 90	2, 2, 2, 2	0
3	A	173/186 (93%)	-0.01	4 (2%) 64 57	2, 2, 2, 2	0
3	B	173/186 (93%)	0.38	10 (5%) 26 21	2, 2, 2, 2	0
All	All	1200/1230 (97%)	-0.10	26 (2%) 65 59	2, 2, 2, 2	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	M	213	CYS	6.7
3	B	170	ASN	6.1
2	H	131	ALA	4.4
2	H	3	GLN	4.0
3	A	170	ASN	3.7
2	I	3	GLN	3.5
3	A	164	THR	3.4
1	M	212	GLU	3.1
2	H	130	SER	3.0
2	H	129	GLY	3.0
3	A	167	LYS	2.9
3	B	162	GLY	2.9
3	B	171	VAL	2.8
3	B	173	LEU	2.8
3	B	134	GLU	2.7
1	L	213	CYS	2.7
1	M	211	ASN	2.7
3	B	168	GLU	2.6
2	H	128	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
2	H	132	ALA	2.4
2	H	42	GLY	2.4
3	B	27	GLY	2.2
3	B	126	GLU	2.2
3	A	165	PHE	2.1
3	B	102	VAL	2.0
3	B	117	ILE	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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.