



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:32 PM UTC

PDB ID : 3D5N / pdb_00003d5n
Title : Crystal structure of the Q97W15_SULSO protein from Sulfolobus solfataricus. NESG target SsR125.
Authors : Vorobiev, S.M.; Chen, Y.; Seetharaman, J.; Lee, D.; Foote, R.E.; Maglaqui, M.; Janjua, H.; Xiao, R.; Acton, T.B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2008-05-16
Resolution : 2.80 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

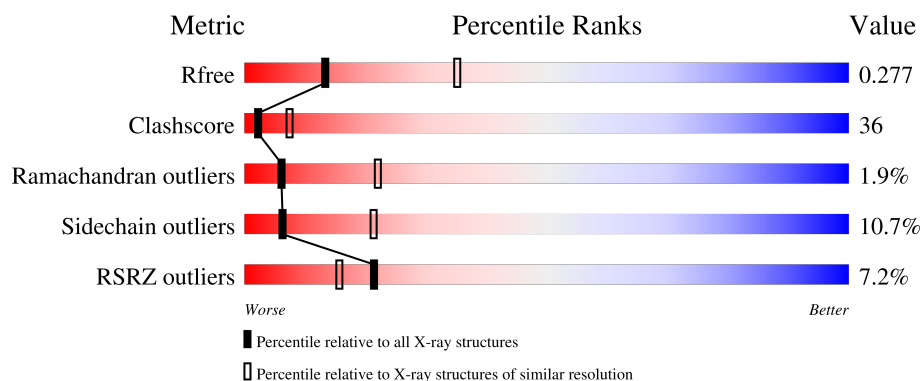
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	197	5% 42% 40% 7% 10%
1	B	197	9% 44% 34% 6% 15%
1	D	197	5% 42% 40% 8% 10%
1	E	197	6% 39% 39% 9% 13%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	197	4%41%41%5%13%
1	G	197	4%35%43%6%16%
1	H	197	7%29%46%13%12%
1	I	197	9%39%37%10%13%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10496 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 Q97W15_SULSO.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	178	Total	C	N	O	S	Se	0	0	0
			1356	882	223	243	3	5			
1	B	167	Total	C	N	O	S	Se	0	0	0
			1268	831	201	229	3	4			
1	D	178	Total	C	N	O	S	Se	0	0	0
			1373	894	226	245	3	5			
1	E	172	Total	C	N	O	S	Se	0	0	0
			1300	848	208	236	3	5			
1	F	172	Total	C	N	O	S	Se	0	0	0
			1292	847	214	223	3	5			
1	G	166	Total	C	N	O	S	Se	0	0	0
			1269	829	207	225	3	5			
1	H	173	Total	C	N	O	S	Se	0	0	0
			1315	855	212	241	3	4			
1	I	171	Total	C	N	O	S	Se	0	0	0
			1290	842	206	235	3	4			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	190	LEU	-	expression tag	UNP Q97W15
A	191	GLU	-	expression tag	UNP Q97W15
A	192	HIS	-	expression tag	UNP Q97W15
A	193	HIS	-	expression tag	UNP Q97W15
A	194	HIS	-	expression tag	UNP Q97W15
A	195	HIS	-	expression tag	UNP Q97W15
A	196	HIS	-	expression tag	UNP Q97W15
A	197	HIS	-	expression tag	UNP Q97W15
B	190	LEU	-	expression tag	UNP Q97W15
B	191	GLU	-	expression tag	UNP Q97W15
B	192	HIS	-	expression tag	UNP Q97W15
B	193	HIS	-	expression tag	UNP Q97W15
B	194	HIS	-	expression tag	UNP Q97W15

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	195	HIS	-	expression tag	UNP Q97W15
B	196	HIS	-	expression tag	UNP Q97W15
B	197	HIS	-	expression tag	UNP Q97W15
D	190	LEU	-	expression tag	UNP Q97W15
D	191	GLU	-	expression tag	UNP Q97W15
D	192	HIS	-	expression tag	UNP Q97W15
D	193	HIS	-	expression tag	UNP Q97W15
D	194	HIS	-	expression tag	UNP Q97W15
D	195	HIS	-	expression tag	UNP Q97W15
D	196	HIS	-	expression tag	UNP Q97W15
D	197	HIS	-	expression tag	UNP Q97W15
E	190	LEU	-	expression tag	UNP Q97W15
E	191	GLU	-	expression tag	UNP Q97W15
E	192	HIS	-	expression tag	UNP Q97W15
E	193	HIS	-	expression tag	UNP Q97W15
E	194	HIS	-	expression tag	UNP Q97W15
E	195	HIS	-	expression tag	UNP Q97W15
E	196	HIS	-	expression tag	UNP Q97W15
E	197	HIS	-	expression tag	UNP Q97W15
F	190	LEU	-	expression tag	UNP Q97W15
F	191	GLU	-	expression tag	UNP Q97W15
F	192	HIS	-	expression tag	UNP Q97W15
F	193	HIS	-	expression tag	UNP Q97W15
F	194	HIS	-	expression tag	UNP Q97W15
F	195	HIS	-	expression tag	UNP Q97W15
F	196	HIS	-	expression tag	UNP Q97W15
F	197	HIS	-	expression tag	UNP Q97W15
G	190	LEU	-	expression tag	UNP Q97W15
G	191	GLU	-	expression tag	UNP Q97W15
G	192	HIS	-	expression tag	UNP Q97W15
G	193	HIS	-	expression tag	UNP Q97W15
G	194	HIS	-	expression tag	UNP Q97W15
G	195	HIS	-	expression tag	UNP Q97W15
G	196	HIS	-	expression tag	UNP Q97W15
G	197	HIS	-	expression tag	UNP Q97W15
H	190	LEU	-	expression tag	UNP Q97W15
H	191	GLU	-	expression tag	UNP Q97W15
H	192	HIS	-	expression tag	UNP Q97W15
H	193	HIS	-	expression tag	UNP Q97W15
H	194	HIS	-	expression tag	UNP Q97W15
H	195	HIS	-	expression tag	UNP Q97W15
H	196	HIS	-	expression tag	UNP Q97W15

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
H	197	HIS	-	expression tag	UNP Q97W15
I	190	LEU	-	expression tag	UNP Q97W15
I	191	GLU	-	expression tag	UNP Q97W15
I	192	HIS	-	expression tag	UNP Q97W15
I	193	HIS	-	expression tag	UNP Q97W15
I	194	HIS	-	expression tag	UNP Q97W15
I	195	HIS	-	expression tag	UNP Q97W15
I	196	HIS	-	expression tag	UNP Q97W15
I	197	HIS	-	expression tag	UNP Q97W15

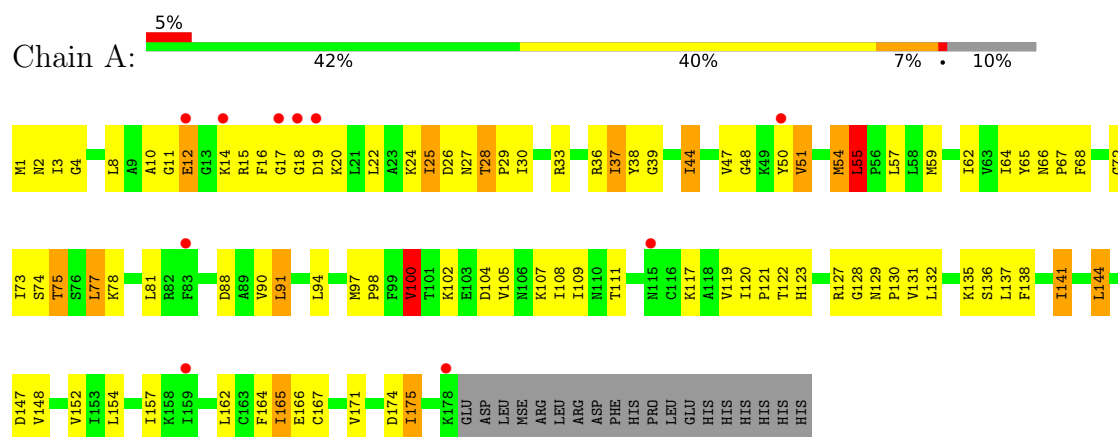
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	3	Total O 3 3	0	0
2	B	4	Total O 4 4	0	0
2	D	6	Total O 6 6	0	0
2	E	4	Total O 4 4	0	0
2	F	4	Total O 4 4	0	0
2	G	2	Total O 2 2	0	0
2	H	5	Total O 5 5	0	0
2	I	5	Total O 5 5	0	0

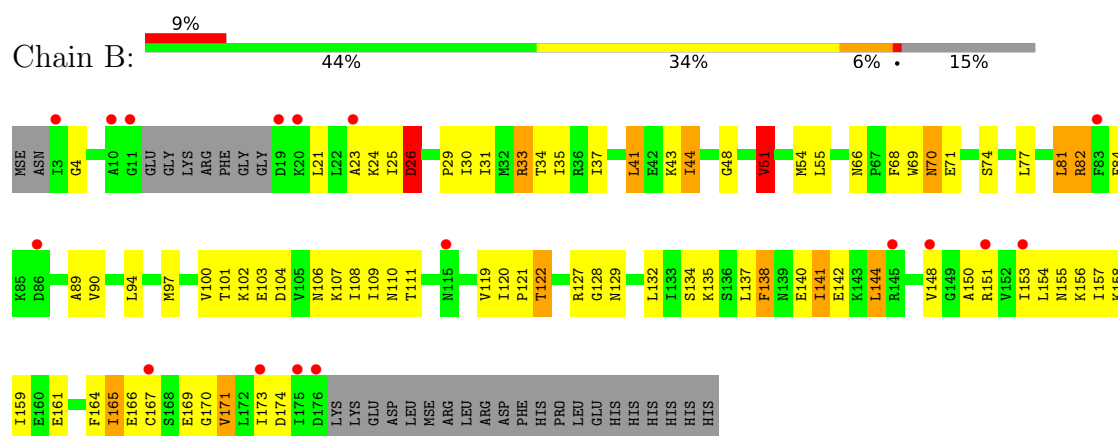
3 Residue-property plots [i](#)

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.

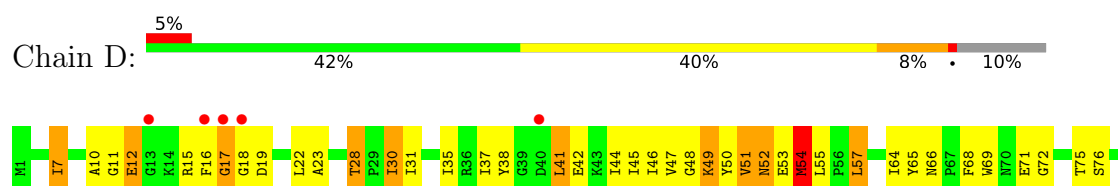
• Molecule 1: Q97W15_SULSO

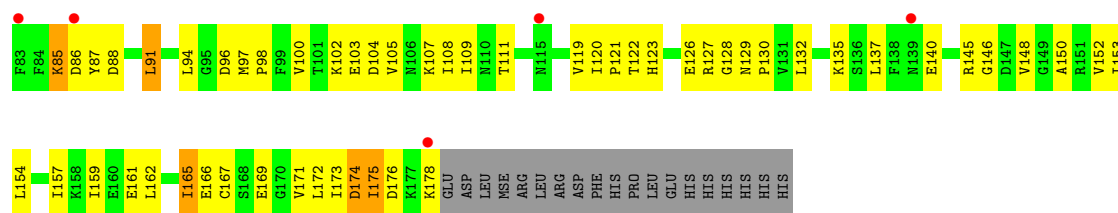


• Molecule 1: Q97W15_SULSO

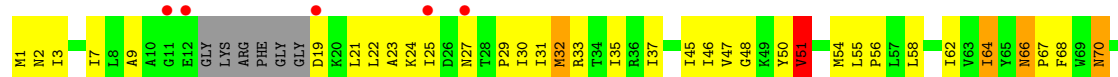


• Molecule 1: Q97W15_SULSO

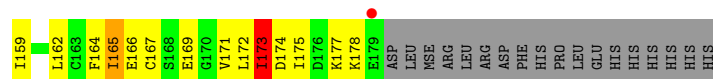
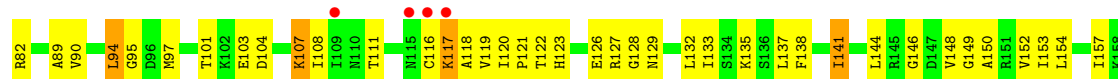
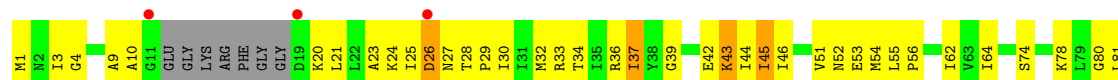




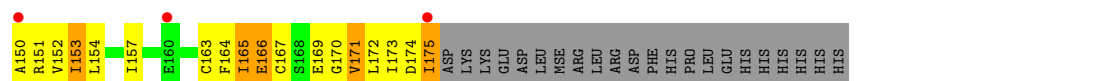
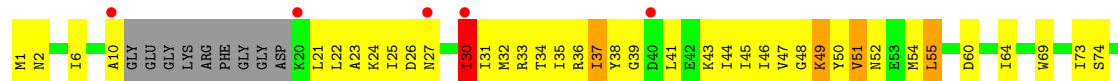
- Molecule 1: Q97W15 SULSO



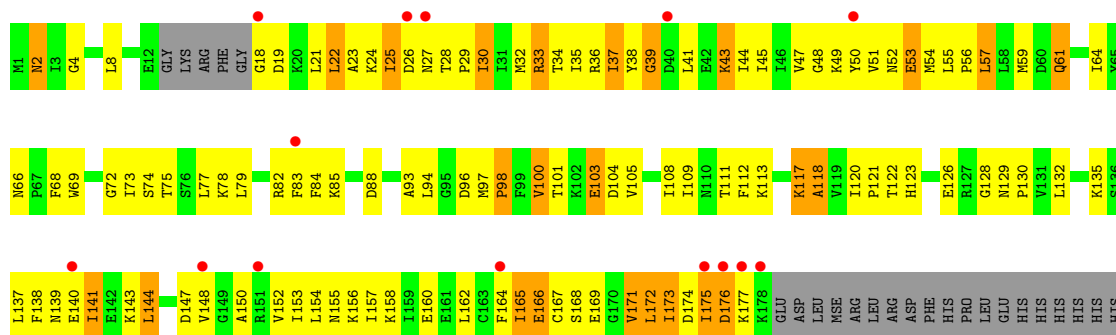
- Molecule 1: Q97W15 SULSO



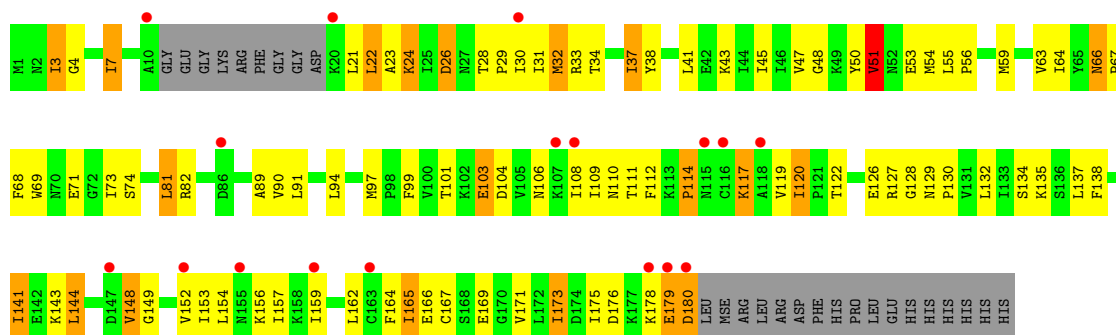
- Molecule 1: Q97W15 SULSO



● Molecule 1: Q97W15_SULSO



● Molecule 1: Q97W15_SULSO



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.96Å 169.23Å 173.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.47 – 2.80 48.47 – 2.80	Depositor EDS
% Data completeness (in resolution range)	86.2 (48.47-2.80) 95.1 (48.47-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 2.77Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.231 , 0.266 0.244 , 0.277	Depositor DCC
R_{free} test set	6108 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	74.0	Xtriage
Anisotropy	0.088	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 68.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.034 for -h,l,k	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10496	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 3.00% 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	0.53	0/1373	1.05	5/1852 (0.3%)
1	B	0.50	0/1283	1.02	6/1734 (0.3%)
1	D	0.50	0/1390	1.03	5/1870 (0.3%)
1	E	0.56	0/1315	1.06	5/1778 (0.3%)
1	F	0.52	1/1307 (0.1%)	1.05	6/1765 (0.3%)
1	G	0.51	0/1284	1.06	6/1736 (0.3%)
1	H	0.58	1/1330 (0.1%)	1.05	4/1797 (0.2%)
1	I	0.49	0/1305	1.04	10/1767 (0.6%)
All	All	0.52	2/10587 (0.0%)	1.04	47/14299 (0.3%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	26	ASP	CA-CB	6.82	1.63	1.53
1	H	177	LYS	CA-CB	5.61	1.61	1.53

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	10	ALA	N-CA-C	-10.06	100.55	113.12
1	A	54	MSE	N-CA-C	8.58	120.25	111.07
1	E	176	ASP	N-CA-C	-8.24	103.01	113.23
1	I	54	MSE	N-CA-C	8.23	120.33	111.36
1	D	17	GLY	N-CA-C	8.12	121.91	112.50
1	E	62	ILE	N-CA-C	-7.45	99.06	108.89
1	E	173	ILE	N-CA-C	7.37	118.90	108.36
1	E	117	LYS	N-CA-C	-7.34	104.32	113.28
1	G	60	ASP	N-CA-C	-6.80	104.60	113.17
1	F	173	ILE	N-CA-C	6.78	118.81	108.71
1	D	150	ALA	N-CA-C	6.77	119.52	111.33
1	H	19	ASP	N-CA-C	-6.71	103.97	111.28
1	B	24	LYS	N-CA-C	6.61	119.18	108.41
1	F	117	LYS	N-CA-C	-6.58	105.21	113.18

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	37	ILE	N-CA-C	-6.52	104.20	113.07
1	E	51	VAL	CB-CA-C	-6.32	103.80	111.88
1	A	14	LYS	N-CA-C	-6.08	105.78	112.72
1	A	55	LEU	N-CA-C	6.00	121.05	113.25
1	F	43	LYS	N-CA-C	5.88	118.01	109.07
1	I	26	ASP	N-CA-C	-5.74	103.48	111.52
1	I	117	LYS	N-CA-C	-5.73	106.29	113.28
1	G	30	ILE	N-CA-C	5.69	117.11	110.62
1	H	117	LYS	N-CA-C	-5.63	105.76	113.30
1	B	51	VAL	CB-CA-C	-5.62	104.77	111.97
1	H	173	ILE	N-CA-C	5.62	116.67	108.58
1	F	116	CYS	N-CA-C	5.58	117.18	109.15
1	F	150	ALA	N-CA-C	5.45	117.98	111.71
1	A	100	VAL	N-CA-C	-5.38	101.89	109.63
1	I	120	ILE	N-CA-C	5.38	113.19	107.76
1	B	33	ARG	N-CA-C	-5.37	105.50	111.36
1	D	54	MSE	N-CA-C	5.32	116.89	111.14
1	A	24	LYS	N-CA-C	5.30	118.30	110.46
1	B	138	PHE	N-CA-C	5.25	116.69	111.07
1	G	117	LYS	N-CA-C	-5.16	107.03	113.38
1	D	52	ASN	N-CA-C	5.13	117.54	111.33
1	D	19	ASP	N-CA-C	-5.12	105.40	111.69
1	B	173	ILE	N-CA-C	5.10	115.82	108.48
1	I	51	VAL	CB-CA-C	-5.08	105.38	111.88
1	I	3	ILE	N-CA-C	5.08	115.89	108.53
1	G	113	LYS	CA-C-N	5.08	124.79	119.82
1	G	113	LYS	C-N-CA	5.08	124.79	119.82
1	I	156	LYS	N-CA-C	-5.07	107.77	114.31
1	H	49	LYS	N-CA-C	5.05	117.18	111.02
1	B	142	GLU	N-CA-C	-5.04	106.40	112.54
1	I	173	ILE	N-CA-C	5.03	116.53	108.89
1	I	50	TYR	N-CA-C	5.01	118.88	112.26
1	I	63	VAL	N-CA-C	5.01	115.06	107.80

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	1356	0	1391	83	0
1	B	1268	0	1300	79	0
1	D	1373	0	1430	94	0
1	E	1300	0	1324	106	0
1	F	1292	0	1335	99	0
1	G	1269	0	1309	101	0
1	H	1315	0	1339	128	0
1	I	1290	0	1305	118	0
2	A	3	0	0	0	0
2	B	4	0	0	1	0
2	D	6	0	0	0	0
2	E	4	0	0	0	0
2	F	4	0	0	0	0
2	G	2	0	0	0	0
2	H	5	0	0	0	0
2	I	5	0	0	0	0
All	All	10496	0	10733	771	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 36.

All (771) 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:G:111:THR:HG21	1:G:165:ILE:HG21	1.20	1.10
1:A:18:GLY:HA3	1:A:50:TYR:HB2	1.31	1.09
1:H:66:ASN:HD21	1:H:75:THR:HG23	1.17	1.09
1:E:68:PHE:HB2	1:E:75:THR:HG21	1.34	1.08
1:G:81:LEU:HD22	1:G:141:ILE:HD11	1.37	1.06
1:E:157:ILE:HD11	1:E:161:GLU:HB2	1.37	1.06
1:I:97:MSE:HE1	1:I:129:ASN:HB2	1.41	1.01
1:E:66:ASN:ND2	1:E:68:PHE:H	1.61	0.97
1:E:70:ASN:HD22	1:E:70:ASN:H	1.03	0.96
1:A:157:ILE:HG13	1:A:162:LEU:HD11	1.48	0.95
1:D:157:ILE:HD11	1:D:162:LEU:HD22	1.50	0.93
1:H:74:SER:HB2	1:H:144:LEU:HD23	1.52	0.92
1:B:70:ASN:H	1:B:70:ASN:HD22	1.18	0.92
1:I:21:LEU:HD23	1:I:30:ILE:HD11	1.51	0.91
1:I:24:LYS:HD3	1:I:24:LYS:H	1.37	0.89

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:66:ASN:ND2	1:H:75:THR:HG23	1.88	0.89
1:H:56:PRO:HA	1:H:59:MSE:CE	2.03	0.88
1:D:66:ASN:ND2	1:D:68:PHE:H	1.71	0.88
1:H:56:PRO:HA	1:H:59:MSE:HE2	1.53	0.87
1:I:66:ASN:ND2	1:I:68:PHE:H	1.74	0.86
1:F:29:PRO:HB2	1:F:32:MSE:HB2	1.57	0.85
1:A:66:ASN:OD1	1:A:75:THR:HG23	1.75	0.85
1:G:101:THR:HG22	1:G:103:GLU:H	1.41	0.85
1:H:111:THR:OG1	1:H:165:ILE:HG21	1.76	0.85
1:E:117:LYS:HB3	1:E:137:LEU:HD11	1.60	0.84
1:E:131:VAL:HG11	1:E:154:LEU:HD11	1.60	0.84
1:F:97:MSE:HE1	1:F:129:ASN:HB2	1.59	0.83
1:G:2:ASN:HB2	1:G:88:ASP:H	1.44	0.83
1:I:97:MSE:HE1	1:I:129:ASN:CB	2.08	0.82
1:E:103:GLU:CD	1:E:103:GLU:H	1.83	0.82
1:H:21:LEU:HB2	1:H:54:MSE:HE1	1.61	0.82
1:A:33:ARG:HD3	1:A:98:PRO:HB2	1.60	0.82
1:B:111:THR:HG21	1:B:165:ILE:HD11	1.59	0.82
1:B:127:ARG:H	1:B:127:ARG:HD2	1.43	0.82
1:G:141:ILE:HG22	1:G:153:ILE:HG21	1.60	0.82
1:I:148:VAL:HG23	1:I:149:GLY:H	1.45	0.82
1:F:94:LEU:HB2	1:F:97:MSE:HE2	1.62	0.81
1:D:22:LEU:HD11	1:D:53:GLU:HB3	1.61	0.80
1:F:101:THR:OG1	1:F:103:GLU:HG2	1.81	0.80
1:E:30:ILE:HG13	1:E:175:ILE:HD13	1.63	0.80
1:E:70:ASN:H	1:E:70:ASN:ND2	1.79	0.80
1:D:128:GLY:HA3	1:D:171:VAL:HG12	1.62	0.80
1:G:144:LEU:HD21	1:G:149:GLY:HA2	1.63	0.79
1:D:85:LYS:HE2	1:D:85:LYS:C	2.08	0.79
1:E:31:ILE:O	1:E:35:ILE:HG12	1.82	0.79
1:D:49:LYS:HG2	1:D:69:TRP:CH2	2.16	0.79
1:B:70:ASN:HD22	1:B:70:ASN:N	1.78	0.78
1:D:45:ILE:HG22	1:D:47:VAL:HG13	1.66	0.78
1:B:66:ASN:ND2	1:B:68:PHE:H	1.81	0.77
1:I:101:THR:HB	1:I:104:ASP:OD1	1.85	0.77
1:I:24:LYS:HD3	1:I:24:LYS:N	1.99	0.77
1:F:157:ILE:HD11	1:F:162:LEU:HD13	1.65	0.76
1:D:102:LYS:HE3	1:I:126:GLU:OE2	1.85	0.76
1:G:24:LYS:HD3	1:G:27:ASN:HA	1.65	0.76
1:D:11:GLY:HA3	1:D:17:GLY:HA2	1.68	0.76
1:G:111:THR:HG21	1:G:165:ILE:CG2	2.11	0.75

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:174:ASP:OD1	1:H:176:ASP:HB3	1.86	0.75
1:F:97:MSE:HE1	1:F:129:ASN:CB	2.16	0.74
1:G:81:LEU:HD22	1:G:141:ILE:CD1	2.15	0.74
1:F:148:VAL:HG23	1:F:149:GLY:H	1.51	0.74
1:I:3:ILE:HD12	1:I:3:ILE:O	1.88	0.73
1:F:175:ILE:HD11	1:I:173:ILE:HD12	1.71	0.73
1:B:70:ASN:H	1:B:70:ASN:ND2	1.86	0.73
1:B:141:ILE:HG22	1:B:153:ILE:HG13	1.70	0.73
1:F:107:LYS:HB3	1:F:165:ILE:HD12	1.70	0.72
1:G:101:THR:HG22	1:G:103:GLU:N	2.03	0.72
1:A:55:LEU:HD13	1:B:55:LEU:HD11	1.72	0.72
1:D:7:ILE:HG12	1:D:45:ILE:HG13	1.71	0.72
1:H:33:ARG:HG3	1:H:33:ARG:HH11	1.54	0.72
1:B:97:MSE:HE1	1:B:129:ASN:CB	2.20	0.71
1:H:44:ILE:HD11	1:H:64:ILE:CD1	2.20	0.71
1:A:25:ILE:O	1:A:26:ASP:HB2	1.91	0.71
1:B:34:THR:O	1:B:37:ILE:HG12	1.90	0.71
1:H:94:LEU:HD22	1:H:97:MSE:HE3	1.73	0.71
1:B:138:PHE:O	1:B:141:ILE:HD13	1.90	0.70
1:H:45:ILE:HD13	1:H:61:GLN:HG2	1.72	0.70
1:A:167:CYS:HB2	1:A:171:VAL:HG21	1.73	0.70
1:B:158:LYS:HB2	1:B:161:GLU:HG3	1.73	0.70
1:A:18:GLY:HA3	1:A:50:TYR:CB	2.16	0.70
1:H:144:LEU:HD12	1:H:153:ILE:HD11	1.74	0.69
1:I:141:ILE:O	1:I:144:LEU:HD22	1.92	0.69
1:D:45:ILE:HG22	1:D:47:VAL:CG1	2.22	0.69
1:G:45:ILE:HG22	1:G:47:VAL:HG13	1.73	0.69
1:D:104:ASP:OD1	1:D:167:CYS:HB3	1.94	0.68
1:D:175:ILE:HD13	1:D:175:ILE:O	1.93	0.68
1:H:94:LEU:HD22	1:H:97:MSE:CE	2.23	0.68
1:E:165:ILE:HD13	1:E:165:ILE:H	1.59	0.68
1:A:131:VAL:HG11	1:A:154:LEU:HD11	1.75	0.68
1:I:101:THR:HG22	1:I:103:GLU:H	1.59	0.68
1:B:122:THR:HG23	1:B:166:GLU:OE2	1.94	0.68
1:I:47:VAL:HG23	1:I:51:VAL:HG13	1.76	0.68
1:D:104:ASP:O	1:D:108:ILE:HG13	1.93	0.67
1:G:22:LEU:HD11	1:G:54:MSE:HB3	1.76	0.67
1:I:28:THR:HB	1:I:33:ARG:HE	1.59	0.67
1:I:22:LEU:H	1:I:22:LEU:HD23	1.60	0.67
1:E:94:LEU:HD12	1:E:97:MSE:HE3	1.77	0.67
1:E:104:ASP:O	1:E:108:ILE:HG13	1.93	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:ILE:HD13	1:A:100:VAL:HG13	1.77	0.67
1:I:34:THR:O	1:I:37:ILE:HG23	1.94	0.67
1:A:47:VAL:HG11	1:A:51:VAL:HA	1.78	0.66
1:A:66:ASN:ND2	1:A:68:PHE:H	1.93	0.66
1:A:111:THR:OG1	1:A:165:ILE:HG21	1.94	0.66
1:E:35:ILE:HD12	1:E:45:ILE:HD11	1.78	0.66
1:E:111:THR:OG1	1:E:165:ILE:HG21	1.94	0.66
1:F:34:THR:O	1:F:37:ILE:HG23	1.95	0.66
1:D:7:ILE:CD1	1:D:45:ILE:HG13	2.25	0.66
1:E:129:ASN:O	1:E:171:VAL:HG13	1.94	0.66
1:A:104:ASP:OD1	1:A:167:CYS:HB3	1.94	0.66
1:F:25:ILE:HD11	1:F:175:ILE:HD12	1.76	0.66
1:G:94:LEU:HB2	1:G:97:MSE:HE2	1.78	0.66
1:B:97:MSE:HE1	1:B:129:ASN:HB2	1.78	0.66
1:D:16:PHE:CE1	1:D:175:ILE:HG21	2.31	0.66
1:E:103:GLU:CD	1:E:103:GLU:N	2.54	0.65
1:G:150:ALA:HA	1:G:153:ILE:HD11	1.78	0.65
1:I:94:LEU:HB2	1:I:97:MSE:HE2	1.78	0.65
1:E:169:GLU:HG2	1:E:173:ILE:HD11	1.79	0.65
1:B:97:MSE:HE3	2:B:302:HOH:O	1.95	0.65
1:D:48:GLY:O	1:D:51:VAL:HG22	1.97	0.65
1:E:121:PRO:HB2	1:E:171:VAL:HG11	1.79	0.65
1:F:174:ASP:HB3	1:I:178:LYS:HE2	1.78	0.65
1:G:34:THR:HG21	1:G:95:GLY:HA2	1.78	0.65
1:G:123:HIS:HA	1:G:166:GLU:OE2	1.97	0.65
1:I:74:SER:HB2	1:I:144:LEU:HD23	1.79	0.65
1:G:120:ILE:HD11	1:G:154:LEU:HD13	1.77	0.65
1:I:22:LEU:H	1:I:22:LEU:CD2	2.09	0.65
1:D:44:ILE:HD12	1:D:64:ILE:HD12	1.80	0.64
1:I:66:ASN:C	1:I:66:ASN:HD22	2.03	0.64
1:E:148:VAL:HG12	1:E:149:GLY:N	2.13	0.64
1:H:33:ARG:HG3	1:H:33:ARG:NH1	2.11	0.64
1:I:175:ILE:HG22	1:I:176:ASP:N	2.13	0.64
1:H:157:ILE:HD11	1:H:162:LEU:HG	1.80	0.64
1:A:17:GLY:O	1:A:50:TYR:HD2	1.81	0.64
1:B:25:ILE:O	1:B:26:ASP:HB2	1.96	0.64
1:F:122:THR:O	1:F:166:GLU:HA	1.96	0.64
1:D:30:ILE:HD11	1:D:96:ASP:HA	1.79	0.63
1:B:141:ILE:O	1:B:144:LEU:HD22	1.98	0.63
1:D:64:ILE:HD13	1:E:83:PHE:HZ	1.63	0.63
1:D:7:ILE:CG1	1:D:45:ILE:HG13	2.29	0.63

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:28:THR:HG23	1:H:33:ARG:HD2	1.80	0.63
1:H:77:LEU:HB3	1:H:141:ILE:HD11	1.81	0.63
1:H:93:ALA:HA	1:H:130:PRO:HB3	1.80	0.62
1:I:7:ILE:HD11	1:I:45:ILE:HG23	1.80	0.62
1:D:10:ALA:HA	1:D:48:GLY:H	1.64	0.62
1:G:51:VAL:HG13	1:G:52:ASN:N	2.13	0.62
1:H:68:PHE:O	1:H:75:THR:HG21	1.99	0.62
1:F:46:ILE:HD11	1:F:80:GLY:HA3	1.82	0.62
1:F:81:LEU:HD22	1:F:141:ILE:HD12	1.81	0.62
1:E:70:ASN:HD22	1:E:70:ASN:N	1.86	0.62
1:B:82:ARG:HH11	1:B:82:ARG:HB3	1.65	0.62
1:H:103:GLU:CD	1:H:103:GLU:H	2.06	0.62
1:B:82:ARG:HB3	1:B:82:ARG:NH1	2.14	0.62
1:A:91:LEU:HD22	1:A:130:PRO:HB2	1.82	0.62
1:H:117:LYS:HB3	1:H:137:LEU:HD21	1.82	0.62
1:E:9:ALA:HB1	1:E:21:LEU:CD1	2.30	0.61
1:E:37:ILE:HD13	1:E:100:VAL:HG13	1.82	0.61
1:E:24:LYS:HE3	1:E:27:ASN:HA	1.82	0.61
1:H:44:ILE:HD11	1:H:64:ILE:HD12	1.80	0.61
1:H:56:PRO:HA	1:H:59:MSE:HE3	1.83	0.61
1:E:48:GLY:O	1:E:51:VAL:HG22	2.00	0.61
1:F:23:ALA:O	1:F:30:ILE:HG13	1.99	0.61
1:H:154:LEU:C	1:H:156:LYS:H	2.08	0.61
1:B:97:MSE:HE1	1:B:129:ASN:HB3	1.82	0.61
1:F:103:GLU:O	1:F:107:LYS:HD3	2.00	0.61
1:G:144:LEU:HG	1:G:148:VAL:HG23	1.83	0.61
1:I:33:ARG:O	1:I:37:ILE:HG22	2.00	0.61
1:E:2:ASN:ND2	1:E:87:TYR:HA	2.16	0.60
1:E:66:ASN:C	1:E:66:ASN:HD22	2.09	0.60
1:I:128:GLY:HA3	1:I:171:VAL:HG12	1.83	0.60
1:A:121:PRO:HB2	1:A:171:VAL:HG11	1.83	0.60
1:E:141:ILE:HG12	1:E:144:LEU:HD22	1.83	0.60
1:F:119:VAL:HB	1:F:132:LEU:HB3	1.84	0.60
1:I:175:ILE:HG22	1:I:176:ASP:H	1.67	0.60
1:G:97:MSE:HE1	1:G:129:ASN:HB2	1.84	0.60
1:A:104:ASP:O	1:A:108:ILE:HG13	2.02	0.59
1:B:48:GLY:O	1:B:51:VAL:HG22	2.02	0.59
1:F:169:GLU:CB	1:I:24:LYS:HE2	2.32	0.59
1:G:25:ILE:HB	1:G:33:ARG:HD2	1.85	0.59
1:D:161:GLU:O	1:D:161:GLU:HG2	2.02	0.59
1:E:46:ILE:HD13	1:E:64:ILE:HG23	1.84	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:82:ARG:HH21	1:G:83:PHE:HD2	1.49	0.59
1:F:141:ILE:HD13	1:F:141:ILE:C	2.27	0.59
1:F:148:VAL:HG23	1:F:149:GLY:N	2.16	0.59
1:G:165:ILE:HD13	1:G:165:ILE:N	2.18	0.59
1:E:1:MSE:HB3	1:E:88:ASP:OD2	2.02	0.59
1:E:66:ASN:ND2	1:E:68:PHE:N	2.43	0.59
1:I:48:GLY:O	1:I:51:VAL:HG22	2.03	0.59
1:A:74:SER:O	1:A:78:LYS:HG3	2.02	0.59
1:D:154:LEU:O	1:D:157:ILE:HG13	2.03	0.58
1:I:66:ASN:ND2	1:I:66:ASN:C	2.61	0.58
1:B:31:ILE:O	1:B:35:ILE:HG13	2.03	0.58
1:B:77:LEU:O	1:B:81:LEU:HB2	2.02	0.58
1:B:121:PRO:HB2	1:B:171:VAL:HG11	1.84	0.58
1:I:26:ASP:HB2	1:I:33:ARG:NH2	2.19	0.58
1:E:66:ASN:HD22	1:E:68:PHE:H	1.46	0.58
1:H:129:ASN:HB3	1:H:130:PRO:HA	1.86	0.58
1:I:3:ILE:HD11	1:I:41:LEU:HD22	1.85	0.58
1:A:164:PHE:O	1:A:165:ILE:HG23	2.04	0.58
1:D:111:THR:OG1	1:D:165:ILE:HG21	2.04	0.58
1:E:47:VAL:CG2	1:E:51:VAL:HG13	2.34	0.58
1:H:141:ILE:HA	1:H:144:LEU:CD1	2.32	0.58
1:A:55:LEU:CD1	1:B:55:LEU:HD11	2.34	0.58
1:B:109:ILE:HD12	1:B:110:ASN:N	2.18	0.58
1:B:129:ASN:O	1:B:171:VAL:HG13	2.04	0.58
1:B:164:PHE:C	1:B:165:ILE:HD13	2.29	0.58
1:H:34:THR:O	1:H:37:ILE:HG12	2.03	0.58
1:G:37:ILE:HD11	1:G:38:TYR:CZ	2.39	0.57
1:G:148:VAL:HB	1:G:152:VAL:HG21	1.86	0.57
1:H:122:THR:HG22	1:H:166:GLU:HG2	1.86	0.57
1:G:33:ARG:O	1:G:37:ILE:HG23	2.05	0.57
1:H:104:ASP:OD1	1:H:167:CYS:HB3	2.04	0.57
1:B:23:ALA:O	1:B:29:PRO:HA	2.05	0.57
1:H:51:VAL:HG13	1:H:52:ASN:N	2.19	0.57
1:I:29:PRO:HB2	1:I:32:MSE:HG3	1.87	0.57
1:A:28:THR:HB	1:I:53:GLU:OE1	2.04	0.57
1:H:120:ILE:HD12	1:H:162:LEU:HD21	1.85	0.57
1:I:81:LEU:HD12	1:I:141:ILE:HD11	1.85	0.57
1:I:129:ASN:O	1:I:171:VAL:HG13	2.04	0.57
1:F:52:ASN:HD21	1:G:55:LEU:HB3	1.70	0.57
1:E:141:ILE:HA	1:E:144:LEU:HD13	1.87	0.57
1:A:11:GLY:HA2	1:A:16:PHE:O	2.05	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:49:LYS:HD2	1:D:50:TYR:CE2	2.41	0.56
1:F:46:ILE:HD11	1:F:80:GLY:CA	2.35	0.56
1:I:97:MSE:CE	1:I:129:ASN:HB2	2.24	0.56
1:D:100:VAL:HG21	1:D:130:PRO:HD3	1.86	0.56
1:F:122:THR:HG22	1:F:165:ILE:C	2.30	0.56
1:B:140:GLU:HG3	1:B:153:ILE:HD12	1.86	0.56
1:D:107:LYS:HG2	1:I:159:ILE:HG12	1.86	0.56
1:H:64:ILE:HD11	1:H:83:PHE:CE2	2.40	0.56
1:I:111:THR:OG1	1:I:165:ILE:HG13	2.05	0.56
1:F:30:ILE:HD13	1:F:175:ILE:HD13	1.88	0.56
1:G:108:ILE:HA	1:G:165:ILE:HD11	1.88	0.56
1:G:117:LYS:HB3	1:G:137:LEU:HD21	1.86	0.56
1:I:24:LYS:H	1:I:24:LYS:CD	2.09	0.56
1:A:171:VAL:HG12	1:A:171:VAL:O	2.06	0.56
1:A:47:VAL:O	1:A:65:TYR:HA	2.06	0.56
1:A:81:LEU:HD22	1:A:141:ILE:HD12	1.87	0.56
1:F:29:PRO:O	1:F:32:MSE:HB3	2.05	0.56
1:F:122:THR:HG23	1:F:166:GLU:HA	1.87	0.56
1:F:23:ALA:HA	1:I:169:GLU:OE1	2.06	0.56
1:F:121:PRO:HB2	1:F:171:VAL:HG11	1.88	0.55
1:H:66:ASN:O	1:H:69:TRP:HD1	1.89	0.55
1:I:26:ASP:HB2	1:I:33:ARG:HH21	1.71	0.55
1:A:122:THR:O	1:A:166:GLU:HA	2.07	0.55
1:A:136:SER:OG	1:A:137:LEU:HD22	2.07	0.55
1:D:121:PRO:HB2	1:D:171:VAL:HG11	1.88	0.55
1:F:141:ILE:HA	1:F:153:ILE:HD11	1.89	0.55
1:I:47:VAL:CG2	1:I:51:VAL:HG13	2.36	0.55
1:E:66:ASN:ND2	1:E:66:ASN:C	2.64	0.55
1:G:51:VAL:HG13	1:G:52:ASN:H	1.71	0.55
1:A:141:ILE:O	1:A:144:LEU:HB2	2.07	0.55
1:D:145:ARG:O	1:D:148:VAL:HG23	2.05	0.55
1:G:1:MSE:HG3	1:G:88:ASP:OD2	2.07	0.55
1:E:47:VAL:HG23	1:E:51:VAL:HG13	1.87	0.55
1:F:94:LEU:HD22	1:F:97:MSE:HE2	1.87	0.55
1:B:69:TRP:HZ2	1:I:71:GLU:HG2	1.72	0.55
1:H:18:GLY:HA2	1:H:50:TYR:CD2	2.40	0.55
1:H:94:LEU:CB	1:H:97:MSE:HE3	2.36	0.55
1:I:144:LEU:HD23	1:I:144:LEU:O	2.07	0.55
1:D:66:ASN:C	1:D:66:ASN:HD22	2.14	0.55
1:H:135:LYS:HA	1:H:138:PHE:CE2	2.42	0.55
1:E:50:TYR:O	1:E:54:MSE:HG3	2.06	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:101:THR:OG1	1:E:103:GLU:HG2	2.07	0.55
1:F:135:LYS:HA	1:F:138:PHE:CD1	2.42	0.55
1:B:97:MSE:HG2	1:B:174:ASP:HB3	1.89	0.54
1:H:148:VAL:HG13	1:H:152:VAL:HG21	1.87	0.54
1:D:7:ILE:CD1	1:D:35:ILE:HG12	2.38	0.54
1:F:1:MSE:HE2	1:F:1:MSE:HA	1.89	0.54
1:H:98:PRO:HG3	1:H:175:ILE:HD11	1.89	0.54
1:I:127:ARG:HD2	1:I:164:PHE:HE1	1.72	0.54
1:A:37:ILE:HG12	1:A:38:TYR:CD1	2.42	0.54
1:E:104:ASP:CG	1:E:167:CYS:HB3	2.32	0.54
1:G:89:ALA:HB2	1:G:134:SER:HA	1.87	0.54
1:H:128:GLY:HA3	1:H:171:VAL:CG1	2.37	0.54
1:A:47:VAL:CG1	1:A:51:VAL:HA	2.38	0.54
1:A:148:VAL:HG13	1:A:152:VAL:HG21	1.87	0.54
1:D:18:GLY:HA3	1:D:50:TYR:HB2	1.88	0.54
1:H:44:ILE:HD11	1:H:64:ILE:HD11	1.90	0.54
1:F:51:VAL:HG13	1:F:52:ASN:N	2.21	0.54
1:I:66:ASN:HD22	1:I:67:PRO:N	2.05	0.54
1:A:128:GLY:HA3	1:A:171:VAL:HG12	1.90	0.54
1:D:122:THR:O	1:D:166:GLU:HA	2.08	0.54
1:I:122:THR:O	1:I:166:GLU:HA	2.07	0.54
1:F:45:ILE:O	1:F:45:ILE:HG13	2.08	0.54
1:H:126:GLU:HB2	1:H:172:LEU:HD21	1.90	0.54
1:A:77:LEU:HD13	1:A:141:ILE:CG1	2.39	0.53
1:E:157:ILE:CG1	1:E:162:LEU:HD13	2.38	0.53
1:I:66:ASN:ND2	1:I:68:PHE:N	2.52	0.53
1:I:141:ILE:HA	1:I:153:ILE:HD11	1.90	0.53
1:G:10:ALA:HB1	1:G:49:LYS:H	1.73	0.53
1:H:162:LEU:HD22	1:H:164:PHE:CZ	2.43	0.53
1:E:157:ILE:HG13	1:E:162:LEU:HD13	1.91	0.53
1:F:81:LEU:CD2	1:F:141:ILE:HD12	2.37	0.53
1:G:44:ILE:HG13	1:G:44:ILE:O	2.08	0.53
1:A:37:ILE:HG12	1:A:38:TYR:CE1	2.44	0.53
1:D:165:ILE:N	1:D:165:ILE:HD13	2.24	0.53
1:F:127:ARG:HD2	1:F:164:PHE:CE1	2.44	0.53
1:G:119:VAL:HB	1:G:132:LEU:HB3	1.89	0.53
1:H:158:LYS:C	1:H:160:GLU:H	2.17	0.53
1:A:72:GLY:O	1:A:75:THR:HG22	2.09	0.53
1:E:29:PRO:HD2	1:E:32:MSE:HG3	1.90	0.53
1:F:74:SER:OG	1:F:146:GLY:O	2.24	0.53
1:F:104:ASP:OD1	1:F:167:CYS:HB3	2.08	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:89:ALA:CB	1:G:134:SER:HA	2.38	0.53
1:B:81:LEU:HD11	1:B:141:ILE:HD11	1.91	0.53
1:G:45:ILE:HG22	1:G:47:VAL:CG1	2.37	0.53
1:G:97:MSE:HG2	1:G:174:ASP:HA	1.90	0.53
1:I:148:VAL:HG23	1:I:149:GLY:N	2.20	0.53
1:A:117:LYS:HB3	1:A:137:LEU:HD21	1.90	0.53
1:E:66:ASN:HD21	1:E:68:PHE:HB2	1.73	0.53
1:F:126:GLU:HB3	1:F:172:LEU:HD11	1.90	0.53
1:I:73:ILE:HG23	1:I:74:SER:N	2.23	0.53
1:H:39:GLY:O	1:H:43:LYS:NZ	2.34	0.52
1:B:94:LEU:HB2	1:B:97:MSE:HE2	1.90	0.52
1:D:7:ILE:HD11	1:D:45:ILE:HG13	1.91	0.52
1:E:30:ILE:HG13	1:E:175:ILE:CD1	2.36	0.52
1:H:38:TYR:O	1:H:41:LEU:HB2	2.08	0.52
1:H:121:PRO:HB2	1:H:171:VAL:HG21	1.92	0.52
1:H:141:ILE:HG12	1:H:144:LEU:HD11	1.92	0.52
1:I:109:ILE:HG13	1:I:110:ASN:N	2.25	0.52
1:A:117:LYS:O	1:A:137:LEU:HD23	2.09	0.52
1:B:167:CYS:HB2	1:B:171:VAL:HG21	1.90	0.52
1:G:128:GLY:HA3	1:G:171:VAL:O	2.10	0.52
1:I:149:GLY:O	1:I:152:VAL:HG22	2.10	0.52
1:A:26:ASP:O	1:A:27:ASN:HB2	2.09	0.52
1:B:111:THR:HG21	1:B:165:ILE:CD1	2.34	0.52
1:H:138:PHE:HA	1:H:141:ILE:HG22	1.91	0.52
1:A:77:LEU:HD13	1:A:141:ILE:HG13	1.91	0.52
1:D:66:ASN:ND2	1:D:66:ASN:C	2.67	0.52
1:E:94:LEU:HD12	1:E:97:MSE:CE	2.39	0.52
1:F:117:LYS:HB3	1:F:137:LEU:HD11	1.92	0.52
1:F:132:LEU:C	1:F:132:LEU:HD13	2.34	0.52
1:H:37:ILE:HD12	1:H:100:VAL:HG22	1.92	0.52
1:D:23:ALA:O	1:D:30:ILE:HG23	2.09	0.52
1:H:32:MSE:HE2	1:H:36:ARG:HH21	1.74	0.52
1:E:141:ILE:HD13	1:E:141:ILE:C	2.35	0.52
1:A:18:GLY:O	1:A:22:LEU:HG	2.10	0.52
1:G:97:MSE:HE1	1:G:129:ASN:CB	2.39	0.52
1:H:111:THR:HG21	1:H:165:ILE:HG23	1.92	0.52
1:D:111:THR:HG22	1:D:111:THR:O	2.09	0.51
1:E:165:ILE:HD13	1:E:165:ILE:N	2.24	0.51
1:H:55:LEU:HD13	1:I:51:VAL:HB	1.91	0.51
1:I:106:ASN:HA	1:I:109:ILE:HG12	1.91	0.51
1:E:119:VAL:HB	1:E:132:LEU:HB3	1.92	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:164:PHE:C	1:A:165:ILE:HD13	2.35	0.51
1:E:171:VAL:HG12	1:E:171:VAL:O	2.10	0.51
1:G:120:ILE:HD11	1:G:154:LEU:CD1	2.40	0.51
1:F:178:LYS:HE3	1:I:94:LEU:CD1	2.40	0.51
1:I:23:ALA:O	1:I:29:PRO:HA	2.11	0.51
1:A:67:PRO:HG2	1:A:68:PHE:CD2	2.46	0.51
1:D:47:VAL:HG23	1:D:47:VAL:O	2.11	0.51
1:E:64:ILE:HD11	1:E:79:LEU:HD21	1.93	0.51
1:F:51:VAL:HG22	1:G:55:LEU:CD1	2.40	0.51
1:G:107:LYS:O	1:G:111:THR:HG23	2.11	0.51
1:H:21:LEU:HD23	1:H:30:ILE:HD11	1.92	0.51
1:E:30:ILE:CG1	1:E:175:ILE:HD13	2.38	0.51
1:H:104:ASP:CG	1:H:167:CYS:HB3	2.36	0.51
1:H:148:VAL:HG13	1:H:152:VAL:CG2	2.41	0.51
1:A:129:ASN:HB3	1:A:130:PRO:HA	1.92	0.51
1:D:38:TYR:O	1:D:41:LEU:HB2	2.10	0.51
1:D:51:VAL:HG12	1:D:55:LEU:HD23	1.92	0.51
1:D:176:ASP:O	1:D:178:LYS:N	2.42	0.51
1:F:36:ARG:HH11	1:F:36:ARG:HG3	1.75	0.51
1:F:51:VAL:HG22	1:G:55:LEU:HD11	1.92	0.51
1:F:127:ARG:HD2	1:F:164:PHE:HE1	1.75	0.51
1:A:44:ILE:HG13	1:A:64:ILE:HD11	1.91	0.51
1:B:128:GLY:HA3	1:B:171:VAL:HG12	1.93	0.51
1:I:171:VAL:HG12	1:I:171:VAL:O	2.10	0.51
1:B:4:GLY:HA3	1:B:84:PHE:CE1	2.46	0.51
1:D:30:ILE:CD1	1:D:96:ASP:HA	2.40	0.51
1:A:94:LEU:HD12	1:A:97:MSE:SE	2.61	0.51
1:E:23:ALA:O	1:E:30:ILE:HG12	2.11	0.51
1:F:127:ARG:HB2	1:I:180:ASP:HB3	1.92	0.51
1:H:94:LEU:HB2	1:H:97:MSE:HE3	1.92	0.51
1:F:141:ILE:HG12	1:F:144:LEU:HD22	1.93	0.50
1:G:94:LEU:HD12	1:G:97:MSE:SE	2.60	0.50
1:B:33:ARG:HG2	1:B:33:ARG:HH11	1.75	0.50
1:D:148:VAL:HG13	1:D:152:VAL:HG21	1.94	0.50
1:G:26:ASP:O	1:G:27:ASN:HB2	2.10	0.50
1:H:112:PHE:O	1:H:113:LYS:HE2	2.11	0.50
1:D:12:GLU:HB3	1:D:69:TRP:O	2.11	0.50
1:F:169:GLU:O	1:F:173:ILE:HG12	2.10	0.50
1:G:30:ILE:C	1:G:30:ILE:HD13	2.35	0.50
1:A:120:ILE:HD11	1:A:154:LEU:HD13	1.93	0.50
1:F:44:ILE:HD12	1:F:64:ILE:HD12	1.92	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:41:LEU:O	1:G:43:LYS:HE2	2.11	0.50
1:G:74:SER:O	1:G:78:LYS:HG3	2.12	0.50
1:H:47:VAL:HG12	1:H:48:GLY:N	2.26	0.50
1:D:71:GLU:HB3	1:D:75:THR:HG21	1.94	0.50
1:E:3:ILE:HD12	1:E:109:ILE:HG23	1.92	0.50
1:G:21:LEU:O	1:G:30:ILE:HG23	2.12	0.50
1:I:117:LYS:O	1:I:137:LEU:HD22	2.11	0.50
1:B:122:THR:O	1:B:166:GLU:HA	2.12	0.50
1:D:157:ILE:C	1:D:157:ILE:HD12	2.37	0.50
1:H:56:PRO:CA	1:H:59:MSE:HE2	2.35	0.50
1:E:129:ASN:HB3	1:E:130:PRO:HA	1.92	0.50
1:E:158:LYS:C	1:E:160:GLU:N	2.69	0.50
1:H:120:ILE:HD11	1:H:154:LEU:HD13	1.93	0.50
1:E:25:ILE:CG2	1:E:33:ARG:HD2	2.42	0.50
1:H:144:LEU:CD1	1:H:153:ILE:HD11	2.42	0.50
1:B:100:VAL:HG12	1:B:101:THR:N	2.27	0.50
1:G:48:GLY:O	1:G:49:LYS:C	2.54	0.50
1:G:101:THR:CG2	1:G:103:GLU:H	2.20	0.50
1:I:7:ILE:HG13	1:I:45:ILE:HA	1.93	0.50
1:D:171:VAL:HG12	1:D:171:VAL:O	2.11	0.49
1:E:157:ILE:HD12	1:E:158:LYS:H	1.77	0.49
1:H:72:GLY:O	1:H:75:THR:HG22	2.12	0.49
1:H:158:LYS:C	1:H:160:GLU:N	2.69	0.49
1:B:106:ASN:O	1:B:109:ILE:HG13	2.12	0.49
1:D:129:ASN:O	1:D:171:VAL:HG13	2.12	0.49
1:F:177:LYS:HG2	1:I:173:ILE:HG22	1.93	0.49
1:F:33:ARG:HG3	1:F:33:ARG:HH11	1.78	0.49
1:H:165:ILE:N	1:H:165:ILE:HD13	2.27	0.49
1:F:74:SER:O	1:F:78:LYS:HG3	2.12	0.49
1:H:25:ILE:O	1:H:26:ASP:HB3	2.12	0.49
1:I:127:ARG:HD2	1:I:164:PHE:CE1	2.46	0.49
1:D:88:ASP:O	1:D:135:LYS:N	2.41	0.49
1:E:3:ILE:CD1	1:E:109:ILE:HG23	2.42	0.49
1:G:44:ILE:HD11	1:G:80:GLY:HA2	1.94	0.49
1:B:4:GLY:O	1:B:90:VAL:HA	2.13	0.49
1:H:94:LEU:HD13	1:H:97:MSE:HE1	1.93	0.49
1:H:113:LYS:HE2	1:H:113:LYS:HA	1.95	0.49
1:I:7:ILE:HG12	1:I:45:ILE:HG13	1.94	0.49
1:B:169:GLU:O	1:B:170:GLY:C	2.55	0.49
1:A:37:ILE:O	1:A:102:LYS:HG2	2.12	0.49
1:I:73:ILE:CG2	1:I:74:SER:N	2.76	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:101:THR:O	1:I:104:ASP:HB2	2.12	0.49
1:D:22:LEU:CD1	1:D:53:GLU:HB3	2.39	0.49
1:D:37:ILE:HD11	1:D:98:PRO:O	2.13	0.49
1:E:137:LEU:N	1:E:137:LEU:HD12	2.27	0.49
1:F:121:PRO:HB2	1:F:171:VAL:CG1	2.43	0.49
1:G:23:ALA:O	1:G:30:ILE:HG22	2.13	0.48
1:G:64:ILE:HD11	1:G:83:PHE:CE1	2.48	0.48
1:B:70:ASN:N	1:B:70:ASN:ND2	2.49	0.48
1:I:120:ILE:HG21	1:I:127:ARG:HB3	1.94	0.48
1:B:66:ASN:HD22	1:B:68:PHE:H	1.59	0.48
1:D:126:GLU:HB2	1:D:172:LEU:HD21	1.95	0.48
1:G:123:HIS:HB3	1:G:172:LEU:CD1	2.44	0.48
1:I:129:ASN:HB3	1:I:130:PRO:HA	1.96	0.48
1:F:52:ASN:ND2	1:G:55:LEU:HB3	2.29	0.48
1:F:81:LEU:HD11	1:F:138:PHE:CD2	2.49	0.48
1:F:20:LYS:HA	1:F:177:LYS:HE3	1.96	0.48
1:G:21:LEU:HD12	1:G:54:MSE:HE1	1.95	0.48
1:I:28:THR:HB	1:I:33:ARG:NE	2.26	0.48
1:B:43:LYS:O	1:B:44:ILE:HD12	2.13	0.48
1:B:41:LEU:O	1:B:43:LYS:HG3	2.14	0.48
1:D:46:ILE:HD13	1:D:64:ILE:HB	1.93	0.48
1:D:94:LEU:HD12	1:D:97:MSE:SE	2.64	0.48
1:F:128:GLY:HA3	1:F:171:VAL:HG12	1.96	0.48
1:F:173:ILE:HG22	1:I:176:ASP:O	2.14	0.48
1:E:140:GLU:HB3	1:E:153:ILE:HD13	1.96	0.48
1:G:2:ASN:HB3	1:G:87:TYR:HA	1.95	0.48
1:E:2:ASN:HD22	1:E:87:TYR:HA	1.79	0.48
1:H:21:LEU:CD2	1:H:30:ILE:HD11	2.43	0.48
1:H:157:ILE:HD11	1:H:162:LEU:CG	2.43	0.48
1:H:137:LEU:HD12	1:H:153:ILE:CG2	2.44	0.47
1:H:154:LEU:C	1:H:156:LYS:N	2.72	0.47
1:I:66:ASN:HD21	1:I:68:PHE:H	1.57	0.47
1:D:85:LYS:HD3	1:E:85:LYS:HE2	1.96	0.47
1:B:25:ILE:HB	1:B:33:ARG:HD3	1.96	0.47
1:E:176:ASP:OD1	1:E:177:LYS:N	2.47	0.47
1:I:22:LEU:CD2	1:I:22:LEU:N	2.77	0.47
1:I:99:PHE:CE2	1:I:175:ILE:HD11	2.49	0.47
1:B:66:ASN:HD22	1:B:66:ASN:C	2.22	0.47
1:B:94:LEU:HG	1:B:97:MSE:HE2	1.96	0.47
1:E:22:LEU:HD11	1:E:54:MSE:HG2	1.95	0.47
1:E:148:VAL:HG13	1:E:152:VAL:CG2	2.45	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:144:LEU:CD1	1:F:153:ILE:HD11	2.45	0.47
1:F:165:ILE:O	1:F:165:ILE:HG23	2.14	0.47
1:D:140:GLU:HB3	1:D:153:ILE:HD13	1.95	0.47
1:B:104:ASP:O	1:B:108:ILE:HG13	2.14	0.47
1:D:22:LEU:HD21	1:D:57:LEU:HD21	1.96	0.47
1:D:41:LEU:HD21	1:D:109:ILE:HD11	1.95	0.47
1:D:85:LYS:HE2	1:D:85:LYS:O	2.14	0.47
1:H:45:ILE:CD1	1:H:61:GLN:HG2	2.42	0.47
1:H:132:LEU:HD23	1:H:132:LEU:C	2.40	0.47
1:A:27:ASN:ND2	1:I:56:PRO:HG2	2.30	0.47
1:D:157:ILE:HD12	1:D:157:ILE:O	2.14	0.47
1:E:9:ALA:HB1	1:E:21:LEU:HD13	1.96	0.47
1:E:167:CYS:HB2	1:E:171:VAL:HG21	1.96	0.47
1:H:2:ASN:N	1:H:88:ASP:OD2	2.47	0.47
1:D:105:VAL:O	1:D:109:ILE:HG13	2.15	0.47
1:D:119:VAL:HB	1:D:132:LEU:HB3	1.97	0.47
1:F:97:MSE:HE1	1:F:129:ASN:HB3	1.95	0.47
1:G:50:TYR:O	1:G:51:VAL:C	2.56	0.47
1:G:101:THR:HB	1:G:104:ASP:CG	2.40	0.47
1:G:169:GLU:O	1:G:173:ILE:HG13	2.14	0.47
1:I:7:ILE:HG13	1:I:7:ILE:O	2.13	0.47
1:A:48:GLY:O	1:A:51:VAL:HG22	2.15	0.47
1:A:57:LEU:HG	1:H:59:MSE:CE	2.45	0.47
1:A:122:THR:HG22	1:A:123:HIS:N	2.30	0.47
1:A:157:ILE:CG1	1:A:162:LEU:HD11	2.34	0.47
1:B:137:LEU:O	1:B:140:GLU:HB3	2.14	0.47
1:D:157:ILE:CD1	1:D:162:LEU:HD22	2.36	0.47
1:F:149:GLY:O	1:F:152:VAL:HG23	2.15	0.47
1:I:132:LEU:HD13	1:I:132:LEU:C	2.40	0.47
1:A:107:LYS:HE3	1:F:159:ILE:HG12	1.97	0.46
1:F:64:ILE:HD13	1:G:83:PHE:HZ	1.80	0.46
1:I:157:ILE:HB	1:I:162:LEU:HD11	1.98	0.46
1:E:158:LYS:C	1:E:160:GLU:H	2.23	0.46
1:F:123:HIS:HB3	1:F:172:LEU:HD23	1.96	0.46
1:H:45:ILE:N	1:H:45:ILE:HD12	2.29	0.46
1:I:171:VAL:O	1:I:171:VAL:CG1	2.63	0.46
1:H:38:TYR:CZ	1:H:93:ALA:HB2	2.50	0.46
1:H:128:GLY:HA3	1:H:171:VAL:HG12	1.97	0.46
1:E:122:THR:O	1:E:166:GLU:HA	2.16	0.46
1:G:25:ILE:O	1:G:26:ASP:HB2	2.14	0.46
1:H:135:LYS:HA	1:H:138:PHE:CD2	2.51	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:135:LYS:HA	1:I:138:PHE:CG	2.51	0.46
1:B:138:PHE:HA	1:B:141:ILE:CD1	2.46	0.46
1:E:117:LYS:HD2	1:E:157:ILE:HD13	1.97	0.46
1:G:104:ASP:CG	1:G:167:CYS:HB3	2.40	0.46
1:I:28:THR:HG22	1:I:33:ARG:HG3	1.97	0.46
1:I:41:LEU:O	1:I:43:LYS:HG3	2.15	0.46
1:A:128:GLY:HA3	1:A:171:VAL:O	2.16	0.46
1:F:157:ILE:HD11	1:F:162:LEU:HD22	1.97	0.46
1:G:32:MSE:O	1:G:36:ARG:HG2	2.15	0.46
1:G:170:GLY:HA2	1:G:173:ILE:HD12	1.98	0.46
1:H:23:ALA:O	1:H:29:PRO:HA	2.16	0.46
1:H:26:ASP:O	1:H:27:ASN:HB2	2.14	0.46
1:A:8:LEU:HD13	1:A:73:ILE:HD11	1.97	0.46
1:A:165:ILE:HD13	1:A:165:ILE:N	2.31	0.46
1:F:82:ARG:NH2	1:G:83:PHE:HD2	2.14	0.46
1:G:123:HIS:CA	1:G:166:GLU:OE2	2.64	0.46
1:G:129:ASN:HB3	1:G:130:PRO:HA	1.97	0.46
1:G:149:GLY:O	1:G:153:ILE:HG12	2.16	0.46
1:D:123:HIS:HB3	1:D:172:LEU:HD13	1.96	0.46
1:E:138:PHE:O	1:E:141:ILE:HG22	2.16	0.46
1:G:35:ILE:HD11	1:G:45:ILE:HG12	1.98	0.46
1:G:154:LEU:O	1:G:157:ILE:HG12	2.16	0.46
1:H:38:TYR:CZ	1:H:93:ALA:CB	2.99	0.46
1:H:51:VAL:HG13	1:H:52:ASN:H	1.81	0.46
1:H:117:LYS:O	1:H:118:ALA:HB2	2.15	0.46
1:D:28:THR:HB	1:F:53:GLU:OE2	2.16	0.45
1:D:85:LYS:HD3	1:E:85:LYS:HZ3	1.81	0.45
1:B:101:THR:HG22	1:B:102:LYS:N	2.31	0.45
1:H:97:MSE:HG2	1:H:173:ILE:O	2.16	0.45
1:H:104:ASP:O	1:H:108:ILE:HG13	2.17	0.45
1:A:12:GLU:HG3	1:A:73:ILE:HB	1.98	0.45
1:B:94:LEU:CB	1:B:97:MSE:HE2	2.47	0.45
1:D:7:ILE:HD13	1:D:35:ILE:HG12	1.97	0.45
1:E:25:ILE:O	1:E:25:ILE:HG23	2.17	0.45
1:F:9:ALA:HB1	1:F:54:MSE:CE	2.47	0.45
1:G:129:ASN:O	1:G:171:VAL:HG22	2.15	0.45
1:A:33:ARG:O	1:A:36:ARG:HB3	2.17	0.45
1:B:135:LYS:HA	1:B:138:PHE:CE2	2.52	0.45
1:E:55:LEU:N	1:E:56:PRO:CD	2.80	0.45
1:F:118:ALA:HB3	1:F:162:LEU:HD13	1.97	0.45
1:G:22:LEU:HD11	1:G:54:MSE:CB	2.45	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:4:GLY:O	1:I:90:VAL:HA	2.17	0.45
1:A:174:ASP:O	1:A:175:ILE:C	2.59	0.45
1:H:138:PHE:O	1:H:139:ASN:C	2.59	0.45
1:B:21:LEU:HA	1:B:30:ILE:HD11	1.99	0.45
1:E:77:LEU:HD13	1:E:141:ILE:HG13	1.98	0.45
1:E:126:GLU:HG3	1:E:172:LEU:HD21	1.98	0.45
1:G:30:ILE:HD13	1:G:31:ILE:N	2.32	0.45
1:H:122:THR:O	1:H:166:GLU:HA	2.15	0.45
1:F:97:MSE:CE	1:F:129:ASN:HB2	2.40	0.45
1:D:157:ILE:HD11	1:D:162:LEU:HD13	1.98	0.45
1:E:148:VAL:HG13	1:E:152:VAL:HG21	1.99	0.45
1:F:24:LYS:H	1:I:169:GLU:CD	2.25	0.45
1:H:83:PHE:HZ	1:I:64:ILE:HD13	1.82	0.45
1:A:105:VAL:O	1:A:109:ILE:HG13	2.17	0.44
1:B:150:ALA:O	1:B:151:ARG:C	2.60	0.44
1:E:70:ASN:ND2	1:E:70:ASN:N	2.53	0.44
1:F:154:LEU:O	1:F:157:ILE:HG12	2.17	0.44
1:H:2:ASN:HD22	1:H:2:ASN:HA	1.56	0.44
1:H:35:ILE:HG23	1:H:43:LYS:HD2	1.98	0.44
1:A:141:ILE:HD13	1:A:141:ILE:C	2.42	0.44
1:D:85:LYS:HE2	1:D:86:ASP:N	2.32	0.44
1:E:121:PRO:HG2	1:E:128:GLY:C	2.43	0.44
1:G:81:LEU:HD22	1:G:141:ILE:CG1	2.47	0.44
1:G:113:LYS:CB	1:G:114:PRO:CD	2.95	0.44
1:H:24:LYS:HD3	1:H:27:ASN:C	2.42	0.44
1:A:19:ASP:CG	1:A:20:LYS:H	2.25	0.44
1:D:51:VAL:HG21	1:D:65:TYR:CE1	2.53	0.44
1:D:157:ILE:HD11	1:D:162:LEU:CD2	2.34	0.44
1:F:44:ILE:HG22	1:F:62:ILE:HB	1.98	0.44
1:G:50:TYR:HB2	1:G:54:MSE:HE3	2.00	0.44
1:A:66:ASN:C	1:A:66:ASN:HD22	2.25	0.44
1:D:120:ILE:HD11	1:D:154:LEU:HD13	1.99	0.44
1:F:39:GLY:O	1:F:43:LYS:NZ	2.48	0.44
1:F:135:LYS:O	1:F:138:PHE:HB2	2.17	0.44
1:H:169:GLU:O	1:H:173:ILE:HG13	2.16	0.44
1:B:103:GLU:HG2	1:B:107:LYS:HE3	1.99	0.44
1:D:16:PHE:CE1	1:D:175:ILE:CG2	3.00	0.44
1:G:174:ASP:O	1:G:175:ILE:HG22	2.18	0.44
1:B:108:ILE:HD11	1:B:121:PRO:HB3	1.99	0.44
1:I:45:ILE:HD12	1:I:45:ILE:N	2.33	0.44
1:I:90:VAL:HG22	1:I:138:PHE:CE1	2.52	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:141:ILE:HA	1:I:153:ILE:CD1	2.47	0.44
1:B:74:SER:HB2	1:B:144:LEU:CD2	2.48	0.44
1:D:46:ILE:HG21	1:D:76:SER:HB3	2.00	0.44
1:E:45:ILE:HD13	1:E:58:LEU:HD22	1.99	0.44
1:H:94:LEU:HD13	1:H:97:MSE:CE	2.47	0.44
1:H:150:ALA:O	1:H:154:LEU:HG	2.18	0.44
1:I:43:LYS:HE3	1:I:43:LYS:HB2	1.87	0.44
1:A:28:THR:HA	1:A:29:PRO:HD3	1.87	0.44
1:A:171:VAL:O	1:A:171:VAL:CG1	2.65	0.44
1:F:178:LYS:HE3	1:I:94:LEU:HD13	2.00	0.44
1:A:119:VAL:HB	1:A:132:LEU:HB3	2.00	0.43
1:D:47:VAL:O	1:D:65:TYR:HA	2.18	0.43
1:E:19:ASP:HA	1:E:22:LEU:HD23	2.00	0.43
1:F:111:THR:O	1:F:111:THR:HG22	2.18	0.43
1:F:122:THR:HA	1:F:126:GLU:O	2.18	0.43
1:H:50:TYR:O	1:H:53:GLU:HG2	2.18	0.43
1:I:7:ILE:CG1	1:I:45:ILE:HG13	2.48	0.43
1:G:122:THR:O	1:G:166:GLU:HA	2.18	0.43
1:G:132:LEU:HD13	1:G:132:LEU:C	2.43	0.43
1:H:78:LYS:HE3	1:H:144:LEU:HD22	2.01	0.43
1:D:22:LEU:CD2	1:D:57:LEU:HD11	2.48	0.43
1:D:49:LYS:HG2	1:D:69:TRP:CZ2	2.53	0.43
1:E:25:ILE:HG23	1:E:33:ARG:HD2	1.99	0.43
1:E:122:THR:HG22	1:E:127:ARG:HB3	2.00	0.43
1:G:51:VAL:CG1	1:G:52:ASN:N	2.80	0.43
1:A:15:ARG:HH12	1:A:147:ASP:CG	2.26	0.43
1:B:119:VAL:HB	1:B:132:LEU:HB3	1.99	0.43
1:E:89:ALA:HB1	1:E:132:LEU:HD13	2.00	0.43
1:I:91:LEU:HD13	1:I:132:LEU:HD23	2.00	0.43
1:I:162:LEU:HD12	1:I:162:LEU:N	2.33	0.43
1:D:174:ASP:OD2	1:D:176:ASP:HB2	2.18	0.43
1:F:108:ILE:CD1	1:F:121:PRO:HG3	2.49	0.43
1:I:175:ILE:CG2	1:I:176:ASP:N	2.82	0.43
1:A:3:ILE:CD1	1:A:109:ILE:HG23	2.48	0.43
1:B:97:MSE:HG2	1:B:174:ASP:CB	2.49	0.43
1:D:91:LEU:HD22	1:D:130:PRO:HB2	2.01	0.43
1:D:175:ILE:HD13	1:D:175:ILE:C	2.43	0.43
1:E:122:THR:HA	1:E:127:ARG:HA	2.01	0.43
1:E:137:LEU:N	1:E:137:LEU:CD1	2.81	0.43
1:F:25:ILE:HG22	1:F:33:ARG:HH12	1.83	0.43
1:G:120:ILE:CD1	1:G:154:LEU:HD13	2.47	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:21:LEU:CB	1:H:54:MSE:HE1	2.40	0.43
1:H:121:PRO:HG2	1:H:128:GLY:C	2.44	0.43
1:A:44:ILE:HD12	1:A:62:ILE:HB	1.99	0.43
1:D:174:ASP:OD1	1:D:175:ILE:N	2.51	0.43
1:E:66:ASN:HD21	1:E:68:PHE:CB	2.32	0.43
1:E:111:THR:HG21	1:E:165:ILE:CG2	2.49	0.43
1:G:2:ASN:HB2	1:G:88:ASP:N	2.23	0.43
1:H:77:LEU:HD13	1:H:141:ILE:HD11	2.01	0.43
1:F:144:LEU:HD11	1:F:153:ILE:HD11	2.01	0.43
1:B:66:ASN:ND2	1:B:66:ASN:C	2.76	0.43
1:D:72:GLY:HA3	1:D:146:GLY:C	2.44	0.43
1:H:101:THR:OG1	1:H:103:GLU:HG2	2.18	0.43
1:H:105:VAL:O	1:H:109:ILE:HG13	2.18	0.43
1:B:120:ILE:HD11	1:B:154:LEU:HD13	2.00	0.43
1:E:141:ILE:HA	1:E:153:ILE:HD11	2.01	0.43
1:G:6:ILE:HG23	1:G:46:ILE:HD13	2.01	0.43
1:H:28:THR:HA	1:H:29:PRO:HD3	1.91	0.43
1:E:31:ILE:HG21	1:E:54:MSE:HE2	2.01	0.42
1:E:128:GLY:HA3	1:E:171:VAL:HG12	2.00	0.42
1:E:141:ILE:HD13	1:E:141:ILE:O	2.18	0.42
1:G:30:ILE:HG12	1:G:98:PRO:HB3	2.01	0.42
1:G:137:LEU:O	1:G:138:PHE:C	2.62	0.42
1:H:30:ILE:CD1	1:H:96:ASP:HA	2.49	0.42
1:H:141:ILE:HA	1:H:144:LEU:HD11	2.01	0.42
1:I:28:THR:CG2	1:I:33:ARG:HG3	2.48	0.42
1:D:111:THR:O	1:D:111:THR:CG2	2.66	0.42
1:I:179:GLU:CD	1:I:179:GLU:H	2.28	0.42
1:D:129:ASN:N	1:D:171:VAL:HG13	2.35	0.42
1:F:4:GLY:O	1:F:90:VAL:HA	2.19	0.42
1:F:101:THR:O	1:F:104:ASP:HB2	2.19	0.42
1:G:123:HIS:HB3	1:G:172:LEU:HD11	2.01	0.42
1:H:33:ARG:NH1	1:H:33:ARG:CG	2.81	0.42
1:D:15:ARG:HG3	1:D:15:ARG:HH11	1.84	0.42
1:F:3:ILE:HA	1:F:89:ALA:O	2.19	0.42
1:G:49:LYS:HA	1:G:69:TRP:CZ2	2.55	0.42
1:H:44:ILE:HG23	1:H:44:ILE:O	2.19	0.42
1:I:141:ILE:C	1:I:143:LYS:H	2.27	0.42
1:A:4:GLY:O	1:A:90:VAL:HA	2.20	0.42
1:B:135:LYS:HA	1:B:138:PHE:CD2	2.55	0.42
1:E:77:LEU:HD13	1:E:141:ILE:CG1	2.50	0.42
1:E:120:ILE:HD11	1:E:154:LEU:HD13	2.01	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:ALA:CB	1:B:134:SER:HA	2.50	0.42
1:G:37:ILE:HG13	1:G:38:TYR:CD1	2.55	0.42
1:G:151:ARG:HG2	1:G:151:ARG:HH11	1.84	0.42
1:A:12:GLU:O	1:A:15:ARG:HB2	2.20	0.42
1:A:44:ILE:HD11	1:A:62:ILE:HG21	2.00	0.42
1:B:81:LEU:CD1	1:B:141:ILE:HD11	2.49	0.42
1:B:164:PHE:CD2	1:B:164:PHE:N	2.87	0.42
1:F:108:ILE:HD13	1:F:121:PRO:HG3	2.01	0.42
1:F:123:HIS:HB3	1:F:172:LEU:CD2	2.50	0.42
1:G:164:PHE:O	1:G:165:ILE:HG23	2.20	0.42
1:H:32:MSE:SE	1:H:57:LEU:HG	2.70	0.42
1:H:78:LYS:O	1:H:82:ARG:HG3	2.19	0.42
1:B:156:LYS:C	1:B:157:ILE:HG13	2.45	0.42
1:H:4:GLY:HA3	1:H:84:PHE:CE1	2.54	0.42
1:D:42:GLU:OE2	1:D:87:TYR:OH	2.36	0.42
1:I:144:LEU:HD23	1:I:144:LEU:C	2.45	0.42
1:I:154:LEU:HD23	1:I:157:ILE:CD1	2.50	0.42
1:A:47:VAL:HG12	1:A:51:VAL:HG13	2.01	0.41
1:A:104:ASP:CG	1:A:167:CYS:HB3	2.44	0.41
1:E:29:PRO:HB2	1:E:32:MSE:HG3	2.02	0.41
1:E:126:GLU:HG3	1:E:172:LEU:CD2	2.50	0.41
1:E:157:ILE:HD12	1:E:158:LYS:N	2.35	0.41
1:F:51:VAL:CG1	1:F:52:ASN:N	2.83	0.41
1:F:89:ALA:HB1	1:F:133:ILE:O	2.19	0.41
1:H:8:LEU:CD1	1:H:73:ILE:HD11	2.49	0.41
1:H:45:ILE:CD1	1:H:45:ILE:N	2.83	0.41
1:H:83:PHE:HD1	1:I:82:ARG:NH2	2.18	0.41
1:H:84:PHE:O	1:H:85:LYS:C	2.63	0.41
1:B:132:LEU:O	1:B:132:LEU:HD13	2.20	0.41
1:B:171:VAL:HG12	1:B:171:VAL:O	2.20	0.41
1:B:158:LYS:HB2	1:B:161:GLU:OE2	2.20	0.41
1:F:141:ILE:HD13	1:F:141:ILE:O	2.20	0.41
1:H:55:LEU:CD1	1:I:55:LEU:HD12	2.50	0.41
1:H:123:HIS:HB3	1:H:172:LEU:HD22	2.02	0.41
1:H:141:ILE:C	1:H:143:LYS:H	2.27	0.41
1:I:3:ILE:HA	1:I:89:ALA:O	2.20	0.41
1:D:169:GLU:HB3	1:D:173:ILE:HD11	2.03	0.41
1:E:66:ASN:HD22	1:E:67:PRO:N	2.18	0.41
1:E:158:LYS:O	1:E:160:GLU:N	2.53	0.41
1:G:171:VAL:O	1:G:171:VAL:CG1	2.69	0.41
1:H:22:LEU:HD23	1:H:54:MSE:HE2	2.01	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:31:ILE:HG23	1:I:32:MSE:N	2.36	0.41
1:I:104:ASP:OD2	1:I:167:CYS:HB3	2.19	0.41
1:A:97:MSE:SE	1:A:129:ASN:HB2	2.71	0.41
1:B:154:LEU:HA	1:B:157:ILE:HD12	2.02	0.41
1:D:94:LEU:HB2	1:D:97:MSE:HG3	2.01	0.41
1:E:91:LEU:HD22	1:E:130:PRO:HB2	2.01	0.41
1:F:120:ILE:HA	1:F:121:PRO:HD3	1.94	0.41
1:G:101:THR:CG2	1:G:102:LYS:N	2.83	0.41
1:F:34:THR:HG21	1:F:95:GLY:HA2	2.03	0.41
1:F:55:LEU:HD22	1:G:55:LEU:HD11	2.01	0.41
1:G:122:THR:HA	1:G:126:GLU:O	2.21	0.41
1:E:97:MSE:HG2	1:E:173:ILE:O	2.20	0.41
1:G:37:ILE:HD12	1:G:37:ILE:O	2.21	0.41
1:H:72:GLY:HA3	1:H:147:ASP:HA	2.01	0.41
1:I:129:ASN:N	1:I:171:VAL:HG13	2.35	0.41
1:A:30:ILE:HG23	1:A:98:PRO:HG3	2.02	0.41
1:E:91:LEU:HD11	1:E:105:VAL:HG13	2.03	0.41
1:E:176:ASP:OD1	1:E:176:ASP:C	2.57	0.41
1:F:56:PRO:HG3	1:G:52:ASN:HD21	1.85	0.41
1:F:174:ASP:OD1	1:F:174:ASP:C	2.64	0.41
1:G:119:VAL:HG22	1:G:163:CYS:HB2	2.03	0.41
1:I:55:LEU:O	1:I:56:PRO:C	2.63	0.41
1:D:44:ILE:HD12	1:D:64:ILE:CD1	2.48	0.41
1:F:26:ASP:O	1:F:27:ASN:HB2	2.21	0.41
1:H:53:GLU:HG2	1:H:53:GLU:H	1.65	0.41
1:H:169:GLU:HG2	1:H:173:ILE:HD11	2.01	0.41
1:I:3:ILE:HD12	1:I:3:ILE:C	2.45	0.41
1:I:37:ILE:HG12	1:I:38:TYR:N	2.36	0.41
1:I:55:LEU:HD23	1:I:55:LEU:HA	1.78	0.41
1:I:104:ASP:O	1:I:108:ILE:HG13	2.20	0.41
1:I:122:THR:HG22	1:I:165:ILE:C	2.46	0.41
1:I:135:LYS:HA	1:I:138:PHE:CD2	2.56	0.41
1:A:66:ASN:ND2	1:A:66:ASN:C	2.77	0.41
1:A:135:LYS:HA	1:A:138:PHE:CD2	2.56	0.41
1:D:31:ILE:O	1:D:35:ILE:HG13	2.21	0.41
1:H:144:LEU:HD13	1:H:144:LEU:H	1.86	0.41
1:I:119:VAL:HB	1:I:132:LEU:HB3	2.02	0.41
1:A:2:ASN:HB2	1:A:88:ASP:OD2	2.21	0.40
1:B:66:ASN:ND2	1:B:68:PHE:N	2.60	0.40
1:B:94:LEU:CG	1:B:97:MSE:HE2	2.50	0.40
1:D:31:ILE:HG21	1:D:54:MSE:HE1	2.03	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:7:ILE:CD1	1:E:35:ILE:HD13	2.51	0.40
1:A:59:MSE:HB2	1:I:69:TRP:CZ2	2.56	0.40
1:G:170:GLY:C	1:G:172:LEU:H	2.28	0.40
1:H:66:ASN:HD21	1:H:75:THR:CG2	2.08	0.40
1:I:74:SER:HB2	1:I:144:LEU:CD2	2.49	0.40
1:A:10:ALA:HA	1:A:48:GLY:H	1.86	0.40
1:A:111:THR:HG21	1:A:165:ILE:HG23	2.02	0.40
1:D:12:GLU:CD	1:D:12:GLU:H	2.29	0.40
1:H:78:LYS:NZ	1:H:144:LEU:O	2.51	0.40
1:B:30:ILE:O	1:B:33:ARG:N	2.54	0.40
1:H:104:ASP:OD2	1:H:168:SER:OG	2.34	0.40
1:H:154:LEU:O	1:H:156:LYS:N	2.55	0.40
1:I:112:PHE:CE1	1:I:134:SER:HB3	2.56	0.40
1:I:135:LYS:C	1:I:137:LEU:H	2.28	0.40
1:B:31:ILE:HG12	1:B:54:MSE:CE	2.51	0.40
1:E:81:LEU:HD22	1:E:141:ILE:HD12	2.04	0.40
1:H:94:LEU:CD2	1:H:97:MSE:HE3	2.47	0.40
1:I:165:ILE:O	1:I:165:ILE: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	176/197 (89%)	155 (88%)	18 (10%)	3 (2%)	7	25
1	B	163/197 (83%)	141 (86%)	19 (12%)	3 (2%)	6	23
1	D	176/197 (89%)	157 (89%)	18 (10%)	1 (1%)	21	51
1	E	168/197 (85%)	150 (89%)	14 (8%)	4 (2%)	4	17
1	F	168/197 (85%)	152 (90%)	16 (10%)	0	100	100
1	G	162/197 (82%)	142 (88%)	15 (9%)	5 (3%)	3	12

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	169/197 (86%)	149 (88%)	14 (8%)	6 (4%)	2	10
1	I	167/197 (85%)	147 (88%)	17 (10%)	3 (2%)	6	23
All	All	1349/1576 (86%)	1193 (88%)	131 (10%)	25 (2%)	6	22

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	85	LYS
1	B	171	VAL
1	D	159	ILE
1	G	49	LYS
1	H	155	ASN
1	H	175	ILE
1	B	26	ASP
1	E	150	ALA
1	G	39	GLY
1	H	118	ALA
1	I	59	MSE
1	A	39	GLY
1	E	144	LEU
1	H	39	GLY
1	H	171	VAL
1	B	148	VAL
1	E	148	VAL
1	G	51	VAL
1	H	25	ILE
1	E	175	ILE
1	G	73	ILE
1	I	114	PRO
1	A	25	ILE
1	A	175	ILE
1	I	148	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	A	144/169 (85%)	128 (89%)	16 (11%)	6	20
1	B	136/169 (80%)	122 (90%)	14 (10%)	7	23
1	D	148/169 (88%)	130 (88%)	18 (12%)	5	16
1	E	139/169 (82%)	123 (88%)	16 (12%)	5	18
1	F	135/169 (80%)	125 (93%)	10 (7%)	13	37
1	G	137/169 (81%)	127 (93%)	10 (7%)	13	38
1	H	141/169 (83%)	121 (86%)	20 (14%)	3	12
1	I	137/169 (81%)	122 (89%)	15 (11%)	6	21
All	All	1117/1352 (83%)	998 (89%)	119 (11%)	6	21

All (119) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	12	GLU
1	A	28	THR
1	A	37	ILE
1	A	44	ILE
1	A	51	VAL
1	A	54	MSE
1	A	55	LEU
1	A	75	THR
1	A	77	LEU
1	A	91	LEU
1	A	100	VAL
1	A	127	ARG
1	A	141	ILE
1	A	144	LEU
1	A	165	ILE
1	B	26	ASP
1	B	41	LEU
1	B	44	ILE
1	B	51	VAL
1	B	70	ASN
1	B	71	GLU
1	B	81	LEU
1	B	82	ARG
1	B	122	THR
1	B	141	ILE
1	B	144	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	155	ASN
1	B	159	ILE
1	B	165	ILE
1	D	7	ILE
1	D	12	GLU
1	D	28	THR
1	D	30	ILE
1	D	41	LEU
1	D	49	LYS
1	D	51	VAL
1	D	52	ASN
1	D	54	MSE
1	D	57	LEU
1	D	85	LYS
1	D	91	LEU
1	D	103	GLU
1	D	127	ARG
1	D	137	LEU
1	D	165	ILE
1	D	174	ASP
1	D	175	ILE
1	E	32	MSE
1	E	51	VAL
1	E	64	ILE
1	E	66	ASN
1	E	70	ASN
1	E	77	LEU
1	E	100	VAL
1	E	103	GLU
1	E	106	ASN
1	E	132	LEU
1	E	141	ILE
1	E	155	ASN
1	E	157	ILE
1	E	162	LEU
1	E	165	ILE
1	E	176	ASP
1	F	21	LEU
1	F	28	THR
1	F	37	ILE
1	F	42	GLU
1	F	45	ILE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	F	94	LEU
1	F	107	LYS
1	F	141	ILE
1	F	165	ILE
1	F	173	ILE
1	G	30	ILE
1	G	55	LEU
1	G	79	LEU
1	G	132	LEU
1	G	139	ASN
1	G	153	ILE
1	G	165	ILE
1	G	166	GLU
1	G	171	VAL
1	G	175	ILE
1	H	2	ASN
1	H	22	LEU
1	H	30	ILE
1	H	33	ARG
1	H	37	ILE
1	H	43	LYS
1	H	53	GLU
1	H	57	LEU
1	H	61	GLN
1	H	79	LEU
1	H	98	PRO
1	H	100	VAL
1	H	103	GLU
1	H	140	GLU
1	H	141	ILE
1	H	144	LEU
1	H	165	ILE
1	H	166	GLU
1	H	172	LEU
1	H	176	ASP
1	I	7	ILE
1	I	22	LEU
1	I	24	LYS
1	I	32	MSE
1	I	37	ILE
1	I	51	VAL
1	I	66	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	I	81	LEU
1	I	103	GLU
1	I	114	PRO
1	I	141	ILE
1	I	144	LEU
1	I	165	ILE
1	I	179	GLU
1	I	180	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	ASN
1	A	123	HIS
1	A	129	ASN
1	A	155	ASN
1	B	27	ASN
1	B	66	ASN
1	B	70	ASN
1	B	106	ASN
1	D	2	ASN
1	D	66	ASN
1	D	106	ASN
1	E	2	ASN
1	E	66	ASN
1	E	70	ASN
1	E	155	ASN
1	F	52	ASN
1	G	52	ASN
1	G	110	ASN
1	H	2	ASN
1	H	106	ASN
1	H	115	ASN
1	I	66	ASN
1	I	70	ASN
1	I	139	ASN

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 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	173/197 (87%)	0.38	10 (5%) 29 22	42, 66, 92, 98	0
1	B	163/197 (82%)	0.66	17 (10%) 11 8	39, 70, 99, 100	0
1	D	173/197 (87%)	0.48	10 (5%) 29 22	45, 69, 94, 100	0
1	E	167/197 (84%)	0.54	12 (7%) 21 16	28, 69, 89, 93	0
1	F	167/197 (84%)	0.58	8 (4%) 35 28	34, 74, 99, 100	0
1	G	161/197 (81%)	0.56	8 (4%) 34 26	45, 70, 90, 96	0
1	H	168/197 (85%)	0.59	14 (8%) 17 12	33, 68, 93, 95	0
1	I	166/197 (84%)	0.72	17 (10%) 12 9	44, 77, 100, 100	0
All	All	1338/1576 (84%)	0.56	96 (7%) 21 16	28, 70, 96, 100	0

All (96) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	18	GLY	7.4
1	B	19	ASP	6.6
1	D	17	GLY	5.4
1	E	176	ASP	5.3
1	F	115	ASN	5.3
1	I	115	ASN	4.9
1	I	20	LYS	4.9
1	H	177	LYS	4.8
1	F	26	ASP	4.6
1	B	3	ILE	4.5
1	D	18	GLY	4.5
1	F	19	ASP	4.3
1	G	10	ALA	4.3
1	D	115	ASN	4.2
1	E	175	ILE	4.1
1	I	10	ALA	4.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	17	GLY	4.0
1	E	177	LYS	3.9
1	B	176	ASP	3.8
1	I	152	VAL	3.8
1	H	40	ASP	3.6
1	I	147	ASP	3.6
1	H	178	LYS	3.6
1	H	176	ASP	3.6
1	I	159	ILE	3.6
1	H	175	ILE	3.4
1	G	27	ASN	3.3
1	F	179	GLU	3.3
1	B	11	GLY	3.3
1	A	115	ASN	3.2
1	A	178	LYS	3.1
1	G	150	ALA	3.1
1	B	83	PHE	3.0
1	I	107	LYS	3.0
1	D	139	ASN	3.0
1	B	175	ILE	2.9
1	D	178	LYS	2.9
1	F	117	LYS	2.9
1	A	19	ASP	2.8
1	A	159	ILE	2.8
1	B	115	ASN	2.8
1	E	11	GLY	2.8
1	I	179	GLU	2.7
1	E	178	LYS	2.7
1	E	12	GLU	2.7
1	I	178	LYS	2.7
1	H	18	GLY	2.7
1	H	164	PHE	2.6
1	F	116	CYS	2.6
1	E	19	ASP	2.6
1	H	140	GLU	2.6
1	B	173	ILE	2.6
1	H	148	VAL	2.6
1	E	115	ASN	2.6
1	H	26	ASP	2.5
1	I	180	ASP	2.5
1	I	163	CYS	2.5
1	E	27	ASN	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	G	20	LYS	2.5
1	H	151	ARG	2.5
1	H	27	ASN	2.5
1	A	14	LYS	2.4
1	F	11	GLY	2.4
1	A	83	PHE	2.4
1	A	50	TYR	2.3
1	B	148	VAL	2.3
1	H	83	PHE	2.3
1	I	116	CYS	2.3
1	B	86	ASP	2.3
1	A	12	GLU	2.3
1	B	167	CYS	2.3
1	I	86	ASP	2.2
1	D	13	GLY	2.2
1	I	30	ILE	2.2
1	I	108	ILE	2.2
1	D	16	PHE	2.2
1	I	155	ASN	2.2
1	B	153	ILE	2.2
1	F	109	ILE	2.2
1	D	40	ASP	2.2
1	G	40	ASP	2.1
1	B	20	LYS	2.1
1	E	164	PHE	2.1
1	E	25	ILE	2.1
1	B	145	ARG	2.1
1	G	160	GLU	2.1
1	G	175	ILE	2.1
1	D	83	PHE	2.1
1	H	50	TYR	2.1
1	E	163	CYS	2.0
1	G	30	ILE	2.0
1	B	10	ALA	2.0
1	B	23	ALA	2.0
1	D	86	ASP	2.0
1	I	118	ALA	2.0
1	B	151	ARG	2.0

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

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.