



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:39 AM GMT

PDB ID : 3BNV
Title : Crystal structure of Cj0977, a sigma28-regulated virulence protein from *Campylobacter jejuni*.
Authors : Yokoyama, T.; Yeo, H.J.
Deposited on : 2007-12-14
Resolution : 2.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

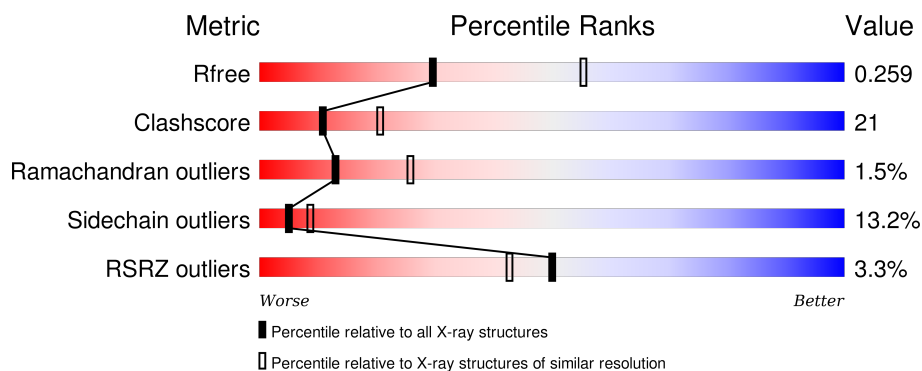
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	152	<div> <div>66%</div> <div>23%</div> <div>7%</div> </div>
1	B	152	<div> <div>59%</div> <div>24%</div> <div>6%</div> <div>10%</div> </div>
1	C	152	<div> <div>58%</div> <div>28%</div> <div>7%</div> <div>7%</div> </div>
1	D	152	<div> <div>69%</div> <div>22%</div> <div>6%</div> </div>
1	E	152	<div> <div>4%</div> <div>59%</div> <div>26%</div> <div>7%</div> <div>7%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	152	<div><div></div><div>4%</div><div>56%</div><div>28%</div><div>8%</div><div>9%</div></div>
1	G	152	<div><div></div><div>11%</div><div>47%</div><div>34%</div><div>• •</div><div>14%</div></div>
1	H	152	<div><div></div><div>2%</div><div>46%</div><div>32%</div><div>7%</div><div>•</div><div>14%</div></div>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 8728 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 Cj0977.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	142	Total	C	N	O	S	Se	0	0	0
			1108	712	175	217	2	2			
1	B	137	Total	C	N	O	S	Se	0	0	0
			1067	687	170	206	2	2			
1	C	142	Total	C	N	O	S	Se	0	0	0
			1091	703	172	212	2	2			
1	D	143	Total	C	N	O	S	Se	0	0	0
			1117	717	177	219	2	2			
1	E	142	Total	C	N	O	S	Se	0	1	0
			1112	715	175	217	2	3			
1	F	139	Total	C	N	O	S	Se	0	1	0
			1096	706	172	213	2	3			
1	G	130	Total	C	N	O	S	Se	0	0	0
			1012	654	160	194	2	2			
1	H	130	Total	C	N	O	S	Se	0	0	0
			1014	654	162	194	2	2			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	23	Total	O	0	0
			23	23		
2	B	18	Total	O	0	0
			18	18		
2	C	13	Total	O	0	0
			13	13		
2	D	19	Total	O	0	0
			19	19		
2	E	9	Total	O	0	0
			9	9		
2	F	13	Total	O	0	0
			13	13		

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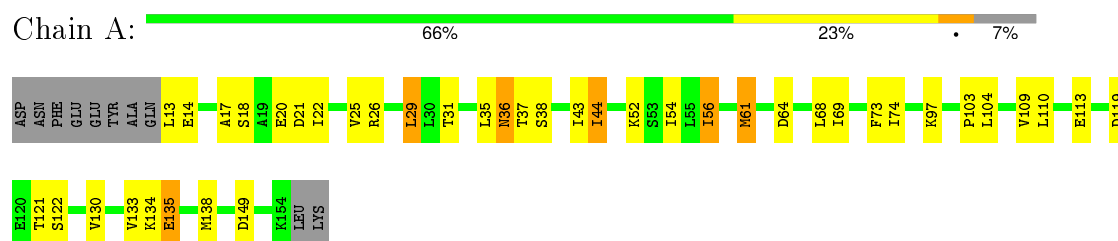
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	5	Total 5	O 5	0	0
2	H	11	Total 11	O 11	0	0

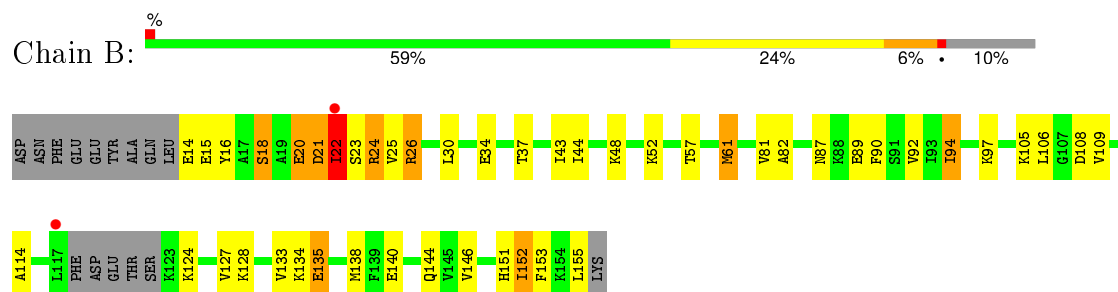
3 Residue-property plots

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.

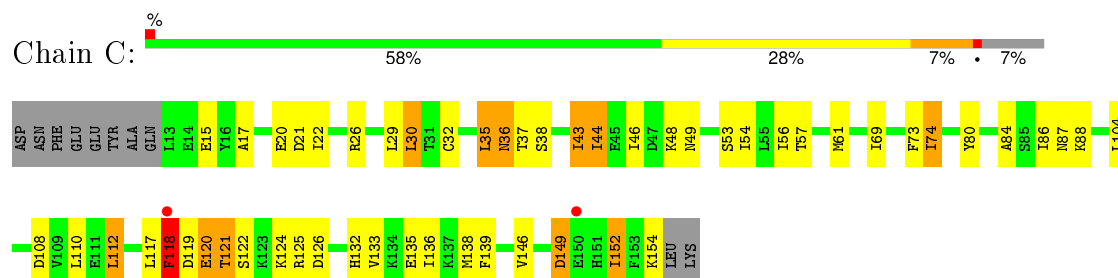
• Molecule 1: Cj0977



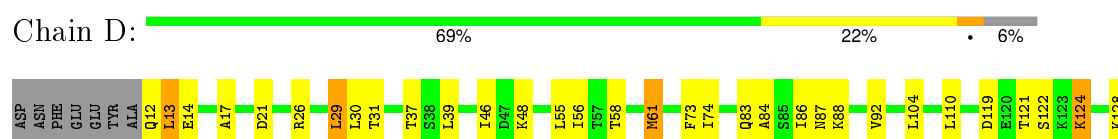
• Molecule 1: Cj0977

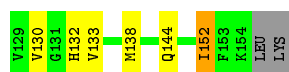


• Molecule 1: Cj0977

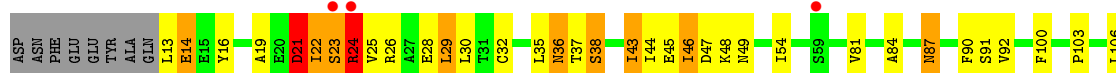


• Molecule 1: Cj0977





• Molecule 1: Cj0977



• Molecule 1: Cj0977



• Molecule 1: Cj0977



• Molecule 1: Cj0977



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	79.88Å 94.53Å 81.82Å 90.00° 99.32° 90.00°	Depositor
Resolution (Å)	37.93 – 2.60 37.92 – 2.60	Depositor EDS
% Data completeness (in resolution range)	96.8 (37.93-2.60) 96.8 (37.92-2.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.213 , 0.269 0.208 , 0.259	Depositor DCC
R_{free} test set	1790 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	44.4	Xtriage
Anisotropy	0.007	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 57.6	EDS
Estimated twinning fraction	0.027 for l,-k,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 35882 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8728	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.44% 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, 0.375 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	A	0.66	0/1123	0.71	0/1513
1	B	0.55	0/1080	0.70	0/1453
1	C	0.98	2/1106 (0.2%)	0.72	2/1492 (0.1%)
1	D	0.68	0/1132	0.74	0/1525
1	E	0.56	0/1131	0.70	0/1523
1	F	1.61	9/1114 (0.8%)	0.80	3/1498 (0.2%)
1	G	0.58	3/1026 (0.3%)	0.65	0/1382
1	H	0.85	5/1028 (0.5%)	0.74	1/1385 (0.1%)
All	All	0.88	19/8740 (0.2%)	0.72	6/11771 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1
1	H	0	1
All	All	0	2

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	28	GLU	CD-OE1	30.20	1.58	1.25
1	F	28	GLU	CD-OE2	23.39	1.51	1.25
1	C	15	GLU	CD-OE1	21.90	1.49	1.25
1	F	120	GLU	CD-OE1	19.18	1.46	1.25
1	F	120	GLU	CD-OE2	17.38	1.44	1.25
1	F	24	ARG	CZ-NH1	15.90	1.53	1.33
1	H	153	PHE	C-O	14.97	1.51	1.23
1	C	15	GLU	CD-OE2	14.86	1.42	1.25
1	H	153	PHE	CE2-CZ	8.38	1.53	1.37
1	H	153	PHE	CG-CD1	7.73	1.50	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	153	PHE	CE1-CZ	7.51	1.51	1.37
1	H	153	PHE	CG-CD2	6.18	1.48	1.38
1	G	120	GLU	CD-OE1	5.89	1.32	1.25
1	G	45	GLU	CD-OE1	5.71	1.31	1.25
1	F	24	ARG	NE-CZ	5.68	1.40	1.33
1	G	120	GLU	CD-OE2	5.64	1.31	1.25
1	F	24	ARG	CZ-NH2	5.30	1.40	1.33
1	F	120	GLU	C-N	5.20	1.46	1.34
1	F	120	GLU	CG-CD	5.04	1.59	1.51

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	24	ARG	NE-CZ-NH1	-9.15	115.72	120.30
1	F	28	GLU	OE1-CD-OE2	7.54	132.35	123.30
1	H	153	PHE	CB-CG-CD2	-6.89	115.97	120.80
1	F	64	ASP	CB-CA-C	-6.20	98.00	110.40
1	C	119	ASP	N-CA-C	5.16	124.92	111.00
1	C	15	GLU	OE1-CD-OE2	5.07	129.38	123.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	118	PHE	Peptide
1	H	153	PHE	Sidechain

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	1108	0	1113	41	0
1	B	1067	0	1081	55	0
1	C	1091	0	1084	45	1
1	D	1117	0	1121	23	0
1	E	1112	0	1118	58	0
1	F	1096	0	1102	49	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	1012	0	1027	48	1
1	H	1014	0	1027	62	0
2	A	23	0	0	2	0
2	B	18	0	0	2	0
2	C	13	0	0	0	0
2	D	19	0	0	2	0
2	E	9	0	0	2	0
2	F	13	0	0	4	0
2	G	5	0	0	1	0
2	H	11	0	0	6	0
All	All	8728	0	8673	365	1

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 21.

All (365) 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:C:43:ILE:CD1	1:C:43:ILE:H	1.60	1.13
1:B:24:ARG:HA	1:B:24:ARG:HH11	1.05	1.09
1:C:43:ILE:HD12	1:C:43:ILE:N	1.68	1.07
1:C:152:ILE:H	1:C:152:ILE:HD13	1.02	1.07
1:B:22:ILE:O	1:B:25:VAL:HG23	1.57	1.04
1:C:152:ILE:CD1	1:C:152:ILE:H	1.74	0.99
1:C:43:ILE:HD12	1:C:43:ILE:H	0.83	0.99
1:C:152:ILE:N	1:C:152:ILE:HD13	1.77	0.97
1:A:14:GLU:OE2	1:A:14:GLU:N	1.97	0.97
1:B:30:LEU:H	1:B:87:ASN:HD21	0.99	0.97
1:E:30:LEU:H	1:E:87:ASN:HD21	1.12	0.96
1:F:14:GLU:HG2	1:F:16:TYR:H	1.30	0.95
1:C:74:ILE:HD11	1:C:139:PHE:CD2	2.02	0.95
1:B:18:SER:HB3	1:B:21:ASP:HB2	1.47	0.94
1:C:30:LEU:H	1:C:87:ASN:HD21	1.10	0.94
1:E:19:ALA:HA	1:E:22:ILE:HD12	1.51	0.92
1:B:24:ARG:HH12	1:B:26:ARG:HB2	1.31	0.92
1:C:124:LYS:HG2	1:C:146:VAL:HG22	1.52	0.90
1:H:61:MSE:HE3	1:H:73:PHE:CB	2.03	0.89
1:F:57:THR:HG22	1:F:61:MSE:HE1	1.55	0.89
1:B:24:ARG:HA	1:B:24:ARG:NH1	1.86	0.89
1:H:74:ILE:HD11	1:H:139:PHE:CG	2.08	0.87
1:A:29:LEU:CD1	1:A:31:THR:HG22	2.04	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:86:ILE:HD13	1:H:145:VAL:HG21	1.57	0.86
1:H:86:ILE:HD13	1:H:145:VAL:CG2	2.05	0.85
1:F:148:THR:HG23	1:F:150:GLU:O	1.77	0.85
1:A:36:ASN:HD22	1:A:36:ASN:C	1.80	0.85
1:E:29:LEU:HD23	1:E:43:ILE:HG12	1.59	0.84
1:D:30:LEU:H	1:D:87:ASN:HD21	1.23	0.83
1:B:21:ASP:O	1:B:23:SER:N	2.10	0.83
1:H:61:MSE:HE3	1:H:73:PHE:HB3	1.61	0.83
1:F:135:GLU:OE2	1:F:135:GLU:N	2.12	0.80
1:E:135:GLU:HA	1:E:135:GLU:OE1	1.81	0.79
1:C:149:ASP:OD1	1:C:149:ASP:N	2.15	0.79
1:C:49:ASN:ND2	1:C:125:ARG:HH22	1.81	0.79
1:B:18:SER:CB	1:B:21:ASP:HB2	2.13	0.78
1:C:74:ILE:HD11	1:C:139:PHE:CG	2.18	0.78
1:E:132:HIS:C	1:E:138[A]:MSE:HE1	2.04	0.78
1:F:22:ILE:O	1:F:25:VAL:HG22	1.84	0.78
1:E:36:ASN:HD22	1:E:36:ASN:C	1.85	0.78
1:C:120:GLU:O	1:C:122:SER:N	2.17	0.78
1:E:133:VAL:N	1:E:138[A]:MSE:HE1	2.00	0.77
1:A:135:GLU:OE1	1:A:135:GLU:N	2.17	0.76
1:B:18:SER:HB3	1:B:21:ASP:CB	2.16	0.75
1:F:104:LEU:HD13	1:F:138[B]:MSE:HE3	1.68	0.75
1:F:74:ILE:HD11	1:F:139:PHE:CD1	2.22	0.75
1:E:30:LEU:H	1:E:87:ASN:ND2	1.85	0.75
1:F:49:ASN:OD1	1:F:125:ARG:NH2	2.15	0.74
1:B:21:ASP:C	1:B:23:SER:H	1.90	0.74
1:H:74:ILE:HD11	1:H:139:PHE:CD1	2.22	0.74
1:F:31:THR:HG21	1:F:83:GLN:HB3	1.70	0.74
1:H:86:ILE:CD1	1:H:145:VAL:HG21	2.17	0.74
1:E:44:ILE:HG21	1:E:54:ILE:HD12	1.70	0.74
1:H:30:LEU:H	1:H:87:ASN:HD21	1.36	0.73
1:E:36:ASN:ND2	1:E:38:SER:H	1.87	0.73
1:F:151:HIS:HE1	1:F:153:PHE:CD1	2.05	0.73
1:H:148:THR:OG1	1:H:150:GLU:O	2.02	0.73
1:C:30:LEU:H	1:C:87:ASN:ND2	1.86	0.72
1:C:43:ILE:HD11	1:C:80:TYR:OH	1.89	0.72
1:B:30:LEU:H	1:B:87:ASN:ND2	1.83	0.72
1:E:14:GLU:HG2	1:E:14:GLU:O	1.88	0.71
1:C:117:LEU:O	1:C:118:PHE:O	2.09	0.71
1:F:42:THR:OG1	2:F:200:HOH:O	2.06	0.71
1:F:31:THR:HG22	1:F:32:CYS:SG	2.32	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:30:LEU:N	1:E:87:ASN:HD21	1.88	0.69
1:A:29:LEU:HD11	1:A:31:THR:HG22	1.74	0.69
1:F:148:THR:CG2	1:F:150:GLU:O	2.40	0.69
1:D:133:VAL:HG23	1:D:138:MSE:HG3	1.74	0.69
1:E:21:ASP:C	1:E:21:ASP:OD2	2.30	0.68
1:A:68:LEU:HD13	1:B:152:ILE:HD11	1.75	0.68
1:H:24:ARG:HG2	1:H:24:ARG:O	1.92	0.68
1:A:36:ASN:ND2	1:A:38:SER:H	1.91	0.68
1:G:86:ILE:HD13	1:G:125:ARG:NH1	2.09	0.68
1:H:26:ARG:CG	1:H:26:ARG:HH11	2.07	0.68
1:G:97:LYS:HD3	1:H:97:LYS:HG2	1.74	0.67
1:G:90:PHE:CD1	1:G:151:HIS:N	2.63	0.67
1:C:61:MSE:HB3	1:C:73:PHE:CD2	2.28	0.67
1:E:24:ARG:HE	1:E:45:GLU:HA	1.60	0.67
1:E:16:TYR:HB2	1:E:54:ILE:CD1	2.25	0.67
1:F:64:ASP:HB3	1:F:66:GLN:H	1.60	0.66
1:G:120:GLU:O	1:G:120:GLU:HG3	1.95	0.66
1:D:30:LEU:H	1:D:87:ASN:ND2	1.92	0.66
1:G:74:ILE:HD11	1:G:139:PHE:HD1	1.61	0.66
1:E:90:PHE:HA	2:E:198:HOH:O	1.95	0.66
1:A:17:ALA:HB2	1:A:44:ILE:CG1	2.26	0.65
1:E:119:ASP:OD1	1:E:122:SER:HB3	1.97	0.65
1:A:17:ALA:HB2	1:A:44:ILE:HG12	1.79	0.65
1:C:152:ILE:CD1	1:C:152:ILE:N	2.44	0.65
1:H:83:GLN:HG3	1:H:93:ILE:HD11	1.79	0.64
1:C:30:LEU:N	1:C:87:ASN:HD21	1.90	0.64
1:E:24:ARG:HG3	1:E:25:VAL:HG23	1.80	0.64
1:H:82:ALA:HB3	1:H:143:ILE:HD13	1.79	0.64
1:B:135:GLU:CA	1:B:135:GLU:OE2	2.46	0.64
1:E:16:TYR:HB2	1:E:54:ILE:HD13	1.80	0.63
1:E:36:ASN:HD22	1:E:38:SER:H	1.46	0.63
1:E:21:ASP:OD2	1:E:22:ILE:N	2.31	0.63
1:E:133:VAL:HG23	1:E:138[A]:MSE:HE2	1.81	0.62
1:F:25:VAL:HG23	1:F:26:ARG:N	2.13	0.62
1:H:82:ALA:CB	1:H:143:ILE:HD13	2.31	0.61
1:A:103:PRO:HG3	1:B:152:ILE:HD13	1.83	0.61
1:D:152:ILE:O	1:D:152:ILE:HG13	1.99	0.61
1:B:135:GLU:HA	1:B:135:GLU:OE2	2.01	0.60
1:G:122:SER:OG	1:G:123:LYS:N	2.35	0.60
1:C:29:LEU:HD11	1:C:43:ILE:HG12	1.82	0.60
1:A:17:ALA:CB	1:A:44:ILE:HG12	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:150:GLU:OE2	1:G:154:LYS:HD3	2.00	0.60
1:E:49:ASN:ND2	1:E:125:ARG:HH22	1.98	0.60
1:E:132:HIS:CA	1:E:138[A]:MSE:HE1	2.32	0.60
1:E:23:SER:O	1:E:25:VAL:N	2.34	0.60
1:D:12:GLN:HG3	1:D:56:ILE:HD13	1.84	0.60
1:H:79:ASN:OD1	1:H:93:ILE:HD13	2.02	0.60
1:H:26:ARG:HH22	1:H:42:THR:CG2	2.14	0.59
1:D:83:GLN:NE2	2:D:202:HOH:O	2.31	0.59
1:G:86:ILE:HG21	1:G:145:VAL:HG11	1.84	0.59
1:E:135:GLU:CA	1:E:135:GLU:OE1	2.48	0.59
1:G:74:ILE:HD11	1:G:139:PHE:CD1	2.37	0.59
1:H:31:THR:HG21	1:H:83:GLN:HB3	1.82	0.59
1:E:49:ASN:HD21	1:E:125:ARG:HH22	1.50	0.59
1:G:43:ILE:HD11	1:G:80:TYR:CE2	2.37	0.59
1:H:151:HIS:O	1:H:153:PHE:N	2.36	0.59
1:D:121:THR:HG22	1:D:122:SER:N	2.17	0.59
1:F:25:VAL:CG2	1:F:26:ARG:N	2.66	0.59
1:B:57:THR:HG22	1:B:61:MSE:HE2	1.84	0.59
1:G:79:ASN:HD22	1:H:72:ALA:HB2	1.67	0.58
1:C:110:LEU:HD12	1:C:132:HIS:O	2.03	0.58
1:A:36:ASN:C	1:A:36:ASN:ND2	2.51	0.58
1:F:152:ILE:O	1:F:152:ILE:CG1	2.51	0.58
1:E:29:LEU:HD12	1:E:29:LEU:O	2.04	0.57
1:E:24:ARG:NE	1:E:45:GLU:HA	2.19	0.57
1:H:86:ILE:CD1	1:H:145:VAL:CG2	2.78	0.57
1:B:24:ARG:HH11	1:B:24:ARG:CA	1.98	0.57
1:A:68:LEU:HD13	1:B:152:ILE:CD1	2.33	0.57
1:E:29:LEU:CD2	1:E:43:ILE:HG12	2.31	0.57
1:D:128:LYS:HD2	2:D:207:HOH:O	2.04	0.57
1:C:43:ILE:HD11	1:C:80:TYR:CZ	2.39	0.56
1:A:36:ASN:HD22	1:A:37:THR:N	2.01	0.56
1:B:22:ILE:CG2	1:B:22:ILE:O	2.54	0.56
1:D:124:LYS:HE3	1:D:144:GLN:OE1	2.04	0.56
1:A:17:ALA:HB2	1:A:44:ILE:HD11	1.87	0.56
1:B:21:ASP:C	1:B:23:SER:N	2.53	0.56
1:E:133:VAL:N	1:E:138[A]:MSE:CE	2.67	0.56
1:H:86:ILE:HD13	1:H:145:VAL:HG23	1.85	0.56
1:G:133:VAL:HB	1:G:138:MSE:HE3	1.88	0.56
1:E:21:ASP:O	1:E:25:VAL:HB	2.06	0.56
1:F:74:ILE:HD13	1:F:100:PHE:HD2	1.70	0.55
1:F:102:ALA:HB3	1:F:138[A]:MSE:HG3	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:45:GLU:OE2	1:G:52:LYS:HD3	2.07	0.55
1:F:152:ILE:O	1:F:152:ILE:HG12	2.06	0.55
1:H:93:ILE:HG12	1:H:143:ILE:CD1	2.36	0.55
1:H:101:TYR:HE2	1:H:140:GLU:HB2	1.72	0.55
1:G:136:ILE:HB	1:G:138:MSE:HE2	1.88	0.55
1:C:17:ALA:HB3	1:C:44:ILE:HG13	1.88	0.55
1:B:44:ILE:HD11	1:B:52:LYS:HD3	1.89	0.55
1:A:17:ALA:HB2	1:A:44:ILE:CD1	2.36	0.55
1:G:44:ILE:HD13	1:G:54:ILE:HD11	1.89	0.54
1:E:152:ILE:O	1:E:152:ILE:HG13	2.06	0.54
1:D:48:LYS:HG2	1:H:118:PHE:CG	2.42	0.54
1:B:30:LEU:N	1:B:87:ASN:HD21	1.84	0.54
1:B:94:ILE:HG22	1:B:144:GLN:HB2	1.90	0.54
1:H:88:LYS:CE	2:H:198:HOH:O	2.56	0.54
1:C:36:ASN:HD22	1:C:36:ASN:C	2.10	0.54
1:C:133:VAL:HG23	1:C:138:MSE:HG3	1.90	0.54
1:H:130:VAL:HG12	2:H:195:HOH:O	2.07	0.53
1:H:61:MSE:HE3	1:H:73:PHE:HB2	1.86	0.53
1:H:26:ARG:HH11	1:H:26:ARG:HG3	1.73	0.53
1:A:74:ILE:HD11	2:A:211:HOH:O	2.08	0.53
1:G:121:THR:HG23	1:G:122:SER:N	2.24	0.53
1:A:61:MSE:HB3	1:A:73:PHE:CD2	2.44	0.53
1:B:43:ILE:N	1:B:43:ILE:HD13	2.24	0.53
1:H:126:ASP:OD1	1:H:144:GLN:HG2	2.08	0.53
1:H:30:LEU:H	1:H:87:ASN:ND2	2.05	0.53
1:C:53:SER:O	1:C:54:ILE:HD12	2.08	0.52
1:G:29:LEU:HD11	1:G:43:ILE:HD12	1.91	0.52
1:G:74:ILE:CD1	1:G:139:PHE:CD1	2.93	0.52
1:C:112:LEU:HD12	1:C:112:LEU:N	2.25	0.52
1:C:49:ASN:ND2	1:C:125:ARG:NH2	2.53	0.52
1:A:68:LEU:CD1	1:B:152:ILE:HD11	2.40	0.51
1:F:152:ILE:HG22	2:F:203:HOH:O	2.10	0.51
1:B:18:SER:CB	1:B:21:ASP:CB	2.83	0.51
1:B:23:SER:C	1:B:25:VAL:H	2.13	0.51
1:F:24:ARG:HG2	1:F:25:VAL:N	2.26	0.51
1:E:23:SER:O	1:E:24:ARG:C	2.48	0.51
1:H:88:LYS:HE2	2:H:198:HOH:O	2.10	0.51
1:C:49:ASN:HD22	1:C:125:ARG:HH22	1.58	0.51
1:G:126:ASP:OD1	1:G:144:GLN:HG3	2.10	0.51
1:B:94:ILE:HG12	2:B:210:HOH:O	2.10	0.51
1:E:103:PRO:HG3	1:F:152:ILE:HD13	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:74:ILE:HD11	1:H:139:PHE:CB	2.42	0.50
1:F:151:HIS:CE1	1:F:153:PHE:CD1	2.93	0.50
1:D:46:ILE:HD11	1:D:84:ALA:O	2.11	0.50
1:D:30:LEU:N	1:D:87:ASN:HD21	2.02	0.50
1:D:110:LEU:HD12	1:D:132:HIS:O	2.10	0.50
1:A:134:LYS:HB2	1:A:135:GLU:OE1	2.12	0.50
1:G:132:HIS:ND1	1:G:137:LYS:HA	2.26	0.50
1:B:24:ARG:NH1	1:B:26:ARG:HB2	2.14	0.50
1:H:26:ARG:CG	1:H:26:ARG:NH1	2.70	0.50
1:A:61:MSE:CE	1:A:69:ILE:CG2	2.90	0.50
1:G:39:LEU:HD21	1:H:39:LEU:HD13	1.93	0.50
1:A:44:ILE:HG22	1:A:52:LYS:HB3	1.93	0.50
1:H:92:VAL:HG13	1:H:152:ILE:HD13	1.94	0.50
1:C:43:ILE:HD11	1:C:80:TYR:CE1	2.47	0.49
1:C:57:THR:HB	1:C:69:ILE:HD11	1.93	0.49
1:H:79:ASN:OD1	1:H:93:ILE:CD1	2.60	0.49
1:G:100:PHE:N	1:G:100:PHE:CD1	2.78	0.49
1:E:19:ALA:CA	1:E:22:ILE:HD12	2.34	0.49
1:A:97:LYS:HD2	1:B:97:LYS:HG3	1.93	0.49
1:E:46:ILE:HD11	1:E:84:ALA:O	2.13	0.49
1:E:16:TYR:HB2	1:E:54:ILE:HD11	1.94	0.49
1:A:25:VAL:HG23	1:A:26:ARG:N	2.27	0.49
1:H:28:GLU:HG2	1:H:28:GLU:O	2.12	0.49
1:E:36:ASN:HD22	1:E:37:THR:N	2.10	0.49
1:E:118:PHE:CD2	1:E:118:PHE:C	2.86	0.49
1:E:21:ASP:O	1:E:25:VAL:N	2.34	0.49
1:G:33:PRO:CG	2:H:202:HOH:O	2.60	0.49
1:G:114:ALA:HB2	1:G:129:VAL:HG22	1.95	0.49
1:G:121:THR:CG2	1:G:122:SER:N	2.76	0.48
1:B:20:GLU:O	1:B:23:SER:N	2.47	0.48
1:H:124:LYS:NZ	1:H:124:LYS:HB3	2.29	0.48
1:B:81:VAL:HG13	1:B:114:ALA:HB2	1.95	0.48
1:B:24:ARG:N	1:B:24:ARG:HD2	2.29	0.48
1:D:104:LEU:HD12	1:D:138:MSE:HE3	1.95	0.48
1:C:20:GLU:O	1:C:22:ILE:N	2.46	0.48
1:G:90:PHE:HD1	1:G:151:HIS:CA	2.27	0.48
1:D:74:ILE:N	1:D:74:ILE:HD12	2.28	0.48
1:G:120:GLU:O	1:G:120:GLU:CG	2.62	0.48
1:E:132:HIS:HA	1:E:138[A]:MSE:HE1	1.96	0.48
1:F:144:GLN:HG3	2:F:202:HOH:O	2.14	0.47
1:G:134:LYS:HB3	1:G:135:GLU:OE1	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:133:VAL:HG23	1:H:138:MSE:HG2	1.95	0.47
1:A:20:GLU:O	1:A:21:ASP:C	2.53	0.47
1:A:36:ASN:HD22	1:A:38:SER:H	1.57	0.47
1:F:24:ARG:O	1:F:27:ALA:N	2.48	0.47
1:F:25:VAL:CG2	1:F:26:ARG:H	2.27	0.47
1:C:117:LEU:HD12	1:C:126:ASP:HB2	1.96	0.47
1:A:64:ASP:HA	1:B:34:GLU:HB2	1.95	0.47
1:B:57:THR:HG22	1:B:61:MSE:CE	2.44	0.47
1:H:26:ARG:NH2	1:H:42:THR:CG2	2.77	0.47
1:G:44:ILE:HG21	1:G:54:ILE:HG13	1.96	0.47
1:B:124:LYS:HG3	1:B:146:VAL:HG22	1.96	0.47
1:G:48:LYS:HG3	1:G:48:LYS:O	2.15	0.47
1:F:71:ASP:O	1:F:74:ILE:HG23	2.14	0.46
1:C:57:THR:HG23	1:C:108:ASP:O	2.15	0.46
1:E:16:TYR:CB	1:E:54:ILE:HD11	2.45	0.46
1:A:29:LEU:HD11	1:A:31:THR:CG2	2.44	0.46
1:H:26:ARG:NH2	2:H:199:HOH:O	2.33	0.46
1:C:57:THR:HB	1:C:69:ILE:CD1	2.45	0.46
1:H:32:CYS:N	1:H:33:PRO:HD3	2.31	0.46
1:F:92:VAL:HG22	1:F:93:ILE:N	2.31	0.46
1:F:54:ILE:HD12	1:F:54:ILE:C	2.35	0.46
1:E:29:LEU:HA	1:E:87:ASN:OD1	2.16	0.46
1:F:32:CYS:N	1:F:33:PRO:CD	2.79	0.46
1:A:17:ALA:CB	1:A:44:ILE:CG1	2.91	0.46
1:H:26:ARG:NH2	1:H:42:THR:HG23	2.31	0.46
1:B:133:VAL:O	1:B:134:LYS:HB2	2.15	0.46
1:H:61:MSE:HE2	1:H:69:ILE:HG23	1.97	0.45
1:B:82:ALA:HA	1:B:127:VAL:HG11	1.99	0.45
1:C:120:GLU:O	1:C:121:THR:C	2.55	0.45
1:A:61:MSE:HE2	1:A:69:ILE:CG2	2.46	0.45
1:B:133:VAL:HG23	1:B:138:MSE:HG3	1.98	0.45
1:B:105:LYS:O	1:B:108:ASP:CG	2.55	0.45
1:B:151:HIS:HE1	1:B:153:PHE:CD1	2.35	0.45
1:B:20:GLU:O	1:B:21:ASP:C	2.55	0.45
1:B:22:ILE:O	1:B:22:ILE:HG23	2.17	0.45
1:F:61:MSE:CE	1:F:69:ILE:CG2	2.94	0.45
1:F:25:VAL:O	1:F:26:ARG:C	2.53	0.45
1:F:61:MSE:HE3	1:F:69:ILE:CG2	2.47	0.45
1:B:135:GLU:N	1:B:135:GLU:OE2	2.50	0.45
1:G:39:LEU:HD23	1:G:73:PHE:HE1	1.82	0.45
1:A:97:LYS:HE2	2:A:213:HOH:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:20:GLU:C	1:C:22:ILE:H	2.20	0.45
1:F:21:ASP:O	1:F:25:VAL:HG13	2.17	0.45
1:A:22:ILE:O	1:A:25:VAL:HG22	2.17	0.44
1:E:100:PHE:CE2	1:E:139:PHE:HE1	2.34	0.44
1:D:86:ILE:HG22	1:D:88:LYS:HG2	1.99	0.44
1:H:101:TYR:CE2	1:H:140:GLU:HB2	2.52	0.44
1:H:61:MSE:HE2	1:H:69:ILE:CG2	2.47	0.44
1:H:124:LYS:NZ	1:H:124:LYS:CB	2.81	0.44
1:A:133:VAL:HG23	1:A:138:MSE:HG3	1.99	0.44
1:B:89:GLU:HG2	1:B:90:PHE:CE2	2.52	0.44
1:A:52:LYS:HG3	1:A:113:GLU:HG2	1.98	0.44
1:F:22:ILE:HG22	1:F:23:SER:N	2.32	0.44
1:A:18:SER:O	1:A:22:ILE:HD12	2.18	0.44
1:D:14:GLU:HA	1:D:17:ALA:O	2.17	0.44
1:F:86:ILE:HD11	1:F:125:ARG:HB2	2.00	0.43
1:G:103:PRO:HD2	2:G:195:HOH:O	2.16	0.43
1:B:16:TYR:CZ	1:B:109:VAL:HG21	2.53	0.43
1:B:22:ILE:HA	1:B:22:ILE:HD13	1.84	0.43
1:H:151:HIS:C	1:H:153:PHE:N	2.72	0.43
1:E:91:SER:N	2:E:198:HOH:O	2.44	0.43
1:G:90:PHE:HD1	1:G:151:HIS:HA	1.84	0.43
1:C:104:LEU:HD12	1:C:138:MSE:HE3	2.00	0.43
1:A:104:LEU:HD21	1:A:110:LEU:HB2	2.01	0.43
1:E:21:ASP:CG	1:E:22:ILE:N	2.71	0.43
1:E:133:VAL:HG23	1:E:138[A]:MSE:CE	2.48	0.43
1:G:64:ASP:OD1	1:G:66:GLN:N	2.39	0.43
1:C:32:CYS:HB3	1:C:35:LEU:HB2	2.01	0.43
1:G:118:PHE:H	1:G:118:PHE:HD1	1.66	0.43
1:G:72:ALA:HB2	1:H:79:ASN:HD22	1.83	0.43
1:G:86:ILE:HD13	1:G:125:ARG:HH11	1.81	0.43
1:G:152:ILE:HD11	1:H:68:LEU:HD13	2.01	0.43
1:D:55:LEU:HD21	1:D:61:MSE:HE1	2.00	0.43
1:H:55:LEU:HD21	1:H:61:MSE:HE1	2.00	0.43
1:C:36:ASN:ND2	1:C:38:SER:H	2.17	0.43
1:G:33:PRO:HG2	2:H:202:HOH:O	2.19	0.43
1:F:61:MSE:CE	1:F:69:ILE:HG21	2.49	0.42
1:E:133:VAL:CG2	1:E:138[A]:MSE:HE2	2.47	0.42
1:A:61:MSE:CE	1:A:69:ILE:HG21	2.48	0.42
1:E:32:CYS:HB3	1:E:35:LEU:HB2	2.01	0.42
1:E:81:VAL:HG13	1:E:114:ALA:HB2	2.01	0.42
1:B:94:ILE:HD13	1:B:94:ILE:HG23	1.78	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:58:THR:O	1:D:61:MSE:HG3	2.19	0.42
1:A:121:THR:OG1	1:A:122:SER:N	2.53	0.42
1:G:32:CYS:HA	1:H:63:ALA:O	2.19	0.42
1:E:44:ILE:HD13	1:E:54:ILE:HD11	2.01	0.42
1:C:36:ASN:HD22	1:C:37:THR:N	2.17	0.42
1:G:152:ILE:HG13	1:G:152:ILE:H	1.56	0.42
1:E:21:ASP:HA	1:E:25:VAL:HG23	2.02	0.42
1:B:152:ILE:HG13	2:B:197:HOH:O	2.20	0.42
1:D:74:ILE:N	1:D:74:ILE:CD1	2.81	0.42
1:B:128:LYS:NZ	1:B:140:GLU:OE2	2.52	0.42
1:E:36:ASN:C	1:E:36:ASN:ND2	2.58	0.42
1:H:26:ARG:HG2	1:H:26:ARG:NH1	2.34	0.42
1:F:144:GLN:CG	2:F:202:HOH:O	2.67	0.42
1:F:77:ALA:HB1	1:F:112:LEU:CD1	2.50	0.42
1:F:24:ARG:O	1:F:25:VAL:C	2.57	0.42
1:A:17:ALA:HB3	1:A:22:ILE:HD11	2.02	0.42
1:G:102:ALA:HB3	1:G:138:MSE:SE	2.69	0.42
1:G:32:CYS:HB3	1:G:35:LEU:HB2	2.01	0.42
1:F:90:PHE:CB	1:F:148:THR:HG22	2.49	0.42
1:H:93:ILE:HG12	1:H:143:ILE:HD12	2.00	0.42
1:A:25:VAL:HG21	1:A:43:ILE:O	2.20	0.41
1:B:89:GLU:HG2	1:B:90:PHE:CD2	2.55	0.41
1:D:61:MSE:HE3	1:D:73:PHE:HB2	2.02	0.41
1:H:52:LYS:HE2	1:H:111:GLU:OE1	2.19	0.41
1:G:45:GLU:HG2	1:G:52:LYS:HB3	2.00	0.41
1:D:13:LEU:HD13	1:D:13:LEU:HA	1.74	0.41
1:F:134:LYS:HB3	1:F:135:GLU:OE2	2.19	0.41
1:H:26:ARG:HH22	1:H:42:THR:HG23	1.84	0.41
1:H:119:ASP:OD2	1:H:122:SER:HB3	2.19	0.41
1:E:48:LYS:O	1:E:49:ASN:HB2	2.21	0.41
1:H:59:SER:O	1:H:62:VAL:HG22	2.21	0.41
1:B:23:SER:C	1:B:25:VAL:N	2.74	0.41
1:F:25:VAL:HG21	1:F:43:ILE:O	2.19	0.41
1:F:74:ILE:HD12	1:F:138[B]:MSE:HE1	2.02	0.41
1:H:37:THR:HB	1:H:41:GLY:O	2.21	0.41
1:F:105:LYS:HB3	1:F:105:LYS:HE3	1.86	0.41
1:B:21:ASP:O	1:B:22:ILE:C	2.58	0.41
1:E:16:TYR:CB	1:E:54:ILE:CD1	2.96	0.41
1:E:46:ILE:HG13	1:E:47:ASP:N	2.36	0.41
1:G:98:CYS:HB2	1:G:100:PHE:CE1	2.56	0.41
1:B:151:HIS:CE1	1:B:153:PHE:CD1	3.09	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:152:ILE:CD1	1:H:68:LEU:HD13	2.51	0.41
1:C:36:ASN:ND2	1:C:36:ASN:C	2.73	0.40
1:C:57:THR:CG2	1:C:69:ILE:HD13	2.51	0.40
1:D:29:LEU:HD22	1:D:31:THR:HG22	2.03	0.40
1:C:46:ILE:HD11	1:C:84:ALA:O	2.20	0.40
1:C:74:ILE:CD1	1:C:139:PHE:CG	2.99	0.40
1:F:26:ARG:O	1:F:29:LEU:HD12	2.22	0.40
1:G:90:PHE:CD1	1:G:151:HIS:CA	3.03	0.40
1:E:25:VAL:HG12	1:E:26:ARG:N	2.36	0.40
1:F:54:ILE:CD1	1:F:56:ILE:HG13	2.51	0.40
1:B:105:LYS:O	1:B:108:ASP:OD2	2.39	0.40
1:A:56:ILE:HG12	1:A:56:ILE:O	2.21	0.40
1:G:72:ALA:CB	1:H:76:ALA:HA	2.51	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:120:GLU:OE1	1:G:150:GLU:OE1[2_647]	2.18	0.02

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	140/152 (92%)	134 (96%)	5 (4%)	1 (1%)	26	51
1	B	133/152 (88%)	128 (96%)	3 (2%)	2 (2%)	13	26
1	C	140/152 (92%)	131 (94%)	5 (4%)	4 (3%)	6	9
1	D	141/152 (93%)	135 (96%)	5 (4%)	1 (1%)	26	51
1	E	141/152 (93%)	130 (92%)	7 (5%)	4 (3%)	6	10
1	F	136/152 (90%)	129 (95%)	5 (4%)	2 (2%)	13	26

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	G	128/152 (84%)	123 (96%)	4 (3%)	1 (1%)	24	46
1	H	128/152 (84%)	123 (96%)	4 (3%)	1 (1%)	24	46
All	All	1087/1216 (89%)	1033 (95%)	38 (4%)	16 (2%)	13	26

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	21	ASP
1	B	22	ILE
1	C	118	PHE
1	C	120	GLU
1	C	121	THR
1	E	21	ASP
1	G	151	HIS
1	H	152	ILE
1	C	21	ASP
1	A	119	ASP
1	D	119	ASP
1	E	24	ARG
1	F	23	SER
1	E	23	SER
1	E	22	ILE
1	F	25	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	121/128 (94%)	109 (90%)	12 (10%)	10	18
1	B	116/128 (91%)	100 (86%)	16 (14%)	4	7
1	C	116/128 (91%)	98 (84%)	18 (16%)	3	5
1	D	122/128 (95%)	111 (91%)	11 (9%)	12	23
1	E	122/128 (95%)	100 (82%)	22 (18%)	2	3

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	121/128 (94%)	107 (88%)	14 (12%)	7	12
1	G	111/128 (87%)	96 (86%)	15 (14%)	5	8
1	H	111/128 (87%)	93 (84%)	18 (16%)	3	5
All	All	940/1024 (92%)	814 (87%)	126 (13%)	5	8

All (126) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LEU
1	A	29	LEU
1	A	35	LEU
1	A	36	ASN
1	A	44	ILE
1	A	54	ILE
1	A	56	ILE
1	A	61	MSE
1	A	109	VAL
1	A	130	VAL
1	A	135	GLU
1	A	149	ASP
1	B	14	GLU
1	B	15	GLU
1	B	18	SER
1	B	20	GLU
1	B	22	ILE
1	B	24	ARG
1	B	26	ARG
1	B	37	THR
1	B	48	LYS
1	B	61	MSE
1	B	92	VAL
1	B	94	ILE
1	B	106	LEU
1	B	135	GLU
1	B	152	ILE
1	B	155	LEU
1	C	26	ARG
1	C	30	LEU
1	C	35	LEU
1	C	36	ASN
1	C	43	ILE

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Mol	Chain	Res	Type
1	C	44	ILE
1	C	48	LYS
1	C	56	ILE
1	C	74	ILE
1	C	86	ILE
1	C	88	LYS
1	C	112	LEU
1	C	118	PHE
1	C	135	GLU
1	C	136	ILE
1	C	149	ASP
1	C	152	ILE
1	C	154	LYS
1	D	13	LEU
1	D	21	ASP
1	D	26	ARG
1	D	29	LEU
1	D	37	THR
1	D	39	LEU
1	D	61	MSE
1	D	92	VAL
1	D	124	LYS
1	D	130	VAL
1	D	152	ILE
1	E	13	LEU
1	E	14	GLU
1	E	21	ASP
1	E	24	ARG
1	E	28	GLU
1	E	29	LEU
1	E	36	ASN
1	E	38	SER
1	E	43	ILE
1	E	46	ILE
1	E	87	ASN
1	E	92	VAL
1	E	106	LEU
1	E	109	VAL
1	E	113	GLU
1	E	118	PHE
1	E	123	LYS
1	E	128	LYS

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Mol	Chain	Res	Type
1	E	138[A]	MSE
1	E	138[B]	MSE
1	E	146	VAL
1	E	149	ASP
1	F	15	GLU
1	F	26	ARG
1	F	29	LEU
1	F	31	THR
1	F	35	LEU
1	F	46	ILE
1	F	74	ILE
1	F	86	ILE
1	F	104	LEU
1	F	109	VAL
1	F	138[A]	MSE
1	F	138[B]	MSE
1	F	148	THR
1	F	154	LYS
1	G	34	GLU
1	G	42	THR
1	G	50	TYR
1	G	66	GLN
1	G	74	ILE
1	G	89	GLU
1	G	92	VAL
1	G	93	ILE
1	G	112	LEU
1	G	117	LEU
1	G	118	PHE
1	G	119	ASP
1	G	120	GLU
1	G	124	LYS
1	G	152	ILE
1	H	24	ARG
1	H	26	ARG
1	H	31	THR
1	H	35	LEU
1	H	48	LYS
1	H	60	GLU
1	H	65	ASP
1	H	74	ILE
1	H	87	ASN

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Mol	Chain	Res	Type
1	H	97	LYS
1	H	109	VAL
1	H	123	LYS
1	H	124	LYS
1	H	135	GLU
1	H	138	MSE
1	H	143	ILE
1	H	146	VAL
1	H	153	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	ASN
1	B	79	ASN
1	B	83	GLN
1	B	87	ASN
1	C	36	ASN
1	C	49	ASN
1	C	87	ASN
1	D	87	ASN
1	E	36	ASN
1	E	49	ASN
1	E	87	ASN
1	F	144	GLN
1	F	151	HIS
1	G	66	GLN
1	G	79	ASN
1	H	87	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no 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, 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	140/152 (92%)	-0.18	0 100 100	30, 33, 39, 44	0
1	B	135/152 (88%)	0.01	2 (1%) 76 71	30, 33, 43, 49	0
1	C	140/152 (92%)	-0.03	2 (1%) 78 74	31, 33, 41, 46	0
1	D	141/152 (92%)	-0.16	0 100 100	27, 33, 39, 46	0
1	E	140/152 (92%)	0.19	6 (4%) 39 31	30, 33, 41, 45	0
1	F	137/152 (90%)	0.06	6 (4%) 38 30	31, 33, 43, 47	0
1	G	128/152 (84%)	0.96	17 (13%) 4 3	31, 33, 37, 39	0
1	H	128/152 (84%)	-0.02	3 (2%) 64 57	31, 33, 37, 45	0
All	All	1089/1216 (89%)	0.10	36 (3%) 50 43	27, 33, 40, 49	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	121	THR	7.1
1	B	22	ILE	4.9
1	H	26	ARG	4.9
1	G	122	SER	4.8
1	G	120	GLU	4.6
1	G	119	ASP	4.5
1	E	24	ARG	4.4
1	G	56	ILE	4.0
1	G	118	PHE	4.0
1	G	44	ILE	3.9
1	G	58	THR	3.6
1	G	106	LEU	3.5
1	E	59	SER	3.3
1	F	27	ALA	3.0
1	E	118	PHE	3.0
1	H	24	ARG	3.0

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Mol	Chain	Res	Type	RSRZ
1	E	119	ASP	2.9
1	F	22	ILE	2.8
1	G	57	THR	2.8
1	G	29	LEU	2.7
1	H	65	ASP	2.7
1	G	50	TYR	2.7
1	E	122	SER	2.6
1	C	150	GLU	2.6
1	F	24	ARG	2.5
1	G	47	ASP	2.5
1	F	23	SER	2.4
1	G	59	SER	2.4
1	F	25	VAL	2.3
1	G	149	ASP	2.3
1	G	150	GLU	2.3
1	E	23	SER	2.3
1	B	117	LEU	2.2
1	G	28	GLU	2.2
1	F	26	ARG	2.2
1	C	118	PHE	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.