Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID	:	2J0O
Title	:	Shigella Flexneri IpaD
Authors	:	Johnson, S.; Roversi, P.; Lea, S.M.
Deposited on	:	2006-08-04
Resolution	:	3.00 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 following versions of software and data (see references ⓘ) were used in the production of this report:

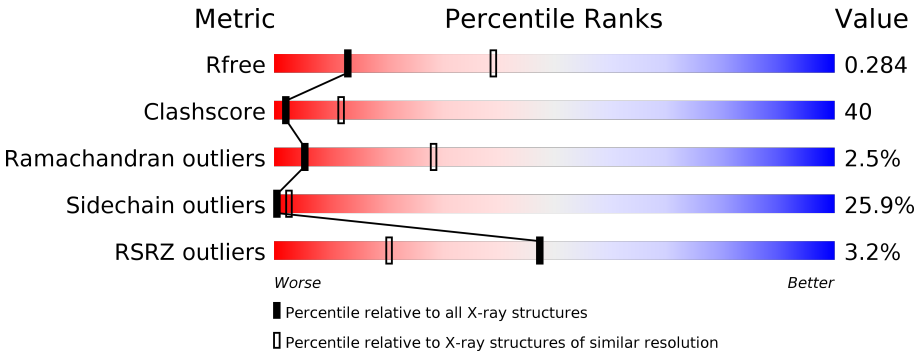
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.00 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}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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 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 <=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	318	3% 35% 39% 15% 11%
1	B	318	3% 36% 40% 14% 11%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4449 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.

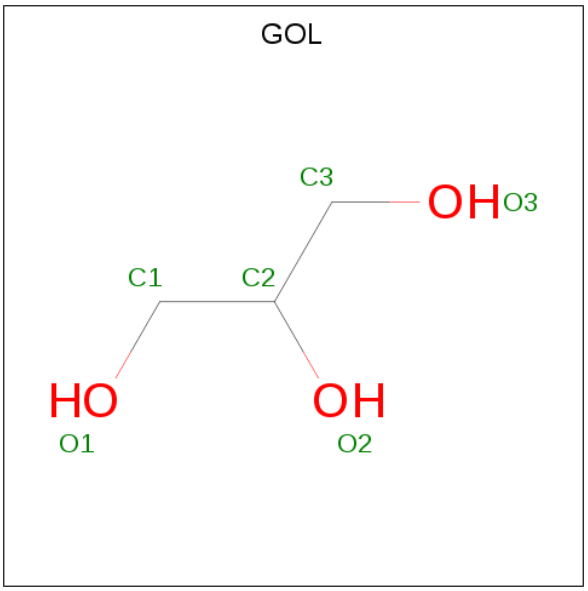
- Molecule 1 is a protein called INVASIN IPAD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	284	Total	C	N	O	S	0	0	0
			2209	1377	377	447	8			
1	B	284	Total	C	N	O	S	0	0	0
			2213	1379	379	447	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	102	HIS	ASN	conflict	UNP P18013
B	102	HIS	ASN	conflict	UNP P18013

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is water.

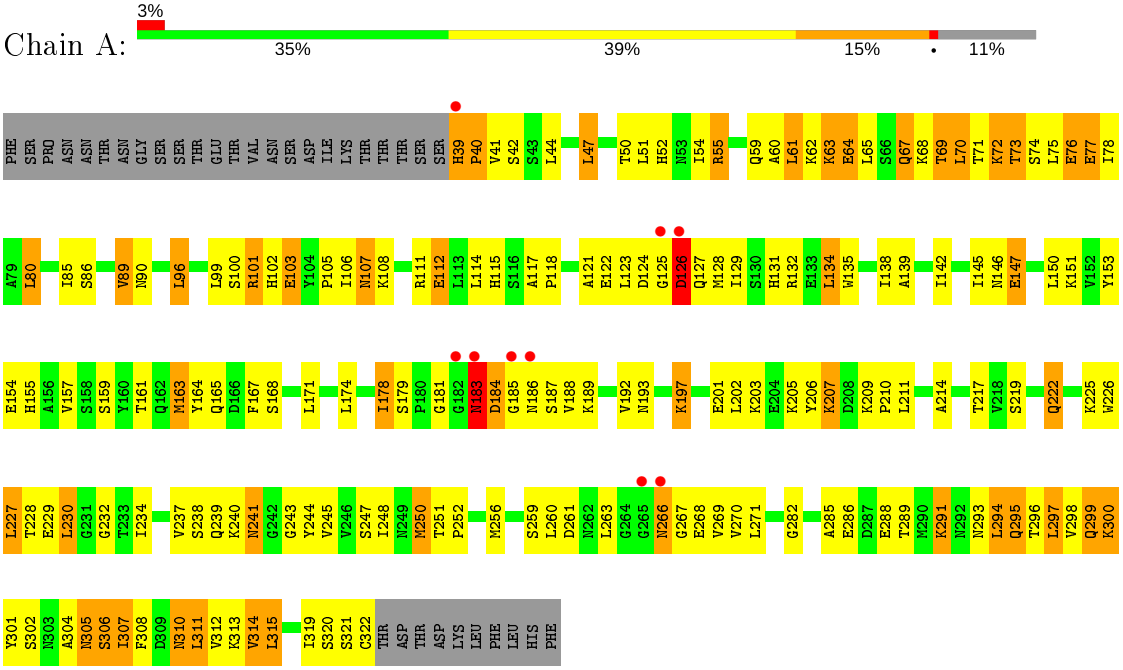
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	17	Total O 17 17	0	0
3	B	4	Total O 4 4	0	0

3 Residue-property plots (i)

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

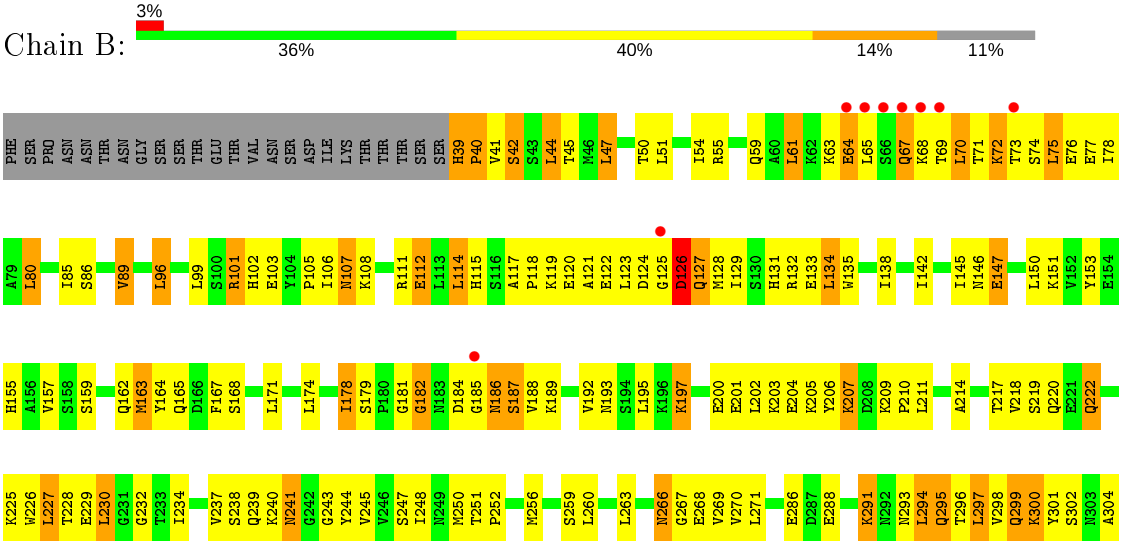
- Molecule 1: INVASIN IPAD

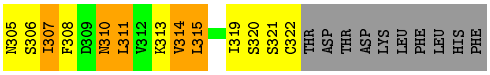
Chain A:



- Molecule 1: INVASIN IPAD

Chain B:




N305
S306
I307
F308
D309
N310
L311
V312
K313
V314
L315
I319
S320
S321
C322
THR
ASP
THR
ASP
LYS
LEU
PHE
LEU
HIS
PHE

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	55.87Å 100.69Å 111.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.00 28.79 – 2.73	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-3.00) 88.8 (28.79-2.73)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.85 (at 2.72Å)	Xtriage
Refinement program	TNT 5.6.1	Depositor
R, R_{free}	0.235 , (Not available) 0.249 , 0.284	Depositor DCC
R_{free} test set	772 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	43.8	Xtriage
Anisotropy	0.440	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 63.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	4449	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 47.91 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 9.2801e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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

Bond lengths and bond angles in the following residue types are not validated in this section: GOL

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	0.22	0/2242	0.39	0/3033
1	B	0.25	0/2247	0.46	0/3040
All	All	0.23	0/4489	0.43	0/6073

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2209	0	2196	175	0
1	B	2213	0	2199	183	0
2	A	6	0	8	2	0
3	A	17	0	0	2	0
3	B	4	0	0	1	0
All	All	4449	0	4403	356	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 40.

All (356) 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:B:182:GLY:HA3	1:B:184:ASP:HA	1.22	1.09
1:B:185:GLY:H	1:B:186:ASN:HA	1.16	1.06
1:B:163:MET:HE2	1:B:248:ILE:HG21	1.36	1.03
1:A:163:MET:HE2	1:A:248:ILE:HG21	1.40	0.98
1:B:147:GLU:HG3	1:B:151:LYS:HE3	1.45	0.98
1:A:147:GLU:HG3	1:A:151:LYS:HE3	1.50	0.93
1:A:227:LEU:HD22	1:A:232:GLY:HA2	1.50	0.93
1:A:39:HIS:HB2	1:A:40:PRO:HD2	1.51	0.93
1:B:185:GLY:N	1:B:186:ASN:HA	1.80	0.92
1:B:145:ILE:HG23	1:B:307:ILE:HD11	1.55	0.89
1:B:227:LEU:HD22	1:B:232:GLY:HA2	1.55	0.88
1:A:145:ILE:HG23	1:A:307:ILE:HD11	1.56	0.87
1:B:127:GLN:HB3	1:B:129:ILE:HG23	1.58	0.86
1:B:39:HIS:HB2	1:B:40:PRO:HD2	1.57	0.85
1:A:96:LEU:HD21	1:A:138:ILE:HG22	1.58	0.85
1:B:147:GLU:HG3	1:B:151:LYS:CE	2.07	0.84
1:B:219:SER:HB3	1:B:222:GLN:HG2	1.60	0.84
1:A:55:ARG:NE	1:A:313:LYS:HE2	1.94	0.82
1:B:96:LEU:HD21	1:B:138:ILE:HG22	1.60	0.82
1:B:182:GLY:HA2	1:B:187:SER:HB2	1.63	0.81
1:B:168:SER:HA	1:B:171:LEU:HD12	1.64	0.80
1:B:142:ILE:HD11	1:B:315:LEU:CD2	2.13	0.79
1:B:105:PRO:HA	1:B:131:HIS:CD2	2.17	0.79
1:A:105:PRO:HA	1:A:131:HIS:CD2	2.18	0.79
1:A:183:ASN:ND2	1:A:185:GLY:N	2.30	0.79
1:A:62:LYS:CE	1:A:305:ASN:HB3	2.12	0.79
1:B:145:ILE:CG2	1:B:307:ILE:HD11	2.12	0.79
1:A:147:GLU:HG3	1:A:151:LYS:CE	2.13	0.78
1:A:85:ILE:O	1:A:89:VAL:HG12	1.84	0.77
1:B:127:GLN:HA	1:B:127:GLN:HE21	1.48	0.77
1:A:145:ILE:CG2	1:A:307:ILE:HD11	2.13	0.77
1:B:85:ILE:O	1:B:89:VAL:HG12	1.85	0.77
1:A:252:PRO:O	1:A:256:MET:HG3	1.85	0.77
1:B:307:ILE:HD13	1:B:308:PHE:N	2.00	0.77
1:B:101:ARG:HG3	1:B:101:ARG:HH21	1.51	0.76
1:A:168:SER:HA	1:A:171:LEU:HD12	1.67	0.76
1:A:307:ILE:HD13	1:A:308:PHE:N	2.01	0.75
1:B:142:ILE:HD11	1:B:315:LEU:HD21	1.67	0.75
1:B:55:ARG:NE	1:B:313:LYS:HE2	2.02	0.74
1:B:108:LYS:HE3	1:B:112:GLU:OE1	1.87	0.74
1:A:76:GLU:O	1:A:80:LEU:HD22	1.86	0.74
1:A:101:ARG:HG3	1:A:101:ARG:HH21	1.52	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:65:LEU:HD21	1:B:301:TYR:CD1	2.24	0.73
1:A:219:SER:HB3	1:A:222:GLN:CG	2.19	0.73
1:B:89:VAL:HG23	1:B:142:ILE:HG22	1.71	0.73
1:A:219:SER:HB3	1:A:222:GLN:HG2	1.72	0.72
1:A:311:LEU:HD22	1:A:315:LEU:HD22	1.71	0.72
1:A:108:LYS:HE3	1:A:112:GLU:OE1	1.89	0.72
1:A:86:SER:HA	1:A:89:VAL:CG1	2.20	0.71
1:B:219:SER:HB3	1:B:222:GLN:CG	2.19	0.71
1:B:182:GLY:HA2	1:B:187:SER:H	1.55	0.71
1:A:55:ARG:CD	1:A:313:LYS:HE2	2.21	0.71
1:A:181:GLY:O	1:A:184:ASP:HB2	1.90	0.70
1:A:64:GLU:OE1	1:A:77:GLU:HB3	1.90	0.70
1:B:182:GLY:CA	1:B:184:ASP:HA	2.12	0.69
1:A:256:MET:CE	1:A:286:GLU:HB2	2.23	0.69
1:B:127:GLN:CA	1:B:127:GLN:HE21	2.06	0.68
1:A:70:LEU:CD1	1:A:294:LEU:HD11	2.24	0.68
1:B:163:MET:HE2	1:B:248:ILE:CG2	2.18	0.68
1:B:72:LYS:HA	1:B:75:LEU:HB2	1.77	0.67
1:B:71:THR:HB	1:B:73:THR:HG22	1.76	0.67
1:A:70:LEU:HD13	1:A:161:THR:OG1	1.96	0.66
1:A:256:MET:HE2	1:A:286:GLU:HB2	1.77	0.66
1:A:142:ILE:HD11	1:A:315:LEU:CD2	2.26	0.66
1:B:96:LEU:HD21	1:B:138:ILE:CG2	2.25	0.66
1:B:256:MET:CE	1:B:286:GLU:HB2	2.26	0.66
1:A:55:ARG:HD3	1:A:313:LYS:HE2	1.77	0.66
1:B:72:LYS:NZ	1:B:162:GLN:HG3	2.11	0.65
1:A:266:ASN:CB	1:A:267:GLY:HA2	2.27	0.65
1:B:147:GLU:HG3	1:B:151:LYS:NZ	2.12	0.65
1:B:252:PRO:O	1:B:256:MET:HG3	1.97	0.64
1:A:250:MET:HA	1:A:250:MET:CE	2.28	0.64
1:B:266:ASN:CB	1:B:267:GLY:HA2	2.26	0.64
1:A:70:LEU:O	1:A:165:GLN:NE2	2.29	0.64
1:A:183:ASN:CG	1:A:184:ASP:HA	2.17	0.64
1:B:250:MET:HA	1:B:250:MET:CE	2.27	0.64
1:B:70:LEU:HB2	1:B:75:LEU:HD13	1.80	0.63
1:A:89:VAL:HG23	1:A:142:ILE:HG22	1.80	0.63
1:A:289:THR:OG1	2:A:1323:GOL:H12	1.98	0.62
1:A:62:LYS:NZ	1:A:306:SER:N	2.48	0.62
1:A:96:LEU:HD21	1:A:138:ILE:CG2	2.30	0.62
1:A:163:MET:HG3	1:A:164:TYR:N	2.15	0.61
1:B:311:LEU:HD22	1:B:315:LEU:HD22	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:127:GLN:HB3	1:B:129:ILE:CG2	2.28	0.61
1:B:182:GLY:HA3	1:B:184:ASP:CA	2.14	0.61
1:B:64:GLU:OE1	1:B:77:GLU:HB3	2.01	0.61
1:A:142:ILE:HD11	1:A:315:LEU:HD21	1.83	0.61
1:A:307:ILE:HD13	1:A:308:PHE:H	1.65	0.60
1:A:217:THR:HB	1:A:243:GLY:HA3	1.83	0.60
1:B:294:LEU:O	1:B:294:LEU:HD12	2.01	0.60
1:A:147:GLU:HG3	1:A:151:LYS:NZ	2.16	0.59
1:A:111:ARG:NH1	1:A:131:HIS:ND1	2.50	0.59
1:B:112:GLU:HA	1:B:125:GLY:O	2.03	0.59
1:A:70:LEU:HB2	1:A:161:THR:HG23	1.84	0.59
1:A:150:LEU:HD11	1:A:308:PHE:CD2	2.38	0.59
1:B:256:MET:HE2	1:B:286:GLU:HB2	1.84	0.59
1:B:146:ASN:HA	1:B:150:LEU:HD12	1.85	0.59
1:B:230:LEU:H	1:B:230:LEU:HD12	1.66	0.59
1:A:225:LYS:O	1:A:228:THR:HG22	2.02	0.59
1:A:62:LYS:NZ	1:A:305:ASN:HB3	2.18	0.58
1:A:62:LYS:HZ1	1:A:306:SER:N	2.01	0.58
1:B:197:LYS:O	1:B:201:GLU:HG3	2.03	0.58
1:B:182:GLY:CA	1:B:187:SER:HB2	2.33	0.58
1:B:192:VAL:HG23	1:B:193:ASN:N	2.18	0.58
1:B:76:GLU:O	1:B:80:LEU:HD22	2.03	0.58
1:A:111:ARG:HH11	1:A:131:HIS:CE1	2.21	0.58
1:A:183:ASN:OD1	1:A:183:ASN:N	2.36	0.58
1:A:183:ASN:ND2	1:A:185:GLY:H	2.01	0.58
1:A:227:LEU:CD2	1:A:232:GLY:HA2	2.30	0.58
1:B:47:LEU:HD22	1:B:51:LEU:HG	1.86	0.58
1:B:310:ASN:O	1:B:314:VAL:HG12	2.04	0.58
1:B:64:GLU:HG3	1:B:74:SER:OG	2.04	0.58
1:A:183:ASN:CB	1:A:184:ASP:HA	2.34	0.58
1:A:192:VAL:HG23	1:A:193:ASN:N	2.19	0.58
1:A:39:HIS:HB2	1:A:40:PRO:CD	2.30	0.58
1:B:86:SER:HA	1:B:89:VAL:CG1	2.34	0.57
1:A:146:ASN:HA	1:A:150:LEU:HD12	1.85	0.57
1:B:150:LEU:HD11	1:B:308:PHE:CD2	2.40	0.57
1:A:153:TYR:O	1:A:157:VAL:HG23	2.05	0.57
1:B:111:ARG:NH1	1:B:131:HIS:ND1	2.52	0.57
1:A:47:LEU:HD22	1:A:51:LEU:HG	1.86	0.57
1:B:178:ILE:HG22	1:B:188:VAL:CG1	2.35	0.57
1:B:142:ILE:HD11	1:B:315:LEU:HD23	1.86	0.56
1:A:250:MET:HA	1:A:250:MET:HE3	1.85	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:PRO:HA	1:A:131:HIS:CG	2.41	0.56
1:B:315:LEU:O	1:B:319:ILE:HG12	2.05	0.56
1:B:189:LYS:HE3	1:B:268:GLU:OE2	2.05	0.56
1:A:184:ASP:OD2	1:A:185:GLY:N	2.39	0.56
1:B:210:PRO:HA	1:B:247:SER:HA	1.87	0.56
1:B:72:LYS:HG3	1:B:72:LYS:O	2.05	0.56
1:B:105:PRO:HA	1:B:131:HIS:CG	2.40	0.56
1:A:106:ILE:HD12	1:A:134:LEU:CD1	2.35	0.56
1:B:111:ARG:HH11	1:B:131:HIS:CE1	2.23	0.56
1:A:294:LEU:O	1:A:294:LEU:HD12	2.05	0.56
1:A:295:GLN:HB3	1:B:299:GLN:OE1	2.06	0.56
1:B:39:HIS:HB2	1:B:40:PRO:CD	2.30	0.56
1:B:96:LEU:HD11	1:B:135:TRP:HE3	1.71	0.55
1:B:71:THR:O	1:B:72:LYS:HB3	2.06	0.55
1:A:163:MET:HE2	1:A:248:ILE:CG2	2.25	0.55
1:B:230:LEU:HA	1:B:300:LYS:HE2	1.89	0.55
1:A:189:LYS:HE2	1:A:268:GLU:OE2	2.07	0.55
1:A:210:PRO:HA	1:A:247:SER:HA	1.88	0.55
1:A:234:ILE:HG12	1:A:293:ASN:CG	2.27	0.55
1:B:225:LYS:O	1:B:228:THR:HG22	2.06	0.55
1:B:307:ILE:HD13	1:B:308:PHE:H	1.69	0.55
1:B:70:LEU:O	1:B:165:GLN:NE2	2.39	0.55
1:A:310:ASN:O	1:A:314:VAL:HG12	2.07	0.54
1:B:72:LYS:HZ1	1:B:162:GLN:HG3	1.72	0.54
1:B:310:ASN:HB2	3:B:2004:HOH:O	2.06	0.54
1:A:40:PRO:HG3	1:A:102:HIS:CG	2.42	0.54
1:A:108:LYS:HE3	1:A:112:GLU:CD	2.28	0.54
1:A:185:GLY:O	1:A:186:ASN:HB2	2.08	0.54
1:B:163:MET:HA	1:B:206:TYR:CZ	2.42	0.54
1:A:239:GLN:HB2	1:A:244:TYR:CE2	2.43	0.54
1:A:61:LEU:HD21	1:A:78:ILE:HG12	1.88	0.54
1:B:108:LYS:HE3	1:B:112:GLU:CD	2.27	0.54
1:B:163:MET:HG3	1:B:164:TYR:N	2.23	0.54
1:B:239:GLN:HB2	1:B:244:TYR:CE2	2.43	0.54
1:B:250:MET:HA	1:B:250:MET:HE3	1.89	0.54
1:A:55:ARG:HG3	1:A:312:VAL:HG23	1.89	0.53
1:B:96:LEU:CD2	1:B:138:ILE:HG22	2.37	0.53
1:B:61:LEU:HD21	1:B:78:ILE:HG12	1.90	0.53
1:A:126:ASP:O	1:A:127:GLN:HB3	2.07	0.53
1:B:115:HIS:CG	1:B:125:GLY:HA3	2.43	0.53
1:A:205:LYS:HG2	1:A:206:TYR:CE2	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163:MET:CE	1:A:248:ILE:HG21	2.27	0.53
1:A:296:THR:O	1:A:299:GLN:HB2	2.09	0.53
1:B:106:ILE:HD12	1:B:134:LEU:CD1	2.39	0.53
1:A:106:ILE:HD12	1:A:134:LEU:HD13	1.91	0.53
1:A:230:LEU:HD12	1:A:230:LEU:H	1.72	0.52
1:B:296:THR:O	1:B:299:GLN:HB2	2.09	0.52
1:B:61:LEU:HD22	1:B:61:LEU:O	2.08	0.52
1:A:127:GLN:O	1:A:127:GLN:HG3	2.10	0.52
1:A:319:ILE:O	1:A:322:CYS:HB2	2.10	0.52
1:A:65:LEU:C	1:A:67:GLN:H	2.14	0.51
1:B:121:ALA:O	1:B:122:GLU:HB2	2.10	0.51
1:A:70:LEU:HB2	1:A:161:THR:CG2	2.40	0.51
1:A:163:MET:HA	1:A:206:TYR:CZ	2.44	0.51
1:B:256:MET:HE3	1:B:286:GLU:HB2	1.91	0.51
1:A:299:GLN:OE1	1:B:295:GLN:HB3	2.10	0.51
1:B:294:LEU:O	1:B:298:VAL:HG23	2.11	0.51
3:A:2016:HOH:O	1:B:313:LYS:HD2	2.10	0.51
1:A:62:LYS:HE2	1:A:305:ASN:HB3	1.92	0.50
1:A:320:SER:C	1:A:322:CYS:H	2.14	0.50
1:A:69:THR:O	1:A:71:THR:N	2.43	0.50
1:A:86:SER:O	1:A:90:ASN:OD1	2.29	0.50
1:A:96:LEU:HD11	1:A:135:TRP:HE3	1.76	0.50
1:A:71:THR:HG22	1:A:72:LYS:N	2.26	0.50
1:A:86:SER:O	1:A:89:VAL:HG13	2.11	0.50
1:B:65:LEU:C	1:B:67:GLN:H	2.15	0.50
1:B:157:VAL:HG22	1:B:297:LEU:HB3	1.93	0.50
1:B:117:ALA:HA	1:B:314:VAL:HG11	1.92	0.50
2:A:1323:GOL:O3	1:B:120:GLU:OE2	2.30	0.50
1:A:304:ALA:O	1:A:307:ILE:HD13	2.11	0.50
1:A:197:LYS:O	1:A:201:GLU:HG3	2.12	0.50
1:A:260:LEU:O	1:A:263:LEU:HB2	2.12	0.50
1:A:71:THR:HG22	1:A:73:THR:H	1.76	0.49
1:A:261:ASP:HB2	3:A:2015:HOH:O	2.13	0.49
1:B:234:ILE:HG12	1:B:293:ASN:CG	2.32	0.49
1:A:256:MET:HE3	1:A:286:GLU:HB2	1.93	0.49
1:B:50:THR:O	1:B:54:ILE:HG13	2.13	0.49
1:A:142:ILE:HD11	1:A:315:LEU:HD23	1.94	0.49
1:B:71:THR:HG22	1:B:73:THR:HB	1.94	0.49
1:B:96:LEU:CD1	1:B:135:TRP:HE3	2.26	0.49
1:B:55:ARG:HE	1:B:313:LYS:HE2	1.76	0.49
1:B:70:LEU:HB2	1:B:75:LEU:CD1	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:LEU:HA	1:A:300:LYS:HE2	1.94	0.49
1:A:52:HIS:CE1	1:A:55:ARG:HH12	2.30	0.49
1:A:178:ILE:HG22	1:A:188:VAL:CG1	2.43	0.49
1:B:163:MET:CE	1:B:250:MET:HE1	2.43	0.49
1:B:320:SER:C	1:B:322:CYS:H	2.16	0.49
1:A:96:LEU:CD1	1:A:135:TRP:HE3	2.26	0.48
1:B:115:HIS:CD2	1:B:125:GLY:H	2.32	0.48
1:A:108:LYS:HE3	1:A:112:GLU:OE2	2.13	0.48
1:A:174:LEU:O	1:A:178:ILE:HG12	2.14	0.48
1:B:192:VAL:CG1	1:B:269:VAL:HG12	2.43	0.48
1:B:230:LEU:HA	1:B:300:LYS:CE	2.44	0.48
1:B:227:LEU:CD2	1:B:232:GLY:HA2	2.36	0.48
1:A:230:LEU:HD13	1:A:230:LEU:C	2.33	0.48
1:A:52:HIS:ND1	1:A:55:ARG:NH1	2.57	0.48
1:B:114:LEU:HB2	1:B:125:GLY:HA2	1.95	0.48
1:B:304:ALA:O	1:B:307:ILE:HD13	2.13	0.48
1:B:319:ILE:O	1:B:322:CYS:HB2	2.13	0.48
1:A:286:GLU:HA	1:A:286:GLU:OE2	2.14	0.48
1:A:86:SER:HA	1:A:89:VAL:HG12	1.93	0.48
1:B:108:LYS:HE3	1:B:112:GLU:OE2	2.13	0.48
1:B:71:THR:CG2	1:B:73:THR:HB	2.44	0.48
1:A:203:LYS:O	1:A:207:LYS:N	2.47	0.48
1:B:230:LEU:C	1:B:230:LEU:HD13	2.33	0.48
1:A:157:VAL:HG21	1:A:301:TYR:HB2	1.96	0.47
1:A:39:HIS:O	1:A:41:VAL:N	2.46	0.47
1:B:153:TYR:O	1:B:157:VAL:HG23	2.14	0.47
1:B:101:ARG:NH2	1:B:101:ARG:HG3	2.25	0.47
1:B:106:ILE:HD12	1:B:134:LEU:HD13	1.96	0.47
1:A:163:MET:CG	1:A:164:TYR:N	2.78	0.47
1:A:50:THR:O	1:A:54:ILE:HG13	2.14	0.47
1:B:127:GLN:HA	1:B:127:GLN:NE2	2.22	0.47
1:B:211:LEU:HD21	1:B:248:ILE:HG12	1.95	0.47
1:B:251:THR:N	1:B:252:PRO:CD	2.78	0.47
1:B:150:LEU:HD11	1:B:308:PHE:CG	2.50	0.47
1:A:121:ALA:O	1:A:122:GLU:HB2	2.14	0.47
1:B:142:ILE:CD1	1:B:315:LEU:HD21	2.41	0.47
1:A:39:HIS:CD2	1:A:41:VAL:HB	2.50	0.47
1:B:260:LEU:O	1:B:263:LEU:HB2	2.15	0.47
1:B:204:GLU:OE1	1:B:207:LYS:HE3	2.15	0.46
1:B:286:GLU:OE2	1:B:286:GLU:HA	2.14	0.46
1:A:101:ARG:CG	1:A:101:ARG:HH21	2.22	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:39:HIS:CD2	1:B:41:VAL:HB	2.50	0.46
1:B:96:LEU:HD11	1:B:135:TRP:CE3	2.48	0.46
1:A:251:THR:N	1:A:252:PRO:CD	2.78	0.46
1:B:315:LEU:HD12	1:B:315:LEU:HA	1.78	0.46
1:A:211:LEU:HD21	1:A:248:ILE:HG12	1.97	0.46
1:A:288:GLU:HA	1:A:291:LYS:HB2	1.96	0.46
1:B:39:HIS:CB	1:B:40:PRO:HD2	2.36	0.46
1:A:184:ASP:OD2	1:A:187:SER:N	2.47	0.46
1:B:203:LYS:O	1:B:207:LYS:N	2.49	0.46
1:B:89:VAL:HG23	1:B:142:ILE:CG2	2.44	0.46
1:A:230:LEU:HA	1:A:300:LYS:CE	2.46	0.46
1:A:150:LEU:HD11	1:A:308:PHE:CG	2.51	0.46
1:B:125:GLY:O	1:B:126:ASP:HB2	2.16	0.46
1:B:185:GLY:N	1:B:186:ASN:CA	2.67	0.46
1:A:73:THR:CG2	1:A:74:SER:N	2.79	0.45
1:B:96:LEU:CD1	1:B:135:TRP:CE3	2.99	0.45
1:B:197:LYS:HA	1:B:197:LYS:HD3	1.46	0.45
1:B:174:LEU:O	1:B:178:ILE:HG12	2.17	0.45
1:B:71:THR:HG22	1:B:73:THR:N	2.32	0.45
1:A:115:HIS:CG	1:A:125:GLY:HA3	2.52	0.45
1:B:40:PRO:HG3	1:B:102:HIS:CG	2.51	0.45
1:B:266:ASN:HB2	1:B:267:GLY:HA2	1.98	0.45
1:B:42:SER:HA	1:B:45:THR:HG22	1.98	0.45
1:A:157:VAL:HG22	1:A:297:LEU:HB3	1.99	0.45
1:A:39:HIS:NE2	1:A:41:VAL:CG2	2.80	0.45
1:B:220:GLN:HB2	1:B:244:TYR:CD1	2.52	0.45
1:B:71:THR:CG2	1:B:73:THR:HG22	2.46	0.45
1:B:86:SER:O	1:B:89:VAL:HG13	2.17	0.45
1:A:55:ARG:HE	1:A:313:LYS:HE2	1.76	0.45
1:A:102:HIS:C	1:A:103:GLU:HG2	2.36	0.45
1:A:70:LEU:HB3	1:A:161:THR:HG21	1.98	0.45
1:B:207:LYS:CB	1:B:207:LYS:NZ	2.80	0.45
1:A:155:HIS:CE1	1:A:214:ALA:HA	2.52	0.44
1:A:163:MET:CE	1:A:250:MET:HE1	2.47	0.44
1:A:250:MET:CA	1:A:250:MET:HE3	2.48	0.44
1:A:266:ASN:HB2	1:A:267:GLY:HA2	1.98	0.44
1:B:72:LYS:HZ3	1:B:162:GLN:HG3	1.81	0.44
1:A:192:VAL:CG2	1:A:193:ASN:N	2.80	0.44
1:A:118:PRO:HG3	1:A:310:ASN:OD1	2.17	0.44
1:A:117:ALA:HA	1:A:314:VAL:HG11	1.98	0.44
1:A:207:LYS:CB	1:A:207:LYS:NZ	2.80	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:LYS:O	1:A:241:ASN:HB2	2.16	0.44
1:A:315:LEU:O	1:A:319:ILE:HG12	2.17	0.44
1:B:118:PRO:HG3	1:B:310:ASN:OD1	2.18	0.44
1:B:288:GLU:HA	1:B:291:LYS:HB2	1.99	0.44
1:A:60:ALA:HA	1:A:63:LYS:NZ	2.33	0.44
1:A:77:GLU:OE1	1:A:77:GLU:HA	2.17	0.44
1:B:192:VAL:CG2	1:B:193:ASN:N	2.80	0.44
1:B:115:HIS:ND1	1:B:125:GLY:HA3	2.34	0.43
1:B:207:LYS:NZ	1:B:207:LYS:HB3	2.33	0.43
1:B:226:TRP:HA	1:B:229:GLU:OE2	2.18	0.43
1:B:163:MET:CG	1:B:164:TYR:N	2.79	0.43
1:A:100:SER:OG	1:A:132:ARG:O	2.33	0.43
1:A:61:LEU:HD23	1:A:61:LEU:HA	1.85	0.43
1:A:207:LYS:NZ	1:A:207:LYS:HB3	2.34	0.43
1:A:226:TRP:HA	1:A:229:GLU:OE2	2.18	0.43
1:A:219:SER:O	1:A:222:GLN:HG3	2.19	0.43
1:A:96:LEU:HD23	1:A:139:ALA:HA	2.01	0.43
1:B:155:HIS:CE1	1:B:214:ALA:HA	2.54	0.43
1:B:240:LYS:O	1:B:241:ASN:HB2	2.18	0.43
1:B:220:GLN:HB2	1:B:244:TYR:CE1	2.54	0.43
1:A:154:GLU:HB2	1:A:301:TYR:CZ	2.54	0.43
1:A:96:LEU:CD1	1:A:135:TRP:CE3	3.01	0.43
1:B:167:PHE:HB2	1:B:202:LEU:HD21	2.00	0.43
1:A:294:LEU:O	1:A:298:VAL:HG23	2.19	0.43
1:B:217:THR:HB	1:B:243:GLY:HA3	2.00	0.42
1:B:294:LEU:HD12	1:B:294:LEU:C	2.40	0.42
1:B:217:THR:O	1:B:218:VAL:HG13	2.19	0.42
1:A:55:ARG:HG3	1:A:312:VAL:CG2	2.49	0.42
1:B:127:GLN:CA	1:B:127:GLN:NE2	2.78	0.42
1:B:70:LEU:H	1:B:70:LEU:HG	1.44	0.42
1:A:101:ARG:NH2	1:A:101:ARG:CG	2.80	0.42
1:B:65:LEU:CD2	1:B:301:TYR:CD1	2.99	0.42
1:A:183:ASN:CB	1:A:184:ASP:CA	2.97	0.42
1:B:44:LEU:HD23	1:B:44:LEU:HA	1.90	0.42
1:B:39:HIS:C	1:B:41:VAL:H	2.22	0.42
1:A:167:PHE:HB2	1:A:202:LEU:HD21	2.00	0.42
1:B:51:LEU:HD23	1:B:51:LEU:HA	1.90	0.42
1:A:101:ARG:NH2	1:A:101:ARG:HG3	2.25	0.42
1:B:114:LEU:HA	1:B:114:LEU:HD12	1.84	0.42
1:B:115:HIS:CE1	1:B:125:GLY:HA3	2.54	0.42
1:B:147:GLU:CG	1:B:151:LYS:NZ	2.81	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:197:LYS:HD3	1:B:200:GLU:OE1	2.20	0.42
1:A:47:LEU:HD22	1:A:47:LEU:O	2.20	0.41
1:B:153:TYR:CE2	1:B:300:LYS:HB3	2.55	0.41
1:A:96:LEU:CD2	1:A:138:ILE:HG22	2.40	0.41
1:A:282:GLY:O	1:A:285:ALA:HB3	2.20	0.41
1:A:96:LEU:HA	1:A:96:LEU:HD12	1.82	0.41
1:B:157:VAL:HG21	1:B:301:TYR:HB2	2.01	0.41
1:A:205:LYS:HE2	1:A:206:TYR:OH	2.20	0.41
1:B:71:THR:CB	1:B:73:THR:HG22	2.46	0.41
1:B:39:HIS:NE2	1:B:41:VAL:CG2	2.84	0.41
1:A:70:LEU:CB	1:A:161:THR:HG21	2.51	0.41
1:A:96:LEU:HD11	1:A:135:TRP:CE3	2.54	0.41
1:A:192:VAL:CG1	1:A:269:VAL:HG12	2.51	0.41
1:B:101:ARG:NH2	1:B:101:ARG:CG	2.80	0.41
1:B:134:LEU:HA	1:B:134:LEU:HD23	1.78	0.41
1:B:71:THR:CG2	1:B:73:THR:CG2	2.99	0.41
1:A:127:GLN:HG3	1:A:129:ILE:CG2	2.51	0.40
1:A:62:LYS:HZ1	1:A:305:ASN:C	2.24	0.40
1:A:80:LEU:N	1:A:80:LEU:HD13	2.36	0.40
1:B:205:LYS:HG2	1:B:206:TYR:CE2	2.56	0.40
1:B:193:ASN:HB2	1:B:267:GLY:HA3	2.01	0.40
1:A:197:LYS:HA	1:A:197:LYS:HD3	1.45	0.40
1:B:132:ARG:HG2	1:B:133:GLU:N	2.36	0.40
1:B:195:LEU:HD23	1:B:260:LEU:HD21	2.03	0.40
1:B:71:THR:HG22	1:B:73:THR:H	1.87	0.40
1:A:147:GLU:CG	1:A:151:LYS:NZ	2.84	0.40
1:A:71:THR:CG2	1:A:72:LYS:N	2.84	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	282/318 (89%)	257 (91%)	18 (6%)	7 (2%)	5 28
1	B	282/318 (89%)	253 (90%)	22 (8%)	7 (2%)	5 28
All	All	564/636 (89%)	510 (90%)	40 (7%)	14 (2%)	5 28

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	126	ASP
1	B	126	ASP
1	B	241	ASN
1	A	70	LEU
1	A	107	ASN
1	A	183	ASN
1	A	241	ASN
1	A	321	SER
1	B	107	ASN
1	B	182	GLY
1	B	321	SER
1	A	40	PRO
1	B	40	PRO
1	B	181	GLY

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	250/284 (88%)	184 (74%)	66 (26%)	0 2
1	B	251/284 (88%)	187 (74%)	64 (26%)	0 3
All	All	501/568 (88%)	371 (74%)	130 (26%)	0 2

All (130) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	HIS
1	A	42	SER

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Mol	Chain	Res	Type
1	A	44	LEU
1	A	47	LEU
1	A	55	ARG
1	A	59	GLN
1	A	61	LEU
1	A	63	LYS
1	A	64	GLU
1	A	67	GLN
1	A	68	LYS
1	A	69	THR
1	A	72	LYS
1	A	73	THR
1	A	75	LEU
1	A	76	GLU
1	A	77	GLU
1	A	80	LEU
1	A	89	VAL
1	A	96	LEU
1	A	99	LEU
1	A	101	ARG
1	A	103	GLU
1	A	107	ASN
1	A	112	GLU
1	A	114	LEU
1	A	123	LEU
1	A	124	ASP
1	A	126	ASP
1	A	128	MET
1	A	134	LEU
1	A	147	GLU
1	A	159	SER
1	A	163	MET
1	A	178	ILE
1	A	179	SER
1	A	183	ASN
1	A	184	ASP
1	A	197	LYS
1	A	207	LYS
1	A	209	LYS
1	A	222	GLN
1	A	227	LEU
1	A	230	LEU

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Mol	Chain	Res	Type
1	A	237	VAL
1	A	238	SER
1	A	245	VAL
1	A	250	MET
1	A	259	SER
1	A	266	ASN
1	A	270	VAL
1	A	271	LEU
1	A	291	LYS
1	A	294	LEU
1	A	295	GLN
1	A	297	LEU
1	A	299	GLN
1	A	300	LYS
1	A	302	SER
1	A	305	ASN
1	A	306	SER
1	A	307	ILE
1	A	310	ASN
1	A	311	LEU
1	A	314	VAL
1	A	315	LEU
1	B	39	HIS
1	B	42	SER
1	B	44	LEU
1	B	47	LEU
1	B	59	GLN
1	B	61	LEU
1	B	63	LYS
1	B	64	GLU
1	B	67	GLN
1	B	68	LYS
1	B	69	THR
1	B	70	LEU
1	B	72	LYS
1	B	75	LEU
1	B	80	LEU
1	B	89	VAL
1	B	96	LEU
1	B	99	LEU
1	B	101	ARG
1	B	103	GLU

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Mol	Chain	Res	Type
1	B	107	ASN
1	B	112	GLU
1	B	114	LEU
1	B	119	LYS
1	B	123	LEU
1	B	124	ASP
1	B	126	ASP
1	B	127	GLN
1	B	128	MET
1	B	134	LEU
1	B	147	GLU
1	B	159	SER
1	B	163	MET
1	B	178	ILE
1	B	179	SER
1	B	186	ASN
1	B	187	SER
1	B	197	LYS
1	B	207	LYS
1	B	209	LYS
1	B	222	GLN
1	B	227	LEU
1	B	230	LEU
1	B	237	VAL
1	B	238	SER
1	B	245	VAL
1	B	259	SER
1	B	266	ASN
1	B	270	VAL
1	B	271	LEU
1	B	291	LYS
1	B	294	LEU
1	B	295	GLN
1	B	297	LEU
1	B	299	GLN
1	B	300	LYS
1	B	302	SER
1	B	305	ASN
1	B	306	SER
1	B	307	ILE
1	B	310	ASN
1	B	311	LEU

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Mol	Chain	Res	Type
1	B	314	VAL
1	B	315	LEU

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

Mol	Chain	Res	Type
1	B	115	HIS
1	B	127	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	GOL	A	1323	-	5,5,5	0.32	0	5,5,5	0.26	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	A	1323	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1323	GOL	2	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	284/318 (89%)	-0.27	9 (3%) 47 20	12, 30, 59, 76	0
1	B	284/318 (89%)	-0.10	9 (3%) 47 20	17, 35, 61, 78	0
All	All	568/636 (89%)	-0.19	18 (3%) 47 20	12, 31, 60, 78	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	69	THR	5.2
1	B	67	GLN	4.9
1	B	125	GLY	4.8
1	B	64	GLU	3.9
1	A	266	ASN	3.2
1	A	183	ASN	3.1
1	A	186	ASN	3.1
1	A	182	GLY	2.9
1	B	185	GLY	2.7
1	A	39	HIS	2.7
1	B	66	SER	2.5
1	B	65	LEU	2.5
1	A	125	GLY	2.4
1	B	68	LYS	2.2
1	A	185	GLY	2.2
1	A	265	GLY	2.1
1	B	73	THR	2.1
1	A	126	ASP	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 carbohydrates in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	GOL	A	1323	6/6	0.82	0.27	34,36,37,37	0

6.5 Other polymers (i)

There are no such residues in this entry.