



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:21 AM GMT

PDB ID : 2WRB
Title : THE STRUCTURE OF INFLUENZA H2 HUMAN SINGAPORE HEMAG-
GLUTININ WITH AVIAN RECEPTOR
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Deposited on : 2009-09-01
Resolution : 3.10 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

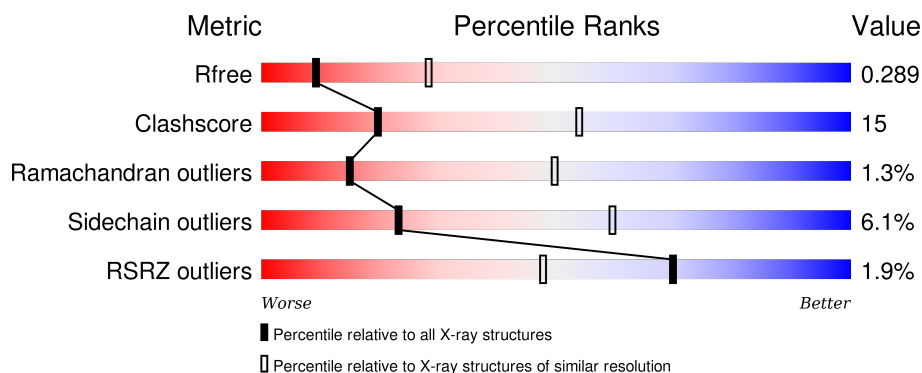
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1114 (3.14-3.06)
Clashscore	102246	1222 (3.14-3.06)
Ramachandran outliers	100387	1174 (3.14-3.06)
Sidechain outliers	100360	1174 (3.14-3.06)
RSRZ outliers	91569	1119 (3.14-3.06)

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	506	<div> <div>2%</div> <div>66%</div> <div>25%</div> <div>• 6%</div> </div>
1	B	506	<div> <div>%</div> <div>66%</div> <div>26%</div> <div>• • •</div> </div>
1	C	506	<div> <div>2%</div> <div>61%</div> <div>23%</div> <div>• 13%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SIA	A	1488	-	-	-	X

2 Entry composition [i](#)

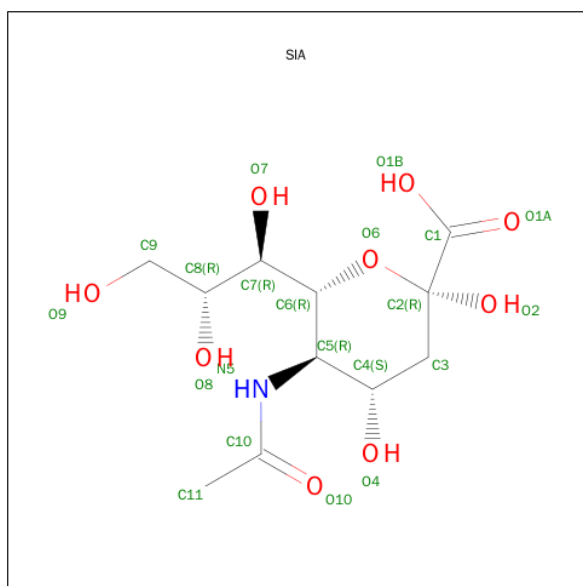
There are 2 unique types of molecules in this entry. The entry contains 11080 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 HEMAGGLUTININ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	474	Total	C	N	O	S	0	0	0
			3736	2347	642	724	23			
1	B	486	Total	C	N	O	S	0	0	0
			3842	2413	659	746	24			
1	C	441	Total	C	N	O	S	0	0	0
			3481	2195	599	667	20			

- Molecule 2 is SUGAR (O-SIALIC ACID) (three-letter code: SIA) (formula: $C_{11}H_{19}NO_9$).

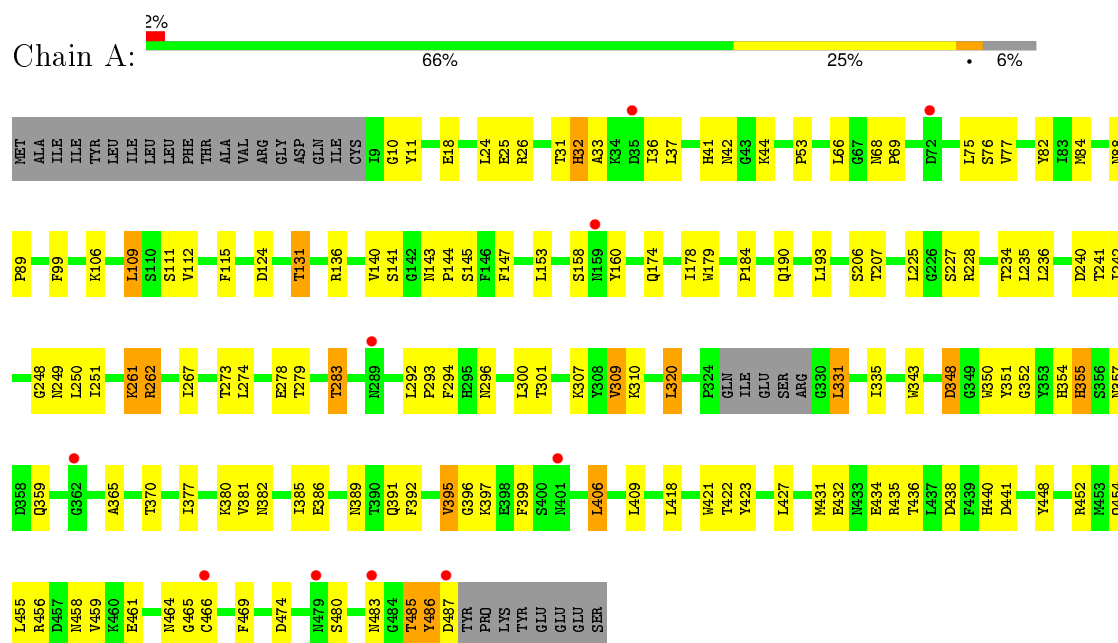


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			21	11	1	9		

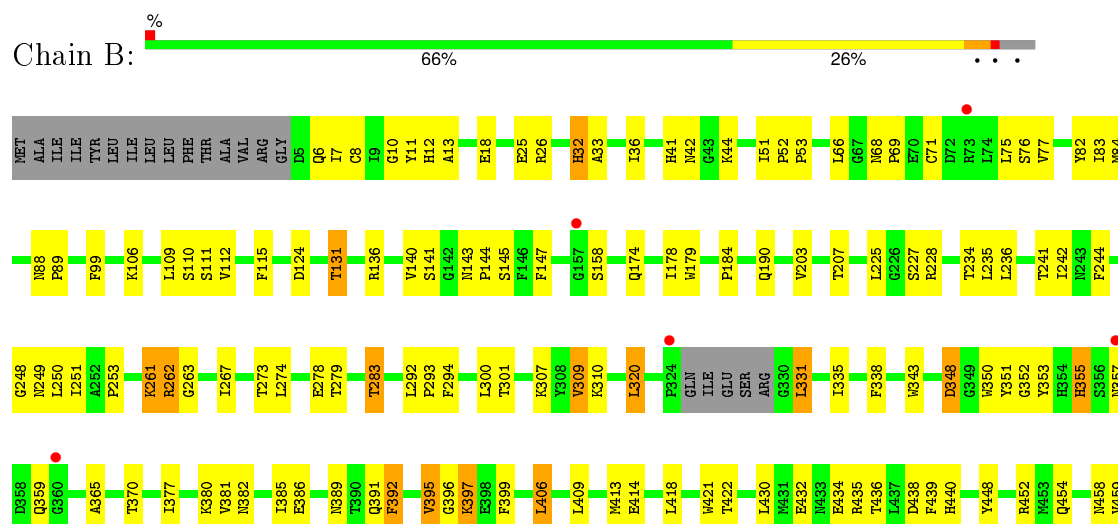
3 Residue-property plots

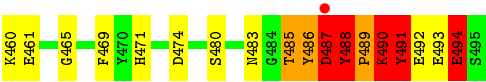
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: HEMAGGLUTININ

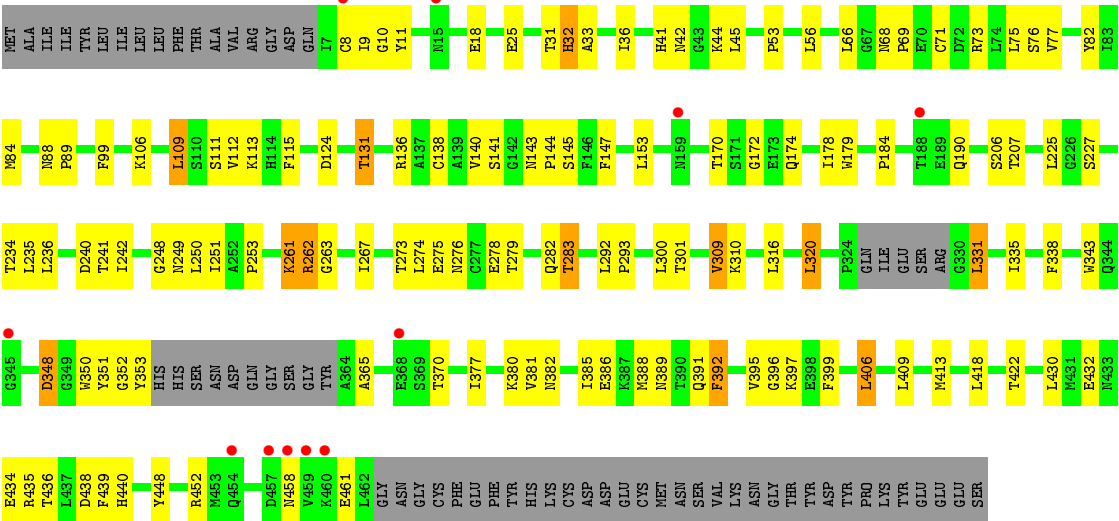


• Molecule 1: HEMAGGLUTININ





● Molecule 1: HEMAGGLUTININ



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.36Å 126.30Å 220.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 3.10 8.00 – 3.10	Depositor EDS
% Data completeness (in resolution range)	98.9 (8.00-3.10) 98.9 (8.00-3.10)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.96 (at 3.08Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.253 , 0.296 0.246 , 0.289	Depositor DCC
R_{free} test set	2007 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	68.5	Xtriage
Anisotropy	0.638	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 59.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Outliers	1 of 40027 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	11080	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.66% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.38	0/3820	0.54	0/5169
1	B	0.47	1/3929 (0.0%)	0.61	4/5316 (0.1%)
1	C	0.43	0/3556	0.56	0/4812
All	All	0.43	1/11305 (0.0%)	0.57	4/15297 (0.0%)

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	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	414	GLU	CG-CD	5.02	1.59	1.51

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	488	TYR	N-CA-C	-6.85	92.50	111.00
1	B	491	TYR	N-CA-C	6.29	127.98	111.00
1	B	488	TYR	CA-CB-CG	-5.40	103.15	113.40
1	B	488	TYR	N-CA-CB	5.14	119.86	110.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	488	TYR	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3736	0	3600	110	0
1	B	3842	0	3694	132	0
1	C	3481	0	3406	95	0
2	A	21	0	17	7	0
All	All	11080	0	10717	325	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 15.

All (325) 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:493:GLU:HA	1:B:494:GLU:HB2	1.38	1.01
1:A:193:LEU:HD21	2:A:1488:SIA:C11	1.92	0.97
1:C:261:LYS:HA	1:C:262:ARG:CB	1.99	0.93
1:C:112:VAL:HG11	1:C:115:PHE:HB2	1.50	0.92
1:B:493:GLU:HA	1:B:494:GLU:CB	1.99	0.92
1:B:261:LYS:HA	1:B:262:ARG:CB	2.00	0.92
1:A:261:LYS:HA	1:A:262:ARG:CB	2.00	0.92
1:A:193:LEU:HD21	2:A:1488:SIA:H111	1.51	0.91
1:A:112:VAL:HG11	1:A:115:PHE:HB2	1.53	0.91
1:A:193:LEU:CD2	2:A:1488:SIA:H111	2.00	0.90
1:B:490:LYS:CA	1:B:491:TYR:HB2	2.02	0.90
1:B:112:VAL:HG11	1:B:115:PHE:HB2	1.54	0.87
2:A:1488:SIA:H113	2:A:1488:SIA:O7	1.73	0.86
1:A:320:LEU:H	1:A:320:LEU:HD23	1.42	0.85
1:B:459:VAL:HG12	1:B:469:PHE:HA	1.57	0.84
1:C:320:LEU:HD23	1:C:320:LEU:H	1.43	0.83
1:A:454:GLN:NE2	1:A:486:TYR:HB3	1.93	0.82
1:B:490:LYS:H	1:B:491:TYR:HB2	1.41	0.82
1:B:320:LEU:HD23	1:B:320:LEU:H	1.44	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:459:VAL:HG12	1:A:469:PHE:HA	1.62	0.81
1:B:493:GLU:CA	1:B:494:GLU:HB2	2.10	0.81
1:B:454:GLN:NE2	1:B:486:TYR:HB3	1.96	0.81
1:B:490:LYS:N	1:B:491:TYR:HB2	1.95	0.80
1:B:488:TYR:HB3	1:B:489:PRO:HD2	1.64	0.80
1:B:261:LYS:HA	1:B:262:ARG:HB3	1.65	0.77
1:C:261:LYS:HA	1:C:262:ARG:HB3	1.65	0.77
1:A:193:LEU:HD21	2:A:1488:SIA:H113	1.68	0.76
1:C:261:LYS:HA	1:C:262:ARG:HB2	1.67	0.74
1:C:348:ASP:HB2	1:C:365:ALA:HB3	1.70	0.73
1:A:348:ASP:HB2	1:A:365:ALA:HB3	1.72	0.72
1:B:348:ASP:HB2	1:B:365:ALA:HB3	1.71	0.72
1:A:184:PRO:HG2	1:A:190:GLN:HE21	1.55	0.71
1:B:261:LYS:HA	1:B:262:ARG:HB2	1.71	0.71
1:B:487:ASP:HA	1:B:488:TYR:HB2	1.73	0.70
1:C:184:PRO:HG2	1:C:190:GLN:HE21	1.56	0.70
1:C:261:LYS:CA	1:C:262:ARG:HB3	2.21	0.69
1:B:261:LYS:CA	1:B:262:ARG:HB3	2.22	0.69
1:B:487:ASP:HA	1:B:488:TYR:CB	2.23	0.68
1:B:261:LYS:CA	1:B:262:ARG:CB	2.72	0.68
1:C:261:LYS:CA	1:C:262:ARG:CB	2.71	0.68
1:B:488:TYR:CD2	1:B:488:TYR:C	2.68	0.67
1:B:184:PRO:HG2	1:B:190:GLN:HE21	1.59	0.66
1:B:492:GLU:HA	1:B:493:GLU:CB	2.24	0.66
1:A:193:LEU:HD11	2:A:1488:SIA:H92	1.76	0.66
1:A:261:LYS:CA	1:A:262:ARG:CB	2.73	0.66
1:B:490:LYS:HA	1:B:491:TYR:HB2	1.76	0.65
1:C:293:PRO:HG3	1:C:385:ILE:HA	1.79	0.65
1:B:488:TYR:HB3	1:B:489:PRO:CD	2.27	0.65
1:A:36:ILE:HA	1:A:292:LEU:HD22	1.78	0.65
1:B:389:ASN:HD21	1:C:310:LYS:HD2	1.63	0.64
1:C:18:GLU:HG2	1:C:33:ALA:HB3	1.81	0.63
1:B:53:PRO:HG2	1:B:84:MET:HE2	1.80	0.63
1:C:53:PRO:HG2	1:C:84:MET:HE2	1.81	0.63
1:B:487:ASP:OD2	1:B:489:PRO:HD2	1.99	0.63
1:C:140:VAL:HG12	1:C:145:SER:HB2	1.81	0.63
1:B:236:LEU:HD22	1:B:242:ILE:HB	1.81	0.63
1:A:331:LEU:HD22	1:A:438:ASP:OD2	2.00	0.62
1:A:18:GLU:HG2	1:A:33:ALA:HB3	1.81	0.62
1:C:66:LEU:O	1:C:147:PHE:HB3	2.00	0.62
1:A:140:VAL:HG12	1:A:145:SER:HB2	1.80	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:36:ILE:HA	1:C:292:LEU:HD22	1.81	0.61
1:C:309:VAL:HG22	1:C:422:THR:HA	1.83	0.61
1:B:36:ILE:HA	1:B:292:LEU:HD22	1.81	0.61
1:B:18:GLU:HG2	1:B:33:ALA:HB3	1.82	0.60
1:B:140:VAL:HG12	1:B:145:SER:HB2	1.82	0.60
1:C:382:ASN:O	1:C:386:GLU:HG2	2.01	0.60
1:C:8:CYS:O	1:C:353:TYR:HA	2.02	0.60
1:A:236:LEU:HD22	1:A:242:ILE:HB	1.83	0.60
1:A:309:VAL:HG22	1:A:422:THR:HA	1.84	0.60
1:A:106:LYS:HB3	1:A:267:ILE:HD11	1.84	0.60
1:A:293:PRO:HG3	1:A:385:ILE:HA	1.84	0.60
1:B:309:VAL:HG22	1:B:422:THR:HA	1.82	0.60
1:A:66:LEU:O	1:A:147:PHE:HB3	2.01	0.60
1:B:392:PHE:CD2	1:B:392:PHE:C	2.75	0.59
1:A:53:PRO:HG2	1:A:84:MET:HE2	1.84	0.59
1:B:106:LYS:HB3	1:B:267:ILE:HD11	1.83	0.59
1:B:293:PRO:HG3	1:B:385:ILE:HA	1.84	0.59
1:B:42:ASN:O	1:B:44:LYS:HG3	2.03	0.59
1:A:99:PHE:HZ	1:A:178:ILE:HD13	1.67	0.59
1:A:184:PRO:HG2	1:A:190:GLN:NE2	2.18	0.58
1:A:310:LYS:HE3	1:A:418:LEU:HD21	1.84	0.58
1:B:389:ASN:ND2	1:C:310:LYS:HD2	2.18	0.58
1:B:99:PHE:HZ	1:B:178:ILE:HD13	1.68	0.58
1:A:179:TRP:HZ3	1:A:234:THR:HG22	1.67	0.58
1:C:184:PRO:HG2	1:C:190:GLN:NE2	2.18	0.58
1:B:452:ARG:HD3	1:B:461:GLU:OE2	2.04	0.58
1:B:320:LEU:N	1:B:320:LEU:HD23	2.17	0.58
1:C:331:LEU:HD22	1:C:438:ASP:OD2	2.04	0.58
1:B:331:LEU:HD22	1:B:438:ASP:OD2	2.04	0.58
1:C:310:LYS:HE3	1:C:418:LEU:HD21	1.86	0.57
1:C:348:ASP:N	1:C:348:ASP:OD1	2.38	0.57
1:A:348:ASP:N	1:A:348:ASP:OD1	2.37	0.57
1:C:392:PHE:C	1:C:392:PHE:CD2	2.77	0.57
1:B:493:GLU:CA	1:B:494:GLU:CB	2.75	0.57
1:A:320:LEU:HD23	1:A:320:LEU:N	2.15	0.56
1:C:99:PHE:HZ	1:C:178:ILE:HD13	1.69	0.56
1:B:66:LEU:O	1:B:147:PHE:HB3	2.06	0.56
1:B:8:CYS:O	1:B:353:TYR:HA	2.06	0.56
1:A:359:GLN:NE2	1:A:474:ASP:HB2	2.21	0.56
1:B:348:ASP:OD1	1:B:348:ASP:N	2.38	0.55
1:C:236:LEU:HD22	1:C:242:ILE:HB	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:488:TYR:HD2	1:B:489:PRO:HD3	1.71	0.55
1:C:42:ASN:O	1:C:44:LYS:HG3	2.07	0.55
1:A:454:GLN:HE21	1:A:486:TYR:HB3	1.69	0.55
1:B:471:HIS:CE1	1:B:490:LYS:HE3	2.42	0.55
1:A:283:THR:HG22	1:A:301:THR:HG22	1.88	0.55
1:C:320:LEU:HD23	1:C:320:LEU:N	2.17	0.54
1:A:42:ASN:O	1:A:44:LYS:HG3	2.07	0.54
1:C:250:LEU:HD12	1:C:251:ILE:N	2.21	0.54
1:B:490:LYS:HD2	1:B:490:LYS:O	2.06	0.54
1:B:184:PRO:HG2	1:B:190:GLN:NE2	2.22	0.54
1:A:348:ASP:HB2	1:A:365:ALA:CB	2.37	0.54
1:A:320:LEU:HB3	1:A:440:HIS:CG	2.42	0.54
1:B:179:TRP:HZ3	1:B:234:THR:HG22	1.72	0.54
1:C:348:ASP:HB2	1:C:365:ALA:CB	2.37	0.54
1:A:359:GLN:HE22	1:A:474:ASP:HB2	1.73	0.54
1:B:382:ASN:O	1:B:386:GLU:HG2	2.07	0.54
1:B:283:THR:HG22	1:B:301:THR:HG22	1.89	0.54
1:A:448:TYR:CE1	1:A:465:GLY:HA2	2.43	0.53
1:B:310:LYS:HE3	1:B:418:LEU:HD21	1.89	0.53
1:B:454:GLN:HE21	1:B:486:TYR:HB3	1.71	0.53
1:A:452:ARG:HD3	1:A:461:GLU:OE2	2.08	0.53
1:A:53:PRO:HG3	1:A:82:TYR:CE2	2.44	0.53
1:B:399:PHE:CE1	1:B:406:LEU:HG	2.43	0.53
1:C:283:THR:HG22	1:C:301:THR:HG22	1.91	0.53
1:A:480:SER:HB2	1:A:486:TYR:H	1.73	0.53
1:B:348:ASP:HB2	1:B:365:ALA:CB	2.37	0.53
1:B:448:TYR:CE1	1:B:465:GLY:HA2	2.43	0.53
1:C:53:PRO:HG3	1:C:82:TYR:CE2	2.44	0.52
1:B:32:HIS:CE1	1:B:350:TRP:HE1	2.28	0.52
1:B:248:GLY:O	1:B:249:ASN:HB2	2.10	0.52
1:A:435:ARG:HH21	1:B:434:GLU:CD	2.13	0.52
1:A:309:VAL:CG2	1:A:422:THR:HA	2.40	0.52
1:C:32:HIS:CE1	1:C:350:TRP:HE1	2.28	0.52
1:B:261:LYS:CB	1:B:262:ARG:HB3	2.40	0.52
1:C:309:VAL:CG2	1:C:422:THR:HA	2.39	0.52
1:A:392:PHE:CD2	1:A:392:PHE:C	2.83	0.52
1:A:193:LEU:CD2	2:A:1488:SIA:C11	2.68	0.51
1:C:88:ASN:N	1:C:89:PRO:HD3	2.25	0.51
1:C:261:LYS:CB	1:C:262:ARG:HB3	2.40	0.51
1:B:262:ARG:HG2	1:B:263:GLY:N	2.25	0.51
1:C:320:LEU:HB3	1:C:440:HIS:CG	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:480:SER:HA	1:B:483:ASN:HB2	1.91	0.51
1:B:480:SER:HB2	1:B:486:TYR:H	1.75	0.51
1:B:68:ASN:OD1	1:B:69:PRO:HD2	2.11	0.51
1:C:377:ILE:HD11	1:C:436:THR:HG23	1.91	0.51
1:C:452:ARG:HD3	1:C:461:GLU:OE2	2.10	0.51
1:B:174:GLN:OE1	1:B:235:LEU:HD13	2.11	0.51
1:B:377:ILE:HD11	1:B:436:THR:HG23	1.93	0.51
1:B:11:TYR:CZ	1:B:335:ILE:HG23	2.46	0.51
1:C:53:PRO:HG2	1:C:84:MET:CE	2.41	0.50
1:B:359:GLN:NE2	1:B:474:ASP:HB2	2.25	0.50
1:B:53:PRO:HG2	1:B:84:MET:CE	2.41	0.50
1:B:309:VAL:CG2	1:B:422:THR:HA	2.41	0.50
1:A:480:SER:HA	1:A:483:ASN:HB2	1.92	0.50
1:C:106:LYS:HB3	1:C:267:ILE:HD11	1.94	0.50
1:C:399:PHE:CE1	1:C:406:LEU:HG	2.47	0.50
1:A:32:HIS:CE1	1:A:350:TRP:HE1	2.29	0.50
1:C:179:TRP:HZ3	1:C:234:THR:HG22	1.76	0.50
1:B:53:PRO:HG3	1:B:82:TYR:CE2	2.47	0.50
1:B:435:ARG:HH21	1:C:434:GLU:CD	2.15	0.50
1:A:131:THR:HG23	1:A:131:THR:O	2.12	0.50
1:C:248:GLY:O	1:C:249:ASN:HB2	2.12	0.50
1:B:320:LEU:HB3	1:B:440:HIS:CG	2.47	0.50
1:C:53:PRO:HD2	1:C:274:LEU:HD22	1.94	0.50
1:B:350:TRP:H	1:B:370:THR:CG2	2.25	0.50
1:B:454:GLN:HE22	1:B:486:TYR:HB3	1.77	0.49
1:A:427:LEU:O	1:A:431:MET:HG3	2.12	0.49
1:C:248:GLY:C	1:C:249:ASN:HD22	2.16	0.49
1:A:480:SER:HB2	1:A:485:THR:HB	1.94	0.49
1:C:241:THR:HG22	1:C:242:ILE:N	2.28	0.49
1:C:350:TRP:H	1:C:370:THR:CG2	2.25	0.49
1:A:381:VAL:HG12	1:A:385:ILE:HD12	1.94	0.49
1:A:357:ASN:HD21	1:A:474:ASP:HA	1.78	0.49
1:A:250:LEU:HD12	1:A:251:ILE:N	2.28	0.48
1:B:10:GLY:HA3	1:B:343:TRP:CH2	2.48	0.48
1:B:359:GLN:HE22	1:B:474:ASP:HB2	1.77	0.48
1:C:262:ARG:HG3	1:C:263:GLY:N	2.29	0.48
1:B:485:THR:O	1:B:486:TYR:O	2.30	0.48
1:C:458:ASN:N	1:C:458:ASN:ND2	2.61	0.48
1:C:300:LEU:HD22	1:C:396:GLY:CA	2.44	0.48
1:C:206:SER:OG	1:C:240:ASP:OD2	2.32	0.48
1:B:241:THR:HG22	1:B:242:ILE:N	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:480:SER:HB2	1:B:485:THR:HB	1.96	0.48
1:A:352:GLY:HA3	1:A:365:ALA:HA	1.96	0.47
1:C:300:LEU:HD22	1:C:396:GLY:HA2	1.96	0.47
1:A:350:TRP:H	1:A:370:THR:CG2	2.27	0.47
1:B:293:PRO:HB2	1:B:294:PHE:CD1	2.50	0.47
1:C:174:GLN:OE1	1:C:235:LEU:HD13	2.13	0.47
1:A:399:PHE:CE1	1:A:406:LEU:HG	2.50	0.47
1:A:174:GLN:OE1	1:A:235:LEU:HD13	2.15	0.47
1:A:75:LEU:O	1:A:76:SER:HB2	2.15	0.47
1:C:225:LEU:HD23	1:C:225:LEU:HA	1.71	0.47
1:B:250:LEU:HD12	1:B:251:ILE:N	2.29	0.47
1:A:227:SER:C	1:A:228:ARG:HG2	2.34	0.47
1:B:140:VAL:HG12	1:B:145:SER:CB	2.44	0.47
1:B:178:ILE:O	1:B:253:PRO:HG3	2.15	0.47
1:A:382:ASN:O	1:A:386:GLU:HG2	2.15	0.47
1:A:68:ASN:OD1	1:A:69:PRO:HD2	2.15	0.46
1:A:458:ASN:ND2	1:A:458:ASN:N	2.63	0.46
1:A:423:TYR:CD1	1:C:388:MET:HA	2.50	0.46
1:A:41:HIS:CD2	1:A:41:HIS:N	2.84	0.46
1:A:248:GLY:C	1:A:249:ASN:HD22	2.18	0.46
1:C:261:LYS:HB2	1:C:262:ARG:HB3	1.98	0.46
1:A:307:LYS:HG2	1:A:421:TRP:CD2	2.49	0.46
1:B:53:PRO:HD2	1:B:274:LEU:HD22	1.98	0.46
1:A:140:VAL:HG12	1:A:145:SER:CB	2.44	0.46
1:C:10:GLY:HA3	1:C:343:TRP:CH2	2.51	0.46
1:A:136:ARG:HG3	1:A:144:PRO:HG3	1.98	0.46
1:C:75:LEU:O	1:C:76:SER:HB2	2.15	0.46
1:B:338:PHE:CD1	1:B:338:PHE:C	2.89	0.46
1:B:488:TYR:CB	1:B:489:PRO:CD	2.94	0.46
1:A:99:PHE:CZ	1:A:178:ILE:HD13	2.51	0.46
1:C:278:GLU:HG3	1:C:279:THR:N	2.31	0.46
1:A:53:PRO:HG2	1:A:84:MET:CE	2.46	0.46
1:B:261:LYS:HB2	1:B:262:ARG:HB3	1.98	0.45
1:C:11:TYR:CZ	1:C:335:ILE:HG23	2.50	0.45
1:B:136:ARG:HG3	1:B:144:PRO:HG3	1.98	0.45
1:B:203:VAL:HG22	1:B:244:PHE:CD2	2.51	0.45
1:A:307:LYS:HG2	1:A:421:TRP:CE2	2.51	0.45
1:A:248:GLY:O	1:A:249:ASN:HB2	2.16	0.45
1:B:83:ILE:HD12	1:B:110:SER:HA	1.99	0.45
1:B:41:HIS:N	1:B:41:HIS:CD2	2.84	0.45
1:A:241:THR:HG22	1:A:242:ILE:N	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:PRO:HD2	1:A:274:LEU:HD22	1.99	0.45
1:B:381:VAL:HG12	1:B:385:ILE:HD12	1.99	0.45
1:A:278:GLU:HG3	1:A:279:THR:N	2.32	0.45
1:C:99:PHE:CZ	1:C:178:ILE:HD13	2.52	0.45
1:C:136:ARG:HG3	1:C:144:PRO:HG3	1.99	0.45
1:B:131:THR:HG23	1:B:131:THR:O	2.16	0.45
1:A:377:ILE:HD11	1:A:436:THR:HG23	1.99	0.45
1:C:352:GLY:HA3	1:C:365:ALA:HA	1.98	0.44
1:C:381:VAL:HG12	1:C:385:ILE:HD12	1.99	0.44
1:B:380:LYS:HE2	1:B:432:GLU:OE2	2.16	0.44
1:A:300:LEU:HD22	1:A:396:GLY:CA	2.47	0.44
1:C:68:ASN:OD1	1:C:69:PRO:HD2	2.17	0.44
1:B:352:GLY:HA3	1:B:365:ALA:HA	1.98	0.44
1:A:293:PRO:HB2	1:A:294:PHE:CD1	2.52	0.44
1:A:225:LEU:HD23	1:A:225:LEU:HA	1.73	0.44
1:C:140:VAL:HG12	1:C:145:SER:CB	2.46	0.44
1:A:53:PRO:HG3	1:A:82:TYR:CZ	2.52	0.44
1:C:41:HIS:CD2	1:C:41:HIS:N	2.85	0.44
1:C:380:LYS:HE2	1:C:432:GLU:OE2	2.17	0.44
1:C:141:SER:C	1:C:143:ASN:H	2.21	0.44
1:A:10:GLY:HA3	1:A:343:TRP:CH2	2.52	0.44
1:B:51:ILE:HA	1:B:52:PRO:HD3	1.87	0.44
1:B:6:GLN:O	1:B:355:HIS:HA	2.18	0.44
1:C:338:PHE:C	1:C:338:PHE:CD1	2.91	0.44
1:A:454:GLN:HE22	1:A:486:TYR:HB3	1.74	0.44
1:A:389:ASN:ND2	1:B:310:LYS:HD2	2.33	0.44
1:C:9:ILE:HG12	1:C:448:TYR:HA	1.99	0.43
1:A:206:SER:OG	1:A:240:ASP:OD2	2.35	0.43
1:A:141:SER:C	1:A:143:ASN:H	2.21	0.43
1:A:389:ASN:HD21	1:B:310:LYS:HD2	1.83	0.43
1:B:490:LYS:HA	1:B:491:TYR:CB	2.47	0.43
1:A:300:LEU:HD22	1:A:396:GLY:HA2	2.00	0.43
1:B:75:LEU:O	1:B:76:SER:HB2	2.18	0.43
1:C:113:LYS:HB2	1:C:262:ARG:HD3	1.99	0.43
1:C:53:PRO:HG3	1:C:82:TYR:CZ	2.53	0.43
1:B:262:ARG:CG	1:B:263:GLY:N	2.81	0.43
1:A:310:LYS:HD2	1:C:389:ASN:HD21	1.83	0.43
1:A:448:TYR:HE1	1:A:465:GLY:HA2	1.84	0.43
1:A:88:ASN:N	1:A:89:PRO:HD3	2.34	0.43
1:C:73:ARG:HB2	1:C:73:ARG:NH1	2.34	0.43
1:B:141:SER:C	1:B:143:ASN:H	2.21	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:434:GLU:CD	1:C:435:ARG:HH21	2.22	0.43
1:C:131:THR:HG23	1:C:131:THR:O	2.17	0.43
1:B:99:PHE:CZ	1:B:178:ILE:HD13	2.51	0.43
1:C:31:THR:HG1	1:C:32:HIS:CE1	2.35	0.43
1:C:436:THR:O	1:C:439:PHE:HB3	2.19	0.43
1:B:300:LEU:HD22	1:B:396:GLY:CA	2.48	0.43
1:B:278:GLU:HG3	1:B:279:THR:N	2.34	0.43
1:B:458:ASN:ND2	1:B:458:ASN:N	2.65	0.42
1:A:406:LEU:HA	1:A:406:LEU:HD12	1.78	0.42
1:A:380:LYS:HE2	1:A:432:GLU:OE2	2.18	0.42
1:B:307:LYS:HG2	1:B:421:TRP:CD2	2.54	0.42
1:B:395:VAL:O	1:B:397:LYS:HE2	2.20	0.42
1:A:31:THR:HG23	1:A:320:LEU:O	2.19	0.42
1:A:485:THR:O	1:A:486:TYR:O	2.37	0.42
1:B:448:TYR:HE1	1:B:465:GLY:HA2	1.85	0.42
1:A:31:THR:HG1	1:A:32:HIS:CE1	2.37	0.42
1:A:160:TYR:CZ	1:A:248:GLY:HA2	2.55	0.42
1:A:395:VAL:O	1:A:397:LYS:HE2	2.20	0.42
1:B:489:PRO:O	1:B:490:LYS:CG	2.67	0.42
1:B:68:ASN:HB3	1:B:71:CYS:SG	2.59	0.42
1:C:68:ASN:HB3	1:C:71:CYS:SG	2.60	0.42
1:B:413:MET:SD	1:C:413:MET:HB2	2.60	0.42
1:B:469:PHE:HB3	1:B:471:HIS:O	2.20	0.41
1:B:377:ILE:HD13	1:B:377:ILE:HA	1.85	0.41
1:B:489:PRO:O	1:B:490:LYS:CB	2.68	0.41
1:C:377:ILE:HA	1:C:377:ILE:HD13	1.85	0.41
1:B:225:LEU:HD23	1:B:225:LEU:HA	1.66	0.41
1:A:335:ILE:HG13	1:A:441:ASP:HA	2.02	0.41
1:A:354:HIS:O	1:A:355:HIS:ND1	2.53	0.41
1:B:494:GLU:HA	1:B:494:GLU:OE1	2.20	0.41
1:B:430:LEU:HD23	1:B:430:LEU:HA	1.89	0.41
1:B:471:HIS:HB2	1:B:492:GLU:OE1	2.20	0.41
1:A:37:LEU:HD21	1:A:296:ASN:ND2	2.34	0.41
1:C:45:LEU:HA	1:C:282:GLN:NE2	2.36	0.41
1:B:227:SER:C	1:B:228:ARG:HG2	2.40	0.41
1:B:488:TYR:HD2	1:B:489:PRO:CD	2.33	0.41
1:B:307:LYS:HG2	1:B:421:TRP:CE2	2.55	0.41
1:A:24:LEU:HD21	1:C:380:LYS:HE3	2.03	0.41
1:C:170:THR:C	1:C:172:GLY:H	2.24	0.41
1:A:109:LEU:HA	1:A:109:LEU:HD12	1.87	0.41
1:B:88:ASN:N	1:B:89:PRO:HD3	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:436:THR:O	1:B:439:PHE:HB3	2.21	0.41
1:A:423:TYR:HD1	1:C:388:MET:HA	1.85	0.41
1:B:300:LEU:HD22	1:B:396:GLY:HA2	2.03	0.41
1:A:11:TYR:CZ	1:A:335:ILE:HG23	2.56	0.41
1:C:430:LEU:HD23	1:C:430:LEU:HA	1.94	0.41
1:C:56:LEU:HD23	1:C:56:LEU:HA	1.82	0.41
1:A:455:LEU:HD13	1:A:459:VAL:HG11	2.03	0.41
1:C:178:ILE:O	1:C:253:PRO:HG3	2.22	0.41
1:B:12:HIS:ND1	1:B:13:ALA:N	2.69	0.41
1:A:32:HIS:HE1	1:A:350:TRP:HE1	1.68	0.40
1:C:138:CYS:O	1:C:145:SER:HB3	2.20	0.40
1:A:456:ARG:HH22	1:B:460:LYS:HE2	1.85	0.40
1:B:357:ASN:HD21	1:B:474:ASP:HA	1.87	0.40
1:A:464:ASN:OD1	1:A:466:CYS:HB2	2.21	0.40
1:A:227:SER:O	1:A:228:ARG:HG2	2.22	0.40
1:A:41:HIS:CD2	1:A:41:HIS:H	2.39	0.40
1:B:52:PRO:HA	1:B:53:PRO:HD3	1.89	0.40
1:C:275:GLU:O	1:C:276:ASN:HB3	2.21	0.40
1:C:109:LEU:HD12	1:C:109:LEU:HA	1.87	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	470/506 (93%)	422 (90%)	43 (9%)	5 (1%)	17 55
1	B	482/506 (95%)	429 (89%)	42 (9%)	11 (2%)	8 35
1	C	435/506 (86%)	396 (91%)	37 (8%)	2 (0%)	34 72
All	All	1387/1518 (91%)	1247 (90%)	122 (9%)	18 (1%)	15 50

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	262	ARG
1	A	486	TYR
1	B	262	ARG
1	B	486	TYR
1	B	488	TYR
1	B	490	LYS
1	B	491	TYR
1	C	262	ARG
1	A	485	THR
1	B	485	THR
1	B	494	GLU
1	A	355	HIS
1	A	395	VAL
1	B	355	HIS
1	C	395	VAL
1	B	395	VAL
1	B	487	ASP
1	B	489	PRO

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	412/443 (93%)	389 (94%)	23 (6%)	26	62
1	B	424/443 (96%)	397 (94%)	27 (6%)	22	57
1	C	385/443 (87%)	361 (94%)	24 (6%)	23	59
All	All	1221/1329 (92%)	1147 (94%)	74 (6%)	23	59

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	GLU
1	A	26	ARG
1	A	32	HIS
1	A	77	VAL
1	A	109	LEU

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Mol	Chain	Res	Type
1	A	111	SER
1	A	124	ASP
1	A	131	THR
1	A	153	LEU
1	A	158	SER
1	A	207	THR
1	A	261	LYS
1	A	273	THR
1	A	283	THR
1	A	309	VAL
1	A	320	LEU
1	A	331	LEU
1	A	348	ASP
1	A	351	TYR
1	A	391	GLN
1	A	406	LEU
1	A	409	LEU
1	A	487	ASP
1	B	7	ILE
1	B	25	GLU
1	B	26	ARG
1	B	32	HIS
1	B	77	VAL
1	B	109	LEU
1	B	111	SER
1	B	124	ASP
1	B	131	THR
1	B	158	SER
1	B	207	THR
1	B	261	LYS
1	B	273	THR
1	B	283	THR
1	B	309	VAL
1	B	320	LEU
1	B	331	LEU
1	B	348	ASP
1	B	351	TYR
1	B	391	GLN
1	B	392	PHE
1	B	397	LYS
1	B	406	LEU
1	B	409	LEU

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Mol	Chain	Res	Type
1	B	487	ASP
1	B	490	LYS
1	B	494	GLU
1	C	25	GLU
1	C	32	HIS
1	C	77	VAL
1	C	109	LEU
1	C	111	SER
1	C	124	ASP
1	C	131	THR
1	C	153	LEU
1	C	207	THR
1	C	227	SER
1	C	261	LYS
1	C	273	THR
1	C	283	THR
1	C	309	VAL
1	C	316	LEU
1	C	320	LEU
1	C	331	LEU
1	C	348	ASP
1	C	351	TYR
1	C	391	GLN
1	C	392	PHE
1	C	397	LYS
1	C	406	LEU
1	C	409	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	129	HIS
1	A	190	GLN
1	A	249	ASN
1	A	359	GLN
1	A	391	GLN
1	A	454	GLN
1	A	458	ASN
1	B	129	HIS
1	B	190	GLN
1	B	195	GLN
1	B	249	ASN

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Mol	Chain	Res	Type
1	B	359	GLN
1	B	391	GLN
1	B	454	GLN
1	B	458	ASN
1	C	128	GLN
1	C	129	HIS
1	C	190	GLN
1	C	195	GLN
1	C	249	ASN
1	C	391	GLN
1	C	424	ASN
1	C	458	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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	SIA	A	1488	1	17,21,21	0.82	0	19,31,31	0.96	2 (10%)

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	SIA	A	1488	1	-	0/14/38/38	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	A	1488	SIA	C7-C6-C5	-2.06	111.20	114.32
2	A	1488	SIA	O6-C6-C7	2.51	111.06	107.26

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1488	SIA	7	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, 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	474/506 (93%)	0.00	10 (2%) 67 44	39, 93, 143, 173	0
1	B	486/506 (96%)	-0.29	6 (1%) 81 64	29, 63, 104, 183	0
1	C	441/506 (87%)	-0.21	11 (2%) 61 37	28, 68, 128, 181	0
All	All	1401/1518 (92%)	-0.17	27 (1%) 70 48	28, 73, 130, 183	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	487	ASP	6.1
1	A	72	ASP	5.1
1	C	457	ASP	4.6
1	A	483	ASN	4.2
1	B	487	ASP	4.0
1	B	357	ASN	3.9
1	C	8	CYS	3.6
1	C	188	THR	3.5
1	C	458	ASN	3.2
1	B	157	GLY	3.2
1	C	454	GLN	3.1
1	B	324	PRO	2.9
1	A	289	ASN	2.9
1	C	15	ASN	2.8
1	C	345	GLY	2.8
1	B	360	GLY	2.7
1	A	466	CYS	2.7
1	C	460	LYS	2.7
1	C	368	GLU	2.6
1	A	159	ASN	2.6
1	C	459	VAL	2.5
1	A	362	GLY	2.5
1	A	401	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	159	ASN	2.2
1	A	479	ASN	2.1
1	A	35	ASP	2.1
1	B	73	ARG	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th 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	LLDF	B-factors(\AA^2)	Q<0.9
2	SIA	A	1488	21/21	0.75	0.34	2.67	98,104,110,112	0

6.5 Other polymers [i](#)

There are no such residues in this entry.