



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:11 AM GMT

PDB ID : 2G1Y
Title : Ketopiperazine-Based Renin Inhibitors: Optimization of the "C" Ring
Authors : Holsworth, D.D.; Jalaiea, M.; Zhanga, E.; Mcconnella, P.
Deposited on : 2006-02-15
Resolution : 2.50 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

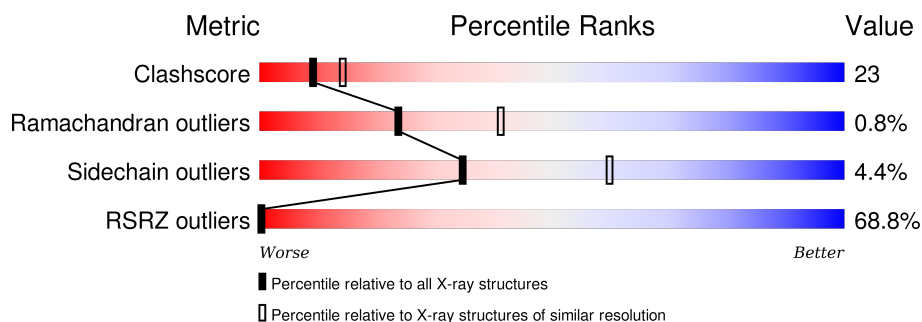
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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	333	
1	B	333	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5550 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 Renin.

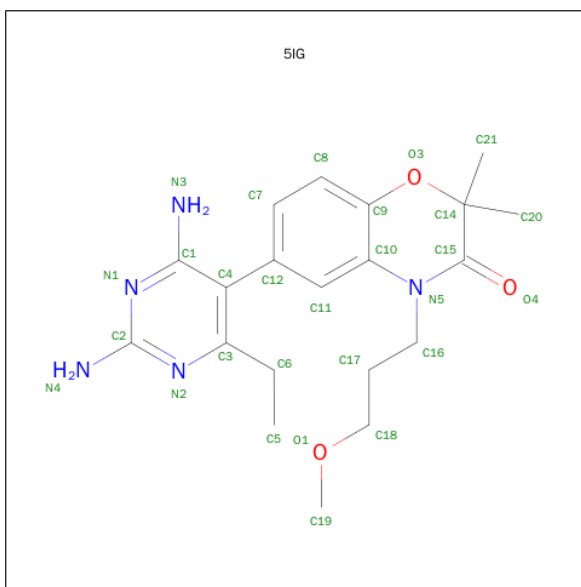
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	333	Total	C	N	O	S	0	0	0
			2572	1642	416	500	14			
1	B	333	Total	C	N	O	S	0	0	0
			2572	1642	416	500	14			

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is 6-(2,4-DIAMINO-6-ETHYLPYRIMIDIN-5-YL)-4-(3-METHOXYPROPYL)-2,2-DIMETHYL-2H-1,4-BENZOXAZIN-3(4H)-ONE (three-letter code: 5IG) (formula: $C_{20}H_{27}N_5O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			28	20	5	3		
3	B	1	Total	C	N	O	0	0
			28	20	5	3		

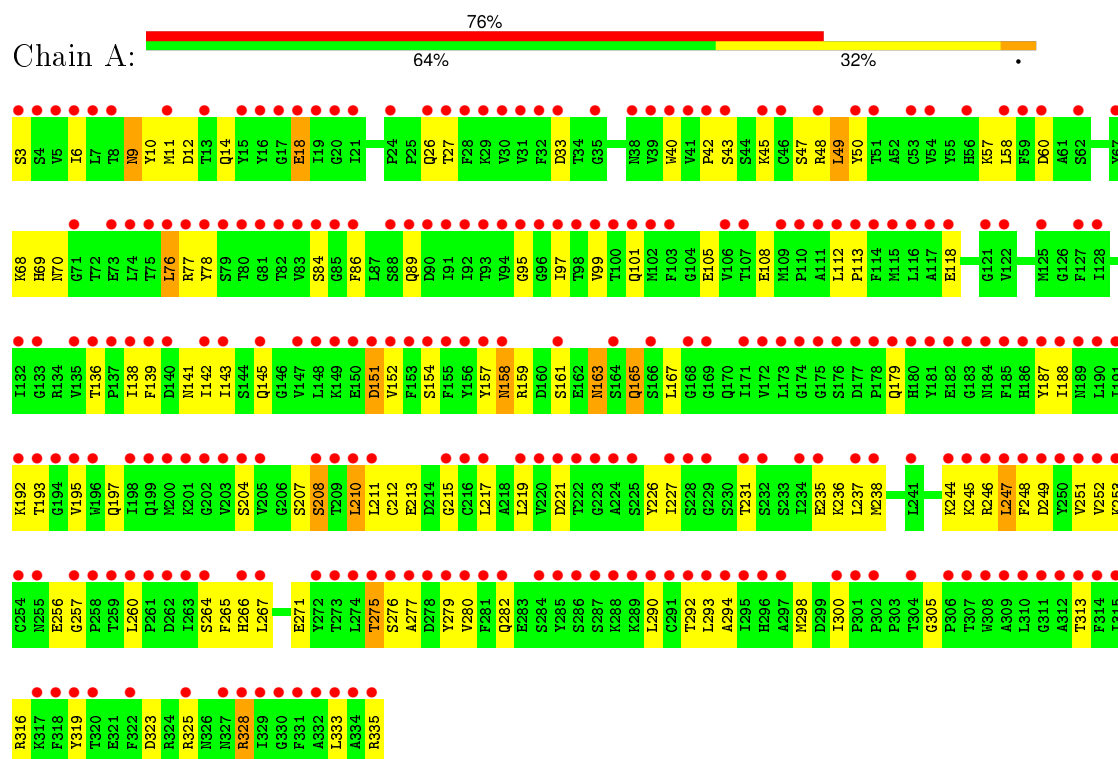
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	198	Total	O	0	0
			198	198		
4	B	124	Total	O	0	0
			124	124		

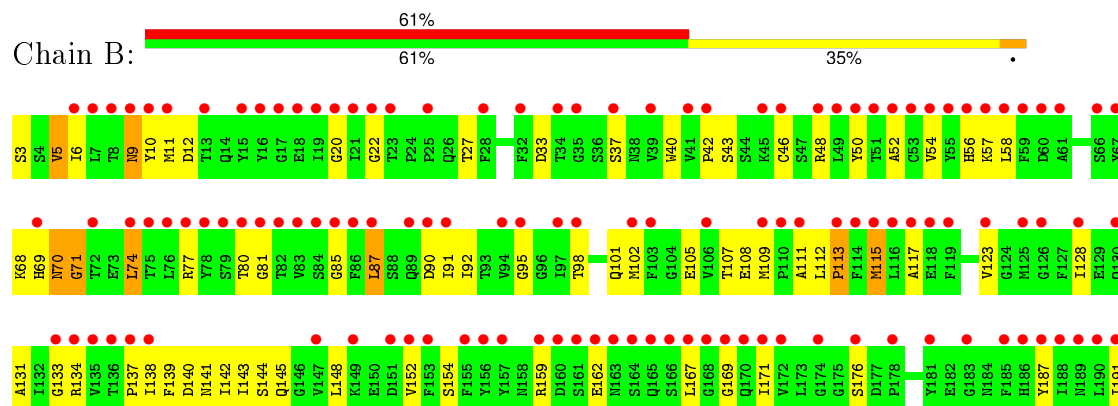
3 Residue-property plots

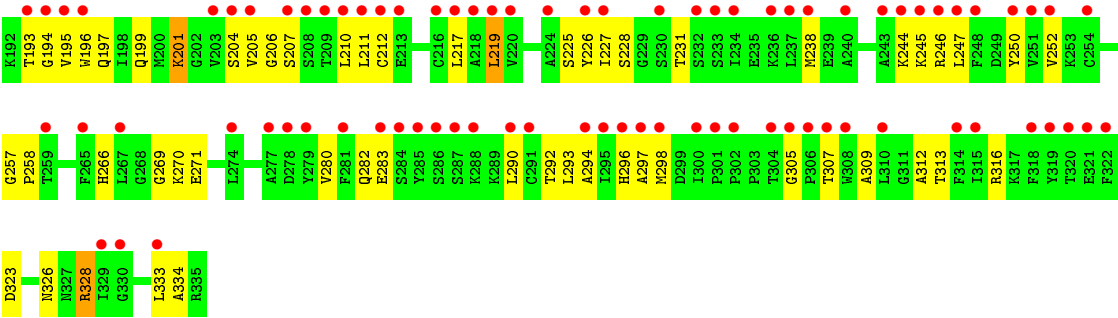
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ($\text{RSRZ} > 2$). 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: Renin



• Molecule 1: Renin





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α , β , γ	141.04Å 141.04Å 141.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 49.86 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.50) 94.2 (49.86-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.69 (at 2.51Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.197 , 0.252 0.229 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	42.0	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 46.0	EDS
Estimated twinning fraction	0.032 for l,-k,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 32300 reflections	Xtriage
F_o, F_c correlation	0.68	EDS
Total number of atoms	5550	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.52% 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

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, 5IG

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	0.33	0/2632	0.66	0/3568
1	B	0.32	0/2632	0.63	0/3568
All	All	0.33	0/5264	0.65	0/7136

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	2572	0	2500	110	0
1	B	2572	0	2500	127	0
2	A	14	0	13	0	0
2	B	14	0	13	5	0
3	A	28	0	27	1	0
3	B	28	0	27	2	0
4	A	198	0	0	18	0
4	B	124	0	0	19	0
All	All	5550	0	5080	237	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 23.

All (237) 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:A:112:LEU:HG	4:A:1158:HOH:O	1.52	1.07
1:A:45:LYS:HD2	4:A:1053:HOH:O	1.62	0.99
1:B:326:ASN:HB2	1:B:328:ARG:HD2	1.54	0.90
1:A:276:SER:O	1:A:280:VAL:HG12	1.74	0.88
1:A:158:ASN:ND2	1:A:159:ARG:H	1.73	0.86
1:B:52:ALA:HA	1:B:115:MET:HE2	1.56	0.85
1:A:152:VAL:HG12	1:A:323:ASP:HA	1.58	0.84
1:B:191:ILE:HG13	4:B:1026:HOH:O	1.78	0.83
1:B:217:LEU:HD22	4:B:1026:HOH:O	1.80	0.81
1:A:26:GLN:HE22	1:A:60:ASP:H	1.29	0.78
1:B:5:VAL:HG13	1:B:171:ILE:HG23	1.66	0.78
1:A:9:ASN:HD22	1:A:10:TYR:N	1.83	0.76
1:B:305:GLY:HA2	1:B:307:THR:HG23	1.66	0.75
1:B:197:GLN:OE1	1:B:307:THR:HG21	1.88	0.74
1:A:68:LYS:HB2	1:A:89:GLN:HB3	1.67	0.74
1:A:9:ASN:HD21	1:A:12:ASP:H	1.33	0.74
1:B:227:ILE:HG13	1:B:313:THR:HB	1.69	0.74
1:A:265:PHE:HE2	4:A:1169:HOH:O	1.72	0.73
1:A:163:ASN:O	1:A:165:GLN:HG2	1.90	0.72
1:B:323:ASP:HB3	1:B:328:ARG:HG2	1.72	0.72
1:A:265:PHE:CE2	4:A:1169:HOH:O	2.43	0.71
1:B:270:LYS:HG2	1:B:271:GLU:H	1.55	0.71
1:A:275:THR:HG21	4:A:1022:HOH:O	1.91	0.71
1:B:9:ASN:HD21	1:B:12:ASP:H	1.37	0.70
1:B:201:LYS:HE2	4:B:999:HOH:O	1.91	0.70
1:B:152:VAL:HG12	1:B:323:ASP:HA	1.74	0.69
1:A:99:VAL:HG12	1:A:145:GLN:OE1	1.93	0.69
1:B:6:ILE:HG23	1:B:167:LEU:CD1	2.23	0.68
1:A:158:ASN:HD22	1:A:159:ARG:H	1.41	0.68
1:B:280:VAL:HG22	1:B:293:LEU:HD22	1.76	0.68
1:B:9:ASN:HD22	1:B:10:TYR:N	1.91	0.68
1:B:298:MET:HE3	4:B:1032:HOH:O	1.93	0.67
1:B:210:LEU:H	1:B:210:LEU:HD23	1.60	0.67
1:A:231:THR:O	1:A:235:GLU:HG2	1.94	0.67
1:B:5:VAL:HG13	1:B:171:ILE:CG2	2.26	0.66
1:B:101:GLN:NE2	1:B:138:ILE:HA	2.09	0.66
1:A:208:SER:O	1:A:210:LEU:HD23	1.96	0.66
1:A:235:GLU:HG3	4:A:1126:HOH:O	1.94	0.66
1:A:207:SER:HB2	4:A:1105:HOH:O	1.95	0.66
1:B:70:ASN:HA	2:B:986:NAG:H81	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:217:LEU:HB2	1:B:307:THR:HG22	1.76	0.65
1:B:217:LEU:HB2	1:B:307:THR:CG2	2.26	0.65
1:B:252:VAL:HG23	4:B:991:HOH:O	1.95	0.65
1:B:257:GLY:HA3	1:B:282:GLN:HE22	1.61	0.65
1:B:50:TYR:OH	1:B:108:GLU:HG2	1.98	0.64
1:B:43:SER:HB2	1:B:105:GLU:HB3	1.78	0.64
1:A:253:LYS:HB2	1:A:256:GLU:OE2	1.99	0.62
1:B:33:ASP:OD1	3:B:886:5IG:N2	2.33	0.62
1:B:42:PRO:HG3	1:B:109:MET:HE1	1.82	0.62
1:B:270:LYS:HG2	1:B:271:GLU:N	2.15	0.61
1:A:27:THR:O	1:A:57:LYS:HG2	1.99	0.61
1:A:298:MET:SD	4:A:1116:HOH:O	2.56	0.61
1:B:9:ASN:C	1:B:9:ASN:HD22	2.03	0.61
1:A:99:VAL:HG11	1:A:142:ILE:HG12	1.82	0.61
1:A:323:ASP:OD1	1:A:328:ARG:HD2	2.00	0.60
1:B:217:LEU:HD12	1:B:307:THR:HG22	1.83	0.60
1:B:143:ILE:HD13	1:B:148:LEU:HD12	1.84	0.59
1:A:163:ASN:H	1:A:163:ASN:HD22	1.51	0.59
1:B:226:TYR:HB3	1:B:294:ALA:O	2.04	0.58
1:A:139:PHE:CE1	1:A:143:ILE:HD11	2.39	0.58
1:A:237:LEU:C	1:A:237:LEU:HD23	2.25	0.57
1:B:176:SER:HB3	4:B:1059:HOH:O	2.02	0.57
1:A:9:ASN:C	1:A:9:ASN:HD22	2.03	0.57
1:B:283:GLU:HB2	4:B:1106:HOH:O	2.04	0.57
1:A:158:ASN:ND2	1:A:159:ARG:N	2.49	0.57
1:B:191:ILE:HD11	1:B:199:GLN:HB2	1.86	0.57
1:A:257:GLY:HA3	1:A:282:GLN:HE22	1.69	0.57
1:B:296:HIS:CE1	4:B:1048:HOH:O	2.57	0.57
1:A:161:SER:OG	1:A:163:ASN:ND2	2.38	0.57
1:B:247:LEU:HA	4:B:1093:HOH:O	2.04	0.56
1:A:245:LYS:HD3	1:A:246:ARG:O	2.06	0.56
1:B:323:ASP:OD1	1:B:328:ARG:HD3	2.06	0.56
1:A:275:THR:HG23	4:A:1097:HOH:O	2.05	0.56
1:B:225:SER:O	1:B:312:ALA:HB3	2.07	0.55
1:A:43:SER:HB2	1:A:105:GLU:HB3	1.88	0.55
1:B:87:LEU:HG	4:B:1105:HOH:O	2.05	0.55
1:A:151:ASP:HB2	1:A:325:ARG:HB2	1.90	0.54
1:A:292:THR:HG21	4:A:1066:HOH:O	2.06	0.54
1:A:6:ILE:HD11	4:A:1130:HOH:O	2.07	0.54
1:A:212:CYS:SG	1:A:212:CYS:O	2.64	0.54
1:B:204:SER:HB3	4:B:1094:HOH:O	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:101:GLN:NE2	1:A:138:ILE:HA	2.22	0.54
1:A:251:VAL:HG21	1:A:290:LEU:HD13	1.89	0.54
1:B:102:MET:SD	4:B:1105:HOH:O	2.58	0.54
1:A:18:GLU:HG3	4:A:1141:HOH:O	2.07	0.54
1:B:139:PHE:O	1:B:143:ILE:HG12	2.08	0.54
1:B:27:THR:O	1:B:57:LYS:HD3	2.08	0.53
1:A:11:MET:O	1:A:12:ASP:HB2	2.08	0.53
1:A:275:THR:CG2	4:A:1097:HOH:O	2.55	0.53
1:A:101:GLN:HE22	1:A:138:ILE:HA	1.73	0.53
1:B:43:SER:O	1:B:58:LEU:HD22	2.08	0.53
1:B:292:THR:HG22	1:B:293:LEU:N	2.24	0.52
1:B:33:ASP:HB2	3:B:886:5IG:H52	1.89	0.52
1:B:141:ASN:O	1:B:144:SER:HB2	2.10	0.52
1:B:210:LEU:HD13	4:B:1004:HOH:O	2.09	0.52
1:B:231:THR:HA	1:B:297:ALA:HB1	1.92	0.52
1:A:9:ASN:HD21	1:A:12:ASP:N	2.03	0.52
1:A:188:ILE:CD1	1:A:267:LEU:HB3	2.40	0.52
1:B:69:HIS:CE1	1:B:71:GLY:H	2.27	0.52
1:A:247:LEU:HD13	1:A:248:PHE:CZ	2.45	0.52
1:A:10:TYR:CD2	1:A:118:GLU:HG3	2.45	0.51
1:B:9:ASN:HD21	1:B:12:ASP:N	2.06	0.51
1:B:85:GLY:HA3	1:B:105:GLU:O	2.10	0.51
1:A:47:SER:OG	1:A:49:LEU:HB2	2.08	0.51
1:A:9:ASN:ND2	1:A:9:ASN:C	2.62	0.51
1:A:246:ARG:HB2	1:A:249:ASP:O	2.10	0.51
1:B:326:ASN:CB	1:B:328:ARG:HD2	2.35	0.51
1:A:45:LYS:HD3	1:A:86:PHE:CZ	2.45	0.51
1:B:187:TYR:HD2	1:B:328:ARG:HG3	1.76	0.51
1:B:48:ARG:HD3	1:B:54:VAL:HG22	1.93	0.51
1:A:42:PRO:HB2	1:A:58:LEU:HD23	1.93	0.50
1:B:9:ASN:C	1:B:9:ASN:ND2	2.64	0.50
1:A:9:ASN:ND2	1:A:12:ASP:H	2.05	0.50
1:B:42:PRO:HG3	1:B:109:MET:CE	2.41	0.50
1:A:163:ASN:H	1:A:163:ASN:ND2	2.10	0.50
1:B:210:LEU:H	1:B:210:LEU:CD2	2.25	0.50
1:A:188:ILE:HD13	1:A:267:LEU:HB3	1.93	0.50
1:A:10:TYR:HB3	1:A:14:GLN:HB2	1.94	0.49
1:B:296:HIS:HE1	4:B:1048:HOH:O	1.93	0.49
1:A:197:GLN:CD	1:A:217:LEU:HD23	2.33	0.49
1:A:300:ILE:O	1:A:305:GLY:HA3	2.13	0.49
1:B:142:ILE:O	1:B:145:GLN:HG2	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:9:ASN:ND2	1:B:12:ASP:H	2.06	0.49
1:A:50:TYR:OH	1:A:108:GLU:HG2	2.13	0.49
1:B:22:GLY:HA2	1:B:90:ASP:OD1	2.13	0.48
1:B:90:ASP:OD1	1:B:91:ILE:N	2.46	0.48
1:A:167:LEU:H	1:A:167:LEU:HD23	1.77	0.48
1:B:217:LEU:HD12	1:B:307:THR:CG2	2.43	0.48
1:A:227:ILE:HG21	4:A:1169:HOH:O	2.13	0.48
1:A:210:LEU:HB2	1:A:236:LYS:NZ	2.28	0.48
1:A:158:ASN:HD22	1:A:159:ARG:N	2.10	0.48
1:B:134:ARG:NH2	2:B:986:NAG:O4	2.46	0.48
1:B:56:HIS:HE1	1:B:117:ALA:O	1.96	0.48
1:A:207:SER:O	1:A:208:SER:HB3	2.14	0.47
1:B:46:CYS:O	1:B:48:ARG:NH1	2.47	0.47
1:B:333:LEU:HD13	1:B:334:ALA:O	2.14	0.47
1:B:134:ARG:HH21	2:B:986:NAG:H61	1.78	0.47
1:A:247:LEU:HD22	1:A:248:PHE:CE2	2.48	0.47
1:A:231:THR:O	1:A:235:GLU:CG	2.60	0.47
1:A:226:TYR:HB3	1:A:294:ALA:O	2.14	0.47
1:A:275:THR:HG22	1:A:277:ALA:N	2.29	0.47
1:B:195:VAL:O	1:B:197:GLN:N	2.47	0.47
1:A:154:SER:HB3	1:A:319:TYR:CE1	2.50	0.47
1:B:245:LYS:HA	1:B:250:TYR:HA	1.97	0.47
1:A:275:THR:CG2	1:A:276:SER:N	2.79	0.46
1:A:213:GLU:C	1:A:215:GLY:H	2.18	0.46
1:A:227:ILE:HG13	1:A:313:THR:HB	1.96	0.46
1:B:70:ASN:HA	2:B:986:NAG:C8	2.45	0.46
1:B:196:TRP:O	1:B:219:LEU:HD23	2.15	0.46
1:B:42:PRO:HB2	1:B:58:LEU:HD23	1.98	0.46
1:A:136:THR:HG21	1:A:141:ASN:ND2	2.31	0.46
1:B:211:LEU:HD22	4:B:1107:HOH:O	2.16	0.46
1:B:154:SER:O	1:B:171:ILE:HA	2.16	0.45
1:B:270:LYS:HG2	4:B:1039:HOH:O	2.16	0.45
1:B:91:ILE:O	1:B:91:ILE:HG13	2.14	0.45
1:B:46:CYS:SG	1:B:107:THR:O	2.75	0.45
1:B:187:TYR:CD2	1:B:328:ARG:HG3	2.51	0.45
1:A:163:ASN:ND2	1:A:163:ASN:N	2.64	0.45
1:B:48:ARG:CD	1:B:54:VAL:HG22	2.46	0.45
1:A:163:ASN:HD22	1:A:163:ASN:N	2.10	0.45
1:B:139:PHE:CE1	1:B:143:ILE:HD11	2.52	0.45
1:A:204:SER:OG	1:A:264:SER:HB2	2.16	0.45
1:B:238:MET:HG3	1:B:250:TYR:CD2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:77:ARG:HG2	1:B:77:ARG:HH11	1.82	0.45
1:A:219:LEU:HD12	1:A:221:ASP:HB2	1.98	0.45
1:A:192:LYS:HG3	1:A:193:THR:O	2.17	0.45
1:A:244:LYS:HE3	4:A:1103:HOH:O	2.17	0.44
1:B:167:LEU:HD11	1:B:169:GLY:O	2.17	0.44
1:B:193:THR:HG22	1:B:194:GLY:N	2.31	0.44
1:B:123:VAL:CG1	1:B:138:ILE:HG21	2.48	0.44
1:B:133:GLY:O	1:B:134:ARG:CG	2.66	0.44
1:B:280:VAL:HG22	1:B:293:LEU:CD2	2.47	0.44
1:B:205:VAL:CG2	1:B:210:LEU:HD21	2.47	0.44
1:B:70:ASN:ND2	2:B:986:NAG:C7	2.81	0.44
1:A:3:SER:HB3	1:A:95:GLY:O	2.17	0.44
1:B:11:MET:O	1:B:12:ASP:HB2	2.17	0.44
1:B:206:GLY:C	4:B:1094:HOH:O	2.56	0.43
1:B:207:SER:N	4:B:1094:HOH:O	2.51	0.43
1:A:33:ASP:OD1	3:A:885:5IG:N2	2.51	0.43
1:B:111:ALA:O	1:B:115:MET:HB2	2.18	0.43
1:B:87:LEU:HD12	1:B:87:LEU:HA	1.87	0.43
1:A:300:ILE:HD12	1:A:300:ILE:N	2.32	0.43
1:A:211:LEU:CD2	1:A:236:LYS:HD3	2.49	0.43
1:B:137:PRO:HD2	1:B:140:ASP:OD2	2.18	0.43
1:A:163:ASN:ND2	1:A:163:ASN:O	2.52	0.43
1:B:139:PHE:CZ	1:B:143:ILE:HD11	2.54	0.43
1:A:237:LEU:O	1:A:237:LEU:HD23	2.18	0.43
1:A:49:LEU:HD12	1:A:49:LEU:HA	1.76	0.43
1:A:219:LEU:CD1	1:A:221:ASP:HB2	2.48	0.43
1:A:10:TYR:CG	1:A:118:GLU:HG3	2.54	0.43
1:B:37:SER:HB2	1:B:131:ALA:HA	2.01	0.43
1:B:187:TYR:N	1:B:187:TYR:CD1	2.87	0.42
1:B:266:HIS:NE2	1:B:269:GLY:HA2	2.33	0.42
1:A:77:ARG:HD3	4:A:1174:HOH:O	2.19	0.42
1:B:238:MET:HG3	1:B:250:TYR:CG	2.55	0.42
1:B:77:ARG:HG2	1:B:77:ARG:NH1	2.34	0.42
1:B:193:THR:CG2	1:B:194:GLY:N	2.82	0.42
1:B:159:ARG:HG2	1:B:316:ARG:HG2	2.02	0.42
1:B:133:GLY:O	1:B:134:ARG:HG2	2.18	0.42
1:A:292:THR:HG22	1:A:293:LEU:N	2.34	0.42
1:B:294:ALA:HB1	1:B:313:THR:OG1	2.20	0.42
1:B:40:TRP:HZ3	1:B:109:MET:HE1	1.84	0.42
1:A:40:TRP:CD1	1:A:76:LEU:HD23	2.54	0.42
1:A:179:GLN:O	1:A:335:ARG:HD2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:210:LEU:N	1:B:210:LEU:HD23	2.32	0.42
1:B:258:PRO:HD3	1:B:282:GLN:HE22	1.85	0.42
1:A:40:TRP:HE1	1:A:78:TYR:HH	1.66	0.42
1:B:80:THR:OG1	1:B:81:GLY:N	2.51	0.42
1:A:260:LEU:HB3	1:A:279:TYR:OH	2.20	0.42
1:A:325:ARG:O	1:A:325:ARG:HD2	2.20	0.42
1:A:238:MET:SD	1:A:293:LEU:HD12	2.60	0.42
1:B:212:CYS:O	1:B:212:CYS:SG	2.78	0.42
1:B:228:SER:OG	1:B:309:ALA:HB3	2.20	0.42
1:A:97:ILE:HG22	1:A:99:VAL:HG13	2.01	0.41
1:A:213:GLU:HB3	4:A:1051:HOH:O	2.19	0.41
1:B:92:ILE:O	1:B:98:THR:HA	2.21	0.41
1:A:275:THR:HG22	1:A:277:ALA:H	1.86	0.41
1:A:187:TYR:CD2	1:A:328:ARG:HD3	2.54	0.41
1:B:128:ILE:HG13	1:B:137:PRO:HD3	2.02	0.41
1:A:266:HIS:CD2	1:A:271:GLU:OE2	2.73	0.41
1:B:52:ALA:HA	1:B:115:MET:CE	2.39	0.41
1:A:84:SER:HB2	1:A:108:GLU:HB2	2.02	0.41
1:B:210:LEU:HB3	4:B:1004:HOH:O	2.21	0.41
1:A:211:LEU:HD22	4:A:1019:HOH:O	2.21	0.41
1:A:167:LEU:HD23	1:A:167:LEU:N	2.36	0.41
1:A:48:ARG:HD3	1:A:48:ARG:HA	1.78	0.41
1:A:237:LEU:CD2	1:A:237:LEU:C	2.90	0.41
1:B:20:GLY:O	1:B:92:ILE:HA	2.20	0.41
1:B:244:LYS:HA	1:B:244:LYS:HD3	1.91	0.41
1:B:68:LYS:HB2	1:B:68:LYS:HE3	1.67	0.40
1:A:69:HIS:CG	1:A:70:ASN:N	2.89	0.40
1:A:157:TYR:O	1:A:316:ARG:HA	2.21	0.40
1:A:195:VAL:HG12	1:A:197:GLN:HB2	2.03	0.40
1:B:112:LEU:HA	1:B:113:PRO:HA	1.63	0.40
1:B:123:VAL:HG13	1:B:138:ILE:HG21	2.03	0.40
1:B:74:LEU:O	1:B:74:LEU:HD12	2.21	0.40
1:B:3:SER:OG	1:B:95:GLY:O	2.36	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	331/333 (99%)	312 (94%)	17 (5%)	2 (1%)	30	50
1	B	331/333 (99%)	310 (94%)	18 (5%)	3 (1%)	21	37
All	All	662/666 (99%)	622 (94%)	35 (5%)	5 (1%)	24	41

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	165	GLN
1	B	70	ASN
1	A	208	SER
1	B	246	ARG
1	B	71	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	284/284 (100%)	270 (95%)	14 (5%)	31	55
1	B	284/284 (100%)	273 (96%)	11 (4%)	39	66
All	All	568/568 (100%)	543 (96%)	25 (4%)	35	60

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ASN

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Mol	Chain	Res	Type
1	A	18	GLU
1	A	49	LEU
1	A	76	LEU
1	A	113	PRO
1	A	151	ASP
1	A	158	ASN
1	A	163	ASN
1	A	210	LEU
1	A	247	LEU
1	A	252	VAL
1	A	275	THR
1	A	328	ARG
1	A	333	LEU
1	B	5	VAL
1	B	9	ASN
1	B	74	LEU
1	B	87	LEU
1	B	113	PRO
1	B	115	MET
1	B	162	GLU
1	B	201	LYS
1	B	219	LEU
1	B	290	LEU
1	B	328	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	26	GLN
1	A	101	GLN
1	A	141	ASN
1	A	158	ASN
1	A	163	ASN
1	A	165	GLN
1	A	189	ASN
1	A	197	GLN
1	A	199	GLN
1	A	266	HIS
1	A	282	GLN
1	B	9	ASN
1	B	26	GLN

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Mol	Chain	Res	Type
1	B	69	HIS
1	B	89	GLN
1	B	101	GLN
1	B	141	ASN
1	B	165	GLN
1	B	189	ASN
1	B	282	GLN
1	B	296	HIS

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](#)

4 ligands are 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	$\# Z > 2$
3	5IG	A	885	-	29,30,30	0.92	0	35,44,44	1.24	3 (8%)
2	NAG	A	985	1	14,14,15	0.49	0	15,19,21	0.83	1 (6%)
3	5IG	B	886	-	29,30,30	1.04	2 (6%)	35,44,44	1.21	4 (11%)
2	NAG	B	986	1	14,14,15	0.62	0	15,19,21	0.58	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
3	5IG	A	885	-	-	0/11/30/30	0/3/3/3
2	NAG	A	985	1	-	0/6/23/26	0/1/1/1
3	5IG	B	886	-	-	0/11/30/30	0/3/3/3
2	NAG	B	986	1	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	886	5IG	C15-N5	2.15	1.39	1.36
3	B	886	5IG	C8-C7	2.26	1.42	1.38

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	885	5IG	C4-C3-N2	-3.64	119.26	123.87
3	B	886	5IG	C4-C3-N2	-3.38	119.59	123.87
3	A	885	5IG	C11-C12-C4	-3.24	115.22	120.57
3	B	886	5IG	C11-C12-C4	-2.91	115.77	120.57
2	A	985	NAG	C2-N2-C7	-2.15	120.28	123.04
3	B	886	5IG	C4-C1-N1	2.03	123.80	122.55
3	B	886	5IG	C7-C12-C4	2.53	125.02	120.78
3	A	885	5IG	C7-C12-C4	2.65	125.22	120.78

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	885	5IG	1	0
3	B	886	5IG	2	0
2	B	986	NAG	5	0

5.7 Other polymers ⓘ

There are no such residues in this entry.







5.8 Polymer linkage issues ⓘ

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	333/333 (100%)	3.06	254 (76%)  	25, 40, 63, 88	0
1	B	333/333 (100%)	2.55	204 (61%)  	28, 46, 71, 89	0
All	All	666/666 (100%)	2.81	458 (68%)  	25, 43, 68, 89	0

All (458) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	175	GLY	9.3
1	A	211	LEU	8.1
1	A	148	LEU	7.8
1	A	147	VAL	7.5
1	B	209	THR	7.0
1	A	38	ASN	6.4
1	A	191	ILE	6.4
1	B	300	ILE	6.3
1	B	134	ARG	6.2
1	A	200	MET	6.2
1	A	99	VAL	6.2
1	B	247	LEU	6.2
1	A	329	ILE	6.2
1	A	93	THR	6.1
1	A	86	PHE	6.1
1	A	219	LEU	6.1
1	B	85	GLY	6.0
1	A	203	VAL	6.0
1	A	188	ILE	5.9
1	A	248	PHE	5.8
1	A	103	PHE	5.8
1	B	308	TRP	5.8
1	A	171	ILE	5.7
1	B	237	LEU	5.7

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Mol	Chain	Res	Type	RSRZ
1	A	193	THR	5.7
1	A	151	ASP	5.7
1	A	78	TYR	5.7
1	A	267	LEU	5.6
1	B	52	ALA	5.6
1	A	138	ILE	5.5
1	A	177	ASP	5.5
1	B	74	LEU	5.4
1	A	19	ILE	5.4
1	A	97	ILE	5.4
1	B	66	SER	5.3
1	B	102	MET	5.3
1	A	155	PHE	5.3
1	A	187	TYR	5.3
1	A	82	THR	5.3
1	B	284	SER	5.3
1	B	22	GLY	5.3
1	A	208	SER	5.2
1	B	195	VAL	5.2
1	A	71	GLY	5.2
1	B	306	PRO	5.1
1	A	205	VAL	5.1
1	A	320	THR	5.1
1	B	304	THR	5.1
1	B	79	SER	5.1
1	A	319	TYR	5.1
1	A	132	ILE	5.0
1	A	81	GLY	5.0
1	A	186	HIS	5.0
1	A	85	GLY	5.0
1	A	290	LEU	4.9
1	B	291	CYS	4.9
1	B	78	TYR	4.9
1	B	294	ALA	4.9
1	A	196	TRP	4.9
1	A	153	PHE	4.8
1	A	40	TRP	4.8
1	A	272	TYR	4.8
1	A	8	THR	4.8
1	A	16	TYR	4.8
1	B	61	ALA	4.7
1	B	310	LEU	4.7

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Mol	Chain	Res	Type	RSRZ
1	A	231	THR	4.7
1	B	19	ILE	4.7
1	B	151	ASP	4.7
1	A	28	PHE	4.7
1	A	308	TRP	4.6
1	A	111	ALA	4.6
1	A	92	ILE	4.6
1	A	190	LEU	4.6
1	B	136	THR	4.5
1	A	247	LEU	4.5
1	B	250	TYR	4.5
1	B	77	ARG	4.5
1	B	53	CYS	4.4
1	B	119	PHE	4.4
1	B	87	LEU	4.4
1	B	167	LEU	4.4
1	B	60	ASP	4.4
1	B	298	MET	4.4
1	B	95	GLY	4.4
1	B	133	GLY	4.4
1	A	74	LEU	4.3
1	B	55	TYR	4.3
1	A	315	ILE	4.3
1	A	109	MET	4.3
1	B	233	SER	4.3
1	A	280	VAL	4.3
1	A	210	LEU	4.2
1	B	267	LEU	4.2
1	A	311	GLY	4.2
1	A	274	LEU	4.2
1	A	156	TYR	4.2
1	A	276	SER	4.2
1	A	21	ILE	4.2
1	B	211	LEU	4.2
1	B	161	SER	4.2
1	B	83	VAL	4.1
1	A	331	PHE	4.1
1	A	198	ILE	4.1
1	B	37	SER	4.1
1	B	166	SER	4.1
1	A	31	VAL	4.1
1	A	258	PRO	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	264	SER	4.1
1	A	300	ILE	4.1
1	A	168	GLY	4.0
1	B	114	PHE	4.0
1	B	34	THR	4.0
1	A	192	LYS	4.0
1	B	274	LEU	4.0
1	A	180	HIS	4.0
1	A	157	TYR	4.0
1	A	263	ILE	4.0
1	B	188	ILE	4.0
1	A	251	VAL	3.9
1	A	281	PHE	3.9
1	B	86	PHE	3.9
1	B	84	SER	3.9
1	A	241	LEU	3.9
1	B	59	PHE	3.9
1	A	142	ILE	3.9
1	A	13	THR	3.9
1	A	254	CYS	3.9
1	B	216	CYS	3.9
1	A	39	VAL	3.9
1	A	195	VAL	3.9
1	B	103	PHE	3.9
1	A	221	ASP	3.8
1	B	251	VAL	3.8
1	B	178	PRO	3.8
1	A	76	LEU	3.8
1	B	21	ILE	3.8
1	A	15	TYR	3.8
1	A	150	GLU	3.8
1	B	9	ASN	3.8
1	A	285	TYR	3.8
1	B	285	TYR	3.8
1	B	295	ILE	3.8
1	A	179	GLN	3.8
1	A	204	SER	3.7
1	B	318	PHE	3.7
1	A	307	THR	3.7
1	A	149	LYS	3.7
1	A	220	VAL	3.7
1	A	334	ALA	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	310	LEU	3.7
1	A	172	VAL	3.7
1	A	327	ASN	3.7
1	A	293	LEU	3.7
1	A	6	ILE	3.7
1	A	295	ILE	3.7
1	B	56	HIS	3.7
1	A	250	TYR	3.6
1	B	319	TYR	3.6
1	A	253	LYS	3.6
1	B	301	PRO	3.6
1	B	333	LEU	3.6
1	A	176	SER	3.6
1	A	277	ALA	3.6
1	A	42	PRO	3.6
1	A	137	PRO	3.6
1	A	174	GLY	3.6
1	B	82	THR	3.6
1	A	18	GLU	3.6
1	A	216	CYS	3.6
1	B	183	GLY	3.6
1	B	219	LEU	3.6
1	A	88	SER	3.6
1	A	136	THR	3.6
1	B	72	THR	3.5
1	A	54	VAL	3.5
1	B	163	ASN	3.5
1	B	13	THR	3.5
1	A	238	MET	3.5
1	B	109	MET	3.5
1	A	106	VAL	3.5
1	B	7	LEU	3.5
1	A	246	ARG	3.5
1	B	307	THR	3.5
1	B	281	PHE	3.5
1	A	252	VAL	3.5
1	A	24	PRO	3.5
1	A	306	PRO	3.5
1	A	75	THR	3.5
1	B	69	HIS	3.5
1	A	30	VAL	3.5
1	A	317	LYS	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	94	VAL	3.5
1	B	128	ILE	3.5
1	A	279	TYR	3.5
1	A	297	ALA	3.5
1	A	318	PHE	3.5
1	A	291	CYS	3.5
1	B	213	GLU	3.5
1	B	76	LEU	3.5
1	B	162	GLU	3.4
1	A	83	VAL	3.4
1	A	27	THR	3.4
1	B	50	TYR	3.4
1	A	91	ILE	3.4
1	A	128	ILE	3.4
1	B	81	GLY	3.4
1	B	196	TRP	3.4
1	B	46	CYS	3.4
1	B	89	GLN	3.4
1	A	135	VAL	3.4
1	B	106	VAL	3.4
1	B	252	VAL	3.4
1	B	165	GLN	3.4
1	B	168	GLY	3.4
1	A	80	THR	3.3
1	A	209	THR	3.3
1	A	62	SER	3.3
1	A	287	SER	3.3
1	B	113	PRO	3.3
1	B	164	SER	3.3
1	B	32	PHE	3.3
1	B	207	SER	3.3
1	B	191	ILE	3.3
1	A	154	SER	3.3
1	B	230	SER	3.3
1	B	330	GLY	3.3
1	B	125	MET	3.3
1	A	17	GLY	3.3
1	A	330	GLY	3.3
1	B	172	VAL	3.3
1	A	224	ALA	3.3
1	A	185	PHE	3.3
1	A	145	GLN	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	121	GLY	3.3
1	B	20	GLY	3.3
1	B	91	ILE	3.3
1	A	235	GLU	3.3
1	A	257	GLY	3.2
1	B	305	GLY	3.2
1	B	149	LYS	3.2
1	A	67	TYR	3.2
1	A	304	THR	3.2
1	B	203	VAL	3.2
1	B	45	LYS	3.2
1	B	75	THR	3.2
1	A	73	GLU	3.2
1	A	166	SER	3.2
1	A	110	PRO	3.2
1	B	18	GLU	3.2
1	A	328	ARG	3.2
1	A	244	LYS	3.2
1	A	314	PHE	3.2
1	B	111	ALA	3.2
1	A	20	GLY	3.2
1	B	279	TYR	3.2
1	B	245	LYS	3.2
1	B	116	LEU	3.2
1	A	286	SER	3.1
1	B	288	LYS	3.1
1	A	282	GLN	3.1
1	B	130	GLN	3.1
1	B	57	LYS	3.1
1	A	260	LEU	3.1
1	B	193	THR	3.1
1	B	123	VAL	3.1
1	A	158	ASN	3.1
1	B	155	PHE	3.1
1	A	118	GLU	3.1
1	A	94	VAL	3.1
1	B	170	GLN	3.0
1	B	181	TYR	3.0
1	A	234	ILE	3.0
1	A	302	PRO	3.0
1	A	296	HIS	3.0
1	A	223	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	254	CYS	3.0
1	B	49	LEU	3.0
1	A	289	LYS	3.0
1	B	208	SER	3.0
1	A	173	LEU	3.0
1	A	322	PHE	3.0
1	A	3	SER	3.0
1	A	261	PRO	3.0
1	B	320	THR	2.9
1	A	288	LYS	2.9
1	A	139	PHE	2.9
1	A	332	ALA	2.9
1	A	77	ARG	2.9
1	A	51	THR	2.9
1	A	225	SER	2.9
1	B	176	SER	2.9
1	A	32	PHE	2.9
1	B	80	THR	2.9
1	A	222	THR	2.9
1	B	97	ILE	2.9
1	B	110	PRO	2.9
1	A	41	VAL	2.9
1	A	95	GLY	2.9
1	B	189	ASN	2.8
1	B	115	MET	2.8
1	A	60	ASP	2.8
1	B	322	PHE	2.8
1	A	182	GLU	2.8
1	A	309	ALA	2.8
1	B	169	GLY	2.8
1	A	127	PHE	2.8
1	B	23	THR	2.8
1	A	245	LYS	2.8
1	B	244	LYS	2.8
1	B	232	SER	2.8
1	B	157	TYR	2.8
1	B	8	THR	2.8
1	A	194	GLY	2.8
1	B	138	ILE	2.8
1	B	126	GLY	2.8
1	A	294	ALA	2.8
1	A	266	HIS	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	133	GLY	2.8
1	B	204	SER	2.8
1	B	234	ILE	2.7
1	A	217	LEU	2.7
1	B	283	GLU	2.7
1	B	6	ILE	2.7
1	B	171	ILE	2.7
1	B	17	GLY	2.7
1	B	236	LYS	2.7
1	A	100	THR	2.7
1	A	125	MET	2.7
1	A	96	GLY	2.7
1	A	169	GLY	2.7
1	A	237	LEU	2.7
1	B	15	TYR	2.7
1	B	246	ARG	2.7
1	A	43	SER	2.7
1	A	232	SER	2.7
1	B	226	TYR	2.7
1	B	329	ILE	2.7
1	A	59	PHE	2.7
1	A	273	THR	2.7
1	A	116	LEU	2.6
1	B	210	LEU	2.6
1	B	217	LEU	2.6
1	B	240	ALA	2.6
1	A	4	SER	2.6
1	B	287	SER	2.6
1	B	205	VAL	2.6
1	B	265	PHE	2.6
1	A	56	HIS	2.6
1	A	58	LEU	2.6
1	A	143	ILE	2.6
1	B	185	PHE	2.6
1	B	16	TYR	2.6
1	A	249	ASP	2.6
1	B	25	PRO	2.6
1	A	114	PHE	2.6
1	B	220	VAL	2.5
1	A	301	PRO	2.5
1	A	259	THR	2.5
1	A	201	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	194	GLY	2.5
1	B	187	TYR	2.5
1	A	90	ASP	2.5
1	B	160	ASP	2.5
1	B	48	ARG	2.5
1	B	290	LEU	2.5
1	A	292	THR	2.5
1	B	238	MET	2.5
1	B	118	GLU	2.5
1	A	140	ASP	2.5
1	B	90	ASP	2.5
1	B	218	ALA	2.5
1	A	164	SER	2.5
1	B	314	PHE	2.5
1	B	41	VAL	2.5
1	B	42	PRO	2.5
1	B	296	HIS	2.5
1	A	325	ARG	2.5
1	B	153	PHE	2.4
1	B	39	VAL	2.4
1	A	112	LEU	2.4
1	A	184	ASN	2.4
1	A	11	MET	2.4
1	B	277	ALA	2.4
1	B	321	GLU	2.4
1	B	10	TYR	2.4
1	A	5	VAL	2.4
1	A	333	LEU	2.4
1	B	212	CYS	2.4
1	B	259	THR	2.4
1	B	67	TYR	2.4
1	A	255	ASN	2.4
1	A	228	SER	2.4
1	A	262	ASP	2.4
1	A	89	GLN	2.4
1	A	215	GLY	2.4
1	A	229	GLY	2.4
1	A	335	ARG	2.4
1	A	278	ASP	2.4
1	B	54	VAL	2.4
1	A	161	SER	2.4
1	A	115	MET	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	224	ALA	2.4
1	B	243	ALA	2.4
1	A	98	THR	2.3
1	B	186	HIS	2.3
1	A	26	GLN	2.3
1	A	152	VAL	2.3
1	A	227	ILE	2.3
1	B	137	PRO	2.3
1	B	156	TYR	2.3
1	A	122	VAL	2.3
1	A	102	MET	2.3
1	A	107	THR	2.3
1	B	315	ILE	2.3
1	A	79	SER	2.3
1	A	181	TYR	2.3
1	A	178	PRO	2.3
1	A	46	CYS	2.2
1	A	189	ASN	2.2
1	B	174	GLY	2.2
1	B	135	VAL	2.2
1	B	248	PHE	2.2
1	A	275	THR	2.2
1	A	313	THR	2.2
1	A	312	ALA	2.2
1	A	35	GLY	2.2
1	A	202	GLY	2.2
1	B	117	ALA	2.2
1	A	33	ASP	2.2
1	B	159	ARG	2.2
1	A	50	TYR	2.2
1	B	286	SER	2.2
1	A	101	GLN	2.2
1	B	58	LEU	2.2
1	B	152	VAL	2.1
1	A	117	ALA	2.1
1	B	11	MET	2.1
1	B	190	LEU	2.1
1	B	51	THR	2.1
1	A	199	GLN	2.1
1	A	29	LYS	2.1
1	A	53	CYS	2.1
1	A	48	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	113	PRO	2.1
1	A	183	GLY	2.1
1	B	302	PRO	2.1
1	A	284	SER	2.1
1	B	227	ILE	2.1
1	A	7	LEU	2.1
1	B	35	GLY	2.1
1	B	98	THR	2.1
1	B	278	ASP	2.0
1	B	297	ALA	2.0
1	A	45	LYS	2.0
1	B	147	VAL	2.0
1	A	84	SER	2.0
1	B	28	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](#)

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(Å ²)	Q<0.9
3	5IG	B	886	28/28	0.70	0.36	0.72	35,39,41,46	0
2	NAG	B	986	14/15	0.53	0.39	0.17	79,83,83,84	0
3	5IG	A	885	28/28	0.79	0.32	-0.09	24,30,36,38	0
2	NAG	A	985	14/15	0.69	0.30	-	48,53,56,58	0

6.5 Other polymers

There are no such residues in this entry.