



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:02 PM GMT

PDB ID : 4KC3
Title : Cytokine/receptor binary complex
Authors : Liu, X.; Wang, X.Q.
Deposited on : 2013-04-24
Resolution : 3.27 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

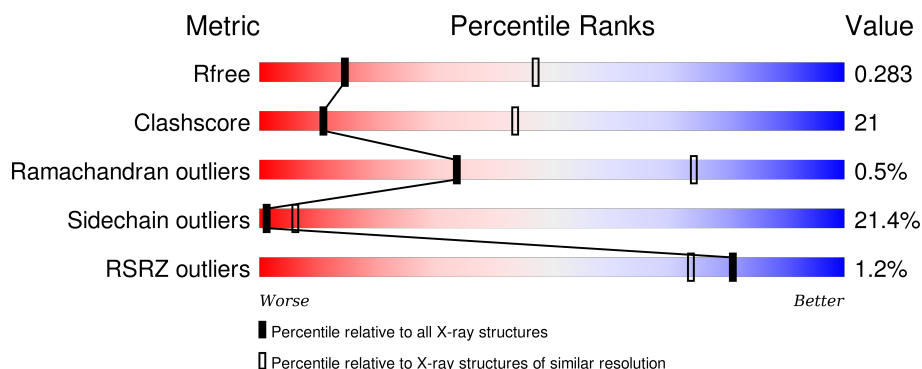
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 3.27 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	1756 (3.36-3.20)
Clashscore	102246	1941 (3.36-3.20)
Ramachandran outliers	100387	1905 (3.36-3.20)
Sidechain outliers	100360	1903 (3.36-3.20)
RSRZ outliers	91569	1764 (3.36-3.20)

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	159	 50% 25% 10% • 14%
2	B	309	 45% 37% 8% • 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3354 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 Interleukin-33.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
1	A	137	Total	C	N	O	S	Se	0	0	0
			1110	712	177	215	3	3			

- Molecule 2 is a protein called Interleukin-1 receptor-like 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	278	Total	C	N	O	S	0	0	0
			2202	1395	378	416	13			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	322	HIS	-	EXPRESSION TAG	UNP Q01638
B	323	HIS	-	EXPRESSION TAG	UNP Q01638
B	324	HIS	-	EXPRESSION TAG	UNP Q01638
B	325	HIS	-	EXPRESSION TAG	UNP Q01638
B	326	HIS	-	EXPRESSION TAG	UNP Q01638
B	327	HIS	-	EXPRESSION TAG	UNP Q01638

- Molecule 3 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C₈H₁₅NO₆).

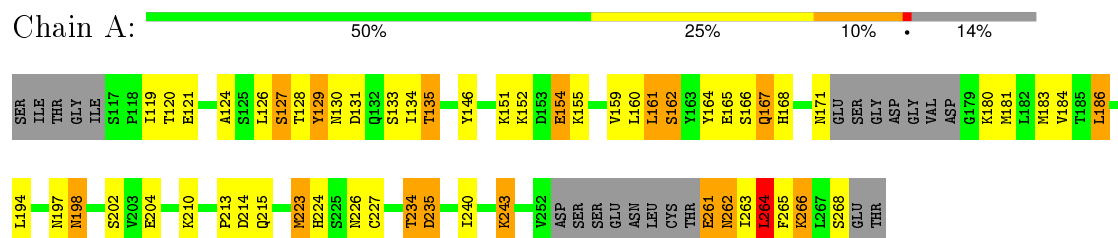


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		

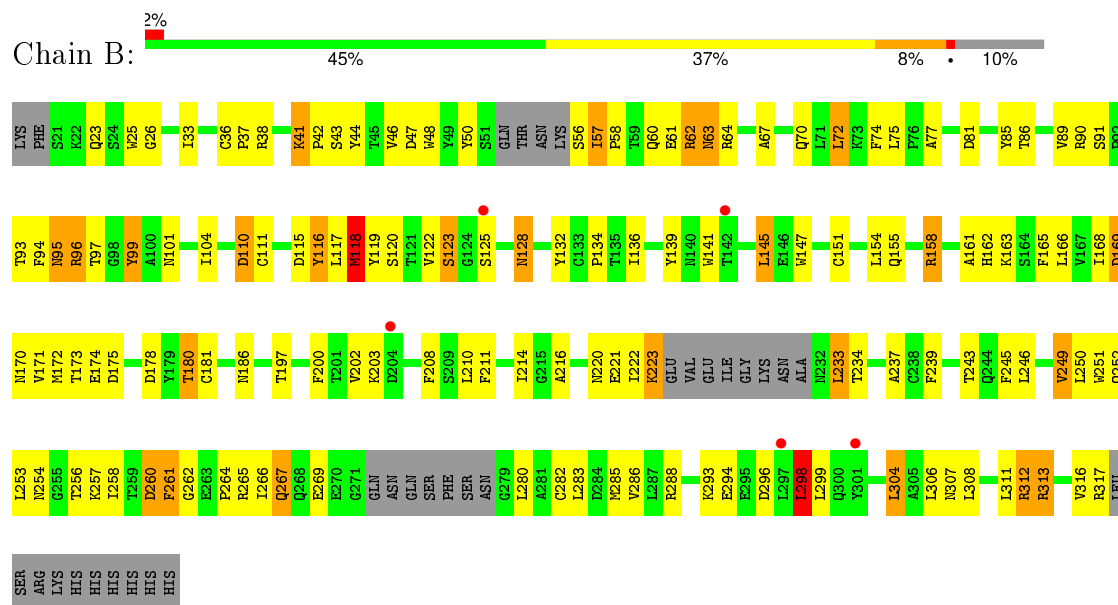
3 Residue-property plots

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: Interleukin-33



• Molecule 2: Interleukin-1 receptor-like 1



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	329.92Å 329.92Å 43.73Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	32.01 – 3.27 32.01 – 3.21	Depositor EDS
% Data completeness (in resolution range)	99.8 (32.01-3.27) 99.7 (32.01-3.21)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.13 (at 3.18Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.238 , 0.283 0.232 , 0.283	Depositor DCC
R_{free} test set	702 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	110.5	Xtriage
Anisotropy	0.010	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 99.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 14817 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3354	wwPDB-VP
Average B, all atoms (Å ²)	133.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.01% 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

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	0.54	0/1132	0.78	2/1523 (0.1%)
2	B	0.53	0/2249	0.79	4/3047 (0.1%)
All	All	0.54	0/3381	0.79	6/4570 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	2

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	264	LEU	CA-CB-CG	5.95	128.98	115.30
2	B	145	LEU	CA-CB-CG	5.82	128.70	115.30
1	A	264	LEU	CB-CG-CD2	5.67	120.64	111.00
2	B	62	ARG	N-CA-C	-5.39	96.44	111.00
2	B	258	ILE	N-CA-C	5.22	125.09	111.00
2	B	61	GLU	N-CA-C	5.19	125.00	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	116	TYR	Peptide
2	B	260	ASP	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	1110	0	1083	45	0
2	B	2202	0	2137	102	0
3	B	42	0	39	3	0
All	All	3354	0	3259	139	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 21.

All (139) 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:129:TYR:H	1:A:264:LEU:HG	1.38	0.86
2:B:117:LEU:HD23	2:B:197:THR:H	1.45	0.81
2:B:155:GLN:O	2:B:158:ARG:NH1	2.16	0.79
1:A:161:LEU:HD12	1:A:186:LEU:HD12	1.66	0.77
2:B:214:ILE:H	2:B:312:ARG:HH12	1.32	0.77
1:A:166:SER:OG	1:A:181:MSE:O	2.04	0.74
2:B:125:SER:HA	2:B:171:VAL:HG12	1.69	0.72
2:B:46:VAL:HG22	2:B:89:VAL:HG22	1.72	0.70
2:B:47:ASP:OD1	2:B:90:ARG:NH1	2.24	0.69
2:B:298:LEU:HD22	2:B:299:LEU:HG	1.73	0.69
2:B:267:GLN:HE22	2:B:288:ARG:HH21	1.41	0.69
2:B:117:LEU:HD23	2:B:197:THR:N	2.07	0.68
2:B:36:CYS:HB3	2:B:70:GLN:HG2	1.77	0.67
2:B:91:SER:OG	2:B:93:THR:N	2.28	0.66
2:B:91:SER:OG	2:B:94:PHE:N	2.29	0.66
2:B:63:ASN:OD1	2:B:63:ASN:N	2.29	0.64
2:B:25:TRP:HE1	2:B:117:LEU:HA	1.62	0.64
1:A:152:LYS:N	2:B:118:MET:O	2.30	0.64
2:B:23:GLN:HG3	2:B:101:ASN:HB3	1.80	0.64
2:B:280:LEU:HD23	2:B:280:LEU:H	1.63	0.63
2:B:77:ALA:HB1	2:B:104:ILE:HG12	1.80	0.62
2:B:25:TRP:HE1	2:B:117:LEU:HD12	1.64	0.60
2:B:253:LEU:HD21	2:B:264:PRO:HG2	1.83	0.60
2:B:25:TRP:NE1	2:B:117:LEU:HD12	2.17	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:123:SER:HB3	2:B:200:PHE:CE1	2.38	0.59
2:B:67:ALA:HB2	2:B:72:LEU:HD23	1.85	0.58
2:B:91:SER:HB3	2:B:94:PHE:CZ	2.38	0.58
2:B:123:SER:OG	2:B:128:ASN:O	2.09	0.57
1:A:235:ASP:OD1	1:A:235:ASP:N	2.38	0.57
1:A:262:ASN:HA	1:A:264:LEU:HD13	1.87	0.56
2:B:158:ARG:HG3	2:B:170:ASN:HD22	1.70	0.56
2:B:118:MET:HG2	2:B:134:PRO:CB	2.35	0.56
2:B:214:ILE:H	2:B:312:ARG:NH1	2.01	0.55
2:B:125:SER:O	2:B:171:VAL:HB	2.07	0.55
1:A:166:SER:HA	1:A:183:MSE:HE2	1.89	0.55
1:A:130:ASN:HA	2:B:132:TYR:CE2	2.41	0.55
1:A:119:ILE:HD13	1:A:164:TYR:HA	1.89	0.55
2:B:94:PHE:CD2	2:B:96:ARG:HD3	2.42	0.54
2:B:25:TRP:CD1	2:B:117:LEU:HD12	2.42	0.54
1:A:240:ILE:O	1:A:263:ILE:HG23	2.08	0.54
2:B:33:ILE:HG12	2:B:139:TYR:CE1	2.43	0.53
2:B:123:SER:HB3	2:B:200:PHE:CZ	2.44	0.53
2:B:298:LEU:HD13	2:B:299:LEU:HD12	1.91	0.52
1:A:152:LYS:CG	2:B:119:TYR:HA	2.39	0.52
2:B:180:THR:HB	2:B:197:THR:OG1	2.08	0.52
2:B:81:ASP:O	2:B:85:TYR:OH	2.20	0.52
2:B:25:TRP:HZ2	2:B:118:MET:H	1.57	0.52
2:B:210:LEU:HG	2:B:211:PHE:H	1.75	0.52
1:A:124:ALA:HB2	2:B:245:PHE:HZ	1.75	0.52
1:A:261:GLU:N	1:A:261:GLU:OE1	2.43	0.51
2:B:125:SER:OG	2:B:171:VAL:O	2.15	0.51
1:A:131:ASP:OD1	1:A:266:LYS:NZ	2.40	0.51
2:B:254:ASN:OD1	2:B:299:LEU:HB3	2.11	0.51
1:A:152:LYS:HG2	2:B:119:TYR:HA	1.91	0.50
2:B:95:ASN:OD1	3:B:401:NAG:O5	2.24	0.50
2:B:186:ASN:HD21	3:B:402:NAG:H4	1.76	0.50
2:B:172:MET:HB2	2:B:174:GLU:HB2	1.93	0.50
2:B:169:ASP:N	2:B:169:ASP:OD2	2.38	0.50
2:B:42:PRO:O	2:B:43:SER:OG	2.28	0.50
2:B:246:LEU:HD12	2:B:308:LEU:HD23	1.94	0.49
1:A:165:GLU:OE1	2:B:313:ARG:NH1	2.44	0.49
2:B:172:MET:O	2:B:202:VAL:HG21	2.12	0.49
1:A:121:GLU:HB2	1:A:160:LEU:HD11	1.95	0.49
1:A:135:THR:HG22	1:A:146:TYR:HB2	1.95	0.49
2:B:155:GLN:N	2:B:155:GLN:OE1	2.45	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:TYR:N	1:A:264:LEU:HG	2.18	0.47
2:B:186:ASN:ND2	3:B:402:NAG:H4	2.29	0.47
2:B:36:CYS:HB2	2:B:48:TRP:CZ2	2.49	0.47
2:B:265:ARG:NH1	2:B:266:ILE:O	2.47	0.47
1:A:128:THR:HG23	1:A:134:ILE:HD11	1.96	0.47
2:B:158:ARG:HG3	2:B:170:ASN:ND2	2.29	0.47
1:A:151:LYS:O	1:A:154:GLU:HB2	2.15	0.47
2:B:214:ILE:HG13	2:B:312:ARG:NH1	2.30	0.46
1:A:162:SER:O	1:A:184:VAL:HA	2.15	0.46
2:B:72:LEU:HD13	2:B:74:PHE:CE2	2.50	0.46
2:B:110:ASP:OD1	2:B:111:CYS:N	2.49	0.46
1:A:262:ASN:CA	1:A:264:LEU:HD13	2.46	0.45
2:B:168:ILE:CG2	2:B:171:VAL:HG23	2.46	0.45
2:B:91:SER:CB	2:B:93:THR:H	2.28	0.45
2:B:239:PHE:HZ	2:B:249:VAL:HG12	1.80	0.45
2:B:74:PHE:O	2:B:75:LEU:HD23	2.16	0.45
2:B:38:ARG:CZ	2:B:96:ARG:HG3	2.46	0.45
1:A:129:TYR:HB3	1:A:264:LEU:HB2	1.98	0.45
2:B:261:PHE:CD2	2:B:262:GLY:N	2.85	0.45
2:B:47:ASP:OD1	2:B:90:ARG:HD2	2.17	0.45
1:A:152:LYS:HE3	1:A:152:LYS:HB3	1.73	0.45
1:A:129:TYR:CG	1:A:264:LEU:HD12	2.52	0.45
1:A:186:LEU:N	1:A:186:LEU:HD13	2.32	0.45
2:B:214:ILE:HG22	2:B:216:ALA:O	2.17	0.45
2:B:38:ARG:NH1	2:B:96:ARG:HG3	2.32	0.44
2:B:254:ASN:OD1	2:B:299:LEU:HD13	2.17	0.44
2:B:136:ILE:HG12	2:B:141:TRP:NE1	2.32	0.44
1:A:130:ASN:HA	2:B:132:TYR:CD2	2.53	0.44
2:B:214:ILE:HA	2:B:237:ALA:HA	1.98	0.44
2:B:33:ILE:HG12	2:B:139:TYR:CD1	2.52	0.44
1:A:186:LEU:HD23	1:A:194:LEU:HD22	1.99	0.44
2:B:158:ARG:HB2	2:B:170:ASN:HB3	1.99	0.44
2:B:293:LYS:N	2:B:296:ASP:OD2	2.33	0.44
2:B:222:ILE:O	2:B:223:LYS:HD3	2.17	0.44
1:A:127:SER:O	1:A:265:PHE:HA	2.16	0.44
1:A:213:PRO:HB2	1:A:215:GLN:OE1	2.18	0.44
1:A:197:ASN:ND2	1:A:204:GLU:OE1	2.41	0.44
2:B:132:TYR:CD1	2:B:165:PHE:HB3	2.53	0.43
1:A:129:TYR:HB3	1:A:264:LEU:CB	2.49	0.43
2:B:122:VAL:HG11	2:B:203:LYS:HZ2	1.84	0.43
2:B:91:SER:HG	2:B:94:PHE:N	2.14	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:48:TRP:O	2:B:57:ILE:HB	2.18	0.43
2:B:210:LEU:O	2:B:307:ASN:ND2	2.52	0.43
1:A:198:ASN:N	1:A:198:ASN:OD1	2.51	0.43
2:B:41:LYS:HB2	2:B:44:TYR:CD2	2.53	0.43
1:A:243:LYS:HA	1:A:243:LYS:HE2	2.01	0.43
1:A:129:TYR:CD2	1:A:264:LEU:HD12	2.54	0.42
2:B:171:VAL:HG22	2:B:175:ASP:HB2	2.00	0.42
1:A:214:ASP:HB3	1:A:234:THR:HG23	2.01	0.42
2:B:116:TYR:N	2:B:116:TYR:CD2	2.85	0.42
2:B:254:ASN:O	2:B:254:ASN:ND2	2.53	0.42
1:A:167:GLN:HA	1:A:180:LYS:HE2	2.02	0.42
2:B:48:TRP:CE3	2:B:72:LEU:HG	2.55	0.42
2:B:266:ILE:HG12	2:B:285:MET:SD	2.60	0.42
2:B:26:GLY:O	2:B:104:ILE:HA	2.19	0.42
2:B:304:LEU:HB3	2:B:311:LEU:HD11	2.03	0.41
1:A:186:LEU:H	1:A:186:LEU:HD22	1.84	0.41
2:B:253:LEU:N	2:B:256:THR:O	2.37	0.41
1:A:167:GLN:CD	1:A:167:GLN:H	2.23	0.41
2:B:99:TYR:N	2:B:99:TYR:CD2	2.88	0.41
2:B:233:LEU:HD12	2:B:234:THR:H	1.84	0.41
1:A:129:TYR:CD2	1:A:130:ASN:N	2.88	0.41
2:B:162:HIS:CE1	2:B:163:LYS:HG3	2.55	0.41
2:B:252:GLN:HG2	2:B:256:THR:N	2.35	0.41
1:A:226:ASN:HD21	2:B:245:PHE:HB2	1.84	0.41
2:B:233:LEU:HD11	2:B:251:TRP:CH2	2.56	0.41
1:A:262:ASN:C	1:A:264:LEU:HD13	2.40	0.41
2:B:161:ALA:HB2	2:B:166:LEU:HD13	2.02	0.41
1:A:223:MSE:N	1:A:223:MSE:SE	3.03	0.41
2:B:118:MET:HG2	2:B:134:PRO:HB3	2.03	0.40
2:B:36:CYS:HA	2:B:37:PRO:HD3	1.88	0.40
1:A:197:ASN:HB3	1:A:202:SER:OG	2.21	0.40
2:B:147:TRP:CE3	2:B:181:CYS:HB3	2.55	0.40
2:B:58:PRO:HG2	2:B:62:ARG:HB3	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	131/159 (82%)	120 (92%)	11 (8%)	0	100	100
2	B	270/309 (87%)	251 (93%)	17 (6%)	2 (1%)	26	69
All	All	401/468 (86%)	371 (92%)	28 (7%)	2 (0%)	34	74

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	118	MET
2	B	298	LEU

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	129/145 (89%)	101 (78%)	28 (22%)	1	6
2	B	240/269 (89%)	189 (79%)	51 (21%)	1	6
All	All	369/414 (89%)	290 (79%)	79 (21%)	1	6

All (79) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	120	THR
1	A	126	LEU
1	A	127	SER
1	A	129	TYR

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Mol	Chain	Res	Type
1	A	133	SER
1	A	135	THR
1	A	154	GLU
1	A	155	LYS
1	A	159	VAL
1	A	161	LEU
1	A	162	SER
1	A	167	GLN
1	A	168	HIS
1	A	171	ASN
1	A	186	LEU
1	A	198	ASN
1	A	210	LYS
1	A	223	MSE
1	A	224	HIS
1	A	227	CYS
1	A	234	THR
1	A	235	ASP
1	A	243	LYS
1	A	261	GLU
1	A	262	ASN
1	A	264	LEU
1	A	266	LYS
1	A	268	SER
2	B	41	LYS
2	B	50	TYR
2	B	56	SER
2	B	57	ILE
2	B	60	GLN
2	B	63	ASN
2	B	64	ARG
2	B	72	LEU
2	B	86	THR
2	B	95	ASN
2	B	96	ARG
2	B	97	THR
2	B	99	TYR
2	B	110	ASP
2	B	115	ASP
2	B	118	MET
2	B	120	SER
2	B	123	SER

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Mol	Chain	Res	Type
2	B	128	ASN
2	B	145	LEU
2	B	151	CYS
2	B	154	LEU
2	B	158	ARG
2	B	169	ASP
2	B	173	THR
2	B	178	ASP
2	B	180	THR
2	B	208	PHE
2	B	220	ASN
2	B	221	GLU
2	B	223	LYS
2	B	233	LEU
2	B	243	THR
2	B	249	VAL
2	B	250	LEU
2	B	257	LYS
2	B	260	ASP
2	B	261	PHE
2	B	267	GLN
2	B	269	GLU
2	B	282	CYS
2	B	283	LEU
2	B	286	VAL
2	B	294	GLU
2	B	298	LEU
2	B	304	LEU
2	B	306	LEU
2	B	312	ARG
2	B	313	ARG
2	B	316	VAL
2	B	317	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	195	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

3 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	NAG	B	401	2	14,14,15	1.05	1 (7%)	15,19,21	1.31	3 (20%)
3	NAG	B	402	2	14,14,15	0.61	0	15,19,21	1.41	2 (13%)
3	NAG	B	403	2	14,14,15	0.50	0	15,19,21	1.18	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	NAG	B	401	2	-	0/6/23/26	0/1/1/1
3	NAG	B	402	2	-	0/6/23/26	0/1/1/1
3	NAG	B	403	2	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	401	NAG	C1-C2	3.06	1.56	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	NAG	O3-C3-C2	-2.37	104.41	109.11
3	B	401	NAG	C3-C4-C5	-2.21	106.35	110.20
3	B	401	NAG	O5-C5-C6	2.01	111.69	107.35
3	B	401	NAG	C2-N2-C7	2.44	126.18	123.04
3	B	402	NAG	C1-O5-C5	2.68	115.64	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	401	NAG	1	0
3	B	402	NAG	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	134/159 (84%)	-0.13	0	100 100	93, 123, 178, 216	0
2	B	278/309 (89%)	-0.05	5 (1%)	71 63	92, 127, 201, 238	0
All	All	412/468 (88%)	-0.08	5 (1%)	81 74	92, 126, 196, 238	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	297	LEU	6.1
2	B	204	ASP	3.3
2	B	142	THR	3.2
2	B	301	TYR	2.3
2	B	125	SER	2.3

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
3	NAG	B	402	14/15	0.94	0.27	0.04	110,110,110,110	0
3	NAG	B	401	14/15	0.85	0.29	-	110,110,110,110	0
3	NAG	B	403	14/15	0.82	0.25	-	110,110,110,110	0

6.5 Other polymers [i](#)

There are no such residues in this entry.