



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:31 PM UTC

PDB ID : 3NQ1 / pdb_00003nq1
Title : Crystal Structure of Tyrosinase from *Bacillus megaterium* in complex with inhibitor kojic acid
Authors : Sendovski, M.; Kanteev, M.; Adir, N.; Fishman, A.
Deposited on : 2010-06-29
Resolution : 2.30 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

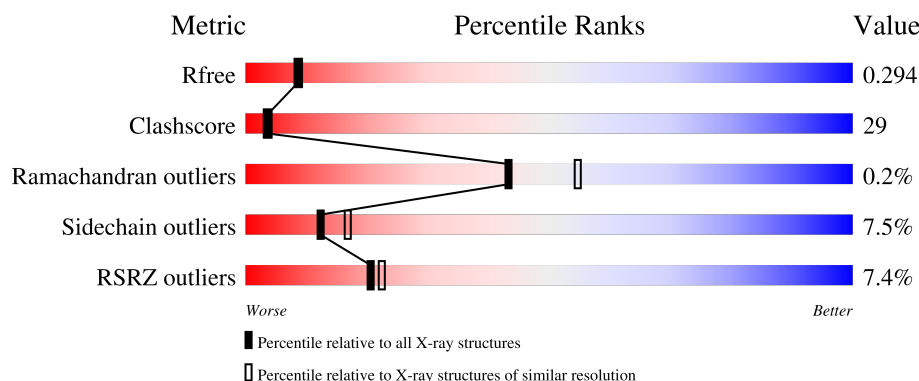
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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 of 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 $\leq 5\%$. 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	303	
1	B	303	

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
4	KOJ	A	1351	-	X	X	-
4	KOJ	B	1351	-	X	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4885 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 Tyrosinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	283	Total	C	N	O	S	0	0	0
			2321	1477	420	416	8			
1	B	283	Total	C	N	O	S	0	0	0
			2321	1477	420	416	8			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	298	HIS	-	expression tag	UNP B2ZB02
A	299	HIS	-	expression tag	UNP B2ZB02
A	300	HIS	-	expression tag	UNP B2ZB02
A	301	HIS	-	expression tag	UNP B2ZB02
A	302	HIS	-	expression tag	UNP B2ZB02
A	303	HIS	-	expression tag	UNP B2ZB02
B	298	HIS	-	expression tag	UNP B2ZB02
B	299	HIS	-	expression tag	UNP B2ZB02
B	300	HIS	-	expression tag	UNP B2ZB02
B	301	HIS	-	expression tag	UNP B2ZB02
B	302	HIS	-	expression tag	UNP B2ZB02
B	303	HIS	-	expression tag	UNP B2ZB02

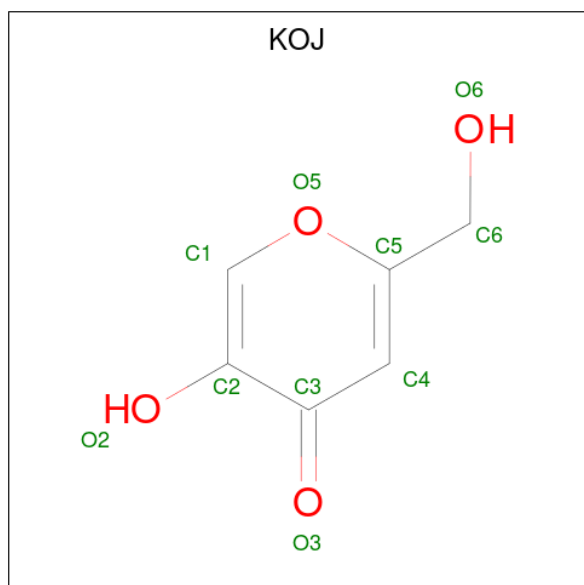
- Molecule 2 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Cu	0	0
			2	2		
2	B	2	Total	Cu	0	0
			2	2		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	8	Total	Zn	0	0
			8	8		
3	B	5	Total	Zn	0	0
			5	5		

- Molecule 4 is 5-HYDROXY-2-(HYDROXYMETHYL)-4H-PYRAN-4-ONE (CCD ID: KOJ) (formula: C₆H₆O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			10	6	4		
4	B	1	Total	C	O	0	0
			10	6	4		

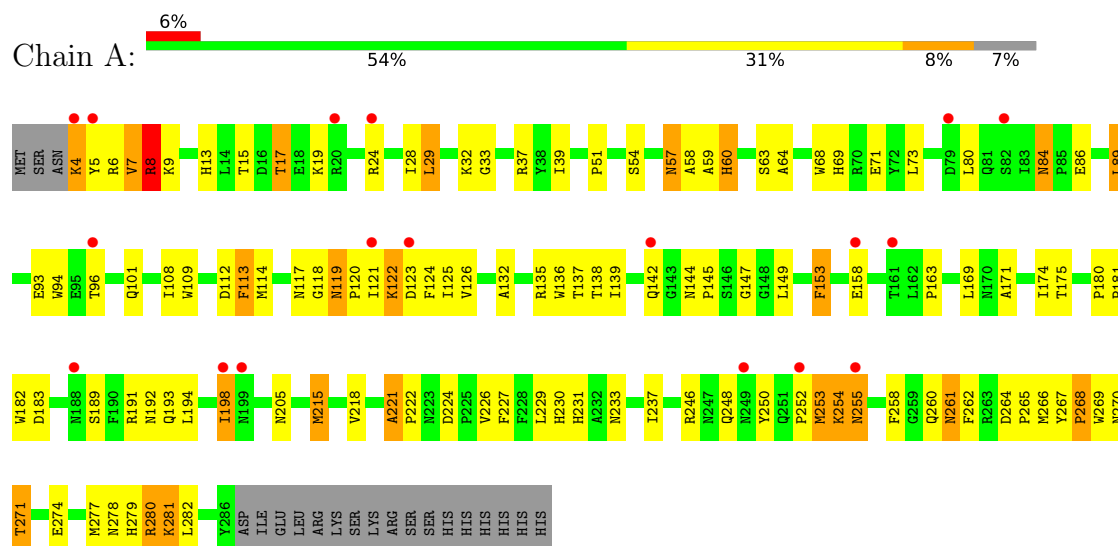
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	102	Total	O	0	0
			102	102		
5	B	104	Total	O	0	0
			104	104		

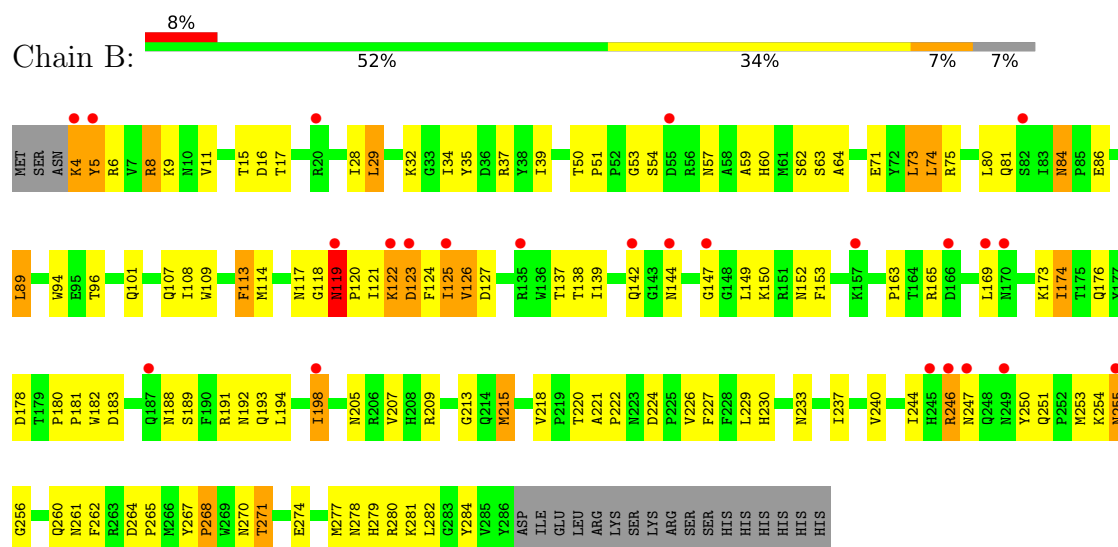
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Tyrosinase



• Molecule 1: Tyrosinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	51.38Å 83.53Å 146.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.09 – 2.30 42.09 – 2.30	Depositor EDS
% Data completeness (in resolution range)	77.1 (42.09-2.30) 83.8 (42.09-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 2.29Å)	Xtriage
Refinement program	REFMAC, PHENIX 1.6.1_357	Depositor
R, R_{free}	0.260 , 0.294 0.264 , 0.294	Depositor DCC
R_{free} test set	1250 reflections (4.35%)	wwPDB-VP
Wilson B-factor (Å ²)	16.4	Xtriage
Anisotropy	1.020	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 45.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	4885	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.32% 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, 0.333 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: ZN, CU, KOJ

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	0.87	13/2399 (0.5%)	1.06	12/3269 (0.4%)
1	B	0.82	11/2399 (0.5%)	1.06	11/3269 (0.3%)
All	All	0.85	24/4798 (0.5%)	1.06	23/6538 (0.4%)

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	137	THR	C-O	-8.05	1.13	1.23
1	B	75	ARG	C-O	-7.36	1.15	1.24
1	A	7	VAL	C-O	-7.28	1.16	1.24
1	B	215	MET	SD-CE	-6.84	1.62	1.79
1	A	215	MET	SD-CE	-6.84	1.62	1.79
1	B	74	LEU	C-O	-6.67	1.16	1.24
1	A	57	ASN	C-O	-6.63	1.15	1.23
1	A	139	ILE	C-O	-6.57	1.17	1.24
1	B	137	THR	C-O	-6.16	1.16	1.23
1	A	125	ILE	C-O	-6.08	1.17	1.24
1	B	15	THR	CA-CB	5.92	1.62	1.53
1	B	73	LEU	C-O	-5.84	1.17	1.24
1	B	247	ASN	C-O	-5.84	1.16	1.24
1	A	126	VAL	C-O	-5.76	1.18	1.24
1	B	138	THR	C-O	-5.74	1.16	1.23
1	A	33	GLY	C-O	-5.69	1.16	1.24
1	A	8	ARG	C-O	-5.38	1.17	1.24
1	B	122	LYS	C-O	-5.33	1.17	1.24
1	A	138	THR	C-O	-5.31	1.17	1.23
1	B	174	ILE	CG1-CD1	-5.14	1.31	1.51
1	B	173	LYS	C-O	-5.13	1.17	1.24
1	A	280	ARG	N-CA	-5.12	1.39	1.46
1	A	136	TRP	C-O	-5.10	1.18	1.23
1	A	58	ALA	C-O	-5.04	1.17	1.24

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	255	ASN	CB-CA-C	-11.95	91.75	111.23
1	B	5	TYR	N-CA-C	11.37	127.27	112.72
1	B	5	TYR	CB-CA-C	-8.93	95.72	110.37
1	B	113	PHE	N-CA-C	7.86	127.55	110.80
1	A	113	PHE	N-CA-C	7.50	126.77	110.80
1	A	261	ASN	CB-CA-C	-7.36	95.11	109.76
1	A	221	ALA	N-CA-C	7.30	125.94	109.81
1	A	57	ASN	N-CA-C	6.39	118.25	109.18
1	B	255	ASN	CB-CA-C	-6.31	97.87	110.42
1	B	63	SER	N-CA-C	6.11	118.72	111.33
1	A	63	SER	N-CA-C	6.09	118.70	111.33
1	B	113	PHE	CB-CA-C	-6.05	98.37	110.42
1	A	218	VAL	CB-CA-C	-5.87	108.12	113.70
1	B	218	VAL	CB-CA-C	-5.87	108.12	113.70
1	A	5	TYR	CB-CA-C	-5.77	100.07	110.23
1	B	119	ASN	CA-C-N	5.38	125.20	119.28
1	B	119	ASN	C-N-CA	5.38	125.20	119.28
1	B	9	LYS	N-CA-C	5.34	118.41	110.14
1	A	5	TYR	N-CA-C	5.25	120.16	113.18
1	A	153	PHE	N-CA-C	5.22	117.75	109.96
1	B	147	GLY	N-CA-C	5.17	121.93	115.21
1	A	9	LYS	N-CA-C	5.14	118.38	110.42
1	A	147	GLY	N-CA-C	5.13	121.88	115.21

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	2321	0	2206	133	1
1	B	2321	0	2206	135	1
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	8	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	5	0	0	0	0
4	A	10	0	4	6	0
4	B	10	0	4	6	0
5	A	102	0	0	16	0
5	B	104	0	0	7	0
All	All	4885	0	4420	261	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 29.

All (261) 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:205:ASN:HB3	4:B:1351:KOJ:H4	1.33	1.09
1:A:215:MET:HE3	1:A:221:ALA:O	1.55	1.07
1:A:215:MET:CE	1:A:221:ALA:O	2.03	1.05
1:A:261:ASN:HB2	1:A:264:ASP:OD2	1.58	1.03
1:B:256:GLY:H	1:B:261:ASN:HD21	1.06	0.97
1:B:122:LYS:HE2	1:B:127:ASP:HB3	1.47	0.96
1:A:205:ASN:OD1	4:A:1351:KOJ:H4	1.66	0.94
4:A:1351:KOJ:H1	5:A:335:HOH:O	1.67	0.93
1:A:261:ASN:HB2	1:A:264:ASP:CG	1.95	0.90
1:A:8:ARG:HG2	1:A:89:LEU:O	1.75	0.86
1:B:205:ASN:CB	4:B:1351:KOJ:H4	2.04	0.86
1:A:278:ASN:HD22	1:A:281:LYS:HD3	1.40	0.86
1:B:57:ASN:HD21	1:B:60:HIS:HD1	1.24	0.86
1:B:246:ARG:HB2	1:B:246:ARG:NH1	1.92	0.85
1:B:215:MET:HE1	1:B:227:PHE:HD2	1.44	0.82
1:B:209:ARG:HG2	4:B:1351:KOJ:H6C2	1.62	0.82
1:A:113:PHE:O	1:A:114:MET:HE2	1.80	0.81
1:B:108:ILE:HD11	1:B:229:LEU:HD23	1.63	0.81
1:A:215:MET:HE1	1:A:227:PHE:HD2	1.44	0.80
1:A:64:ALA:H	1:A:260:GLN:HE22	1.29	0.80
1:B:125:ILE:HG22	1:B:125:ILE:O	1.82	0.80
1:B:64:ALA:H	1:B:260:GLN:HE22	1.29	0.79
1:B:122:LYS:O	1:B:123:ASP:HB2	1.81	0.77
1:A:57:ASN:OD1	1:A:60:HIS:HD2	1.67	0.77
1:B:256:GLY:N	1:B:261:ASN:HD21	1.81	0.77
1:A:4:LYS:HG2	1:A:280:ARG:HD3	1.68	0.76
1:B:8:ARG:HG2	1:B:89:LEU:O	1.86	0.76
1:B:205:ASN:HB3	4:B:1351:KOJ:C4	2.15	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:LYS:HG3	1:A:280:ARG:HG2	1.69	0.74
1:B:122:LYS:HE2	1:B:127:ASP:CB	2.17	0.74
1:A:93:GLU:OE1	1:A:96:THR:HG21	1.88	0.74
1:A:205:ASN:CG	4:A:1351:KOJ:H4	2.13	0.73
1:A:119:ASN:C	1:A:119:ASN:HD22	1.95	0.73
1:A:254:LYS:O	1:A:255:ASN:HB2	1.86	0.73
1:B:11:VAL:HG11	1:B:108:ILE:HD12	1.69	0.73
1:B:246:ARG:HB2	1:B:246:ARG:CZ	2.17	0.73
1:A:64:ALA:H	1:A:260:GLN:NE2	1.86	0.72
1:B:278:ASN:HD22	1:B:281:LYS:HD3	1.53	0.72
1:B:122:LYS:CE	1:B:127:ASP:HB3	2.20	0.71
1:B:246:ARG:CZ	1:B:246:ARG:CB	2.67	0.71
1:B:113:PHE:O	1:B:114:MET:HE2	1.89	0.71
1:B:215:MET:HE1	1:B:227:PHE:CD2	2.25	0.71
1:B:119:ASN:C	1:B:119:ASN:HD22	1.95	0.71
1:A:119:ASN:HD22	1:A:120:PRO:CD	2.04	0.71
1:B:198:ILE:N	1:B:198:ILE:HD12	2.07	0.70
1:B:119:ASN:HD22	1:B:120:PRO:N	1.90	0.70
1:A:215:MET:HE1	1:A:227:PHE:CD2	2.25	0.69
1:A:13:HIS:HB3	5:A:304:HOH:O	1.92	0.69
1:A:198:ILE:HD12	1:A:198:ILE:N	2.07	0.69
1:B:193:GLN:HA	1:B:198:ILE:HD13	1.75	0.68
1:A:119:ASN:ND2	1:A:120:PRO:HD2	2.08	0.68
1:B:94:TRP:CE3	1:B:163:PRO:HG2	2.29	0.68
1:B:119:ASN:ND2	1:B:121:ILE:H	1.91	0.67
1:B:278:ASN:ND2	1:B:281:LYS:HD3	2.09	0.67
1:A:94:TRP:CE3	1:A:163:PRO:HG2	2.29	0.66
1:A:4:LYS:HG2	1:A:280:ARG:CD	2.26	0.66
1:A:37:ARG:NH2	1:B:270:ASN:ND2	2.44	0.66
1:A:37:ARG:NH2	1:B:270:ASN:HD22	1.94	0.66
1:A:193:GLN:HA	1:A:198:ILE:HD13	1.76	0.66
1:A:278:ASN:ND2	1:A:281:LYS:HD3	2.09	0.66
1:A:119:ASN:HD22	1:A:120:PRO:HD2	1.58	0.66
1:A:215:MET:CE	1:A:227:PHE:HB2	2.26	0.66
1:B:209:ARG:CG	4:B:1351:KOJ:H6C2	2.25	0.65
1:A:198:ILE:N	1:A:198:ILE:CD1	2.60	0.65
1:B:198:ILE:N	1:B:198:ILE:CD1	2.60	0.65
1:B:221:ALA:HB3	1:B:222:PRO:HD3	1.79	0.65
1:B:4:LYS:HG3	1:B:280:ARG:HG3	1.80	0.64
1:A:4:LYS:CG	1:A:280:ARG:CG	2.76	0.64
1:B:64:ALA:H	1:B:260:GLN:NE2	1.96	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:ARG:HB2	5:A:328:HOH:O	1.97	0.63
1:B:215:MET:CE	1:B:227:PHE:HB2	2.27	0.63
1:A:221:ALA:HB3	1:A:222:PRO:HD3	1.79	0.63
1:A:182:TRP:CE3	1:A:252:PRO:HD3	2.34	0.63
1:B:233:ASN:O	1:B:237:ILE:HG12	1.99	0.62
1:A:215:MET:HE2	1:A:221:ALA:O	1.99	0.62
1:A:193:GLN:HG2	1:A:198:ILE:HG12	1.81	0.62
1:A:233:ASN:O	1:A:237:ILE:HG12	1.99	0.62
1:A:123:ASP:OD1	1:A:123:ASP:O	2.18	0.61
1:B:5:TYR:HB2	5:B:356:HOH:O	1.99	0.61
1:B:57:ASN:ND2	1:B:60:HIS:H	1.98	0.61
1:B:215:MET:HE3	1:B:221:ALA:O	2.00	0.61
1:B:193:GLN:HG2	1:B:198:ILE:HG12	1.81	0.61
1:A:119:ASN:HD22	1:A:120:PRO:N	1.99	0.60
1:B:198:ILE:CD1	1:B:198:ILE:H	2.14	0.60
1:B:125:ILE:O	1:B:126:VAL:C	2.44	0.60
1:A:278:ASN:HD22	1:A:281:LYS:HB2	1.67	0.60
1:A:261:ASN:CB	1:A:264:ASP:OD2	2.41	0.60
1:B:176:GLN:NE2	1:B:251:GLN:HE22	2.01	0.59
1:A:198:ILE:CD1	1:A:198:ILE:H	2.14	0.59
1:A:57:ASN:OD1	1:A:60:HIS:CD2	2.53	0.59
1:B:215:MET:HA	1:B:215:MET:HE2	1.85	0.59
1:A:123:ASP:O	1:A:124:PHE:HB2	2.03	0.59
1:B:84:ASN:C	1:B:84:ASN:HD22	2.11	0.58
1:A:84:ASN:C	1:A:84:ASN:HD22	2.12	0.58
1:A:117:ASN:ND2	1:A:118:GLY:H	2.02	0.58
1:B:198:ILE:HG13	5:B:354:HOH:O	2.03	0.57
1:A:254:LYS:HA	1:A:261:ASN:ND2	2.18	0.57
1:A:215:MET:HE2	1:A:215:MET:HA	1.85	0.57
1:A:4:LYS:HG3	1:A:280:ARG:CG	2.34	0.57
1:A:6:ARG:NH2	1:A:282:LEU:O	2.37	0.57
1:B:117:ASN:ND2	1:B:118:GLY:H	2.02	0.56
1:A:119:ASN:C	1:A:119:ASN:ND2	2.64	0.56
1:B:119:ASN:CB	1:B:122:LYS:HD3	2.37	0.55
1:B:205:ASN:CG	4:B:1351:KOJ:H4	2.31	0.55
1:B:142:GLN:HE21	1:B:144:ASN:HD22	1.54	0.55
1:A:4:LYS:CG	1:A:280:ARG:HG2	2.36	0.55
1:B:119:ASN:C	1:B:119:ASN:ND2	2.64	0.55
1:B:119:ASN:HB3	1:B:122:LYS:HD3	1.89	0.55
1:A:253:MET:O	1:A:254:LYS:HG2	2.07	0.55
1:B:4:LYS:HG3	1:B:280:ARG:CG	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:6:ARG:HD2	1:B:81:GLN:OE1	2.07	0.54
1:A:142:GLN:HE21	1:A:144:ASN:HD22	1.54	0.54
1:B:35:TYR:CE2	1:B:39:ILE:HD11	2.42	0.54
1:A:215:MET:HE1	1:A:227:PHE:HB2	1.89	0.54
1:B:4:LYS:CG	1:B:280:ARG:HG3	2.37	0.54
1:A:15:THR:OG1	1:A:17:THR:HG23	2.08	0.54
1:A:94:TRP:HE1	1:A:230:HIS:HD2	1.56	0.54
1:A:181:PRO:HG2	1:A:183:ASP:OD2	2.08	0.54
1:B:181:PRO:HG2	1:B:183:ASP:OD2	2.08	0.54
1:A:253:MET:HA	1:A:262:PHE:HB2	1.90	0.54
1:A:253:MET:C	1:A:254:LYS:CG	2.80	0.54
1:A:142:GLN:HE21	1:A:144:ASN:ND2	2.07	0.53
1:B:261:ASN:HB2	1:B:264:ASP:OD2	2.08	0.53
1:A:224:ASP:CG	1:A:226:VAL:HG22	2.34	0.53
1:B:39:ILE:HD13	1:B:222:PRO:HB2	1.90	0.53
1:A:39:ILE:HD13	1:A:222:PRO:HB2	1.91	0.53
1:B:94:TRP:HE1	1:B:230:HIS:HD2	1.56	0.53
1:B:215:MET:HE1	1:B:227:PHE:HB2	1.89	0.53
1:B:4:LYS:CG	1:B:280:ARG:CG	2.87	0.53
1:A:84:ASN:HD22	1:A:86:GLU:H	1.56	0.53
1:B:240:VAL:O	1:B:244:ILE:HG12	2.09	0.53
1:A:253:MET:C	1:A:254:LYS:HG3	2.33	0.52
1:A:205:ASN:CG	4:A:1351:KOJ:H6C1	2.35	0.52
1:B:142:GLN:HE21	1:B:144:ASN:ND2	2.07	0.52
1:B:224:ASP:CG	1:B:226:VAL:HG22	2.34	0.52
1:A:119:ASN:ND2	1:A:121:ILE:H	2.07	0.52
1:B:84:ASN:ND2	1:B:86:GLU:H	2.08	0.52
1:B:246:ARG:NH1	1:B:246:ARG:CB	2.66	0.52
1:A:132:ALA:N	5:A:328:HOH:O	2.41	0.52
1:B:119:ASN:CG	1:B:122:LYS:HD3	2.35	0.52
1:A:267:TYR:CD1	1:A:268:PRO:HA	2.45	0.52
1:B:215:MET:CE	1:B:221:ALA:O	2.57	0.52
1:B:84:ASN:HD22	1:B:86:GLU:H	1.57	0.51
1:B:11:VAL:HG21	1:B:108:ILE:CD1	2.40	0.51
1:B:267:TYR:CD1	1:B:268:PRO:HA	2.45	0.51
1:B:57:ASN:ND2	1:B:60:HIS:HD1	2.00	0.51
1:B:57:ASN:HD22	1:B:59:ALA:H	1.58	0.51
1:A:37:ARG:HH21	1:B:270:ASN:ND2	2.08	0.51
1:A:171:ALA:O	1:A:174:ILE:HG12	2.11	0.51
1:A:122:LYS:O	1:A:123:ASP:HB3	2.11	0.51
1:B:119:ASN:OD1	1:B:122:LYS:HD3	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:254:LYS:O	1:B:255:ASN:HB2	2.12	0.50
1:A:101:GLN:HE21	1:A:101:GLN:HA	1.76	0.50
1:A:269:TRP:HB3	5:A:314:HOH:O	2.11	0.50
1:B:11:VAL:HG21	1:B:108:ILE:HD13	1.92	0.50
1:A:93:GLU:HG2	1:A:96:THR:HG23	1.92	0.50
1:A:224:ASP:OD1	1:A:226:VAL:HG22	2.12	0.50
1:B:101:GLN:HE21	1:B:101:GLN:HA	1.76	0.50
1:A:205:ASN:HB3	4:A:1351:KOJ:C4	2.42	0.49
1:A:93:GLU:OE1	1:A:96:THR:CG2	2.57	0.49
1:B:101:GLN:HA	1:B:101:GLN:NE2	2.27	0.49
1:B:224:ASP:OD1	1:B:226:VAL:HG22	2.12	0.49
1:A:101:GLN:HA	1:A:101:GLN:NE2	2.27	0.49
1:A:215:MET:HE3	1:A:227:PHE:HB2	1.95	0.49
1:A:262:PHE:CZ	1:A:274:GLU:HG3	2.48	0.49
1:A:265:PRO:HB3	1:A:270:ASN:HA	1.93	0.49
1:B:254:LYS:HD3	1:B:255:ASN:OD1	2.13	0.49
1:A:6:ARG:CZ	5:A:378:HOH:O	2.61	0.48
1:B:265:PRO:HB3	1:B:270:ASN:HA	1.93	0.48
1:B:117:ASN:HD22	1:B:118:GLY:H	1.61	0.48
1:A:253:MET:O	1:A:254:LYS:CG	2.61	0.48
1:B:62:SER:HB3	5:B:386:HOH:O	2.13	0.48
1:A:29:LEU:HD22	1:A:80:LEU:HD23	1.95	0.48
1:B:220:THR:HA	5:B:322:HOH:O	2.14	0.48
1:A:84:ASN:ND2	1:A:86:GLU:H	2.11	0.47
1:B:174:ILE:CD1	1:B:193:GLN:HG3	2.44	0.47
1:B:29:LEU:HD22	1:B:80:LEU:HD23	1.95	0.47
1:B:4:LYS:N	1:B:4:LYS:HD2	2.29	0.47
1:B:51:PRO:O	1:B:54:SER:HB3	2.15	0.47
1:A:191:ARG:C	1:A:191:ARG:HD2	2.40	0.47
1:A:144:ASN:HA	5:A:346:HOH:O	2.14	0.47
1:A:37:ARG:HH22	1:B:270:ASN:HD22	1.63	0.47
1:A:117:ASN:HD22	1:A:118:GLY:H	1.61	0.47
1:B:107:GLN:HG3	5:B:333:HOH:O	2.15	0.47
1:A:51:PRO:O	1:A:54:SER:HB3	2.14	0.47
1:B:191:ARG:HD2	1:B:191:ARG:C	2.40	0.46
1:B:124:PHE:CE2	1:B:152:ASN:ND2	2.83	0.46
1:B:178:ASP:HA	1:B:189:SER:OG	2.16	0.46
1:B:215:MET:HE3	1:B:227:PHE:HB2	1.95	0.46
1:A:8:ARG:CG	1:A:89:LEU:O	2.55	0.46
1:A:231:HIS:ND1	5:A:355:HOH:O	2.31	0.46
1:B:35:TYR:HE2	1:B:39:ILE:HD11	1.77	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:256:GLY:H	1:B:261:ASN:ND2	1.91	0.46
1:B:11:VAL:HG11	1:B:108:ILE:CD1	2.40	0.46
1:A:270:ASN:HD22	1:B:37:ARG:NH2	2.14	0.45
1:B:119:ASN:HA	1:B:120:PRO:HD2	1.86	0.45
1:A:135:ARG:NE	5:A:328:HOH:O	2.49	0.45
1:A:270:ASN:ND2	1:B:37:ARG:NH2	2.64	0.45
1:B:123:ASP:O	1:B:124:PHE:HB2	2.16	0.45
1:B:188:ASN:HB3	5:B:397:HOH:O	2.16	0.45
1:A:69:HIS:O	1:A:73:LEU:HG	2.17	0.45
1:A:117:ASN:HD21	1:A:153:PHE:H	1.65	0.44
1:A:271:THR:HB	5:A:364:HOH:O	2.17	0.44
1:B:180:PRO:HA	1:B:182:TRP:N	2.33	0.44
1:A:182:TRP:CZ2	1:A:252:PRO:HB3	2.53	0.44
1:B:279:HIS:HB2	1:B:284:TYR:CZ	2.52	0.44
1:A:59:ALA:HB2	1:A:68:TRP:CD2	2.53	0.43
1:A:253:MET:HA	1:A:262:PHE:CB	2.47	0.43
1:A:253:MET:HB3	1:A:253:MET:HE3	1.80	0.43
1:B:28:ILE:CG2	1:B:32:LYS:HE2	2.48	0.43
1:B:262:PHE:CZ	1:B:274:GLU:HG3	2.52	0.43
1:B:278:ASN:HD22	1:B:281:LYS:HB2	1.83	0.43
1:B:279:HIS:HA	1:B:282:LEU:HD12	1.99	0.43
1:A:180:PRO:HA	1:A:182:TRP:N	2.33	0.43
1:A:231:HIS:HB3	5:A:355:HOH:O	2.17	0.43
1:B:117:ASN:HD21	1:B:153:PHE:H	1.65	0.43
1:B:250:TYR:HB2	1:B:277:MET:CE	2.49	0.43
1:A:258:PHE:CZ	1:B:34:ILE:CD1	3.02	0.43
1:A:250:TYR:HB2	1:A:277:MET:CE	2.49	0.43
4:A:1351:KOJ:C6	5:A:356:HOH:O	2.66	0.43
1:B:246:ARG:HH11	1:B:246:ARG:CG	2.32	0.43
1:A:246:ARG:HG3	1:A:246:ARG:HH11	1.84	0.43
1:A:145:PRO:HD3	1:B:53:GLY:N	2.34	0.43
1:B:253:MET:HE1	1:B:262:PHE:CD2	2.54	0.43
1:A:28:ILE:CG2	1:A:32:LYS:HE2	2.48	0.43
1:A:123:ASP:O	1:A:123:ASP:CG	2.62	0.43
1:A:180:PRO:HA	1:A:182:TRP:H	1.83	0.43
1:A:254:LYS:CA	1:A:261:ASN:ND2	2.81	0.42
1:A:278:ASN:HD22	1:A:281:LYS:CD	2.21	0.42
1:B:50:THR:HA	1:B:51:PRO:HA	1.89	0.42
1:B:96:THR:HG22	1:B:165:ARG:CZ	2.49	0.42
1:B:119:ASN:HB3	1:B:122:LYS:HB2	2.00	0.42
1:B:124:PHE:CE2	1:B:152:ASN:CG	2.98	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:144:ASN:HA	5:B:323:HOH:O	2.19	0.42
1:B:109:TRP:CZ3	1:B:226:VAL:HG11	2.55	0.42
1:B:215:MET:HE2	1:B:221:ALA:HB1	2.01	0.42
1:A:109:TRP:CZ3	1:A:226:VAL:HG11	2.55	0.42
1:A:278:ASN:ND2	1:A:281:LYS:HB2	2.33	0.42
1:A:279:HIS:HA	1:A:282:LEU:HD12	2.00	0.42
1:B:96:THR:HG22	1:B:165:ARG:NH1	2.35	0.42
1:B:180:PRO:HA	1:B:182:TRP:H	1.84	0.42
1:A:19:LYS:HG2	1:A:113:PHE:CE1	2.55	0.41
1:A:71:GLU:HG2	1:A:271:THR:HG21	2.03	0.41
1:A:158:GLU:HB2	5:A:391:HOH:O	2.20	0.41
1:A:101:GLN:HB3	5:A:372:HOH:O	2.20	0.41
1:B:192:ASN:C	1:B:198:ILE:CD1	2.94	0.41
1:A:258:PHE:HZ	1:B:34:ILE:HD11	1.85	0.41
1:B:6:ARG:HH22	1:B:74:LEU:CD1	2.34	0.41
1:B:207:VAL:HB	1:B:230:HIS:CE1	2.56	0.41
1:A:175:THR:O	1:A:248:GLN:HG3	2.21	0.41
1:A:215:MET:CE	1:A:221:ALA:C	2.87	0.41
1:A:215:MET:HE2	1:A:221:ALA:HB1	2.01	0.41
1:B:73:LEU:HD22	1:B:89:LEU:HD11	2.02	0.41
1:A:24:ARG:HG3	5:A:383:HOH:O	2.21	0.41
1:A:108:ILE:HD11	1:A:229:LEU:HD23	2.02	0.41
1:A:192:ASN:C	1:A:198:ILE:CD1	2.94	0.41
1:B:150:LYS:O	1:B:213:GLY:HA3	2.21	0.41
1:A:39:ILE:CD1	1:A:222:PRO:HB2	2.51	0.40
1:A:266:MET:O	1:A:267:TYR:C	2.64	0.40
1:B:71:GLU:HG2	1:B:271:THR:HG21	2.03	0.40
1:A:13:HIS:CB	5:A:304:HOH:O	2.59	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:ASP:OD2	1:B:16:ASP:OD1[2_455]	2.16	0.04

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	281/303 (93%)	272 (97%)	9 (3%)	0	100	100
1	B	281/303 (93%)	270 (96%)	10 (4%)	1 (0%)	30	38
All	All	562/606 (93%)	542 (96%)	19 (3%)	1 (0%)	43	55

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	126	VAL

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	246/266 (92%)	226 (92%)	20 (8%)	11	14
1	B	246/266 (92%)	229 (93%)	17 (7%)	14	20
All	All	492/532 (92%)	455 (92%)	37 (8%)	12	17

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LYS
1	A	7	VAL
1	A	8	ARG
1	A	17	THR
1	A	29	LEU

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Mol	Chain	Res	Type
1	A	60	HIS
1	A	84	ASN
1	A	89	LEU
1	A	119	ASN
1	A	122	LYS
1	A	149	LEU
1	A	169	LEU
1	A	189	SER
1	A	194	LEU
1	A	198	ILE
1	A	253	MET
1	A	254	LYS
1	A	268	PRO
1	A	271	THR
1	A	281	LYS
1	B	4	LYS
1	B	8	ARG
1	B	17	THR
1	B	29	LEU
1	B	84	ASN
1	B	89	LEU
1	B	119	ASN
1	B	123	ASP
1	B	125	ILE
1	B	139	ILE
1	B	149	LEU
1	B	169	LEU
1	B	194	LEU
1	B	198	ILE
1	B	246	ARG
1	B	268	PRO
1	B	271	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	HIS
1	A	84	ASN
1	A	101	GLN
1	A	117	ASN
1	A	119	ASN
1	A	142	GLN

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Mol	Chain	Res	Type
1	A	144	ASN
1	A	152	ASN
1	A	176	GLN
1	A	187	GLN
1	A	202	GLN
1	A	205	ASN
1	A	230	HIS
1	A	248	GLN
1	A	255	ASN
1	A	260	GLN
1	A	261	ASN
1	A	270	ASN
1	A	278	ASN
1	B	57	ASN
1	B	84	ASN
1	B	101	GLN
1	B	117	ASN
1	B	119	ASN
1	B	142	GLN
1	B	144	ASN
1	B	152	ASN
1	B	176	GLN
1	B	202	GLN
1	B	205	ASN
1	B	230	HIS
1	B	251	GLN
1	B	260	GLN
1	B	261	ASN
1	B	270	ASN
1	B	278	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 17 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
4	KOJ	A	1351	-	10,10,10	3.21	6 (60%)	12,13,13	2.79	6 (50%)
4	KOJ	B	1351	-	10,10,10	3.21	6 (60%)	12,13,13	2.77	6 (50%)

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
4	KOJ	A	1351	-	-	1/2/2/2	0/1/1/1
4	KOJ	B	1351	-	-	1/2/2/2	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1351	KOJ	O5-C5	-5.25	1.25	1.35
4	B	1351	KOJ	O5-C5	-5.16	1.25	1.35
4	B	1351	KOJ	C1-C2	4.88	1.38	1.34
4	A	1351	KOJ	C1-C2	4.80	1.38	1.34
4	B	1351	KOJ	C2-C3	-4.02	1.35	1.47
4	A	1351	KOJ	C2-C3	-3.99	1.35	1.47
4	A	1351	KOJ	O5-C1	3.71	1.44	1.37
4	B	1351	KOJ	O5-C1	3.71	1.44	1.37
4	B	1351	KOJ	O3-C3	3.02	1.32	1.24
4	A	1351	KOJ	O3-C3	3.00	1.32	1.24
4	A	1351	KOJ	O6-C6	-2.58	1.33	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1351	KOJ	O6-C6	-2.58	1.33	1.41

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1351	KOJ	O5-C1-C2	-5.72	119.26	122.75
4	B	1351	KOJ	O5-C1-C2	-5.55	119.36	122.75
4	B	1351	KOJ	O5-C5-C6	4.42	118.26	111.06
4	A	1351	KOJ	O5-C5-C6	4.39	118.20	111.06
4	A	1351	KOJ	C6-C5-C4	-3.51	121.43	126.78
4	B	1351	KOJ	C6-C5-C4	-3.48	121.47	126.78
4	A	1351	KOJ	O6-C6-C5	2.72	118.77	111.24
4	B	1351	KOJ	O6-C6-C5	2.69	118.69	111.24
4	A	1351	KOJ	O2-C2-C1	2.58	123.78	120.15
4	B	1351	KOJ	O2-C2-C1	2.56	123.75	120.15
4	A	1351	KOJ	C1-O5-C5	2.26	121.67	119.33
4	B	1351	KOJ	C1-O5-C5	2.25	121.66	119.33

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1351	KOJ	C4-C5-C6-O6
4	B	1351	KOJ	C4-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1351	KOJ	6	0
4	B	1351	KOJ	6	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	283/303 (93%)	0.70	18 (6%) 25 27	8, 18, 34, 52	0
1	B	283/303 (93%)	0.79	24 (8%) 16 18	11, 18, 38, 64	0
All	All	566/606 (93%)	0.74	42 (7%) 20 22	8, 18, 37, 64	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	5	TYR	6.1
1	B	5	TYR	4.6
1	A	20	ARG	4.3
1	B	255	ASN	3.8
1	B	119	ASN	3.4
1	A	24	ARG	3.2
1	B	142	GLN	3.2
1	B	147	GLY	3.1
1	A	142	GLN	3.1
1	B	157	LYS	3.1
1	B	125	ILE	3.0
1	A	4	LYS	2.9
1	B	4	LYS	2.8
1	A	82	SER	2.7
1	A	96	THR	2.7
1	A	255	ASN	2.7
1	A	199	ASN	2.7
1	B	249	ASN	2.6
1	B	82	SER	2.6
1	B	246	ARG	2.6
1	B	166	ASP	2.6
1	A	158	GLU	2.5
1	B	20	ARG	2.5
1	B	55	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	198	ILE	2.4
1	A	161	THR	2.4
1	A	121	ILE	2.4
1	A	249	ASN	2.4
1	B	123	ASP	2.4
1	B	247	ASN	2.3
1	A	123	ASP	2.2
1	A	188	ASN	2.2
1	A	252	PRO	2.2
1	B	169	LEU	2.2
1	A	79	ASP	2.1
1	B	144	ASN	2.1
1	A	198	ILE	2.1
1	B	187	GLN	2.1
1	B	170	ASN	2.0
1	B	135	ARG	2.0
1	B	122	LYS	2.0
1	B	245	HIS	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 oligosaccharides 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, 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	B-factors($\text{\AA}^2$)	Q<0.9
3	ZN	B	506	1/1	0.68	0.19	77,77,77,77	0
3	ZN	A	508	1/1	0.76	0.12	64,64,64,64	0
3	ZN	A	506	1/1	0.76	0.11	27,27,27,27	0
4	KOJ	B	1351	10/10	0.77	0.18	14,32,36,43	0
4	KOJ	A	1351	10/10	0.82	0.21	15,30,36,47	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ZN	B	504	1/1	0.83	0.11	60,60,60,60	0
3	ZN	B	505	1/1	0.86	0.09	30,30,30,30	0
3	ZN	A	505	1/1	0.86	0.10	42,42,42,42	0
3	ZN	A	510	1/1	0.87	0.10	43,43,43,43	0
3	ZN	A	503	1/1	0.89	0.07	22,22,22,22	0
3	ZN	A	504	1/1	0.90	0.08	57,57,57,57	0
3	ZN	B	503	1/1	0.93	0.07	18,18,18,18	0
3	ZN	A	507	1/1	0.93	0.08	48,48,48,48	0
2	CU	B	501	1/1	0.97	0.03	14,14,14,14	0
3	ZN	B	507	1/1	0.97	0.05	34,34,34,34	0
3	ZN	A	509	1/1	0.97	0.05	41,41,41,41	0
2	CU	A	501	1/1	0.97	0.06	22,22,22,22	0
2	CU	B	502	1/1	0.99	0.02	16,16,16,16	0
2	CU	A	502	1/1	1.00	0.02	14,14,14,14	0

6.5 Other polymers [i](#)

There are no such residues in this entry.