



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 09:47 pm GMT

PDB ID : 1BT2  
Title : CATECHOL OXIDASE FROM IPOMOEA BATATAS (SWEET POTATOES) IN THE REDUCED CU(I)-CU(I) STATE  
Authors : Klabunde, T.; Eicken, C.; Sacchettini, J.C.; Krebs, B.  
Deposited on : 1998-09-02  
Resolution : 2.70 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

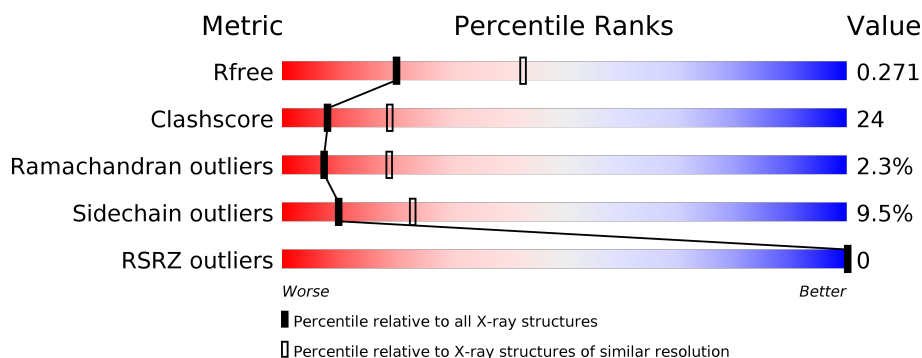
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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))
$R_{free}$	100719	2259 (2.70-2.70)
Clashscore	112137	2590 (2.70-2.70)
Ramachandran outliers	110173	2550 (2.70-2.70)
Sidechain outliers	110143	2550 (2.70-2.70)
RSRZ outliers	101464	2275 (2.70-2.70)

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  $\geq 3$ , 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	345	
1	B	345	

## 2 Entry composition [i](#)

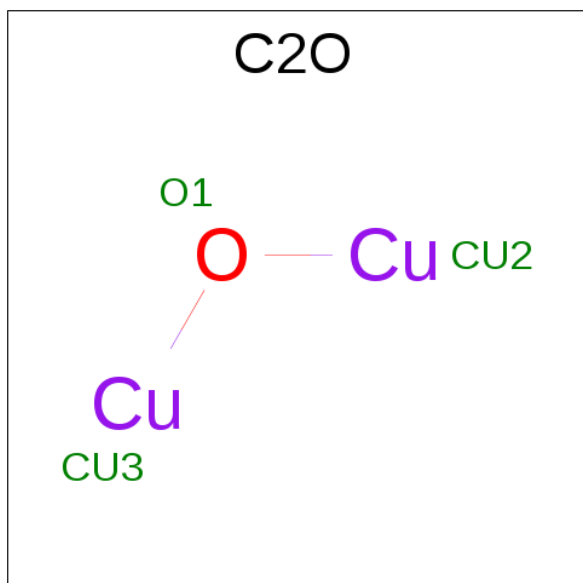
There are 3 unique types of molecules in this entry. The entry contains 5512 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PROTEIN (CATECHOL OXIDASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	336	Total	C	N	O	S	0	0	0
			2666	1703	452	496	15			
1	B	336	Total	C	N	O	S	0	0	0
			2666	1703	452	496	15			

- Molecule 2 is CU-O-CU LINKAGE (three-letter code: C2O) (formula:  $\text{Cu}_2\text{O}$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Cu	O	0	0
			3	2	1		
2	B	1	Total	Cu	O	0	0
			3	2	1		

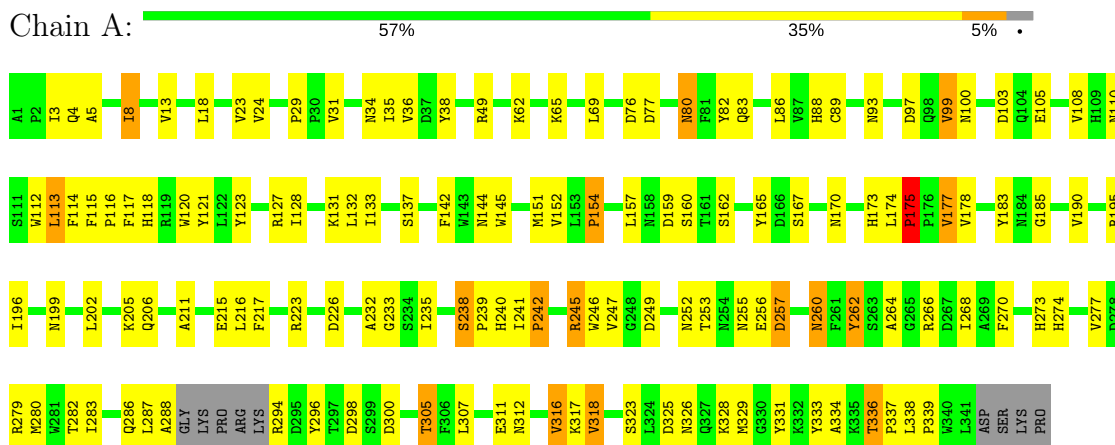
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	85	Total 85	O 85	0	0
3	B	89	Total 89	O 89	0	0

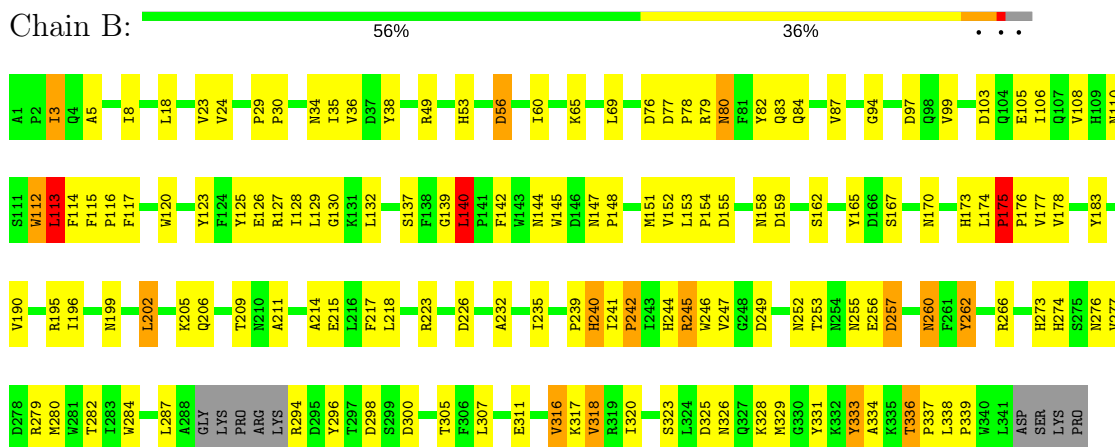
### 3 Residue-property plots

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



#### • Molecule 1: PROTEIN (CATECHOL OXIDASE)



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	45.25Å 162.60Å 51.49Å 90.00° 95.30° 90.00°	Depositor
Resolution (Å)	8.00 – 2.70 7.99 – 2.70	Depositor EDS
% Data completeness (in resolution range)	88.3 (8.00-2.70) 88.3 (7.99-2.70)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	10.23 (at 2.70Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, $R_{free}$	0.184 , 0.278 0.184 , 0.271	Depositor DCC
$R_{free}$ test set	847 reflections (4.92%)	DCC
Wilson B-factor (Å <sup>2</sup> )	27.8	Xtriage
Anisotropy	0.241	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 62.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	5512	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.50% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: C2O

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	0.69	0/2750	0.89	2/3758 (0.1%)
1	B	0.65	0/2750	0.88	2/3758 (0.1%)
All	All	0.67	0/5500	0.88	4/7516 (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
1	A	0	1
1	B	0	2
All	All	0	3

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	113	LEU	CA-CB-CG	6.30	129.80	115.30
1	B	140	LEU	CA-CB-CG	6.05	129.22	115.30
1	A	238	SER	N-CA-C	5.09	124.73	111.00
1	B	113	LEU	CA-CB-CG	5.05	126.91	115.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	331	TYR	Sidechain
1	B	331	TYR	Sidechain

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Mol	Chain	Res	Type	Group
1	B	333	TYR	Sidechain

## 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	2666	0	2525	122	0
1	B	2666	0	2525	124	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	85	0	0	12	0
3	B	89	0	0	10	0
All	All	5512	0	5050	246	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 24.

All (246) 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:337:PRO:O	1:A:339:PRO:HD3	1.81	0.81
1:A:131:LYS:HE3	1:A:312:ASN:ND2	1.97	0.80
1:B:223:ARG:NH1	1:B:300:ASP:HB2	1.98	0.79
1:B:145:TRP:HE1	1:B:273:HIS:HD2	1.29	0.79
1:B:174:LEU:HB3	1:B:175:PRO:HD2	1.63	0.78
1:B:337:PRO:O	1:B:339:PRO:HD3	1.82	0.78
1:A:88:HIS:HB3	3:A:517:HOH:O	1.83	0.77
1:B:80:ASN:HD22	1:B:83:GLN:H	1.29	0.77
1:A:145:TRP:HE1	1:A:273:HIS:HD2	1.32	0.76
1:B:8:ILE:HG23	1:B:29:PRO:HG2	1.72	0.71
1:A:174:LEU:HB3	1:A:175:PRO:HD2	1.72	0.71
1:A:131:LYS:HE3	1:A:312:ASN:HD21	1.56	0.70
1:A:223:ARG:NH1	1:A:300:ASP:HB2	2.08	0.69
1:A:120:TRP:NE1	1:A:323:SER:HB3	2.07	0.69
1:A:62:LYS:HG2	1:A:133:ILE:HD12	1.75	0.68
1:B:151:MET:HE1	1:B:196:ILE:HA	1.75	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:145:TRP:NE1	1:B:273:HIS:HD2	1.92	0.68
1:B:152:VAL:HG13	1:B:175:PRO:O	1.93	0.67
1:A:152:VAL:HG13	1:A:175:PRO:O	1.94	0.67
1:B:174:LEU:O	1:B:177:VAL:HG12	1.95	0.66
1:A:100:ASN:N	3:A:560:HOH:O	2.27	0.66
1:A:145:TRP:NE1	1:A:273:HIS:HD2	1.94	0.66
1:A:279:ARG:HD2	3:A:529:HOH:O	1.97	0.65
1:A:185:GLY:HA2	3:A:567:HOH:O	1.96	0.65
1:B:108:VAL:HG11	1:B:117:PHE:CD2	2.31	0.65
1:A:246:TRP:HH2	3:A:523:HOH:O	1.78	0.64
1:B:151:MET:CE	1:B:196:ILE:HA	2.28	0.63
1:A:80:ASN:HD22	1:A:83:GLN:H	1.44	0.63
1:A:123:TYR:O	1:A:127:ARG:HG3	1.98	0.62
1:B:279:ARG:HG3	1:B:333:TYR:CG	2.34	0.62
1:B:120:TRP:CE2	1:B:318:VAL:HG13	2.35	0.62
1:B:214:ALA:O	1:B:218:LEU:HB2	1.99	0.62
1:A:245:ARG:HB3	1:A:245:ARG:HH11	1.65	0.61
1:B:260:ASN:HD22	1:B:262:TYR:H	1.48	0.61
1:A:151:MET:CE	1:A:196:ILE:HA	2.29	0.61
1:A:178:VAL:HG22	1:A:195:ARG:HD2	1.83	0.61
1:A:65:LYS:CE	1:A:69:LEU:HD21	2.31	0.61
1:A:142:PHE:HZ	1:A:336:THR:HG21	1.66	0.61
1:A:288:ALA:HB2	1:A:294:ARG:HD2	1.82	0.60
1:B:108:VAL:HG11	1:B:117:PHE:CG	2.36	0.60
1:B:284:TRP:HA	1:B:287:LEU:HD12	1.81	0.60
1:B:142:PHE:HZ	1:B:336:THR:HG21	1.65	0.60
1:B:211:ALA:HB1	1:B:217:PHE:HB2	1.81	0.60
1:B:36:VAL:HG23	1:B:316:VAL:HG23	1.84	0.59
1:B:56:ASP:O	1:B:60:ILE:HG13	2.03	0.59
1:B:130:GLY:HA3	3:B:582:HOH:O	2.02	0.59
1:B:49:ARG:HD2	1:B:333:TYR:CZ	2.38	0.59
1:A:38:TYR:HB2	1:A:316:VAL:HG22	1.84	0.59
1:A:211:ALA:HB1	1:A:217:PHE:HB2	1.84	0.59
1:A:151:MET:HE1	1:A:196:ILE:HA	1.84	0.58
1:B:128:ILE:HG12	1:B:311:GLU:HA	1.84	0.58
1:B:151:MET:HE3	1:B:199:ASN:HD22	1.68	0.58
1:B:178:VAL:HG22	1:B:195:ARG:HD2	1.85	0.58
1:B:206:GLN:HG3	1:B:239:PRO:HD3	1.84	0.58
1:B:246:TRP:HH2	3:B:527:HOH:O	1.86	0.58
1:A:215:GLU:HA	1:A:296:TYR:OH	2.03	0.58
1:A:38:TYR:HD1	1:A:316:VAL:HG21	1.69	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:77:ASP:O	1:B:83:GLN:HG3	2.03	0.58
1:B:245:ARG:HB3	1:B:245:ARG:HH11	1.68	0.57
1:B:257:ASP:OD1	1:B:266:ARG:HB2	2.04	0.57
1:B:334:ALA:O	1:B:336:THR:HG22	2.04	0.57
1:A:49:ARG:HD2	1:A:333:TYR:CZ	2.39	0.57
1:B:120:TRP:NE1	1:B:323:SER:HB3	2.18	0.57
1:B:158:ASN:HA	3:B:510:HOH:O	2.04	0.57
1:B:38:TYR:HD1	1:B:316:VAL:HG21	1.69	0.57
1:B:223:ARG:HH12	1:B:300:ASP:HB2	1.66	0.57
1:A:174:LEU:O	1:A:177:VAL:HG12	2.04	0.57
1:A:97:ASP:HA	1:A:105:GLU:HA	1.87	0.56
1:A:65:LYS:O	1:A:69:LEU:HD23	2.05	0.56
1:A:120:TRP:CE2	1:A:318:VAL:HG13	2.40	0.56
1:B:65:LYS:CE	1:B:69:LEU:HD21	2.34	0.56
1:A:260:ASN:ND2	1:A:262:TYR:HD2	2.04	0.56
1:B:23:VAL:HG23	1:B:24:VAL:HG13	1.86	0.56
1:A:5:ALA:H	1:A:305:THR:CG2	2.19	0.56
1:A:5:ALA:H	1:A:305:THR:HG23	1.71	0.55
1:A:108:VAL:HG11	1:A:117:PHE:CG	2.40	0.55
1:B:305:THR:HG21	1:B:317:LYS:HE2	1.87	0.55
1:A:159:ASP:O	1:A:165:TYR:HB2	2.06	0.55
1:A:249:ASP:HB3	1:A:252:ASN:ND2	2.21	0.55
1:A:38:TYR:CD1	1:A:316:VAL:HG21	2.41	0.55
1:B:159:ASP:O	1:B:165:TYR:HB2	2.07	0.55
1:A:35:ILE:HD13	1:A:307:LEU:HD11	1.88	0.55
1:A:8:ILE:HG23	1:A:29:PRO:HG2	1.89	0.55
1:B:8:ILE:HD12	1:B:29:PRO:HG3	1.89	0.55
1:B:80:ASN:ND2	1:B:83:GLN:H	2.02	0.55
1:A:151:MET:HE3	1:A:199:ASN:HD22	1.72	0.54
1:A:65:LYS:HE3	1:A:69:LEU:HD21	1.88	0.54
1:B:127:ARG:HD3	1:B:311:GLU:OE1	2.07	0.54
1:B:49:ARG:NH2	3:B:562:HOH:O	2.39	0.54
1:B:260:ASN:ND2	1:B:262:TYR:H	2.05	0.54
1:B:173:HIS:ND1	1:B:246:TRP:O	2.40	0.54
1:B:142:PHE:CD2	1:B:333:TYR:HD2	2.25	0.54
1:A:257:ASP:OD1	1:A:266:ARG:HB2	2.07	0.53
1:B:65:LYS:O	1:B:69:LEU:HD23	2.08	0.53
1:A:279:ARG:HG3	1:A:333:TYR:CG	2.43	0.53
1:B:8:ILE:HD11	1:B:307:LEU:HD22	1.90	0.53
1:B:277:VAL:O	1:B:280:MET:HB2	2.08	0.53
1:B:69:LEU:HB3	1:B:132:LEU:HD13	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:VAL:HG23	1:A:316:VAL:HG23	1.92	0.52
1:B:223:ARG:O	1:B:226:ASP:HB2	2.08	0.52
1:A:305:THR:HG21	1:A:317:LYS:HE2	1.89	0.52
1:A:80:ASN:ND2	1:A:83:GLN:H	2.08	0.52
1:B:202:LEU:CD1	1:B:206:GLN:HG2	2.39	0.52
1:B:249:ASP:HB3	1:B:252:ASN:ND2	2.25	0.52
1:A:108:VAL:HG11	1:A:117:PHE:CD2	2.44	0.52
1:A:206:GLN:HG3	1:A:239:PRO:HD3	1.90	0.52
1:A:89:CYS:O	1:A:93:ASN:HB2	2.10	0.52
1:B:125:TYR:HD2	1:B:140:LEU:HD21	1.74	0.52
1:B:279:ARG:HD2	3:B:532:HOH:O	2.10	0.52
1:B:35:ILE:HD13	1:B:307:LEU:HD11	1.93	0.51
1:B:38:TYR:CD1	1:B:316:VAL:HG21	2.45	0.51
1:A:205:LYS:HG2	1:A:206:GLN:NE2	2.24	0.51
1:A:274:HIS:HA	1:A:277:VAL:HB	1.93	0.51
1:A:65:LYS:NZ	1:A:69:LEU:HD21	2.26	0.51
1:B:276:ASN:ND2	1:B:279:ARG:HE	2.09	0.51
1:B:215:GLU:HA	1:B:296:TYR:OH	2.10	0.51
1:A:264:ALA:HB1	1:A:270:PHE:CD1	2.47	0.50
1:A:260:ASN:HD22	1:A:262:TYR:H	1.59	0.50
1:B:110:ASN:HA	1:B:232:ALA:HA	1.94	0.50
1:A:249:ASP:HB3	1:A:252:ASN:HD22	1.76	0.50
1:B:38:TYR:HB2	1:B:316:VAL:HG22	1.94	0.50
1:A:8:ILE:HD12	1:A:29:PRO:HG3	1.93	0.50
1:B:5:ALA:H	1:B:305:THR:HG23	1.77	0.50
1:A:170:ASN:ND2	1:A:173:HIS:CE1	2.80	0.49
1:B:65:LYS:NZ	1:B:69:LEU:HD21	2.27	0.49
1:A:23:VAL:HG23	1:A:24:VAL:HG13	1.93	0.49
1:A:247:VAL:O	1:A:247:VAL:HG12	2.13	0.49
1:B:120:TRP:CD2	1:B:318:VAL:CG1	2.96	0.49
1:A:283:ILE:O	1:A:286:GLN:HB3	2.13	0.49
1:B:129:LEU:CD1	1:B:140:LEU:HD22	2.43	0.49
1:A:173:HIS:HB3	1:A:246:TRP:CZ2	2.47	0.48
1:A:334:ALA:O	1:A:336:THR:HG22	2.13	0.48
1:B:5:ALA:HB1	1:B:307:LEU:HB2	1.95	0.48
1:B:147:ASN:O	1:B:151:MET:HG3	2.13	0.48
1:A:110:ASN:HA	1:A:232:ALA:HA	1.96	0.48
1:A:8:ILE:HD12	1:A:29:PRO:CG	2.43	0.48
1:A:38:TYR:HB2	1:A:316:VAL:CG2	2.43	0.48
1:B:260:ASN:ND2	1:B:262:TYR:HD2	2.11	0.48
1:B:5:ALA:H	1:B:305:THR:CG2	2.27	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:202:LEU:HD12	1:B:206:GLN:HG2	1.94	0.48
1:A:8:ILE:CG2	1:A:31:VAL:HA	2.43	0.48
1:B:241:ILE:O	1:B:244:HIS:HB2	2.14	0.47
1:B:247:VAL:HG12	1:B:247:VAL:O	2.14	0.47
1:A:326:ASN:ND2	3:A:572:HOH:O	2.44	0.47
1:A:260:ASN:ND2	1:A:262:TYR:H	2.12	0.47
1:A:223:ARG:HH12	1:A:300:ASP:HB2	1.77	0.47
1:A:325:ASP:HB3	1:A:328:LYS:HB2	1.97	0.47
1:B:274:HIS:HA	1:B:277:VAL:HB	1.96	0.47
1:B:325:ASP:HB3	1:B:328:LYS:HB2	1.97	0.47
1:A:173:HIS:ND1	1:A:246:TRP:O	2.46	0.47
1:B:183:TYR:HD2	1:B:245:ARG:HB2	1.79	0.47
1:B:126:GLU:OE2	1:B:139:GLY:HA2	2.15	0.47
1:B:94:GLY:HA2	3:B:508:HOH:O	2.15	0.47
1:B:18:LEU:HD21	1:B:24:VAL:O	2.15	0.47
1:B:173:HIS:HB3	1:B:246:TRP:CZ2	2.49	0.47
1:B:120:TRP:CG	1:B:318:VAL:HG11	2.49	0.47
1:A:241:ILE:HD11	3:A:505:HOH:O	2.14	0.47
1:B:29:PRO:HD3	3:B:528:HOH:O	2.15	0.47
1:A:151:MET:HE1	1:A:199:ASN:HB2	1.97	0.46
1:A:223:ARG:O	1:A:226:ASP:HB2	2.15	0.46
1:A:5:ALA:HB3	1:A:317:LYS:HE3	1.97	0.46
1:B:125:TYR:HD2	1:B:140:LEU:CD2	2.28	0.46
1:A:268:ILE:HA	1:A:268:ILE:HD12	1.77	0.46
1:A:323:SER:HA	1:A:329:MET:SD	2.55	0.46
1:A:233:GLY:HA3	3:A:515:HOH:O	2.14	0.46
1:A:277:VAL:O	1:A:280:MET:HB2	2.16	0.46
1:A:205:LYS:HG2	1:A:206:GLN:HE21	1.80	0.46
1:B:175:PRO:O	1:B:177:VAL:N	2.49	0.46
1:B:240:HIS:HE2	1:B:274:HIS:CE1	2.34	0.46
1:B:8:ILE:HD12	1:B:29:PRO:CG	2.45	0.46
1:A:18:LEU:HD21	1:A:24:VAL:O	2.16	0.45
1:A:288:ALA:HB2	1:A:294:ARG:CD	2.46	0.45
1:A:80:ASN:ND2	1:A:80:ASN:C	2.70	0.45
1:B:30:PRO:HD2	1:B:78:PRO:HB3	1.98	0.45
1:B:87:VAL:HA	3:B:547:HOH:O	2.16	0.45
1:A:128:ILE:HG12	1:A:311:GLU:HA	1.98	0.45
1:B:53:HIS:HE1	3:B:549:HOH:O	1.99	0.45
1:A:80:ASN:ND2	1:A:82:TYR:H	2.13	0.45
1:B:125:TYR:CD2	1:B:140:LEU:HD21	2.52	0.45
1:A:77:ASP:O	1:A:83:GLN:HG3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:97:ASP:HA	1:B:105:GLU:HA	1.99	0.45
1:B:120:TRP:CD2	1:B:318:VAL:HG11	2.52	0.45
1:A:238:SER:O	1:A:242:PRO:HD2	2.17	0.44
1:B:337:PRO:C	1:B:339:PRO:HD3	2.37	0.44
1:A:233:GLY:N	3:A:521:HOH:O	2.50	0.44
1:B:338:LEU:HD22	1:B:338:LEU:N	2.32	0.44
1:A:175:PRO:O	1:A:177:VAL:N	2.50	0.44
1:A:157:LEU:HD22	1:A:268:ILE:CG2	2.47	0.44
1:B:170:ASN:ND2	1:B:173:HIS:CE1	2.86	0.44
1:B:205:LYS:HA	1:B:209:THR:HB	2.00	0.44
1:A:123:TYR:HA	3:A:553:HOH:O	2.16	0.43
1:A:235:ILE:HA	1:A:239:PRO:HD2	1.99	0.43
1:A:132:LEU:HD23	1:A:132:LEU:HA	1.77	0.43
1:A:235:ILE:HD13	1:A:235:ILE:HA	1.86	0.43
1:B:80:ASN:C	1:B:80:ASN:ND2	2.71	0.43
1:B:249:ASP:HB3	1:B:252:ASN:HD22	1.83	0.43
1:A:183:TYR:HD2	1:A:245:ARG:HB2	1.82	0.43
1:A:77:ASP:O	1:A:80:ASN:HB3	2.19	0.43
1:A:318:VAL:O	1:A:318:VAL:HG12	2.17	0.43
1:A:8:ILE:HG22	1:A:31:VAL:HA	2.01	0.43
1:A:337:PRO:C	1:A:339:PRO:HD3	2.38	0.43
1:B:115:PHE:HB2	1:B:116:PRO:HD3	2.00	0.43
1:B:323:SER:HA	1:B:329:MET:SD	2.59	0.43
1:B:80:ASN:ND2	1:B:82:TYR:H	2.15	0.43
1:A:249:ASP:O	1:A:255:ASN:HA	2.19	0.42
1:B:260:ASN:HD22	1:B:262:TYR:N	2.13	0.42
1:B:65:LYS:HE3	1:B:69:LEU:HD21	2.00	0.42
1:A:99:VAL:HG12	1:A:100:ASN:N	2.34	0.42
1:B:120:TRP:CD2	1:B:318:VAL:HG13	2.54	0.42
1:B:241:ILE:HB	1:B:242:PRO:HD3	2.01	0.42
1:B:69:LEU:N	1:B:69:LEU:HD22	2.34	0.42
1:B:112:TRP:CD1	1:B:113:LEU:HD13	2.54	0.42
1:B:152:VAL:HG12	1:B:153:LEU:O	2.19	0.42
1:B:49:ARG:NH1	1:B:333:TYR:OH	2.50	0.42
1:A:233:GLY:HA3	3:A:521:HOH:O	2.20	0.42
1:A:120:TRP:CD2	1:A:318:VAL:CG1	3.03	0.42
1:A:86:LEU:HA	1:A:86:LEU:HD23	1.81	0.42
1:B:235:ILE:HA	1:B:239:PRO:HD2	2.01	0.42
1:A:151:MET:HE2	1:A:196:ILE:HA	2.00	0.42
1:B:155:ASP:HA	1:B:175:PRO:HG3	2.00	0.42
1:B:240:HIS:HD2	1:B:277:VAL:HG21	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116:PRO:HB3	1:B:320:ILE:HG23	2.03	0.41
1:B:79:ARG:HA	1:B:84:GLN:HE21	1.85	0.41
1:A:131:LYS:CE	1:A:312:ASN:HD21	2.29	0.41
1:B:249:ASP:O	1:B:255:ASN:HA	2.20	0.41
1:B:129:LEU:HD12	1:B:140:LEU:HD22	2.01	0.41
1:B:148:PRO:HA	1:B:151:MET:SD	2.61	0.41
1:A:110:ASN:HA	1:A:232:ALA:CA	2.50	0.41
1:B:123:TYR:O	1:B:127:ARG:HG3	2.21	0.41
1:A:282:THR:HG23	1:A:326:ASN:HD21	1.84	0.41
1:A:13:VAL:HA	3:A:546:HOH:O	2.20	0.41
1:A:120:TRP:CG	1:A:318:VAL:HG11	2.56	0.41
1:A:120:TRP:CD2	1:A:318:VAL:HG13	2.56	0.41
1:B:3:ILE:HD13	1:B:106:ILE:HG23	2.03	0.41
1:A:238:SER:O	1:A:242:PRO:HG2	2.21	0.41
1:B:175:PRO:HB2	1:B:176:PRO:CD	2.51	0.41
1:A:216:LEU:HA	1:A:216:LEU:HD23	1.90	0.40
1:A:4:GLN:HA	1:A:305:THR:HG22	2.03	0.40
1:A:338:LEU:HD22	1:A:338:LEU:N	2.36	0.40
1:B:282:THR:HG23	1:B:326:ASN:HD21	1.85	0.40
1:A:118:HIS:CD2	1:A:274:HIS:CE1	3.09	0.40
1:B:223:ARG:NH1	1:B:300:ASP:CB	2.77	0.40
1:A:88:HIS:HB2	1:A:121:TYR:OH	2.21	0.40
1:B:53:HIS:HB3	3:B:552:HOH:O	2.21	0.40
1:A:115:PHE:HB2	1:A:116:PRO:HD3	2.02	0.40
1:A:80:ASN:ND2	1:A:82:TYR:N	2.69	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	332/345 (96%)	299 (90%)	24 (7%)	9 (3%)	<b>6</b> <b>15</b>

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	332/345 (96%)	304 (92%)	22 (7%)	6 (2%)	10	25
All	All	664/690 (96%)	603 (91%)	46 (7%)	15 (2%)	7	19

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	99	VAL
1	A	175	PRO
1	B	99	VAL
1	B	175	PRO
1	A	114	PHE
1	A	257	ASP
1	A	287	LEU
1	B	114	PHE
1	B	257	ASP
1	A	240	HIS
1	A	256	GLU
1	B	240	HIS
1	B	256	GLU
1	A	8	ILE
1	A	154	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	285/293 (97%)	258 (90%)	27 (10%)	10	23
1	B	285/293 (97%)	258 (90%)	27 (10%)	10	23
All	All	570/586 (97%)	516 (90%)	54 (10%)	10	23

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ILE

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Mol	Chain	Res	Type
1	A	34	ASN
1	A	76	ASP
1	A	80	ASN
1	A	103	ASP
1	A	112	TRP
1	A	113	LEU
1	A	137	SER
1	A	144	ASN
1	A	154	PRO
1	A	160	SER
1	A	162	SER
1	A	167	SER
1	A	175	PRO
1	A	177	VAL
1	A	190	VAL
1	A	202	LEU
1	A	242	PRO
1	A	245	ARG
1	A	253	THR
1	A	260	ASN
1	A	262	TYR
1	A	298	ASP
1	A	305	THR
1	A	316	VAL
1	A	318	VAL
1	A	336	THR
1	B	3	ILE
1	B	34	ASN
1	B	56	ASP
1	B	76	ASP
1	B	80	ASN
1	B	103	ASP
1	B	112	TRP
1	B	113	LEU
1	B	137	SER
1	B	140	LEU
1	B	144	ASN
1	B	154	PRO
1	B	162	SER
1	B	167	SER
1	B	175	PRO
1	B	190	VAL

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Mol	Chain	Res	Type
1	B	202	LEU
1	B	242	PRO
1	B	245	ARG
1	B	253	THR
1	B	260	ASN
1	B	262	TYR
1	B	294	ARG
1	B	298	ASP
1	B	316	VAL
1	B	318	VAL
1	B	336	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	ASN
1	A	93	ASN
1	A	199	ASN
1	A	206	GLN
1	A	210	ASN
1	A	260	ASN
1	A	273	HIS
1	A	276	ASN
1	A	303	ASN
1	A	312	ASN
1	A	326	ASN
1	B	53	HIS
1	B	80	ASN
1	B	93	ASN
1	B	199	ASN
1	B	206	GLN
1	B	210	ASN
1	B	260	ASN
1	B	273	HIS
1	B	276	ASN
1	B	326	ASN

### 5.3.3 RNA ⓘ

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](#)

2 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	$\# Z  > 2$
2	C2O	A	500	1	0,2,2	0.00	-	0,1,1	0.00	-
2	C2O	B	500	1	0,2,2	0.00	-	0,1,1	0.00	-

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	C2O	A	500	1	-	0/0/0/0	0/0/0/0
2	C2O	B	500	1	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	336/345 (97%)	-0.81	0 100 100	4, 16, 29, 42	0
1	B	336/345 (97%)	-0.87	0 100 100	4, 16, 29, 41	0
All	All	672/690 (97%)	-0.84	0 100 100	4, 16, 29, 42	0

There are no RSRZ outliers to report.

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	C2O	A	500	3/3	0.99	0.05	-3.06	17,17,19,22	0
2	C2O	B	500	3/3	0.99	0.04	-4.71	18,18,19,25	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.