



Full wwPDB X-ray Structure Validation Report

Jul 13, 2014 – 04:36 AM EDT

PDB ID : 4LUY
Title : Crystal structure of CdALR mutant K 271 T
Authors : Asojo, O.A.
Deposited on : 2013-07-25
Resolution : 2.60 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

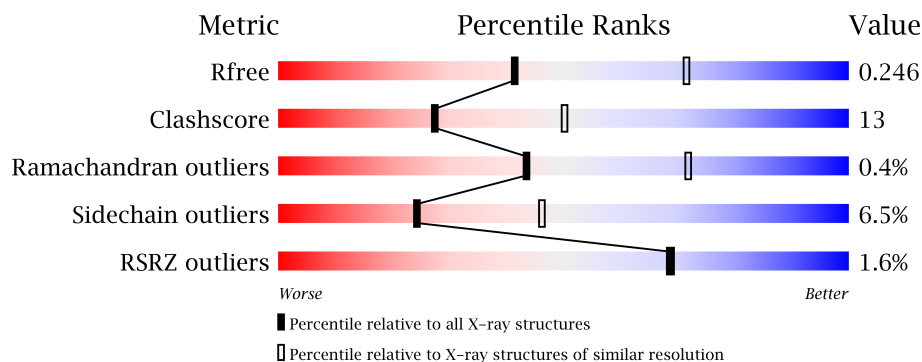
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : **FAILED**
Xtriage (Phenix) : dev-1439
EDS : stable23161
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable23161

1 Overall quality at a glance

The reported resolution of this entry is 2.60 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}	66092	1718 (2.60-2.60)
Clashscore	79885	2154 (2.60-2.60)
Ramachandran outliers	78287	2113 (2.60-2.60)
Sidechain outliers	78261	2113 (2.60-2.60)
RSRZ outliers	66119	1718 (2.60-2.60)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	385	
1	B	385	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6158 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Alanine racemase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	383	Total	C	N	O	P	S	0	0	0
			3036	1927	505	587	1	16			
1	B	383	Total	C	N	O	P	S	0	0	0
			3036	1927	505	587	1	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	271	THR	LYS	ENGINEERED MUTATION	UNP Q180W0
B	271	THR	LYS	ENGINEERED MUTATION	UNP Q180W0

- Molecule 2 is water.

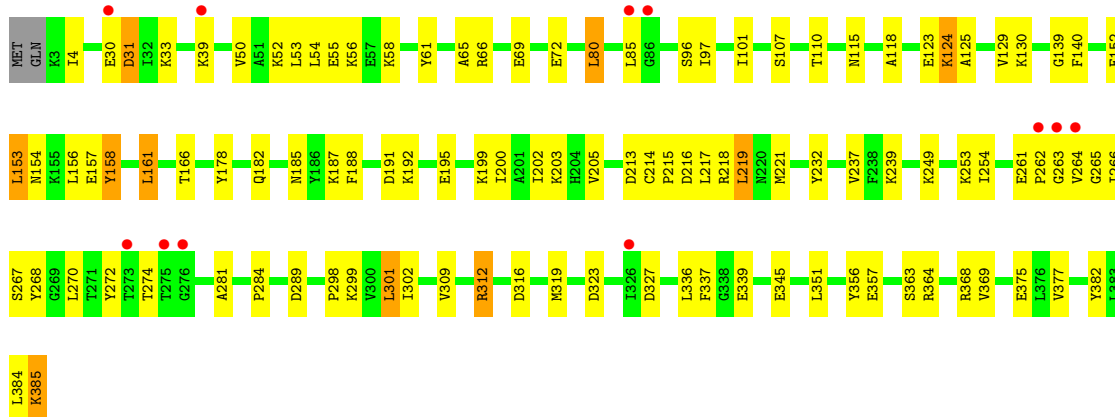
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	44	Total	O	0	0
			44	44		
2	B	42	Total	O	0	0
			42	42		

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.

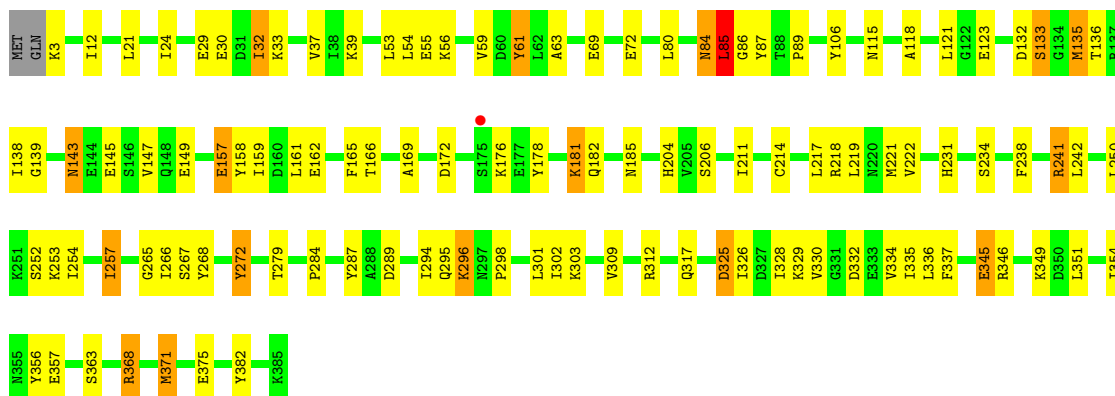
- Molecule 1: Alanine racemase

Chain A: 



- Molecule 1: Alanine racemase

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.97Å 154.33Å 55.77Å 90.00° 113.63° 90.00°	Depositor
Resolution (Å)	28.83 – 2.60 28.83 – 2.60	Depositor EDS
% Data completeness (in resolution range)	97.1 (28.83-2.60) 97.2 (28.83-2.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.40 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.8.0069	Depositor
R, R_{free}	0.181 , 0.252 0.182 , 0.246	Depositor DCC
R_{free} test set	1182 reflections (5.39%)	DCC
Wilson B-factor (Å ²)	41.8	Xtriage
Anisotropy	0.014	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 29.1	EDS
Estimated twinning fraction	0.044 for h,-k,-h-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 23092 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6158	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.86% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	1.69	8/3041 (0.3%)	0.92	4/4102 (0.1%)
1	B	1.72	11/3041 (0.4%)	0.90	2/4102 (0.0%)
All	All	1.71	19/6082 (0.3%)	0.91	6/8204 (0.1%)

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	357	GLU	CD-OE1	-6.26	1.18	1.25
1	B	69	GLU	CD-OE1	-6.09	1.19	1.25
1	B	162	GLU	CD-OE2	-5.87	1.19	1.25
1	B	72	GLU	CD-OE1	-5.79	1.19	1.25
1	B	345	GLU	CD-OE2	-5.77	1.19	1.25
1	A	357	GLU	CD-OE1	-5.69	1.19	1.25
1	A	345	GLU	CD-OE2	-5.61	1.19	1.25
1	B	72	GLU	CD-OE2	-5.55	1.19	1.25
1	B	357	GLU	CD-OE2	-5.52	1.19	1.25
1	A	72	GLU	CD-OE1	-5.50	1.19	1.25
1	A	72	GLU	CD-OE2	-5.33	1.19	1.25
1	A	382	TYR	CE1-CZ	-5.32	1.31	1.38
1	B	272	TYR	CE1-CZ	-5.28	1.31	1.38
1	B	317	GLN	C-O	-5.21	1.13	1.23
1	A	357	GLU	CD-OE2	-5.08	1.20	1.25
1	A	69	GLU	CD-OE2	-5.07	1.20	1.25
1	B	287	TYR	CE1-CZ	-5.05	1.31	1.38
1	A	364	ARG	CZ-NH2	-5.05	1.26	1.33
1	B	382	TYR	CE1-CZ	-5.01	1.32	1.38

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	158	TYR	N-CA-CB	-9.71	93.12	110.60
1	A	153	LEU	CA-CB-CG	6.71	130.72	115.30
1	A	158	TYR	N-CA-C	-5.33	96.60	111.00
1	B	332	ASP	CB-CG-OD1	5.24	123.01	118.30
1	A	356	TYR	CA-CB-CG	5.04	122.98	113.40
1	B	85	LEU	CA-CB-CG	5.02	126.84	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3036	0	3084	78	0
1	B	3036	0	3085	80	0
2	A	44	0	0	2	0
2	B	42	0	0	2	0
All	All	6158	0	6169	158	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 13.

All (158) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:115:ASN:ND2	1:B:157:GLU:O	1.83	1.10
1:A:264:VAL:H	1:A:274:THR:HG22	1.18	1.08
1:B:295:GLN:O	1:B:296:LYS:HB2	1.49	1.06
1:A:52:LYS:O	1:A:56:LYS:HD2	1.65	0.96
1:A:261:GLU:O	1:A:274:THR:HG21	1.65	0.95
1:A:52:LYS:O	1:A:56:LYS:CD	2.19	0.89
1:B:149:GLU:OE1	2:B:422:HOH:O	1.88	0.88
1:A:214:CYS:HB3	1:A:217:LEU:HD12	1.56	0.86
1:A:214:CYS:HB3	1:A:217:LEU:CD1	2.12	0.78
1:B:118:ALA:HB1	1:B:123:GLU:O	1.84	0.77
1:A:195:GLU:OE2	2:A:529:HOH:O	2.04	0.75

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:214:CYS:HB3	1:B:217:LEU:HD12	1.70	0.72
1:A:298:PRO:HB2	1:A:309:VAL:HB	1.71	0.72
1:A:55:GLU:HG2	1:A:80:LEU:HD22	1.72	0.72
1:A:214:CYS:CB	1:A:217:LEU:HD12	2.20	0.72
1:B:135:MET:O	1:B:136:THR:OG1	2.08	0.71
1:A:262:PRO:HA	1:A:274:THR:HG23	1.71	0.70
1:A:266:ILE:HD12	1:A:272:TYR:CD2	2.27	0.69
1:A:107:SER:OG	1:A:110:THR:HG23	1.91	0.69
1:B:61:TYR:HB2	1:B:221:MET:HE2	1.73	0.69
1:B:135:MET:HE1	1:B:169:ALA:HA	1.74	0.68
1:A:53:LEU:C	1:A:53:LEU:HD23	2.14	0.68
1:B:211:ILE:HD11	1:B:222:VAL:HB	1.75	0.68
1:A:312:ARG:HG2	1:A:312:ARG:HH11	1.59	0.66
1:A:52:LYS:O	1:A:56:LYS:HD3	1.95	0.66
1:B:133:SER:OG	1:B:166:THR:OG1	2.11	0.66
1:A:266:ILE:HD12	1:A:272:TYR:HD2	1.60	0.66
1:A:261:GLU:O	1:A:274:THR:CG2	2.43	0.65
1:A:53:LEU:O	1:A:53:LEU:HD23	1.97	0.65
1:B:61:TYR:HB2	1:B:221:MET:CE	2.26	0.65
1:B:143:ASN:OD1	1:B:145:GLU:N	2.28	0.64
1:A:264:VAL:N	1:A:274:THR:HG22	2.03	0.64
1:B:84:ASN:ND2	1:B:86:GLY:H	1.96	0.63
1:B:84:ASN:C	1:B:84:ASN:HD22	2.00	0.63
1:B:84:ASN:HD22	1:B:85:LEU:N	1.98	0.62
1:A:263:GLY:N	1:A:274:THR:O	2.26	0.62
1:A:267:SER:OG	1:A:268:TYR:N	2.29	0.62
1:B:161:LEU:O	1:B:161:LEU:HG	1.98	0.62
1:B:39:LLP:H2'3	1:B:85:LEU:HG	1.81	0.62
1:A:336:LEU:HD22	1:A:336:LEU:H	1.65	0.62
1:B:157:GLU:O	1:B:158:TYR:HB2	2.00	0.61
1:A:336:LEU:HD22	1:A:336:LEU:N	2.15	0.61
1:B:325:ASP:OD1	1:B:325:ASP:N	2.31	0.61
1:A:213:ASP:C	1:A:215:PRO:HD3	2.21	0.61
1:B:214:CYS:CB	1:B:217:LEU:HD12	2.30	0.61
1:B:61:TYR:HD2	1:B:221:MET:HE1	1.67	0.60
1:B:106:TYR:HB3	1:B:139:GLY:HA2	1.84	0.59
1:A:312:ARG:HH11	1:A:312:ARG:CG	2.17	0.58
1:A:375:GLU:O	1:A:377:VAL:HG13	2.04	0.58
1:B:55:GLU:HG2	1:B:80:LEU:HG	1.86	0.57
1:B:181:LYS:O	1:B:185:ASN:ND2	2.37	0.57
1:B:178:TYR:O	1:B:182:GLN:HG3	2.05	0.56
1:A:213:ASP:O	1:A:215:PRO:HD3	2.05	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:371:MET:HA	1:B:375:GLU:O	2.06	0.56
1:A:50:VAL:O	1:A:54:LEU:HG	2.07	0.55
1:B:295:GLN:O	1:B:296:LYS:CB	2.36	0.55
1:B:63:ALA:HB1	1:B:85:LEU:HD22	1.89	0.54
1:B:87:TYR:CE2	1:B:89:PRO:HD3	2.43	0.54
1:B:84:ASN:C	1:B:84:ASN:ND2	2.59	0.54
1:B:298:PRO:HG2	1:B:309:VAL:HB	1.90	0.54
1:A:185:ASN:O	1:A:188:PHE:HB3	2.08	0.54
1:B:115:ASN:ND2	1:B:158:TYR:HB2	2.22	0.53
1:B:135:MET:C	1:B:136:THR:OG1	2.47	0.53
1:A:249:LYS:NZ	1:A:339:GLU:OE2	2.42	0.53
1:B:165:PHE:HB3	1:B:204:HIS:CE1	2.44	0.52
1:B:325:ASP:C	1:B:326:ILE:HG23	2.29	0.52
1:B:268:TYR:O	1:B:312:ARG:HD2	2.10	0.52
1:B:32:ILE:N	1:B:32:ILE:CD1	2.73	0.52
1:A:203:LYS:HB3	1:A:219:LEU:HD23	1.90	0.52
1:B:250:LEU:HG	1:B:336:LEU:HD23	1.91	0.52
1:A:384:LEU:O	1:A:385:LYS:HB2	2.09	0.52
1:A:30:GLU:HA	1:A:30:GLU:OE1	2.09	0.51
1:A:154:ASN:HB2	1:A:161:LEU:HD12	1.92	0.50
1:B:157:GLU:O	1:B:158:TYR:CB	2.57	0.50
1:B:115:ASN:HD22	1:B:159:ILE:HG12	1.76	0.50
1:B:267:SER:OG	1:B:268:TYR:N	2.42	0.50
1:B:257:ILE:HD11	1:B:279:THR:CG2	2.42	0.50
1:A:263:GLY:N	1:A:274:THR:CG2	2.74	0.50
1:B:30:GLU:H	1:B:30:GLU:CD	2.12	0.49
1:B:337:PHE:C	1:B:337:PHE:CD1	2.85	0.49
1:B:185:ASN:HD22	1:B:185:ASN:H	1.59	0.49
1:B:253:LYS:O	1:B:284:PRO:HD2	2.12	0.49
1:A:214:CYS:N	1:A:215:PRO:HD3	2.28	0.49
1:B:214:CYS:HB3	1:B:217:LEU:CD1	2.42	0.49
1:A:124:LYS:HE3	1:A:157:GLU:O	2.13	0.48
1:A:385:LYS:HA	1:A:385:LYS:HD2	1.66	0.48
1:B:61:TYR:CD2	1:B:221:MET:HE1	2.48	0.48
1:A:124:LYS:CE	1:A:157:GLU:O	2.61	0.48
1:A:253:LYS:O	1:A:284:PRO:HD2	2.14	0.48
1:A:4:ILE:HD11	1:A:369:VAL:HG11	1.96	0.47
1:B:294:ILE:HG22	1:B:294:ILE:O	2.14	0.47
1:B:143:ASN:O	1:B:147:VAL:HG23	2.15	0.47
1:B:302:ILE:HG12	1:B:334:VAL:HG22	1.96	0.47
1:A:97:ILE:HG12	1:A:125:ALA:HB2	1.97	0.47
1:A:130:KCX:HE2	1:A:166:THR:HA	1.97	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:132:ASP:OD2	1:B:136:THR:HA	2.15	0.46
1:A:39:LLP:C2	1:A:85:LEU:HD13	2.45	0.46
1:A:254:ILE:HD13	1:A:281:ALA:HB1	1.97	0.46
1:B:33:LYS:HB3	1:B:221:MET:HE3	1.98	0.46
1:B:238:PHE:HB3	1:B:241:ARG:HG3	1.98	0.46
1:A:33:LYS:CB	1:A:221:MET:HG3	2.46	0.45
1:A:65:ALA:C	1:A:66:ARG:HG2	2.36	0.45
1:B:32:ILE:N	1:B:32:ILE:HD12	2.31	0.45
1:A:96:SER:HB3	1:A:101:ILE:HB	1.96	0.45
1:A:263:GLY:N	1:A:274:THR:HG23	2.30	0.45
1:B:266:ILE:O	1:B:267:SER:HB3	2.17	0.45
1:B:39:LLP:C2'	1:B:85:LEU:HG	2.46	0.45
1:A:192:LYS:HD3	1:A:192:LYS:HA	1.74	0.45
1:A:166:THR:HG22	1:A:205:VAL:HG12	1.98	0.45
1:A:337:PHE:C	1:A:337:PHE:CD1	2.90	0.45
1:B:301:LEU:HB3	1:B:335:ILE:HB	1.98	0.45
1:A:115:ASN:OD1	1:A:158:TYR:HB2	2.16	0.45
1:A:268:TYR:O	1:A:312:ARG:HD2	2.16	0.45
1:B:325:ASP:C	1:B:326:ILE:CG2	2.85	0.44
1:A:53:LEU:CD2	1:A:53:LEU:C	2.85	0.44
1:B:303:LYS:HB2	1:B:326:ILE:HD12	2.00	0.44
1:B:37:VAL:HG11	1:B:39:LLP:HE3	2.00	0.44
1:B:21:LEU:O	1:B:24:ILE:HB	2.18	0.44
1:A:312:ARG:NH1	1:A:312:ARG:CG	2.79	0.44
1:B:354:ILE:HD12	1:B:356:TYR:CD1	2.53	0.44
1:A:39:LLP:H2'3	1:A:85:LEU:HD13	2.00	0.43
1:A:266:ILE:O	1:A:267:SER:HB3	2.18	0.43
1:A:153:LEU:O	1:A:156:LEU:HD12	2.18	0.43
1:B:54:LEU:HD22	1:B:59:VAL:HG11	2.00	0.43
1:A:129:VAL:HG13	1:A:140:PHE:CE2	2.54	0.43
1:B:206:SER:HB2	1:B:222:VAL:HG12	2.00	0.42
1:A:39:LLP:C2'	1:A:85:LEU:HD13	2.49	0.42
1:B:231:HIS:HD2	1:B:345:GLU:OE2	2.01	0.42
1:A:130:KCX:HB2	1:A:139:GLY:HA2	2.01	0.42
1:B:218:ARG:NH1	2:B:411:HOH:O	2.52	0.42
1:B:3:LYS:HD2	1:B:3:LYS:HA	1.62	0.42
1:A:178:TYR:O	1:A:182:GLN:HG3	2.19	0.42
1:B:218:ARG:O	1:B:219:LEU:HB2	2.18	0.42
1:A:202:ILE:HG21	1:A:221:MET:CE	2.50	0.42
1:A:289:ASP:HA	1:A:363:SER:HB2	2.01	0.42
1:A:263:GLY:N	1:A:274:THR:HG22	2.34	0.42
1:A:263:GLY:H	1:A:274:THR:HG23	1.85	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:301:LEU:C	1:A:302:ILE:HG13	2.40	0.42
1:B:257:ILE:HD11	1:B:279:THR:HG22	2.02	0.42
1:A:187:LYS:NZ	1:A:191:ASP:OD2	2.46	0.41
1:A:232:TYR:CD2	1:A:239:LYS:HE2	2.55	0.41
1:A:265:GLY:HA3	1:A:270:LEU:HD22	2.02	0.41
1:B:326:ILE:HG13	1:B:328:ILE:HG23	2.01	0.41
1:A:118:ALA:HB1	1:A:123:GLU:O	2.20	0.41
1:B:12:ILE:CD1	1:B:368:ARG:HG2	2.51	0.41
1:A:80:LEU:HD12	1:A:80:LEU:HA	1.84	0.41
1:B:265:GLY:HA2	1:B:272:TYR:O	2.21	0.41
1:B:172:ASP:HB2	1:B:234:SER:HB3	2.02	0.41
1:B:29:GLU:HB2	1:B:32:ILE:HD13	2.02	0.41
1:B:53:LEU:HD23	1:B:53:LEU:C	2.41	0.41
1:A:218:ARG:HG2	1:A:218:ARG:H	1.74	0.40
1:B:289:ASP:HA	1:B:363:SER:OG	2.21	0.40
1:A:262:PRO:CA	1:A:274:THR:HG23	2.47	0.40
1:B:138:ILE:HD13	1:B:138:ILE:HG21	1.85	0.40
1:A:216:ASP:N	1:A:216:ASP:OD1	2.48	0.40
1:A:219:LEU:HD12	2:A:525:HOH:O	2.21	0.40
1:B:169:ALA:H	1:B:182:GLN:HE22	1.68	0.40
1:B:254:ILE:HG22	1:B:330:VAL:HA	2.04	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	379/385 (98%)	365 (96%)	12 (3%)	2 (0%)	38	67
1	B	379/385 (98%)	369 (97%)	9 (2%)	1 (0%)	50	77
All	All	758/770 (98%)	734 (97%)	21 (3%)	3 (0%)	43	72

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	296	LYS
1	A	200	ILE
1	A	31	ASP

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	330/332 (99%)	310 (94%)	20 (6%)	26	50
1	B	330/332 (99%)	307 (93%)	23 (7%)	21	41
All	All	660/664 (99%)	617 (94%)	43 (6%)	24	46

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	ASP
1	A	58	LYS
1	A	61	TYR
1	A	80	LEU
1	A	124	LYS
1	A	152	GLU
1	A	161	LEU
1	A	199	LYS
1	A	219	LEU
1	A	237	VAL
1	A	299	LYS
1	A	301	LEU
1	A	312	ARG
1	A	316	ASP
1	A	319	MET
1	A	323	ASP
1	A	327	ASP
1	A	351	LEU
1	A	368	ARG
1	A	385	LYS
1	B	32	ILE
1	B	56	LYS
1	B	61	TYR

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Mol	Chain	Res	Type
1	B	84	ASN
1	B	85	LEU
1	B	121	LEU
1	B	133	SER
1	B	135	MET
1	B	143	ASN
1	B	157	GLU
1	B	176	LYS
1	B	181	LYS
1	B	241	ARG
1	B	242	LEU
1	B	252	SER
1	B	257	ILE
1	B	325	ASP
1	B	329	LYS
1	B	346	ARG
1	B	349	LYS
1	B	351	LEU
1	B	368	ARG
1	B	371	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	84	ASN
1	B	141	GLN
1	B	182	GLN
1	B	185	ASN
1	B	231	HIS
1	B	259	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

Mogul failed to run properly - this section will therefore be empty.

5.5 Carbohydrates ⓘ

Mogul failed to run properly - this section will therefore be empty.

5.6 Ligand geometry ⓘ

Mogul failed to run properly - this section will therefore be empty.

5.7 Other polymers ⓘ

Mogul failed to run properly - this section will therefore be empty.

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	383/385 (99%)	-0.15	11 (2%) 49 46	26, 39, 58, 79	0
1	B	383/385 (99%)	-0.27	1 (0%) 91 92	23, 39, 59, 67	0
All	All	766/770 (99%)	-0.21	12 (1%) 68 69	23, 39, 59, 79	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	85	LEU	3.6
1	A	30	GLU	2.9
1	A	275	THR	2.6
1	A	276	GLY	2.6
1	A	262	PRO	2.5
1	B	175	SER	2.4
1	A	273	THR	2.4
1	A	39	LLP	2.3
1	A	86	GLY	2.1
1	A	326	ILE	2.1
1	A	263	GLY	2.1
1	A	264	VAL	2.0

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

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	KCX	A	130	12/13	0.24	1.24	29,31,42,42	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	KCX	B	130	12/13	0.21	0.85	42,47,57,61	0
1	LLP	B	39	24/25	0.24	0.84	30,35,38,38	0
1	LLP	A	39	24/25	0.27	0.83	29,33,37,42	0

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.