



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:06 PM UTC

PDB ID : 1CMX / pdb_00001cmx
Title : STRUCTURAL BASIS FOR THE SPECIFICITY OF UBIQUITIN C-
TERMINAL HYDROLASES
Authors : Johnston, S.C.; Riddle, S.M.; Cohen, R.E.; Hill, C.P.
Deposited on : 1999-05-12
Resolution : 2.25 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

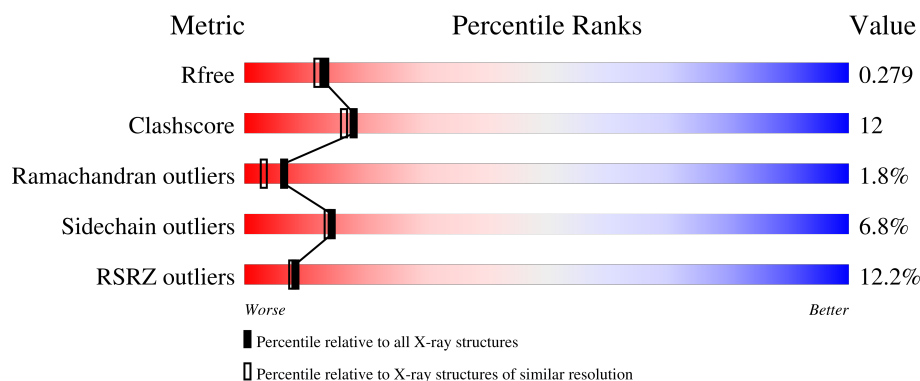
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	235	
1	C	235	
2	B	76	
2	D	76	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4142 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 (UBIQUITIN YUH1-UBAL).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	10	0	0
			1684	1072	274	335	3			
1	C	210	Total	C	N	O	S	24	0	0
			1649	1052	266	328	3			

- Molecule 2 is a protein called PROTEIN (UBIQUITIN YUH1-UBAL).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	76	Total	C	N	O	S	10	0	0
			601	378	105	117	1			
2	D	17	Total	C	N	O		6	0	0
			133	83	29	21				

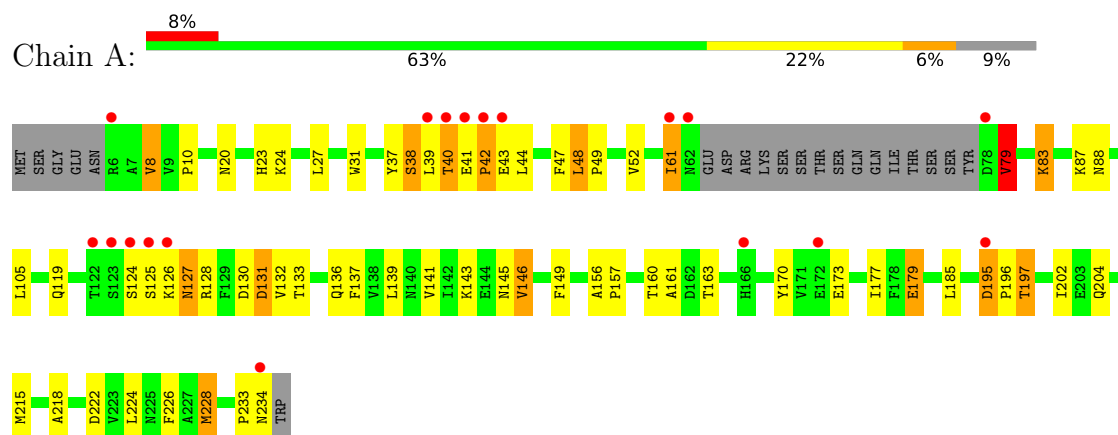
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	38	Total	O	0	0
			38	38		
3	B	16	Total	O	0	0
			16	16		
3	C	18	Total	O	0	0
			18	18		
3	D	3	Total	O	0	0
			3	3		

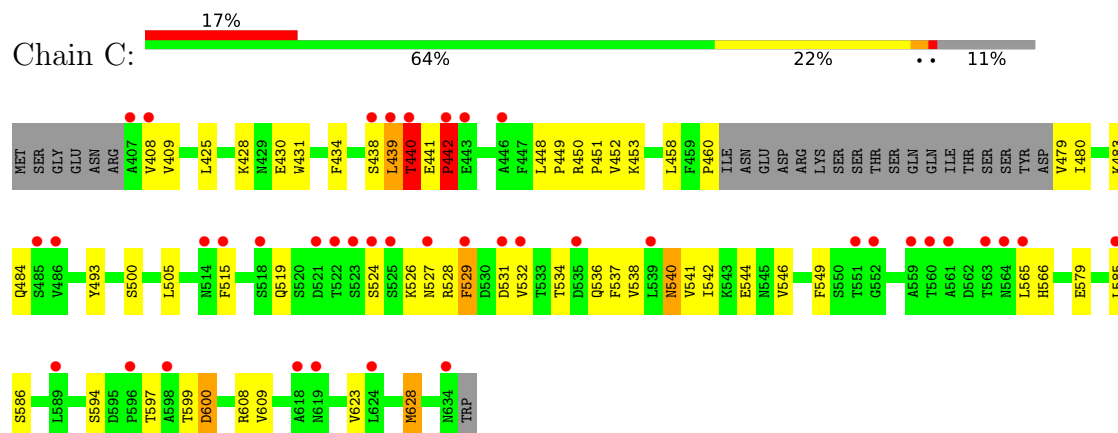
3 Residue-property plots [i](#)

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.

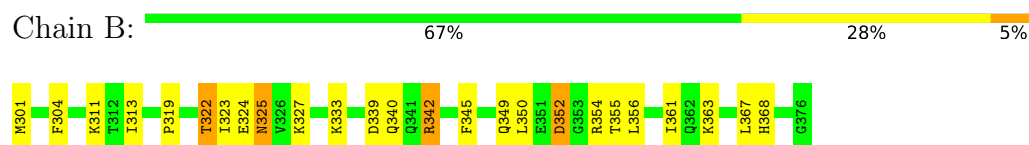
- Molecule 1: PROTEIN (UBIQUITIN YUH1-UBAL)



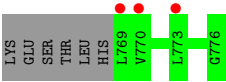
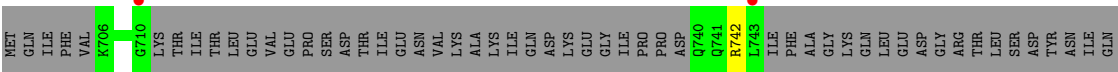
- Molecule 1: PROTEIN (UBIQUITIN YUH1-UBAL)



- Molecule 2: PROTEIN (UBIQUITIN YUH1-UBAL)



- Molecule 2: PROTEIN (UBIQUITIN YUH1-UBAL)



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	199.30Å 199.30Å 36.80Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.25 30.00 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.0 (30.00-2.25) 98.9 (30.00-2.25)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.29 (at 2.22Å)	Xtriage
Refinement program	X-PLOR 98.0	Depositor
R, R_{free}	0.248 , 0.285 0.242 , 0.279	Depositor DCC
R_{free} test set	1262 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	46.1	Xtriage
Anisotropy	0.344	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 48.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.022 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4142	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% 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: GLZ

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	0.89	1/1719 (0.1%)	1.19	12/2340 (0.5%)
1	C	0.81	2/1684 (0.1%)	1.16	10/2293 (0.4%)
2	B	1.10	1/603 (0.2%)	1.16	2/811 (0.2%)
2	D	0.77	0/126	0.96	1/163 (0.6%)
All	All	0.89	4/4132 (0.1%)	1.17	25/5607 (0.4%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	322	THR	CA-CB	7.67	1.66	1.53
1	C	609	VAL	CA-CB	5.67	1.61	1.54
1	A	228	MET	SD-CE	-5.29	1.66	1.79
1	C	628	MET	SD-CE	-5.04	1.67	1.79

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	439	LEU	N-CA-C	9.95	123.24	109.11
1	C	532	VAL	N-CA-C	-9.34	102.80	111.67
1	A	48	LEU	CA-C-N	7.70	127.69	119.76
1	A	48	LEU	C-N-CA	7.70	127.69	119.76
1	C	440	THR	N-CA-CB	7.49	123.14	110.49
1	C	546	VAL	N-CA-C	6.74	117.44	110.36
1	A	79	VAL	CB-CA-C	-6.66	102.00	111.40
1	A	132	VAL	N-CA-C	-6.63	102.76	112.04
1	A	131	ASP	N-CA-C	6.59	118.51	110.41
1	A	47	PHE	N-CA-C	-6.47	104.97	112.92
1	C	600	ASP	N-CA-C	6.34	116.44	108.45
1	A	179	GLU	N-CA-C	-6.27	98.68	108.90
1	A	146	VAL	N-CA-C	5.75	119.10	111.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	352	ASP	N-CA-C	5.68	122.90	110.80
1	C	529	PHE	N-CA-C	5.65	117.95	108.23
1	C	599	THR	N-CA-C	-5.58	105.20	111.28
1	A	197	THR	N-CA-C	-5.54	106.31	112.57
1	C	579	GLU	N-CA-C	-5.46	100.28	109.07
1	C	597	THR	N-CA-C	-5.35	106.70	113.18
1	A	161	ALA	N-CA-C	5.21	118.71	109.96
2	B	325	ASN	N-CA-C	-5.12	105.70	111.28
1	C	442	PRO	N-CA-C	5.11	123.00	112.47
2	D	742	ARG	CD-NE-CZ	-5.05	117.33	124.40
1	A	195	ASP	CA-C-N	5.04	126.14	119.84
1	A	195	ASP	C-N-CA	5.04	126.14	119.84

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	1684	0	1634	50	0
1	C	1649	0	1600	39	0
2	B	601	0	625	15	0
2	D	133	0	151	0	0
3	A	38	0	0	0	0
3	B	16	0	0	2	0
3	C	18	0	0	1	0
3	D	3	0	0	0	0
All	All	4142	0	4010	97	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 12.

All (97) 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:C:439:LEU:HD11	1:C:628:MET:HE1	1.45	0.97
1:C:439:LEU:O	1:C:440:THR:HG23	1.65	0.95
1:A:39:LEU:HD22	1:A:228:MET:HE1	1.60	0.83
1:A:83:LYS:HZ1	1:A:185:LEU:HD22	1.45	0.81
1:A:41:GLU:HG2	1:A:44:LEU:HD12	1.62	0.80
1:A:83:LYS:NZ	1:A:185:LEU:HD22	1.98	0.79
1:A:48:LEU:HD21	1:A:202:ILE:CD1	2.18	0.73
1:A:40:THR:HG23	1:C:440:THR:HB	1.71	0.73
1:C:458:LEU:HD11	1:C:565:LEU:HB3	1.71	0.72
1:A:42:PRO:HG2	1:A:43:GLU:H	1.56	0.71
1:C:529:PHE:HB3	1:C:534:THR:HG21	1.73	0.70
1:C:538:VAL:O	1:C:542:ILE:HG13	1.91	0.69
1:A:224:LEU:HD21	2:B:311:LYS:HD3	1.76	0.67
1:C:515:PHE:HZ	1:C:534:THR:HG22	1.61	0.67
2:B:301:MET:HE3	2:B:319:PRO:HG3	1.77	0.66
2:B:361:ILE:HD13	2:B:367:LEU:HD21	1.79	0.63
1:A:48:LEU:HD21	1:A:202:ILE:HD13	1.83	0.59
2:B:301:MET:HE1	3:B:838:HOH:O	2.02	0.59
1:C:439:LEU:CD1	1:C:628:MET:HE1	2.29	0.59
1:A:79:VAL:HG12	1:A:173:GLU:HG3	1.85	0.59
1:C:428:LYS:HB3	1:C:430:GLU:OE1	2.03	0.58
1:C:439:LEU:O	1:C:440:THR:CG2	2.48	0.58
1:A:10:PRO:HD3	1:A:88:ASN:ND2	2.19	0.58
1:A:37:TYR:HE1	1:A:44:LEU:HD22	1.69	0.56
1:C:441:GLU:CD	1:C:442:PRO:HD2	2.31	0.56
1:A:31:TRP:CE2	1:A:105:LEU:HD22	2.41	0.56
1:A:8:VAL:CG1	1:A:149:PHE:CD2	2.89	0.56
1:A:40:THR:HG22	1:A:40:THR:O	2.05	0.56
2:B:323:ILE:HB	2:B:352:ASP:HA	1.89	0.55
1:A:10:PRO:HD3	1:A:88:ASN:HD22	1.71	0.54
1:A:40:THR:HG21	1:C:440:THR:HG22	1.88	0.54
1:C:434:PHE:CD1	1:C:448:LEU:HD22	2.43	0.54
1:A:8:VAL:HG11	1:A:149:PHE:CE2	2.43	0.53
1:C:434:PHE:HZ	1:C:452:VAL:HG22	1.72	0.53
1:A:24:LYS:HD3	1:A:145:ASN:OD1	2.08	0.53
1:A:40:THR:HG22	1:C:608:ARG:NH1	2.23	0.53
1:A:39:LEU:HD22	1:A:228:MET:CE	2.34	0.53
1:C:458:LEU:HD12	1:C:566:HIS:O	2.08	0.53
2:B:325:ASN:ND2	3:B:871:HOH:O	2.34	0.53
1:C:408:VAL:HG11	1:C:549:PHE:CD2	2.44	0.52
1:C:408:VAL:HG12	1:C:409:VAL:N	2.25	0.52
2:B:345:PHE:HB3	2:B:350:LEU:HD21	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:ILE:HD11	1:A:204:GLN:OE1	2.09	0.52
1:A:8:VAL:HG11	1:A:149:PHE:CD2	2.45	0.51
1:A:218:ALA:HB2	1:A:226:PHE:CE1	2.46	0.51
1:C:519:GLN:OE1	1:C:529:PHE:CE1	2.64	0.51
1:A:170:TYR:CE2	1:A:179:GLU:HG3	2.47	0.50
1:C:408:VAL:HG12	1:C:409:VAL:H	1.77	0.50
1:C:537:PHE:O	1:C:541:VAL:HG23	2.10	0.50
1:A:44:LEU:HD11	2:B:368:HIS:NE2	2.27	0.49
1:C:540:ASN:O	1:C:544:GLU:HG3	2.11	0.49
2:B:322:THR:HA	2:B:355:THR:HA	1.93	0.49
1:C:450:ARG:NE	1:C:600:ASP:OD2	2.44	0.49
1:A:38:SER:HB3	1:A:41:GLU:HB2	1.95	0.49
1:A:38:SER:HB3	1:A:41:GLU:CB	2.42	0.49
1:A:119:GLN:NE2	1:A:133:THR:HG21	2.28	0.48
1:C:480:ILE:HD12	1:C:500:SER:O	2.13	0.48
1:A:177:ILE:HD11	1:A:204:GLN:CD	2.39	0.48
1:A:141:VAL:O	1:A:145:ASN:ND2	2.37	0.47
2:B:339:ASP:OD1	2:B:340:GLN:HG3	2.15	0.47
1:C:431:TRP:CE2	1:C:505:LEU:HD22	2.50	0.47
1:A:23:HIS:HD2	1:A:27:LEU:O	1.96	0.47
1:A:139:LEU:O	1:A:143:LYS:HG3	2.15	0.47
1:C:425:LEU:HD21	1:C:541:VAL:HG11	1.97	0.46
1:A:215:MET:O	1:A:218:ALA:HB3	2.14	0.46
1:A:52:VAL:HG21	1:A:202:ILE:HD12	1.98	0.46
1:C:519:GLN:OE1	1:C:529:PHE:CD1	2.69	0.45
1:A:48:LEU:HA	1:A:49:PRO:HD3	1.73	0.45
1:A:79:VAL:O	1:A:127:ASN:HB2	2.16	0.45
1:C:448:LEU:HA	1:C:449:PRO:HD3	1.83	0.45
1:C:549:PHE:HB2	3:C:824:HOH:O	2.17	0.45
1:A:42:PRO:CG	1:A:43:GLU:H	2.29	0.44
1:A:119:GLN:HE22	1:A:133:THR:HG21	1.80	0.44
2:B:356:LEU:HD22	2:B:361:ILE:HD12	1.99	0.44
1:A:87:LYS:NZ	1:A:160:THR:O	2.50	0.44
1:C:458:LEU:HA	1:C:566:HIS:O	2.17	0.44
2:B:333:LYS:HB3	2:B:333:LYS:HE2	1.69	0.44
1:C:451:PRO:CG	1:C:453:LYS:HE2	2.48	0.43
1:C:529:PHE:CB	1:C:534:THR:HG21	2.46	0.43
1:A:39:LEU:HD22	1:A:228:MET:SD	2.59	0.43
1:A:37:TYR:CE1	1:A:44:LEU:HD22	2.52	0.42
1:A:156:ALA:HA	1:A:157:PRO:HD3	1.85	0.42
1:A:137:PHE:O	1:A:141:VAL:HG23	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:304:PHE:HA	2:B:313:ILE:O	2.20	0.42
1:A:233:PRO:O	1:A:234:ASN:HB3	2.20	0.42
1:C:460:PRO:HG2	1:C:623:VAL:HG11	2.02	0.41
1:C:450:ARG:HH11	1:C:600:ASP:CG	2.28	0.41
1:C:515:PHE:CZ	1:C:534:THR:HG22	2.48	0.41
1:A:41:GLU:HG2	1:A:44:LEU:CD1	2.43	0.41
1:A:83:LYS:NZ	1:A:185:LEU:CD2	2.76	0.41
1:A:195:ASP:HA	1:A:196:PRO:HD2	1.72	0.41
2:B:342:ARG:HD3	2:B:349:GLN:OE1	2.21	0.41
1:A:40:THR:CG2	1:C:608:ARG:CZ	2.99	0.41
2:B:323:ILE:HG22	2:B:327:LYS:HE3	2.01	0.41
1:A:40:THR:OG1	1:C:440:THR:HG21	2.20	0.40
1:C:408:VAL:HG11	1:C:549:PHE:CE2	2.56	0.40
1:C:484:GLN:HA	1:C:493:TYR:CZ	2.56	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	210/235 (89%)	191 (91%)	14 (7%)	5 (2%)	4	2
1	C	206/235 (88%)	193 (94%)	9 (4%)	4 (2%)	6	3
2	B	74/76 (97%)	72 (97%)	2 (3%)	0	100	100
2	D	11/76 (14%)	9 (82%)	2 (18%)	0	100	100
All	All	501/622 (80%)	465 (93%)	27 (5%)	9 (2%)	6	3

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	40	THR

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Mol	Chain	Res	Type
1	C	440	THR
1	C	524	SER
1	C	527	ASN
1	A	61	ILE
1	A	124	SER
1	A	127	ASN
1	A	42	PRO
1	C	442	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	189/209 (90%)	173 (92%)	16 (8%)	10	8
1	C	185/209 (88%)	174 (94%)	11 (6%)	18	18
2	B	68/68 (100%)	64 (94%)	4 (6%)	18	18
2	D	14/68 (21%)	14 (100%)	0	100	100
All	All	456/554 (82%)	425 (93%)	31 (7%)	14	14

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	VAL
1	A	20	ASN
1	A	38	SER
1	A	61	ILE
1	A	79	VAL
1	A	83	LYS
1	A	125	SER
1	A	126	LYS
1	A	128	ARG
1	A	130	ASP
1	A	131	ASP
1	A	136	GLN
1	A	146	VAL

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Mol	Chain	Res	Type
1	A	163	THR
1	A	197	THR
1	A	222	ASP
2	B	324	GLU
2	B	342	ARG
2	B	354	ARG
2	B	363	LYS
1	C	438	SER
1	C	479	VAL
1	C	483	LYS
1	C	526	LYS
1	C	528	ARG
1	C	531	ASP
1	C	536	GLN
1	C	540	ASN
1	C	585	LEU
1	C	586	SER
1	C	594	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	119	GLN
2	B	340	GLN
1	C	420	ASN
1	C	423	HIS
1	C	519	GLN
1	C	617	ASN
2	D	740	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	GLZ	B	376	2,1	3,3,3	0.84	0	1,2,2	0.53	0
2	GLZ	D	776	2,1	3,3,3	0.78	0	1,2,2	0.71	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
2	GLZ	B	376	2,1	-	0/0/1/1	-
2	GLZ	D	776	2,1	-	0/0/1/1	-

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.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	214/235 (91%)	0.45	18 (8%) 17 15	23, 48, 92, 100	2 (0%)
1	C	210/235 (89%)	0.97	40 (19%) 3 3	25, 56, 100, 100	5 (2%)
2	B	75/76 (98%)	-0.11	0 100 100	16, 38, 57, 63	2 (2%)
2	D	16/76 (21%)	1.48	5 (31%) 1 0	39, 62, 86, 92	1 (6%)
All	All	515/622 (82%)	0.61	63 (12%) 8 8	16, 49, 98, 100	10 (1%)

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	522	THR	6.3
1	A	39	LEU	6.0
1	A	40	THR	5.8
1	C	552	GLY	5.2
1	A	123	SER	5.2
1	C	521	ASP	4.4
1	C	440	THR	4.2
1	C	524	SER	4.1
2	D	743	LEU	3.9
1	C	563	THR	3.8
1	C	439	LEU	3.7
1	C	560	THR	3.6
1	C	551	THR	3.5
1	C	618	ALA	3.3
1	C	446	ALA	3.2
1	A	78	ASP	3.2
1	C	559	ALA	3.1
1	C	408	VAL	3.1
1	C	539	LEU	3.1
1	A	172	GLU	3.0
1	C	525	SER	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	125	SER	2.9
1	C	527	ASN	2.9
1	C	565	LEU	2.8
1	C	529	PHE	2.8
1	C	523	SER	2.8
1	C	407	ALA	2.8
1	C	564	ASN	2.7
1	A	124	SER	2.7
1	C	518	SER	2.7
1	C	485	SER	2.7
1	C	486	VAL	2.6
1	C	585	LEU	2.6
1	C	598	ALA	2.5
1	A	62	ASN	2.5
1	A	122	THR	2.4
1	A	126	LYS	2.4
1	C	514	ASN	2.4
1	A	42	PRO	2.3
1	C	589	LEU	2.3
1	C	596	PRO	2.3
2	D	710	GLY	2.3
1	C	532	VAL	2.3
1	C	438	SER	2.3
1	C	535	ASP	2.3
1	A	195	ASP	2.2
1	C	442	PRO	2.2
1	A	43	GLU	2.2
1	C	561	ALA	2.2
1	C	531	ASP	2.2
2	D	773	LEU	2.1
1	C	443	GLU	2.1
1	C	634	ASN	2.1
1	A	166	HIS	2.1
1	C	515	PHE	2.1
1	A	41	GLU	2.1
1	C	624	LEU	2.1
2	D	770	VAL	2.1
1	A	6	ARG	2.1
2	D	769	LEU	2.0
1	A	234	ASN	2.0
1	C	619	ASN	2.0
1	A	61	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
2	GLZ	D	776	4/4	0.84	0.11	57,61,62,66	0
2	GLZ	B	376	4/4	0.95	0.07	26,32,32,32	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.