



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:06 PM UTC

PDB ID : 5ZQ3 / pdb_00005zq3
Title : PDE-Ubiquitin
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Deposited on : 2018-04-17
Resolution : 2.18 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
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

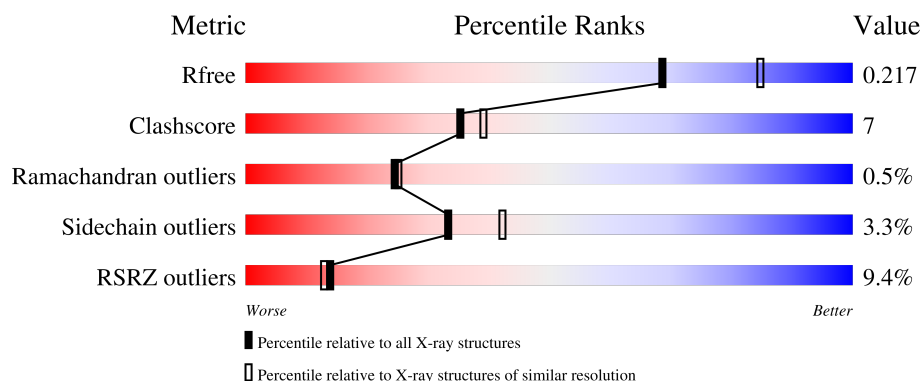
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.18 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	8975 (2.20-2.16)
Clashscore	190562	9786 (2.20-2.16)
Ramachandran outliers	187476	9664 (2.20-2.16)
Sidechain outliers	187428	9664 (2.20-2.16)
RSRZ outliers	180081	8979 (2.20-2.16)

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	377	6% 77% 14% • 7%
1	B	377	4% 82% 10% • 7%
2	C	78	5% 76% 18% • 5%
2	D	78	51% 56% 38% 5%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7209 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 Side.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	349	Total	C	N	O	S	0	0	0
			2839	1797	492	543	7			
1	B	350	Total	C	N	O	S	0	0	0
			2844	1799	493	545	7			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	221	MET	-	initiating methionine	UNP Q6RCR1
A	590	LEU	-	expression tag	UNP Q6RCR1
A	591	GLU	-	expression tag	UNP Q6RCR1
A	592	HIS	-	expression tag	UNP Q6RCR1
A	593	HIS	-	expression tag	UNP Q6RCR1
A	594	HIS	-	expression tag	UNP Q6RCR1
A	595	HIS	-	expression tag	UNP Q6RCR1
A	596	HIS	-	expression tag	UNP Q6RCR1
A	597	HIS	-	expression tag	UNP Q6RCR1
B	221	MET	-	initiating methionine	UNP Q6RCR1
B	590	LEU	-	expression tag	UNP Q6RCR1
B	591	GLU	-	expression tag	UNP Q6RCR1
B	592	HIS	-	expression tag	UNP Q6RCR1
B	593	HIS	-	expression tag	UNP Q6RCR1
B	594	HIS	-	expression tag	UNP Q6RCR1
B	595	HIS	-	expression tag	UNP Q6RCR1
B	596	HIS	-	expression tag	UNP Q6RCR1
B	597	HIS	-	expression tag	UNP Q6RCR1

- Molecule 2 is a protein called Ubiquitin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	74	Total	C	N	O	S	0	0	0
			593	374	103	115	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	74	Total	C	N	O	S	0	0	0
			593	374	103	115	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	GLY	-	expression tag	UNP P0CG48
C	0	SER	-	expression tag	UNP P0CG48
D	-1	GLY	-	expression tag	UNP P0CG48
D	0	SER	-	expression tag	UNP P0CG48

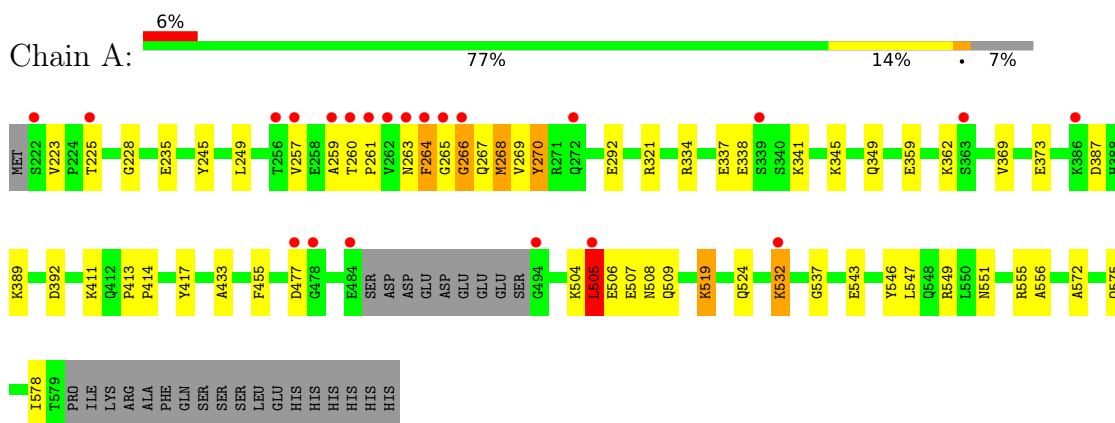
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	124	Total	O	0	0
			124	124		
3	B	188	Total	O	0	0
			188	188		
3	C	18	Total	O	0	0
			18	18		
3	D	10	Total	O	0	0
			10	10		

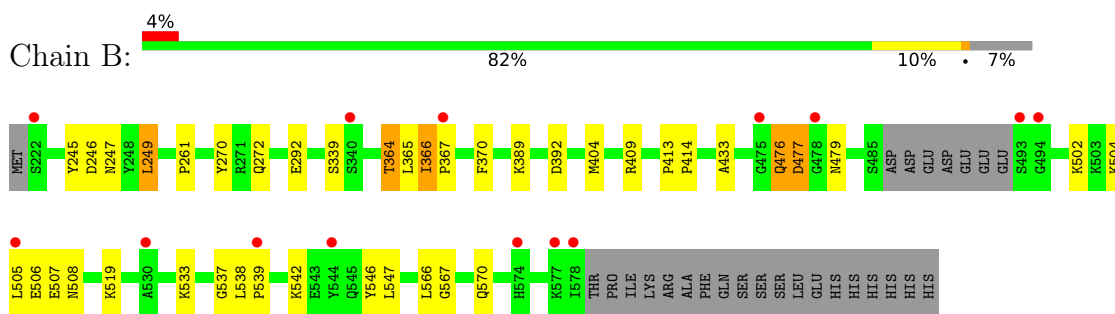
3 Residue-property plots

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.

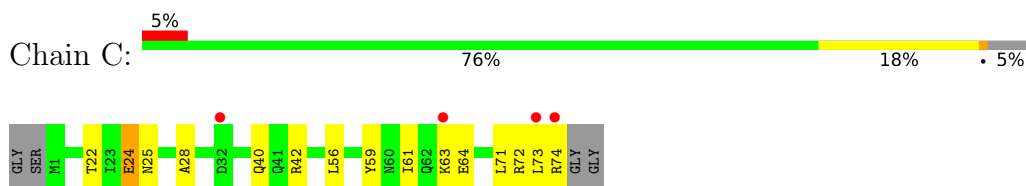
• Molecule 1: SidE



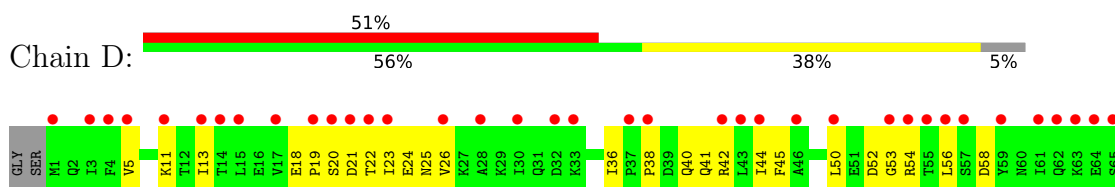
• Molecule 1: SidE

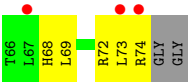


• Molecule 2: Ubiquitin



• Molecule 2: Ubiquitin





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	96.63Å 55.89Å 99.75Å 90.00° 110.50° 90.00°	Depositor
Resolution (Å)	49.11 – 2.18 49.11 – 2.18	Depositor EDS
% Data completeness (in resolution range)	92.5 (49.11-2.18) 92.5 (49.11-2.18)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.18Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.188 , 0.213 0.193 , 0.217	Depositor DCC
R_{free} test set	2405 reflections (4.59%)	wwPDB-VP
Wilson B-factor (Å ²)	28.3	Xtriage
Anisotropy	0.051	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 50.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.014 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7209	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.92% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.40	0/2904	0.76	4/3921 (0.1%)
1	B	0.53	3/2909 (0.1%)	0.72	3/3927 (0.1%)
2	C	0.37	0/599	0.72	0/806
2	D	0.54	0/599	0.93	5/806 (0.6%)
All	All	0.47	3/7011 (0.0%)	0.76	12/9460 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	366	ILE	C-O	-6.48	1.18	1.23
1	B	246	ASP	C-O	-5.16	1.18	1.24
1	B	249	LEU	C-O	-5.00	1.17	1.24

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	365	LEU	N-CA-C	8.35	120.46	111.36
1	A	228	GLY	N-CA-C	-6.91	105.70	115.30
2	D	53	GLY	N-CA-C	-6.61	105.66	114.25
1	A	268	MET	N-CA-C	6.47	120.05	109.76
1	B	247	ASN	N-CA-C	6.20	121.00	113.38
1	B	477	ASP	N-CA-C	-5.79	101.71	110.28
1	A	505	LEU	N-CA-C	5.65	119.18	110.42
2	D	56	LEU	N-CA-C	5.27	116.72	110.97
1	A	270	TYR	N-CA-C	5.22	116.78	111.14
2	D	18	GLU	CA-C-N	5.22	126.36	119.84
2	D	18	GLU	C-N-CA	5.22	126.36	119.84
2	D	21	ASP	N-CA-C	-5.11	102.48	110.10

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	2839	0	2742	51	0
1	B	2844	0	2745	21	0
2	C	593	0	623	11	0
2	D	593	0	623	28	0
3	A	124	0	0	2	0
3	B	188	0	0	4	0
3	C	18	0	0	1	0
3	D	10	0	0	1	0
All	All	7209	0	6733	100	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 7.

All (100) 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:264:PHE:CE2	2:D:73:LEU:CD1	2.56	0.89
2:D:40:GLN:NE2	2:D:74:ARG:HB3	1.88	0.87
2:D:38:PRO:HA	2:D:41:GLN:HE21	1.42	0.81
2:D:42:ARG:HH12	2:D:72:ARG:NH1	1.80	0.79
1:A:264:PHE:CE2	2:D:73:LEU:HD13	2.19	0.77
1:A:264:PHE:CZ	2:D:73:LEU:CD1	2.68	0.77
2:D:40:GLN:HE21	2:D:74:ARG:HB3	1.50	0.74
1:A:264:PHE:CZ	2:D:73:LEU:HD12	2.25	0.72
2:C:73:LEU:HG	2:C:74:ARG:HG2	1.73	0.71
1:A:264:PHE:CB	1:A:556:ALA:HB2	2.20	0.70
2:D:5:VAL:HB	2:D:13:ILE:HB	1.74	0.68
1:A:264:PHE:C	1:A:266:GLY:H	2.02	0.67
2:D:22:THR:OG1	2:D:25:ASN:ND2	2.27	0.67
1:B:272:GLN:OE1	2:C:72:ARG:NH2	2.29	0.65
1:A:264:PHE:CD1	1:A:269:VAL:HG21	2.31	0.65
1:B:477:ASP:N	1:B:477:ASP:OD1	2.28	0.64
1:A:551:ASN:HD21	2:D:74:ARG:H	1.46	0.64
1:A:506:GLU:HG2	1:A:507:GLU:N	2.12	0.63
2:D:42:ARG:NH1	2:D:72:ARG:NH1	2.47	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:519:LYS:HA	1:A:524:GLN:HE21	1.63	0.62
2:D:38:PRO:HA	2:D:41:GLN:NE2	2.15	0.62
1:A:264:PHE:CD1	1:A:269:VAL:CG2	2.83	0.61
1:A:549:ARG:NH1	1:B:570:GLN:OE1	2.33	0.61
1:B:364:THR:HG21	3:B:780:HOH:O	1.99	0.61
1:A:264:PHE:O	1:A:266:GLY:N	2.34	0.60
2:C:42:ARG:CZ	3:C:104:HOH:O	2.50	0.59
1:B:507:GLU:O	1:B:508:ASN:HB2	2.00	0.59
1:A:263:ASN:OD1	1:A:268:MET:SD	2.64	0.56
1:A:264:PHE:HB2	1:A:556:ALA:HB2	1.88	0.56
2:D:42:ARG:HH12	2:D:72:ARG:HH12	1.50	0.55
1:A:264:PHE:C	1:A:266:GLY:N	2.64	0.55
1:A:506:GLU:HG2	1:A:507:GLU:O	2.06	0.55
1:B:538:LEU:HD12	1:B:539:PRO:HD2	1.89	0.55
1:A:506:GLU:HB3	1:A:509:GLN:CD	2.33	0.54
1:A:373:GLU:H	1:A:373:GLU:CD	2.16	0.54
1:A:345:LYS:O	1:A:349:GLN:HG3	2.07	0.54
1:B:476:GLN:HG3	1:B:477:ASP:N	2.23	0.53
1:A:264:PHE:HD1	1:A:269:VAL:CG2	2.21	0.53
1:A:334:ARG:HH21	1:A:337:GLU:HA	1.74	0.53
1:A:387:ASP:HB3	1:A:389:LYS:CE	2.39	0.53
2:D:23:ILE:HA	2:D:26:VAL:HG12	1.90	0.53
2:C:56:LEU:HD22	2:C:61:ILE:HD12	1.91	0.52
2:D:24:GLU:CD	2:D:52:ASP:HB3	2.35	0.51
1:A:264:PHE:CE2	2:D:73:LEU:HD11	2.43	0.51
1:A:505:LEU:HD12	1:A:505:LEU:N	2.26	0.51
2:D:40:GLN:NE2	2:D:74:ARG:CB	2.70	0.50
1:B:537:GLY:HA2	1:B:546:TYR:CZ	2.47	0.50
1:A:257:VAL:HG21	1:A:338:GLU:OE1	2.11	0.50
2:D:68:HIS:ND1	3:D:101:HOH:O	2.25	0.50
2:C:24:GLU:O	2:C:28:ALA:N	2.35	0.49
1:A:455:PHE:CD1	1:A:532:LYS:HE3	2.48	0.49
1:B:542:LYS:N	1:B:542:LYS:HD3	2.28	0.49
2:D:44:ILE:HB	2:D:68:HIS:HB2	1.95	0.49
1:A:321:ARG:HD2	1:A:369:VAL:HG13	1.96	0.48
1:B:477:ASP:O	1:B:479:ASN:N	2.43	0.48
2:D:42:ARG:NH1	2:D:72:ARG:HH12	2.10	0.48
2:C:59:TYR:HB2	2:C:61:ILE:HG13	1.96	0.48
1:A:264:PHE:CE2	1:A:555:ARG:HG2	2.49	0.47
1:B:366:ILE:HA	1:B:367:PRO:HA	1.68	0.47
1:A:261:PRO:HG3	1:A:270:TYR:HD1	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:54:ARG:NH1	2:D:58:ASP:HB3	2.30	0.47
1:A:387:ASP:O	1:A:389:LYS:HE2	2.15	0.46
2:C:63:LYS:HB3	2:C:63:LYS:HE3	1.70	0.46
2:D:41:GLN:HG2	2:D:69:LEU:HD11	1.97	0.46
1:A:572:ALA:HA	3:B:626:HOH:O	2.15	0.46
2:D:20:SER:O	2:D:20:SER:OG	2.24	0.46
1:A:389:LYS:HB3	1:A:392:ASP:CG	2.41	0.45
1:B:292:GLU:HB3	1:B:433:ALA:HB1	1.98	0.45
1:A:292:GLU:HB3	1:A:433:ALA:HB1	1.98	0.45
1:A:359:GLU:OE2	1:A:362:LYS:HE2	2.17	0.45
2:D:45:PHE:HB3	2:D:50:LEU:HD21	1.98	0.45
1:A:575:GLN:NE2	3:A:616:HOH:O	2.50	0.45
1:B:245:TYR:HA	1:B:249:LEU:HB2	1.98	0.45
2:D:54:ARG:CZ	2:D:58:ASP:HB3	2.47	0.45
1:A:387:ASP:HB3	1:A:389:LYS:HE2	1.98	0.45
1:B:547:LEU:HD22	2:C:71:LEU:HB3	1.99	0.44
1:A:551:ASN:ND2	2:D:73:LEU:HD22	2.33	0.44
1:A:245:TYR:HA	1:A:249:LEU:HB2	2.00	0.43
2:C:22:THR:OG1	2:C:25:ASN:OD1	2.23	0.43
2:C:24:GLU:CD	2:C:24:GLU:N	2.76	0.43
1:A:264:PHE:HB3	1:A:556:ALA:HB2	1.99	0.43
1:A:264:PHE:N	1:A:267:GLN:O	2.52	0.43
1:B:261:PRO:HB3	1:B:270:TYR:CD1	2.53	0.43
1:A:413:PRO:HA	1:A:414:PRO:HD3	1.89	0.43
1:A:519:LYS:HE2	1:A:519:LYS:HB3	1.89	0.42
1:B:389:LYS:HB3	1:B:392:ASP:HB2	2.01	0.42
1:A:259:ALA:O	1:A:261:PRO:HD3	2.20	0.42
1:A:235:GLU:HG3	3:A:709:HOH:O	2.19	0.42
1:B:413:PRO:HA	1:B:414:PRO:HD3	1.82	0.41
1:B:567:GLY:HA2	1:B:570:GLN:HG2	2.02	0.41
1:A:477:ASP:N	1:A:477:ASP:OD1	2.53	0.41
1:A:411:LYS:HB3	1:A:417:TYR:CD2	2.56	0.41
1:A:387:ASP:HB3	1:A:389:LYS:HE3	2.02	0.41
1:A:519:LYS:HA	1:A:524:GLN:NE2	2.33	0.41
1:B:409:ARG:NE	3:B:607:HOH:O	2.37	0.41
2:C:40:GLN:HE21	2:C:73:LEU:HD12	1.85	0.41
1:A:537:GLY:HA2	1:A:546:TYR:CZ	2.55	0.40
1:B:519:LYS:HE3	3:B:769:HOH:O	2.20	0.40
1:A:543:GLU:HG2	2:D:11:LYS:CD	2.51	0.40
1:B:367:PRO:HA	1:B:370:PHE:O	2.22	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	345/377 (92%)	331 (96%)	11 (3%)	3 (1%)	14	12
1	B	346/377 (92%)	336 (97%)	10 (3%)	0	100	100
2	C	72/78 (92%)	72 (100%)	0	0	100	100
2	D	72/78 (92%)	68 (94%)	3 (4%)	1 (1%)	9	6
All	All	835/910 (92%)	807 (97%)	24 (3%)	4 (0%)	24	25

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	264	PHE
1	A	265	GLY
2	D	19	PRO
1	A	266	GLY

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	300/327 (92%)	289 (96%)	11 (4%)	30	38
1	B	301/327 (92%)	291 (97%)	10 (3%)	33	42
2	C	68/69 (99%)	66 (97%)	2 (3%)	37	47
2	D	68/69 (99%)	67 (98%)	1 (2%)	57	70
All	All	737/792 (93%)	713 (97%)	24 (3%)	33	42

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	223	VAL
1	A	225	THR
1	A	260	THR
1	A	341	LYS
1	A	504	LYS
1	A	505	LEU
1	A	508	ASN
1	A	519	LYS
1	A	532	LYS
1	A	547	LEU
1	A	578	ILE
1	B	339	SER
1	B	364	THR
1	B	404	MET
1	B	476	GLN
1	B	502	LYS
1	B	504	LYS
1	B	505	LEU
1	B	506	GLU
1	B	533	LYS
1	B	566	LEU
2	C	24	GLU
2	C	64	GLU
2	D	36	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	247	ASN
1	A	412	GLN
1	A	463	HIS
1	A	508	ASN
1	A	524	GLN
1	A	551	ASN
1	A	570	GLN
1	B	443	GLN
2	D	25	ASN

5.3.3 RNA [i](#)

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 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 [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	349/377 (92%)	0.40	22 (6%) 26 25	17, 46, 91, 122	0
1	B	350/377 (92%)	0.05	14 (4%) 42 41	16, 36, 78, 114	0
2	C	74/78 (94%)	0.61	4 (5%) 31 30	30, 56, 82, 93	0
2	D	74/78 (94%)	2.12	40 (54%) 0 0	58, 95, 123, 128	0
All	All	847/910 (93%)	0.43	80 (9%) 14 13	16, 44, 99, 128	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	265	GLY	11.2
1	A	264	PHE	8.2
2	D	28	ALA	4.6
2	D	23	ILE	4.4
2	D	56	LEU	4.3
2	D	17	VAL	4.2
2	D	1	MET	4.1
2	D	61	ILE	3.9
1	A	494	GLY	3.9
2	D	50	LEU	3.6
2	C	73	LEU	3.6
2	D	13	ILE	3.5
2	D	74	ARG	3.5
2	D	43	LEU	3.4
1	A	263	ASN	3.3
2	D	20	SER	3.3
2	D	65	SER	3.3
2	D	57	SER	3.2
2	D	73	LEU	3.1
2	D	11	LYS	3.1
1	B	530	ALA	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	367	PRO	3.1
2	D	15	LEU	3.1
1	B	539	PRO	3.0
1	B	222	SER	3.0
2	D	19	PRO	2.9
1	A	363	SER	2.9
1	B	505	LEU	2.8
2	D	37	PRO	2.8
1	B	493	SER	2.8
1	A	261	PRO	2.8
1	B	475	GLY	2.7
1	A	386	LYS	2.7
2	D	14	THR	2.7
1	A	266	GLY	2.7
2	D	26	VAL	2.6
1	B	577	LYS	2.6
2	D	59	TYR	2.6
1	B	578	ILE	2.6
2	D	67	LEU	2.6
1	B	340	SER	2.5
2	D	4	PHE	2.5
1	A	257	VAL	2.5
1	A	256	THR	2.5
1	A	477	ASP	2.5
1	A	225	THR	2.5
1	A	339	SER	2.5
2	D	33	LYS	2.5
1	A	484	GLU	2.4
2	D	63	LYS	2.4
2	D	21	ASP	2.4
1	A	262	VAL	2.4
1	A	478	GLY	2.3
1	B	574	HIS	2.3
2	D	38	PRO	2.3
2	C	74	ARG	2.3
1	A	505	LEU	2.3
1	B	544	TYR	2.3
2	D	55	THR	2.2
2	D	53	GLY	2.2
1	A	222	SER	2.2
2	D	42	ARG	2.2
2	D	3	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	532	LYS	2.2
1	B	478	GLY	2.2
1	A	259	ALA	2.2
2	D	22	THR	2.2
2	D	30	ILE	2.1
2	D	62	GLN	2.1
2	D	32	ASP	2.1
1	A	260	THR	2.1
1	A	272	GLN	2.1
2	D	44	ILE	2.1
2	D	54	ARG	2.1
2	C	32	ASP	2.1
2	D	5	VAL	2.1
2	C	63	LYS	2.1
2	D	46	ALA	2.1
2	D	64	GLU	2.0
1	B	494	GLY	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.