



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 09:15 AM UTC

PDB ID : 9WQU / pdb_00009wqu
Title : b-b' domain fragment of ER-60 (ERp57) under microgravity
Authors : Furubayashi, N.; Shinohara, Y.; Inaka, K.; Kaito, S.; Kobayashi, Y.; Kamo, M.; Urade, R.
Deposited on : 2025-09-11
Resolution : 1.80 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

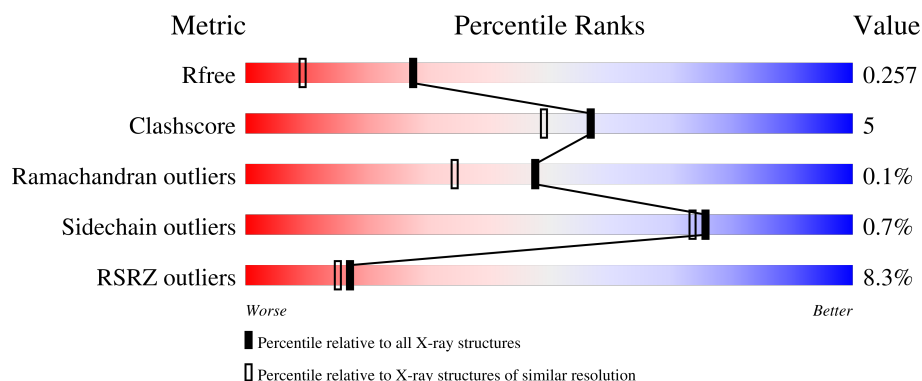
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	252	6% 80% 12% 8%
1	B	252	6% 79% 13% 7%
1	C	252	11% 79% 12% 8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6220 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 disulfide-isomerase A3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	233	Total	C	N	O	S	0	0	0
			1892	1204	320	362	6			
1	B	234	Total	C	N	O	S	0	0	0
			1898	1207	321	364	6			
1	C	232	Total	C	N	O	S	0	1	0
			1886	1200	318	362	6			

There are 27 discrepancies between the modelled and reference sequences:

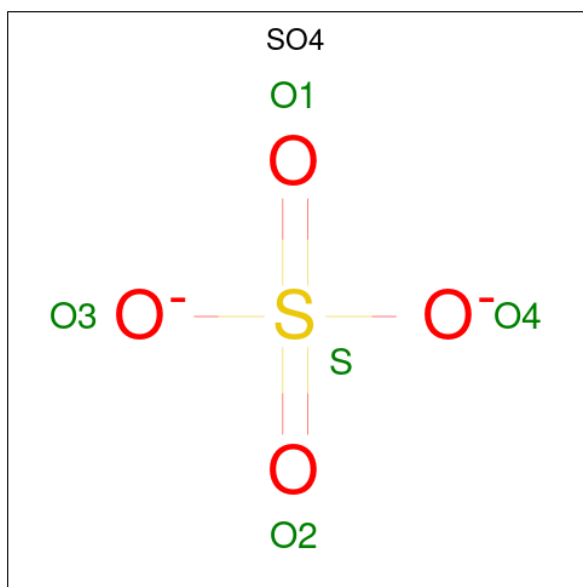
Chain	Residue	Modelled	Actual	Comment	Reference
A	129	GLY	-	expression tag	UNP P30101
A	130	PRO	-	expression tag	UNP P30101
A	131	LEU	-	expression tag	UNP P30101
A	132	GLY	-	expression tag	UNP P30101
A	133	SER	-	expression tag	UNP P30101
A	377	ALA	-	cloning artifact	UNP P30101
A	378	ALA	-	cloning artifact	UNP P30101
A	379	ALA	-	cloning artifact	UNP P30101
A	380	SER	-	cloning artifact	UNP P30101
B	129	GLY	-	expression tag	UNP P30101
B	130	PRO	-	expression tag	UNP P30101
B	131	LEU	-	expression tag	UNP P30101
B	132	GLY	-	expression tag	UNP P30101
B	133	SER	-	expression tag	UNP P30101
B	377	ALA	-	cloning artifact	UNP P30101
B	378	ALA	-	cloning artifact	UNP P30101
B	379	ALA	-	cloning artifact	UNP P30101
B	380	SER	-	cloning artifact	UNP P30101
C	129	GLY	-	expression tag	UNP P30101
C	130	PRO	-	expression tag	UNP P30101
C	131	LEU	-	expression tag	UNP P30101
C	132	GLY	-	expression tag	UNP P30101
C	133	SER	-	expression tag	UNP P30101

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Chain	Residue	Modelled	Actual	Comment	Reference
C	377	ALA	-	cloning artifact	UNP P30101
C	378	ALA	-	cloning artifact	UNP P30101
C	379	ALA	-	cloning artifact	UNP P30101
C	380	SER	-	cloning artifact	UNP P30101

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

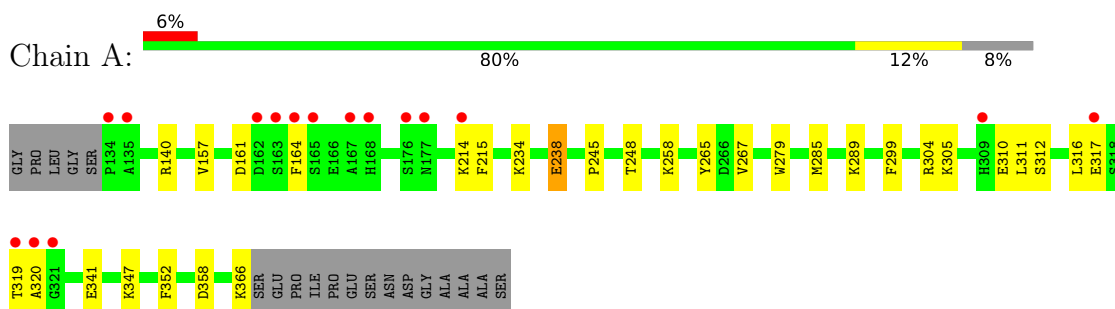
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	179	Total	O	0	0
			179	179		
3	B	185	Total	O	0	0
			185	185		
3	C	170	Total	O	0	0
			170	170		

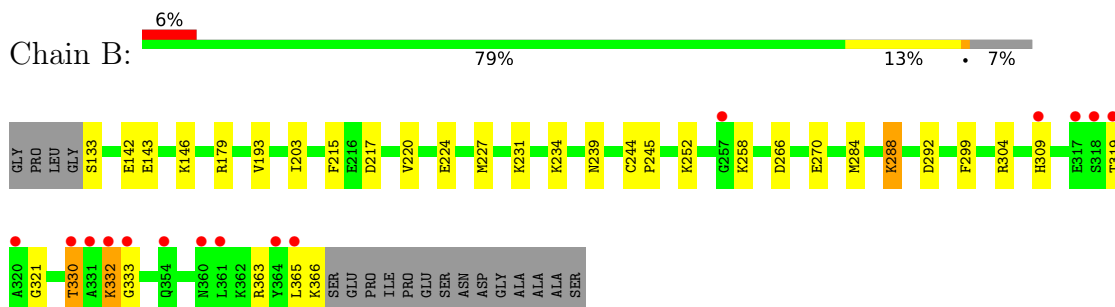
3 Residue-property plots [i](#)

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.

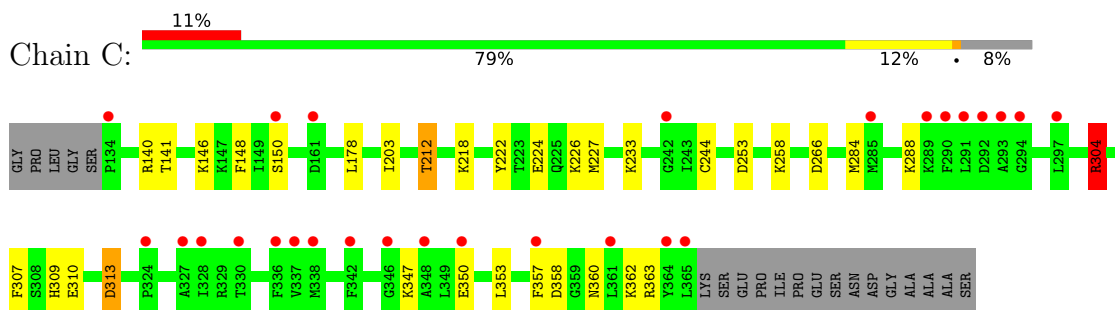
• Molecule 1: Protein disulfide-isomerase A3



• Molecule 1: Protein disulfide-isomerase A3



• Molecule 1: Protein disulfide-isomerase A3



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.31Å 62.90Å 97.15Å 90.00° 98.64° 90.00°	Depositor
Resolution (Å)	47.78 – 1.80 47.78 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.6 (47.78-1.80) 99.7 (47.78-1.80)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.8.0430	Depositor
R, R_{free}	0.219 , 0.263 0.214 , 0.257	Depositor DCC
R_{free} test set	4065 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	31.8	Xtriage
Anisotropy	0.015	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 32.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6220	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.30% 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: SO4

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	0.66	0/1932	1.34	15/2592 (0.6%)
1	B	0.63	0/1938	1.34	10/2601 (0.4%)
1	C	0.65	0/1929	1.36	13/2589 (0.5%)
All	All	0.65	0/5799	1.35	38/7782 (0.5%)

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	C	0	1

There are no bond length outliers.

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	146	LYS	CB-CA-C	8.94	124.91	110.88
1	A	310	GLU	CB-CG-CD	8.69	127.36	112.60
1	B	217	ASP	CB-CA-C	8.13	123.22	109.80
1	B	143	GLU	CB-CG-CD	8.02	126.24	112.60
1	C	304	ARG	CD-NE-CZ	7.47	134.86	124.40
1	B	215	PHE	CA-CB-CG	-7.38	106.42	113.80
1	C	141	THR	CA-CB-OG1	-7.37	98.54	109.60
1	A	358	ASP	CA-CB-CG	6.76	119.36	112.60
1	A	161	ASP	CB-CA-C	6.76	123.25	109.67
1	B	258	LYS	CB-CA-C	6.59	125.01	112.43
1	B	193	VAL	N-CA-CB	6.46	118.11	110.55
1	A	238	GLU	CB-CG-CD	6.38	123.45	112.60
1	B	288	LYS	CB-CG-CD	6.28	125.74	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	140	ARG	CD-NE-CZ	6.02	132.82	124.40
1	C	212	THR	CA-CB-OG1	-5.76	100.96	109.60
1	B	142	GLU	N-CA-CB	5.68	118.58	110.06
1	A	215	PHE	CA-CB-CG	-5.67	108.13	113.80
1	A	164	PHE	CA-CB-CG	-5.65	108.15	113.80
1	C	307	PHE	CA-CB-CG	-5.56	108.24	113.80
1	A	317	GLU	CB-CG-CD	-5.55	103.17	112.60
1	A	157	VAL	N-CA-CB	5.53	118.83	111.64
1	C	357	PHE	CB-CA-C	5.50	119.93	110.79
1	B	330	THR	CA-CB-OG1	-5.48	101.38	109.60
1	B	252	LYS	N-CA-CB	5.48	118.66	110.22
1	C	313	ASP	CA-CB-CG	5.47	118.07	112.60
1	C	304	ARG	CB-CG-CD	5.44	123.81	111.30
1	C	146	LYS	CB-CG-CD	5.43	123.78	111.30
1	C	309	HIS	CA-CB-CG	-5.35	108.45	113.80
1	C	148	PHE	CA-CB-CG	-5.26	108.53	113.80
1	A	352	PHE	CA-CB-CG	5.26	119.06	113.80
1	B	224	GLU	CB-CG-CD	-5.20	103.75	112.60
1	A	317	GLU	N-CA-CB	-5.14	102.55	110.56
1	C	146	LYS	N-CA-CB	-5.13	102.57	110.01
1	A	214	LYS	CB-CA-C	-5.10	99.66	110.31
1	A	248	THR	CA-CB-OG1	-5.07	102.00	109.60
1	C	350	GLU	CB-CA-C	5.03	119.14	110.79
1	A	316	LEU	CA-C-N	-5.01	113.61	122.74
1	A	316	LEU	C-N-CA	-5.01	113.61	122.74

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	304	ARG	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	1892	0	1842	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1898	0	1846	16	0
1	C	1886	0	1834	22	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	179	0	0	3	0
3	B	185	0	0	6	0
3	C	170	0	0	7	0
All	All	6220	0	5522	51	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 5.

All (51) 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:203:ILE:HD11	1:C:227:MET:HE3	1.47	0.94
1:A:289:LYS:HD3	3:A:635:HOH:O	1.76	0.84
1:B:363:ARG:HE	1:B:365:LEU:HD21	1.45	0.81
1:B:266:ASP:HB3	3:B:644:HOH:O	1.80	0.81
1:C:212:THR:HG22	3:C:458:HOH:O	1.81	0.81
1:B:203:ILE:HD11	1:B:227:MET:HE3	1.62	0.79
1:A:304:ARG:HG3	1:A:311:LEU:HD11	1.65	0.78
1:B:366:LYS:HE2	3:B:640:HOH:O	1.91	0.69
1:C:362:LYS:HG2	3:C:532:HOH:O	1.95	0.66
1:B:288:LYS:HE3	1:B:292:ASP:OD2	1.97	0.64
1:A:304:ARG:HG3	1:A:311:LEU:CD1	2.28	0.63
1:C:258:LYS:HG3	3:C:465:HOH:O	1.99	0.62
1:A:347:LYS:HE2	3:A:620:HOH:O	2.01	0.60
1:C:244:CYS:HB3	1:C:284:MET:HG2	1.83	0.58
1:C:266:ASP:OD1	1:C:304:ARG:HD3	2.04	0.57
1:B:245:PRO:HD2	1:B:299:PHE:O	2.05	0.56
1:A:319:THR:HG22	1:C:288:LYS:HE3	1.87	0.55
1:B:146:LYS:HE2	3:B:671:HOH:O	2.06	0.55
1:C:150:SER:OG	1:C:150:SER:O	2.19	0.55
1:C:203:ILE:HD11	1:C:227:MET:CE	2.30	0.54
1:C:203:ILE:CD1	1:C:227:MET:HE3	2.31	0.53
1:B:330:THR:HB	1:B:332:LYS:HD2	1.91	0.52
1:C:140:ARG:HH11	1:C:140:ARG:HG3	1.75	0.52
1:B:133:SER:OG	1:B:179:ARG:NH2	2.43	0.51
1:B:304:ARG:CZ	1:B:321:GLY:HA2	2.42	0.49
1:C:358:ASP:HB2	1:C:360:ASN:OD1	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:305:LYS:HE3	1:A:320:ALA:HA	1.95	0.48
1:C:253:ASP:HA	3:C:489:HOH:O	2.12	0.48
1:B:244:CYS:HB2	1:B:284:MET:HG2	1.95	0.48
1:C:347:LYS:NZ	3:C:406:HOH:O	2.48	0.47
1:C:258:LYS:HA	1:C:258:LYS:HD2	1.63	0.46
1:C:310:GLU:O	1:C:313:ASP:HB2	2.16	0.46
1:A:265:TYR:CZ	1:A:267:VAL:HG22	2.51	0.45
1:B:234:LYS:HE3	3:B:602:HOH:O	2.17	0.45
1:B:220:VAL:HG11	1:B:239:ASN:HB3	2.01	0.43
1:C:222:TYR:CE1	1:C:224:GLU:HB2	2.54	0.42
1:C:363:ARG:HG2	3:C:532:HOH:O	2.18	0.42
1:A:258:LYS:HB3	1:A:258:LYS:HE2	1.57	0.42
1:B:203:ILE:HD11	1:B:227:MET:CE	2.40	0.42
1:C:178:LEU:HD11	1:C:233:LYS:HB2	2.02	0.42
1:B:231:LYS:NZ	3:B:505:HOH:O	2.52	0.41
1:C:353:LEU:HD23	1:C:353:LEU:HA	1.94	0.41
1:A:265:TYR:CE2	1:A:279:TRP:HB2	2.56	0.41
1:B:309:HIS:CD2	3:B:651:HOH:O	2.72	0.41
1:C:218:LYS:O	1:C:218:LYS:HG2	2.21	0.41
1:C:226:LYS:HD2	3:C:494:HOH:O	2.20	0.41
1:A:285:MET:HE2	1:A:285:MET:HB3	2.01	0.41
1:A:234:LYS:HE2	3:A:535:HOH:O	2.20	0.40
1:A:234:LYS:HE3	1:A:238:GLU:OE2	2.21	0.40
1:A:245:PRO:HD2	1:A:299:PHE:O	2.21	0.40
1:A:312:SER:O	1:A:366:LYS:HD2	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	231/252 (92%)	225 (97%)	6 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	232/252 (92%)	222 (96%)	9 (4%)	1 (0%)	30	19
1	C	231/252 (92%)	224 (97%)	7 (3%)	0	100	100
All	All	694/756 (92%)	671 (97%)	22 (3%)	1 (0%)	48	34

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	333	GLY

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	203/216 (94%)	202 (100%)	1 (0%)	81	80
1	B	204/216 (94%)	201 (98%)	3 (2%)	57	49
1	C	203/216 (94%)	203 (100%)	0	100	100
All	All	610/648 (94%)	606 (99%)	4 (1%)	76	73

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	341	GLU
1	B	270	GLU
1	B	319	THR
1	B	332	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 oligosaccharides 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	SO4	A	401	-	4,4,4	0.28	0	6,6,6	0.12	0
2	SO4	B	401	-	4,4,4	0.33	0	6,6,6	0.10	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

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	233/252 (92%)	0.53	16 (6%) 23 21	18, 32, 51, 78	1 (0%)
1	B	234/252 (92%)	0.43	15 (6%) 25 24	21, 32, 57, 76	1 (0%)
1	C	232/252 (92%)	0.90	27 (11%) 9 8	20, 34, 53, 62	1 (0%)
All	All	699/756 (92%)	0.62	58 (8%) 17 15	18, 33, 53, 78	3 (0%)

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	320	ALA	4.5
1	C	365	LEU	4.5
1	B	331	ALA	4.4
1	C	290	PHE	4.3
1	A	134	PRO	4.1
1	B	309	HIS	3.9
1	B	333	GLY	3.7
1	C	291	LEU	3.6
1	C	285	MET	3.5
1	C	327	ALA	3.5
1	C	348	ALA	3.1
1	B	365	LEU	3.0
1	B	332	LYS	3.0
1	A	176	SER	3.0
1	B	318	SER	2.9
1	A	321	GLY	2.9
1	C	289	LYS	2.9
1	B	361	LEU	2.9
1	C	357	PHE	2.8
1	A	164	PHE	2.8
1	C	336	PHE	2.8
1	B	360	ASN	2.7
1	C	328	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	294	GLY	2.6
1	C	297	LEU	2.6
1	C	161	ASP	2.6
1	C	293	ALA	2.6
1	C	150	SER	2.6
1	B	354	GLN	2.6
1	B	320	ALA	2.6
1	B	257	GLY	2.5
1	A	135	ALA	2.5
1	A	317	GLU	2.5
1	A	319	THR	2.5
1	B	319	THR	2.5
1	C	134	PRO	2.5
1	C	364	TYR	2.4
1	A	167	ALA	2.4
1	C	361	LEU	2.4
1	C	292	ASP	2.4
1	A	177	ASN	2.3
1	A	309	HIS	2.3
1	C	330	THR	2.3
1	B	364	TYR	2.3
1	A	163	SER	2.3
1	B	317	GLU	2.3
1	C	324	PRO	2.2
1	C	350	GLU	2.2
1	A	168	HIS	2.2
1	B	330	THR	2.2
1	C	338	MET	2.2
1	A	162	ASP	2.2
1	C	346	GLY	2.2
1	C	342	PHE	2.1
1	C	337	VAL	2.1
1	C	242	GLY	2.0
1	A	214	LYS	2.0
1	A	165	SER	2.0

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

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
2	SO4	A	401	5/5	0.89	0.10	45,54,63,70	0
2	SO4	B	401	5/5	0.94	0.10	41,51,56,57	0

6.5 Other polymers [i](#)

There are no such residues in this entry.