



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 08:39 AM UTC

PDB ID : 8XWX / pdb_00008xwx
Title : Crystal structure of FIS1-BAP31 complex from human
Authors : Nguyen, M.D.; Bong, S.M.; Lee, B.I.
Deposited on : 2024-01-17
Resolution : 2.69 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

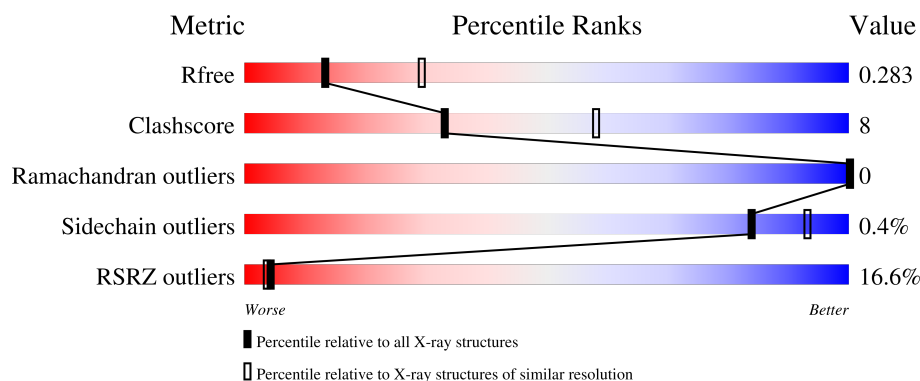
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.69 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (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 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	D	72	10% 79% 11% 10%
1	E	72	24% 69% 11% 19%
1	F	72	8% 78% 14% 8%
1	G	72	14% 64% 15% 21%
2	A	125	4% 81% 15% .

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Mol	Chain	Length	Quality of chain
2	B	125	14% 73% 22% 5%
2	C	125	30% 68% 25% 6%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4763 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 B-cell receptor-associated protein 31.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	58	Total	C	N	O		0	0	0
			399	247	73	79				
1	G	57	Total	C	N	O		0	0	0
			368	226	68	74				
1	D	65	Total	C	N	O	S	0	0	0
			518	318	93	106	1			
1	F	66	Total	C	N	O	S	0	0	0
			523	321	94	107	1			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	166	GLY	-	expression tag	UNP P51572
E	167	SER	-	expression tag	UNP P51572
E	234	SER	-	expression tag	UNP P51572
E	235	ALA	-	expression tag	UNP P51572
E	236	ALA	-	expression tag	UNP P51572
E	237	ALA	-	expression tag	UNP P51572
G	166	GLY	-	expression tag	UNP P51572
G	167	SER	-	expression tag	UNP P51572
G	234	SER	-	expression tag	UNP P51572
G	235	ALA	-	expression tag	UNP P51572
G	236	ALA	-	expression tag	UNP P51572
G	237	ALA	-	expression tag	UNP P51572
D	166	GLY	-	expression tag	UNP P51572
D	167	SER	-	expression tag	UNP P51572
D	234	SER	-	expression tag	UNP P51572
D	235	ALA	-	expression tag	UNP P51572
D	236	ALA	-	expression tag	UNP P51572
D	237	ALA	-	expression tag	UNP P51572
F	166	GLY	-	expression tag	UNP P51572
F	167	SER	-	expression tag	UNP P51572
F	234	SER	-	expression tag	UNP P51572

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Chain	Residue	Modelled	Actual	Comment	Reference
F	235	ALA	-	expression tag	UNP P51572
F	236	ALA	-	expression tag	UNP P51572
F	237	ALA	-	expression tag	UNP P51572

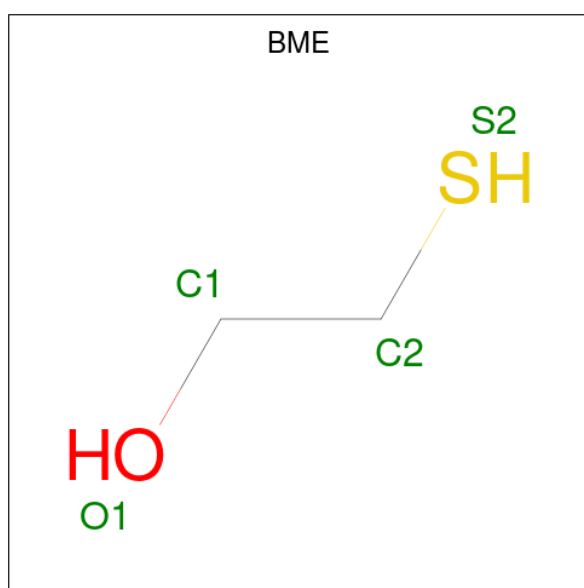
- Molecule 2 is a protein called Mitochondrial fission 1 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	120	Total	C	N	O	S	0	0	0
			988	632	166	187	3			
2	C	117	Total	C	N	O	S	0	0	0
			953	613	158	180	2			
2	B	119	Total	C	N	O	S	0	0	0
			980	627	165	186	2			

There are 6 discrepancies between the modelled and reference sequences:

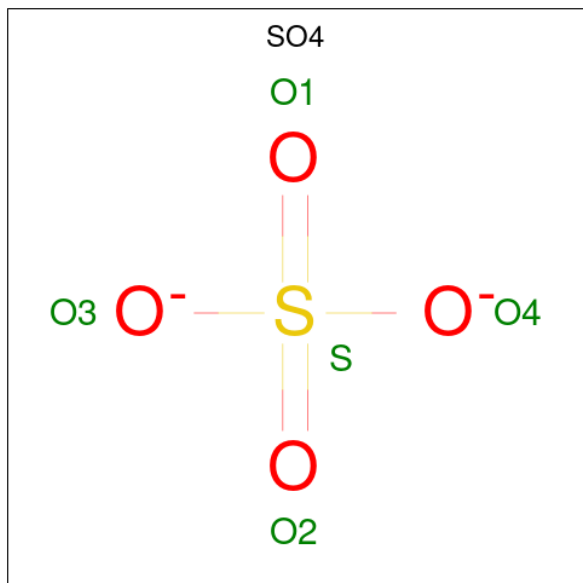
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q9Y3D6
A	0	SER	-	expression tag	UNP Q9Y3D6
C	-1	GLY	-	expression tag	UNP Q9Y3D6
C	0	SER	-	expression tag	UNP Q9Y3D6
B	-1	GLY	-	expression tag	UNP Q9Y3D6
B	0	SER	-	expression tag	UNP Q9Y3D6

- Molecule 3 is BETA-MERCAPTOETHANOL (CCD ID: BME) (formula: C₂H₆OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	S	0	0
			4	2	1	1		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	O	S	0	0
			5	4	1		

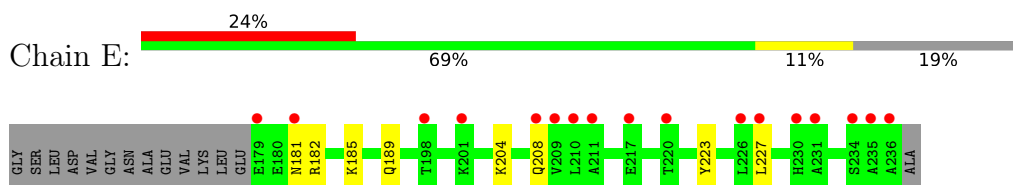
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	10	Total	O	0	0
			10	10		
5	D	3	Total	O	0	0
			3	3		
5	F	3	Total	O	0	0
			3	3		
5	C	3	Total	O	0	0
			3	3		
5	B	6	Total	O	0	0
			6	6		

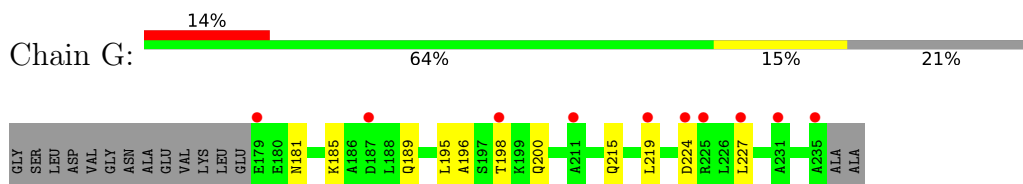
3 Residue-property plots [i](#)

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.

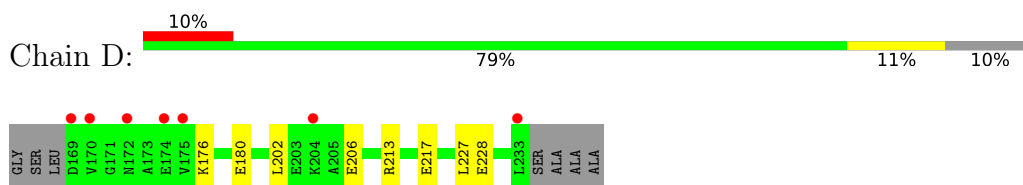
- Molecule 1: B-cell receptor-associated protein 31



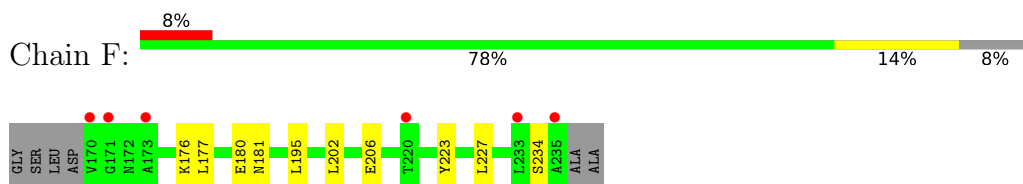
- Molecule 1: B-cell receptor-associated protein 31



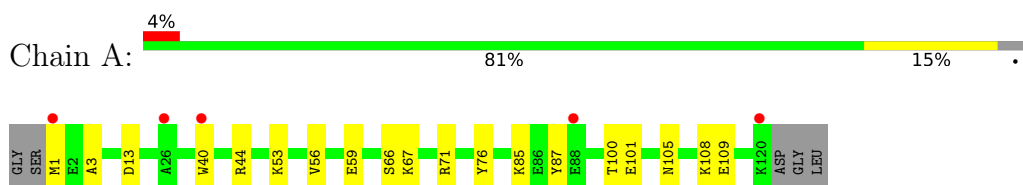
- Molecule 1: B-cell receptor-associated protein 31



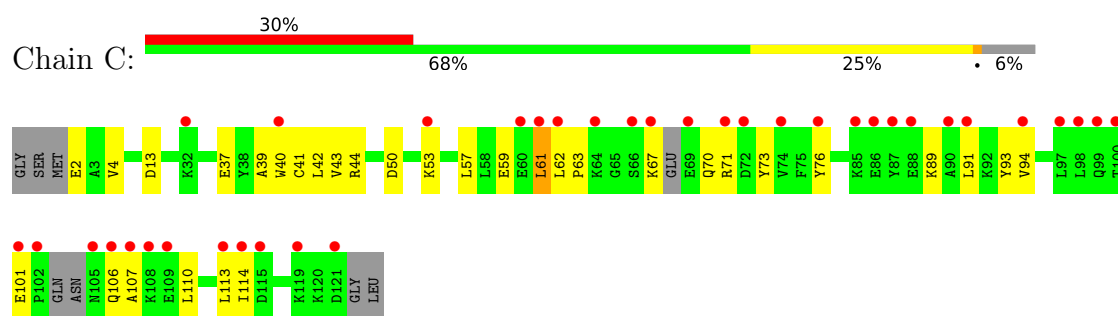
- Molecule 1: B-cell receptor-associated protein 31



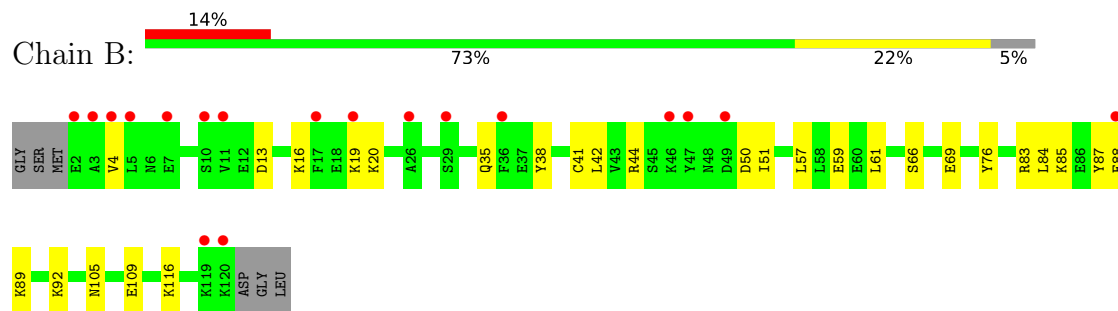
- Molecule 2: Mitochondrial fission 1 protein



- Molecule 2: Mitochondrial fission 1 protein



- Molecule 2: Mitochondrial fission 1 protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	104.63Å 237.27Å 90.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.96 – 2.69 40.96 – 2.69	Depositor EDS
% Data completeness (in resolution range)	85.8 (40.96-2.69) 85.8 (40.96-2.69)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.80 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.21rc1_5156	Depositor
R, R_{free}	0.254 , 0.287 (Not available) , 0.283	Depositor DCC
R_{free} test set	2000 reflections (6.29%)	wwPDB-VP
Wilson B-factor (Å ²)	36.5	Xtriage
Anisotropy	0.238	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 63.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4763	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 4.88% 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: BME, 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	D	0.27	0/519	0.41	0/689
1	E	0.26	0/399	0.29	0/537
1	F	0.27	0/524	0.42	0/696
1	G	0.21	0/368	0.30	0/500
2	A	0.29	0/1002	0.41	0/1340
2	B	0.26	0/994	0.47	0/1330
2	C	0.26	0/965	0.46	0/1289
All	All	0.27	0/4771	0.42	0/6381

There are no bond length outliers.

There are no bond angle outliers.

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	D	518	0	529	5	0
1	E	399	0	376	4	0
1	F	523	0	534	8	0
1	G	368	0	309	7	0
2	A	988	0	1016	11	1
2	B	980	0	1004	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	953	0	973	26	0
3	A	4	0	6	0	0
4	C	5	0	0	0	0
5	A	10	0	0	0	0
5	B	6	0	0	0	0
5	C	3	0	0	0	0
5	D	3	0	0	0	0
5	F	3	0	0	0	0
All	All	4763	0	4747	74	1

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 8.

All (74) 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)
2:A:40:TRP:HE1	2:A:44:ARG:HH21	1.23	0.86
2:B:38:TYR:O	2:B:42:LEU:HD23	1.85	0.76
2:B:35:GLN:HG3	2:B:57:LEU:HD23	1.74	0.69
2:C:76:TYR:OH	2:B:13:ASP:OD1	2.11	0.66
2:C:40:TRP:HE1	2:C:44:ARG:HH11	1.44	0.65
2:C:110:LEU:O	2:C:114:ILE:HG13	1.98	0.63
2:B:16:LYS:HD3	2:B:19:LYS:HD2	1.80	0.63
1:E:182:ARG:HA	1:E:185:LYS:HD3	1.81	0.61
1:F:176:LYS:O	1:F:180:GLU:HG2	2.01	0.61
2:C:101:GLU:OE2	2:B:16:LYS:NZ	2.35	0.59
2:C:44:ARG:NH1	2:B:13:ASP:OD2	2.32	0.58
1:D:202:LEU:O	1:D:206:GLU:HG3	2.04	0.58
2:C:61:LEU:HB3	2:C:73:TYR:CE1	2.38	0.57
1:E:223:TYR:O	1:E:227:LEU:HG	2.03	0.57
2:A:1:MET:HG3	2:A:3:ALA:H	1.69	0.57
2:C:42:LEU:HD13	2:C:53:LYS:HB3	1.87	0.56
2:B:66:SER:H	2:B:69:GLU:HB2	1.71	0.55
2:C:113:LEU:HD11	2:B:4:VAL:HG12	1.88	0.55
1:G:196:ALA:O	1:G:200:GLN:HG2	2.07	0.54
1:F:202:LEU:O	1:F:206:GLU:HG3	2.07	0.54
2:C:89:LYS:O	2:C:93:TYR:HD1	1.89	0.54
2:B:44:ARG:HA	2:B:83:ARG:NH1	2.24	0.53
1:E:204:LYS:O	1:E:208:GLN:HG2	2.09	0.53
1:G:219:LEU:HD23	1:F:223:TYR:OH	2.10	0.52
1:D:176:LYS:O	1:D:180:GLU:HG2	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:62:LEU:O	2:C:70:GLN:NE2	2.26	0.51
2:A:67:LYS:O	2:A:71:ARG:HD2	2.11	0.50
2:A:100:THR:HG23	2:A:101:GLU:HG3	1.93	0.50
2:B:38:TYR:CE1	2:B:42:LEU:HD21	2.47	0.50
2:B:89:LYS:HA	2:B:92:LYS:HE2	1.93	0.50
2:C:39:ALA:HB2	2:C:57:LEU:HB2	1.93	0.48
2:C:91:LEU:O	2:C:94:VAL:HG12	2.14	0.47
1:D:213:ARG:O	1:D:217:GLU:HG2	2.15	0.47
2:B:20:LYS:HE3	2:B:20:LYS:HB2	1.72	0.46
2:C:40:TRP:HA	2:C:43:VAL:HG22	1.97	0.46
1:G:185:LYS:O	1:G:189:GLN:HG3	2.16	0.46
2:A:66:SER:OG	2:A:67:LYS:N	2.47	0.46
2:B:13:ASP:HB3	2:B:41:CYS:SG	2.56	0.45
2:C:67:LYS:HE3	2:C:71:ARG:HH12	1.81	0.45
2:C:2:GLU:HG3	2:C:4:VAL:H	1.81	0.45
2:C:40:TRP:HE1	2:C:44:ARG:NH1	2.13	0.45
2:B:88:GLU:H	2:B:88:GLU:CD	2.25	0.45
2:C:94:VAL:HG23	2:C:107:ALA:HB1	1.99	0.44
1:G:195:LEU:O	1:G:198:THR:HG22	2.17	0.44
2:B:105:ASN:O	2:B:109:GLU:HG3	2.17	0.44
2:A:105:ASN:HA	2:A:108:LYS:HD3	1.99	0.44
1:G:224:ASP:HA	1:G:227:LEU:HD12	1.99	0.44
2:C:106:GLN:HG3	2:C:107:ALA:H	1.82	0.44
2:A:53:LYS:HB2	2:A:53:LYS:HE3	1.74	0.44
1:F:195:LEU:HD23	1:F:195:LEU:HA	1.82	0.44
2:B:51:ILE:HB	2:B:84:LEU:HD21	1.99	0.44
2:B:116:LYS:HB2	2:B:116:LYS:HE2	1.84	0.44
2:A:56:VAL:HG11	1:D:228:GLU:HG2	2.00	0.43
2:A:105:ASN:O	2:A:109:GLU:HG3	2.18	0.43
2:A:59:GLU:OE1	1:D:227:LEU:HD22	2.18	0.43
1:F:176:LYS:HE3	1:F:176:LYS:HB2	1.64	0.43
2:C:13:ASP:OD1	2:B:76:TYR:OH	2.25	0.43
2:A:85:LYS:HA	2:A:87:TYR:CE1	2.55	0.42
1:F:177:LEU:O	1:F:181:ASN:HB2	2.19	0.42
1:F:234:SER:O	2:B:89:LYS:NZ	2.51	0.42
2:C:42:LEU:HB3	2:C:50:ASP:O	2.18	0.42
1:G:215:GLN:O	1:G:219:LEU:HD12	2.20	0.42
2:C:2:GLU:HG3	2:C:4:VAL:N	2.34	0.42
1:G:181:ASN:O	1:G:185:LYS:HG3	2.20	0.42
2:C:91:LEU:HD12	2:C:91:LEU:HA	1.92	0.42
2:B:42:LEU:HB3	2:B:50:ASP:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:37:GLU:HA	2:C:40:TRP:HE3	1.85	0.41
1:F:227:LEU:HD21	2:B:59:GLU:O	2.21	0.41
2:B:85:LYS:HA	2:B:87:TYR:CE1	2.56	0.41
2:C:59:GLU:O	2:C:63:PRO:HD3	2.21	0.41
2:B:61:LEU:HD23	2:B:61:LEU:HA	1.88	0.41
2:C:13:ASP:HB3	2:C:41:CYS:SG	2.61	0.40
1:E:181:ASN:OD1	1:E:181:ASN:N	2.54	0.40
2:C:94:VAL:CG2	2:C:107:ALA:HB1	2.52	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:13:ASP:OD1	2:A:76:TYR:OH[3_656]	2.00	0.20

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	D	63/72 (88%)	63 (100%)	0	0	100	100
1	E	56/72 (78%)	55 (98%)	1 (2%)	0	100	100
1	F	64/72 (89%)	64 (100%)	0	0	100	100
1	G	55/72 (76%)	55 (100%)	0	0	100	100
2	A	118/125 (94%)	114 (97%)	4 (3%)	0	100	100
2	B	117/125 (94%)	115 (98%)	2 (2%)	0	100	100
2	C	111/125 (89%)	109 (98%)	2 (2%)	0	100	100
All	All	584/663 (88%)	575 (98%)	9 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	D	55/60 (92%)	55 (100%)	0	100	100
1	E	34/60 (57%)	33 (97%)	1 (3%)	37	67
1	F	55/60 (92%)	55 (100%)	0	100	100
1	G	26/60 (43%)	26 (100%)	0	100	100
2	A	107/110 (97%)	107 (100%)	0	100	100
2	B	106/110 (96%)	106 (100%)	0	100	100
2	C	102/110 (93%)	101 (99%)	1 (1%)	68	86
All	All	485/570 (85%)	483 (100%)	2 (0%)	84	93

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	189	GLN
2	C	61	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	208	GLN
1	D	172	ASN
1	D	189	GLN
2	C	106	GLN
2	B	106	GLN

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	$\# Z > 2$
3	BME	A	201	-	3,3,3	0.43	0	2,2,2	0.48	0
4	SO4	C	201	-	4,4,4	0.67	0	6,6,6	0.09	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
3	BME	A	201	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	201	BME	O1-C1-C2-S2

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	D	65/72 (90%)	0.97	7 (10%) 11 9	38, 50, 78, 96	0
1	E	58/72 (80%)	1.50	17 (29%) 1 1	34, 50, 76, 82	0
1	F	66/72 (91%)	0.88	6 (9%) 15 12	39, 53, 74, 82	0
1	G	57/72 (79%)	1.03	10 (17%) 4 3	39, 55, 81, 86	0
2	A	120/125 (96%)	0.18	5 (4%) 40 37	14, 31, 63, 79	0
2	B	119/125 (95%)	0.92	18 (15%) 5 4	23, 53, 96, 102	0
2	C	117/125 (93%)	1.64	37 (31%) 1 1	32, 68, 104, 111	0
All	All	602/663 (90%)	0.98	100 (16%) 4 4	14, 51, 93, 111	0

All (100) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	74	VAL	6.6
1	E	235	ALA	5.4
2	C	99	GLN	5.1
2	C	98	LEU	4.8
2	C	121	ASP	4.7
2	C	107	ALA	4.5
1	E	201	LYS	4.4
1	F	173	ALA	4.4
1	E	217	GLU	4.3
1	G	235	ALA	4.2
2	C	106	GLN	4.1
1	D	174	GLU	4.0
1	E	236	ALA	3.9
1	D	169	ASP	3.9
2	C	97	LEU	3.9
2	C	102	PRO	3.8
1	G	227	LEU	3.8

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Mol	Chain	Res	Type	RSRZ
1	F	170	VAL	3.8
2	C	87	TYR	3.8
2	C	90	ALA	3.8
2	B	2	GLU	3.6
1	D	172	ASN	3.5
1	G	219	LEU	3.5
2	C	53	LYS	3.4
2	C	88	GLU	3.4
1	E	210	LEU	3.3
2	B	4	VAL	3.3
1	D	204	LYS	3.3
2	C	105	ASN	3.3
1	E	227	LEU	3.2
2	B	47	TYR	3.2
2	C	62	LEU	3.1
1	E	231	ALA	3.1
2	C	94	VAL	3.1
2	A	88	GLU	3.0
2	A	1	MET	3.0
2	B	26	ALA	3.0
1	E	230	HIS	3.0
1	F	235	ALA	3.0
2	B	3	ALA	2.9
2	B	17	PHE	2.9
1	E	234	SER	2.9
1	G	198	THR	2.9
2	C	64	LYS	2.9
2	C	119	LYS	2.8
1	E	209	VAL	2.8
2	B	7	GLU	2.8
1	G	187	ASP	2.8
2	C	76	TYR	2.7
1	D	233	LEU	2.7
2	B	10	SER	2.7
2	C	71	ARG	2.7
2	C	66	SER	2.7
2	C	108	LYS	2.7
1	D	170	VAL	2.7
2	C	67	LYS	2.6
2	B	19	LYS	2.6
2	C	72	ASP	2.5
2	C	85	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	E	179	GLU	2.5
2	B	36	PHE	2.5
1	E	220	THR	2.4
2	B	5	LEU	2.4
1	G	225	ARG	2.4
2	C	115	ASP	2.4
2	B	119	LYS	2.4
2	C	114	ILE	2.4
2	C	86	GLU	2.4
2	C	40	TRP	2.4
2	B	11	VAL	2.4
2	B	88	GLU	2.3
2	A	120	LYS	2.3
2	B	46	LYS	2.3
2	A	40	TRP	2.3
1	D	175	VAL	2.3
2	C	61	LEU	2.3
1	E	198	THR	2.3
2	C	101	GLU	2.3
1	E	211	ALA	2.3
1	F	233	LEU	2.3
1	F	220	THR	2.3
1	G	179	GLU	2.3
2	C	91	LEU	2.3
2	A	26	ALA	2.3
2	B	49	ASP	2.2
1	F	171	GLY	2.2
2	B	29	SER	2.2
1	E	226	LEU	2.2
2	C	69	GLU	2.2
1	G	224	ASP	2.2
1	E	208	GLN	2.1
1	G	211	ALA	2.1
2	C	32	LYS	2.1
2	B	120	LYS	2.1
2	C	113	LEU	2.1
2	C	109	GLU	2.1
1	E	181	ASN	2.1
1	G	231	ALA	2.1
2	C	60	GLU	2.1
2	C	100	THR	2.1

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](#)

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
4	SO4	C	201	5/5	0.83	0.17	59,59,90,99	0
3	BME	A	201	4/4	0.94	0.25	41,44,45,63	0

6.5 Other polymers [i](#)

There are no such residues in this entry.