



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:11 AM UTC

PDB ID : 6EN9 / pdb_00006en9
Title : E. coli Hydrogenase-2 (hydrogen reduced form)
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Deposited on : 2017-10-04
Resolution : 1.50 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

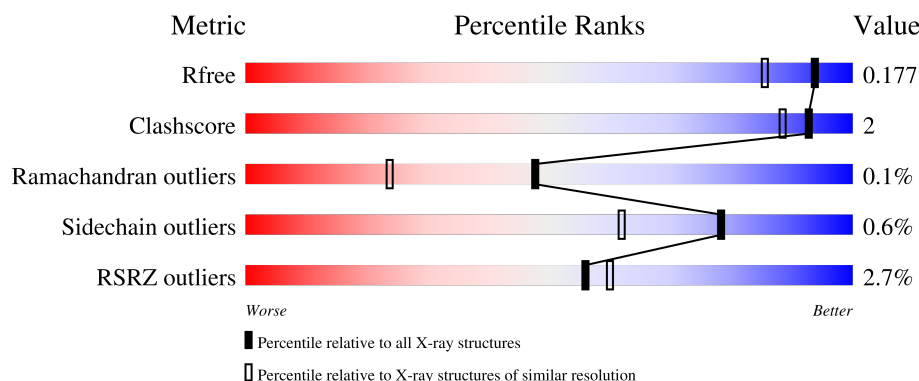
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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	S	298	
1	T	298	
2	L	567	
2	M	567	

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 13927 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 Hydrogenase-2 small chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	S	267	Total	C	N	O	S	0	4	0
			2059	1304	364	378	13			
1	T	267	Total	C	N	O	S	0	5	0
			2059	1306	362	378	13			

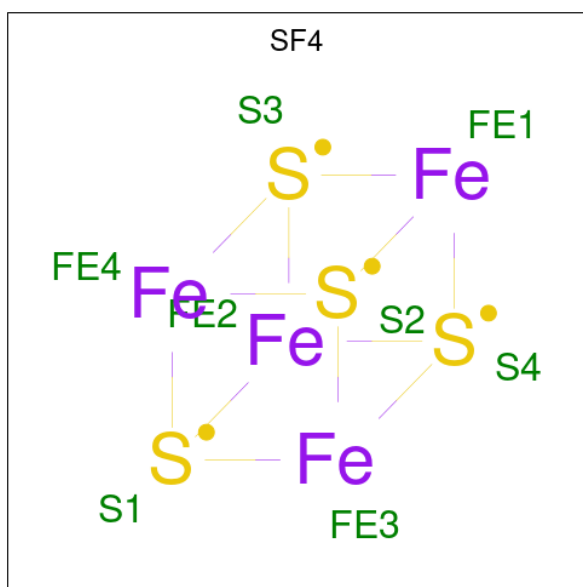
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
S	294	HIS	ASN	conflict	UNP P69741
S	295	HIS	-	expression tag	UNP P69741
S	296	HIS	-	expression tag	UNP P69741
S	297	HIS	-	expression tag	UNP P69741
S	298	HIS	-	expression tag	UNP P69741
S	299	HIS	-	expression tag	UNP P69741
T	294	HIS	ASN	conflict	UNP P69741
T	295	HIS	-	expression tag	UNP P69741
T	296	HIS	-	expression tag	UNP P69741
T	297	HIS	-	expression tag	UNP P69741
T	298	HIS	-	expression tag	UNP P69741
T	299	HIS	-	expression tag	UNP P69741

- Molecule 2 is a protein called Hydrogenase-2 large chain.

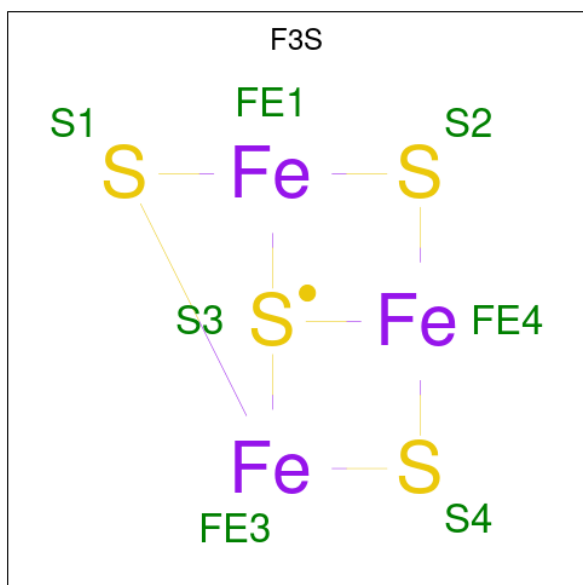
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	551	Total	C	N	O	S	0	6	0
			4319	2748	741	812	18			
2	M	551	Total	C	N	O	S	0	3	0
			4303	2738	740	807	18			

- Molecule 3 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄).



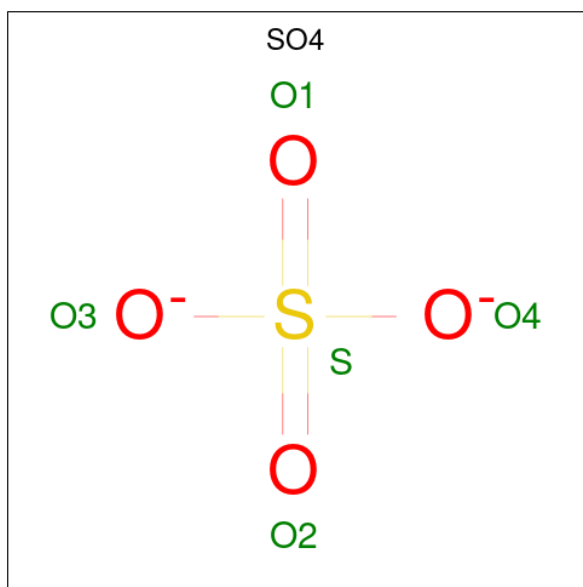
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	S	1	Total	Fe	S	0	0
			8	4	4		
3	S	1	Total	Fe	S	0	0
			8	4	4		
3	T	1	Total	Fe	S	0	0
			8	4	4		
3	T	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 4 is FE3-S4 CLUSTER (CCD ID: F3S) (formula: Fe_3S_4).



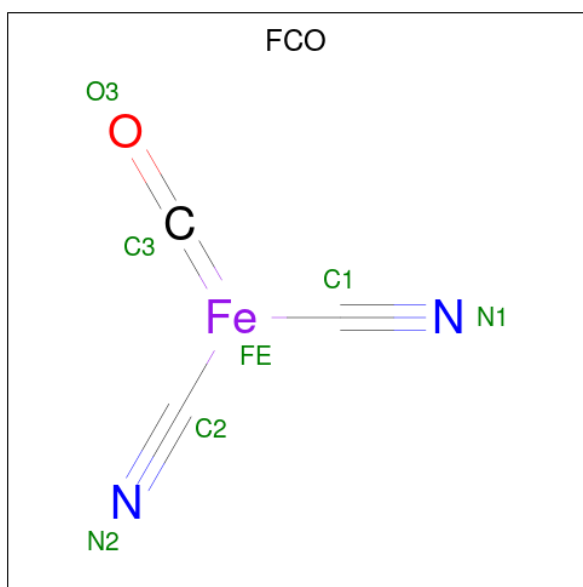
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	S	1	Total	Fe	S	0	0
			7	3	4		
4	T	1	Total	Fe	S	0	0
			7	3	4		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 6 is CARBONMONOXIDE-(DICYANO) IRON (CCD ID: FCO) (formula: C₃FeN₂O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	L	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		
6	M	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		

- Molecule 7 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	L	1	Total	Ni	0	0
			1	1		
7	M	1	Total	Ni	0	0
			1	1		

- Molecule 8 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	L	1	Total	Mg	0	0
			1	1		
8	M	1	Total	Mg	0	0
			1	1		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	S	189	Total	O	0	0
			189	189		

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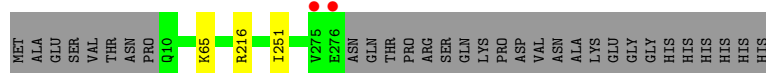
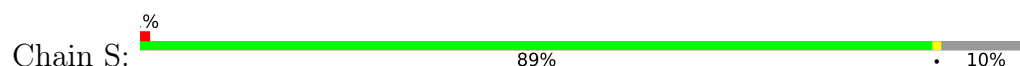
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	L	408	Total 408	O 408	0	0
9	T	163	Total 163	O 163	0	0
9	M	353	Total 353	O 353	0	0

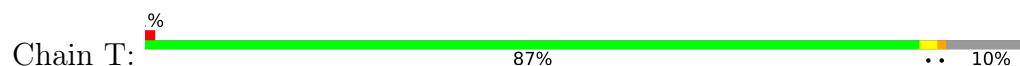
3 Residue-property plots [i](#)

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.

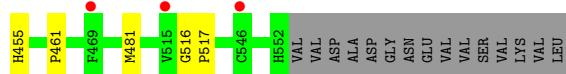
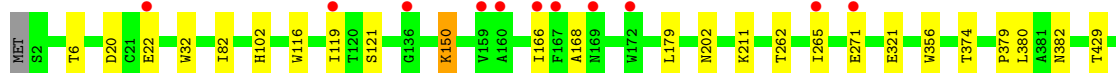
- Molecule 1: Hydrogenase-2 small chain



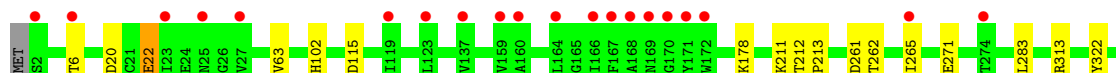
- Molecule 1: Hydrogenase-2 small chain



- Molecule 2: Hydrogenase-2 large chain



- Molecule 2: Hydrogenase-2 large chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	99.60Å 100.54Å 168.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	86.35 – 1.50 86.35 – 1.50	Depositor EDS
% Data completeness (in resolution range)	99.1 (86.35-1.50) 99.1 (86.35-1.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.75 (at 1.50Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.152 , 0.175 0.154 , 0.177	Depositor DCC
R_{free} test set	13461 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	16.3	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 34.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	13927	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.36% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SF4, NI, SO4, FCO, F3S, MG

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	S	1.05	2/2122 (0.1%)	1.01	0/2888
1	T	1.15	2/2131 (0.1%)	1.09	2/2900 (0.1%)
2	L	1.10	8/4441 (0.2%)	1.01	1/6054 (0.0%)
2	M	1.11	4/4419 (0.1%)	1.06	2/6024 (0.0%)
All	All	1.11	16/13113 (0.1%)	1.04	5/17866 (0.0%)

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	121	SER	C-O	-6.33	1.16	1.24
2	M	545	PRO	C-N	-6.01	1.25	1.33
2	L	271	GLU	CD-OE1	-5.89	1.14	1.25
2	M	283	LEU	C-O	-5.84	1.16	1.23
1	T	15[A]	ILE	CA-C	5.79	1.60	1.52
1	T	15[B]	ILE	CA-C	5.79	1.60	1.52
2	L	179	LEU	C-O	-5.72	1.16	1.23
2	M	350	ILE	CA-C	5.65	1.57	1.53
2	M	22	GLU	CA-C	-5.60	1.45	1.52
2	L	380	LEU	N-CA	5.58	1.53	1.46
2	L	168	ALA	N-CA	5.53	1.52	1.46
2	L	321	GLU	C-O	-5.43	1.17	1.24
1	S	216[A]	ARG	CA-C	5.38	1.60	1.52
1	S	216[B]	ARG	CA-C	5.38	1.60	1.52
2	L	119	ILE	C-O	-5.25	1.17	1.24
2	L	429	THR	N-CA	-5.06	1.40	1.46

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	395[A]	GLN	CA-C-O	6.21	127.55	120.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	395[B]	GLN	CA-C-O	6.21	127.55	120.90
1	T	169	ILE	N-CA-C	6.06	116.24	110.42
1	T	23	THR	CB-CA-C	5.58	115.35	109.83
2	L	271	GLU	CB-CG-CD	5.08	121.23	112.60

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	S	2059	0	1987	1	0
1	T	2059	0	1995	3	0
2	L	4319	0	4263	20	0
2	M	4303	0	4246	15	0
3	S	16	0	0	0	0
3	T	16	0	0	0	0
4	S	7	0	0	0	0
4	T	7	0	0	0	0
5	S	5	0	0	0	0
5	T	5	0	0	0	0
6	L	7	0	0	0	0
6	M	7	0	0	0	0
7	L	1	0	0	0	0
7	M	1	0	0	0	0
8	L	1	0	0	0	0
8	M	1	0	0	0	0
9	L	408	0	0	7	1
9	M	353	0	0	7	0
9	S	189	0	0	2	0
9	T	163	0	0	0	0
All	All	13927	0	12491	38	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 2.

All (38) 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:M:6:THR:HB	9:M:983:HOH:O	1.67	0.93
2:L:356:TRP:HH2	9:L:1016:HOH:O	1.71	0.72
2:L:22:GLU:OE1	9:L:701:HOH:O	2.10	0.70
2:M:455:HIS:HD2	9:M:984:HOH:O	1.78	0.66
2:M:22:GLU:HG3	9:M:1017:HOH:O	2.03	0.59
2:M:6:THR:HG23	2:M:20:ASP:OD1	2.03	0.58
2:L:150:LYS:HD3	9:M:1037:HOH:O	2.04	0.58
2:L:455:HIS:HD2	9:L:1050:HOH:O	1.89	0.55
2:M:261:ASP:O	2:M:265:ILE:HG23	2.08	0.53
2:L:262:THR:O	2:L:265:ILE:HG22	2.07	0.53
2:M:481:MET:HE3	9:M:705:HOH:O	2.08	0.52
2:L:150:LYS:HE2	9:M:1037:HOH:O	2.09	0.52
2:M:379:PRO:HB2	9:M:772:HOH:O	2.09	0.52
9:S:548:HOH:O	2:L:166:ILE:HG22	2.11	0.50
2:M:313:ARG:HD3	2:M:322:TYR:CE2	2.47	0.50
1:T:87:LYS:HE3	1:T:133:PRO:O	2.11	0.50
2:L:379:PRO:HB2	9:L:921:HOH:O	2.10	0.50
1:S:65:LYS:HE2	9:S:402:HOH:O	2.13	0.49
2:L:6[A]:THR:HG23	2:L:20:ASP:OD1	2.13	0.48
2:M:178:LYS:HE2	2:M:271:GLU:OE2	2.13	0.48
2:L:116:TRP:CG	2:L:265:ILE:HD11	2.49	0.47
2:L:374[B]:THR:HG21	2:L:516:GLY:HA2	1.96	0.47
2:L:356:TRP:CH2	9:L:1016:HOH:O	2.54	0.47
1:T:15[B]:ILE:CD1	1:T:69:LEU:HD23	2.44	0.47
2:L:481:MET:HE3	9:L:742:HOH:O	2.15	0.46
2:L:6[A]:THR:HG21	9:L:852:HOH:O	2.17	0.44
2:L:374[B]:THR:CG2	2:L:517:PRO:HD3	2.47	0.44
2:M:537:VAL:CG1	2:M:541:HIS:CE1	3.01	0.43
2:L:356:TRP:CD1	2:L:356:TRP:C	2.97	0.43
2:M:212:THR:HA	2:M:213:PRO:C	2.44	0.42
2:M:115:ASP:HB3	2:M:538:ARG:HG2	2.02	0.41
2:L:82:ILE:HD11	2:L:461:PRO:HB3	2.02	0.41
2:L:150:LYS:HB3	2:L:150:LYS:HE3	1.91	0.41
1:T:23:THR:HB	2:M:63:VAL:HG23	2.03	0.41
2:M:262:THR:O	2:M:265:ILE:HG12	2.21	0.41
2:L:22:GLU:HG3	2:L:32:TRP:CZ3	2.56	0.40
2:L:202:ASN:HD22	2:L:202:ASN:HA	1.74	0.40
2:M:468:GLU:HA	2:M:488:ILE:O	2.21	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)
9:L:704:HOH:O	9:L:769:HOH:O[3_755]	1.13	1.07

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	S	269/298 (90%)	260 (97%)	9 (3%)	0	100	100
1	T	270/298 (91%)	261 (97%)	9 (3%)	0	100	100
2	L	555/567 (98%)	532 (96%)	22 (4%)	1 (0%)	43	22
2	M	552/567 (97%)	532 (96%)	19 (3%)	1 (0%)	43	22
All	All	1646/1730 (95%)	1585 (96%)	59 (4%)	2 (0%)	48	24

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	L	211	LYS
2	M	211	LYS

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	S	216/239 (90%)	215 (100%)	1 (0%)	81	66
1	T	217/239 (91%)	216 (100%)	1 (0%)	81	66
2	L	471/479 (98%)	467 (99%)	4 (1%)	73	54

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	M	468/479 (98%)	464 (99%)	4 (1%)	70	49
All	All	1372/1436 (96%)	1362 (99%)	10 (1%)	78	57

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	S	251	ILE
2	L	102	HIS
2	L	150	LYS
2	L	382[A]	ASN
2	L	382[B]	ASN
1	T	251	ILE
2	M	102	HIS
2	M	382[A]	ASN
2	M	382[B]	ASN
2	M	414	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	S	20	GLN
1	S	75	GLN
1	S	266	HIS
2	L	3	GLN
2	L	202	ASN
2	L	411	ASN
2	L	449	ASN
2	L	455	HIS
1	T	60	GLN
1	T	75	GLN
1	T	274	ASN
2	M	47	ASN
2	M	406	GLN
2	M	411	ASN
2	M	455	HIS

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](#)

Of 14 ligands modelled in this entry, 4 are monoatomic - leaving 10 for Mogul analysis.

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	$\# Z > 2$
6	FCO	L	601	2	0,6,6	-	-	-		
3	SF4	T	303	1	0,12,12	-	-	-		
3	SF4	S	301	1	0,12,12	-	-	-		
6	FCO	M	601	2	0,6,6	-	-	-		
4	F3S	T	302	1	0,9,9	-	-	-		
4	F3S	S	302	1	0,9,9	-	-	-		
5	SO4	T	304	-	4,4,4	0.61	0	6,6,6	0.14	0
5	SO4	S	304	-	4,4,4	0.65	0	6,6,6	0.56	0
3	SF4	S	303	1	0,12,12	-	-	-		
3	SF4	T	301	1	0,12,12	-	-	-		

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	SF4	T	303	1	-	-	0/6/5/5
3	SF4	S	301	1	-	-	0/6/5/5
4	F3S	T	302	1	-	-	0/3/3/3
4	F3S	S	302	1	-	-	0/3/3/3
3	SF4	S	303	1	-	-	0/6/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SF4	T	301	1	-	-	0/6/5/5

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	S	267/298 (89%)	-0.15	2 (0%) 84 87	7, 16, 28, 43	4 (1%)
1	T	267/298 (89%)	0.12	3 (1%) 78 81	10, 18, 34, 47	5 (1%)
2	L	551/567 (97%)	-0.14	14 (2%) 58 62	9, 16, 30, 41	6 (1%)
2	M	551/567 (97%)	0.17	25 (4%) 38 42	10, 19, 35, 56	3 (0%)
All	All	1636/1730 (94%)	0.01	44 (2%) 56 60	7, 17, 33, 56	18 (1%)

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	M	265	ILE	5.0
2	M	172	TRP	3.9
2	L	265	ILE	3.7
2	L	172	TRP	3.5
2	M	2	SER	3.5
2	M	160	ALA	3.3
2	L	119	ILE	3.2
2	L	167	PHE	3.1
2	L	515	VAL	3.1
2	M	119	ILE	2.9
2	M	166	ILE	2.9
2	M	469	PHE	2.9
2	M	413	LEU	2.8
2	M	515	VAL	2.8
2	M	6	THR	2.8
2	M	27	VAL	2.7
2	M	23	ILE	2.7
2	L	166	ILE	2.6
2	L	160	ALA	2.6
2	L	159	VAL	2.6
1	T	276	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
2	M	274	THR	2.5
2	M	167	PHE	2.5
1	T	145	GLY	2.5
2	L	22	GLU	2.5
2	M	123	LEU	2.4
2	L	546	CYS	2.4
1	T	89	ASN	2.4
2	M	531	ASN	2.4
2	M	159	VAL	2.3
1	S	276	GLU	2.3
2	L	271	GLU	2.3
2	M	395[A]	GLN	2.3
2	M	168	ALA	2.2
2	M	171	TYR	2.2
2	M	169	ASN	2.2
2	L	136	GLY	2.1
2	M	137	VAL	2.1
2	M	170	GLY	2.1
2	M	25	ASN	2.1
2	L	469	PHE	2.1
1	S	275	VAL	2.0
2	M	164	LEU	2.0
2	L	169	ASN	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](#)

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.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	SO4	T	304	5/5	0.75	0.13	64,71,72,75	0
5	SO4	S	304	5/5	0.89	0.11	41,44,45,48	0
3	SF4	S	301	8/8	0.99	0.03	14,14,14,14	0
3	SF4	T	303	8/8	0.99	0.03	14,14,14,14	0
7	NI	M	602	1/1	0.99	0.04	15,15,15,15	0
8	MG	M	603	1/1	0.99	0.03	12,12,12,12	0
3	SF4	T	301	8/8	1.00	0.03	14,15,15,15	0
3	SF4	S	303	8/8	1.00	0.01	11,12,12,12	0
6	FCO	L	601	7/7	1.00	0.02	11,12,12,13	0
6	FCO	M	601	7/7	1.00	0.04	12,13,15,15	0
7	NI	L	602	1/1	1.00	0.02	13,13,13,13	0
4	F3S	S	302	7/7	1.00	0.02	12,12,13,13	0
8	MG	L	603	1/1	1.00	0.04	9,9,9,9	0
4	F3S	T	302	7/7	1.00	0.02	13,13,13,14	0

6.5 Other polymers [i](#)

There are no such residues in this entry.