



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:59 PM GMT

PDB ID : 4UCW
Title : Structure of the T18V small subunit mutant of D. fructosovorans NiFe- hydrogenase
Authors : Abou-Hamdan, A.; Ceccaldi, P.; Lebrette, H.; Gutierrez-Sanz, O.; Richaud, P.; Cournac, L.; Guigliarelli, B.; deLacey, A.L.; Leger, C.; Volbeda, A.; Burlat, B.; Dementin, S.
Deposited on : 2014-12-04
Resolution : 2.30 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

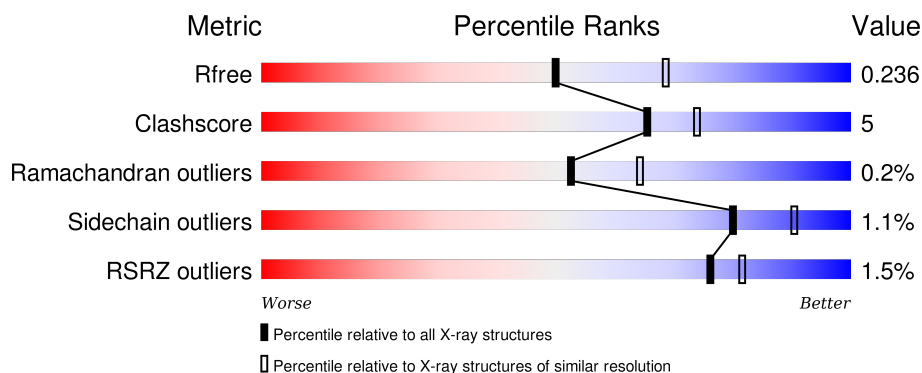
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	264	<div> <div></div> <div>88%12%</div> </div>
1	B	264	<div> <div></div> <div>87%11%</div> </div>
1	C	264	<div> <div>8%</div> <div>80%19%</div> </div>
2	Q	563	<div> <div></div> <div>86%11%</div> </div>
2	R	563	<div> <div>%</div> <div>85%12%</div> </div>

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Mol	Chain	Length	Quality of chain
2	S	563	 2% 84% 13%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	A	271	-	-	-	X
3	GOL	Q	1562	-	-	-	X

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 19144 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 (NIFE) SMALL SUBUNIT HYDA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	1	0
			1981	1261	331	374	15			
1	B	260	Total	C	N	O	S	0	2	0
			1968	1253	327	373	15			
1	C	261	Total	C	N	O	S	0	1	0
			1970	1255	330	370	15			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	18	VAL	THR	ENGINEERED MUTATION	UNP E1K248
B	18	VAL	THR	ENGINEERED MUTATION	UNP E1K248
C	18	VAL	THR	ENGINEERED MUTATION	UNP E1K248

- Molecule 2 is a protein called NICKEL-DEPENDENT HYDROGENASE LARGE SUB-UNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	Q	544	Total	C	N	O	S	0	5	0
			4183	2665	724	772	22			
2	R	545	Total	C	N	O	S	0	1	0
			4174	2659	724	769	22			
2	S	545	Total	C	N	O	S	0	1	0
			4174	2657	725	770	22			

There are 45 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Q	-13	ALA	-	EXPRESSION TAG	UNP E1K247
Q	-12	SER	-	EXPRESSION TAG	UNP E1K247
Q	-11	TRP	-	EXPRESSION TAG	UNP E1K247
Q	-10	SER	-	EXPRESSION TAG	UNP E1K247

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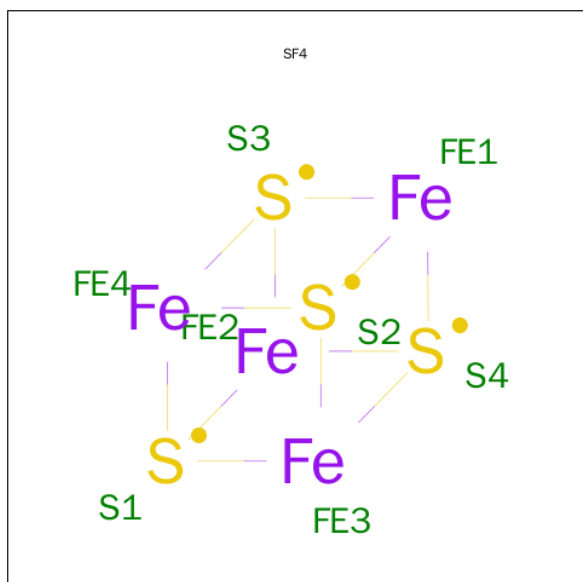
Chain	Residue	Modelled	Actual	Comment	Reference
Q	-9	HIS	-	EXPRESSION TAG	UNP E1K247
Q	-8	PRO	-	EXPRESSION TAG	UNP E1K247
Q	-7	GLN	-	EXPRESSION TAG	UNP E1K247
Q	-6	PHE	-	EXPRESSION TAG	UNP E1K247
Q	-5	GLU	-	EXPRESSION TAG	UNP E1K247
Q	-4	LYS	-	EXPRESSION TAG	UNP E1K247
Q	-3	GLY	-	EXPRESSION TAG	UNP E1K247
Q	-2	ALA	-	EXPRESSION TAG	UNP E1K247
Q	-1	SER	-	EXPRESSION TAG	UNP E1K247
Q	0	GLY	-	EXPRESSION TAG	UNP E1K247
Q	1	ALA	-	EXPRESSION TAG	UNP E1K247
R	-13	ALA	-	EXPRESSION TAG	UNP E1K247
R	-12	SER	-	EXPRESSION TAG	UNP E1K247
R	-11	TRP	-	EXPRESSION TAG	UNP E1K247
R	-10	SER	-	EXPRESSION TAG	UNP E1K247
R	-9	HIS	-	EXPRESSION TAG	UNP E1K247
R	-8	PRO	-	EXPRESSION TAG	UNP E1K247
R	-7	GLN	-	EXPRESSION TAG	UNP E1K247
R	-6	PHE	-	EXPRESSION TAG	UNP E1K247
R	-5	GLU	-	EXPRESSION TAG	UNP E1K247
R	-4	LYS	-	EXPRESSION TAG	UNP E1K247
R	-3	GLY	-	EXPRESSION TAG	UNP E1K247
R	-2	ALA	-	EXPRESSION TAG	UNP E1K247
R	-1	SER	-	EXPRESSION TAG	UNP E1K247
R	0	GLY	-	EXPRESSION TAG	UNP E1K247
R	1	ALA	-	EXPRESSION TAG	UNP E1K247
S	-13	ALA	-	EXPRESSION TAG	UNP E1K247
S	-12	SER	-	EXPRESSION TAG	UNP E1K247
S	-11	TRP	-	EXPRESSION TAG	UNP E1K247
S	-10	SER	-	EXPRESSION TAG	UNP E1K247
S	-9	HIS	-	EXPRESSION TAG	UNP E1K247
S	-8	PRO	-	EXPRESSION TAG	UNP E1K247
S	-7	GLN	-	EXPRESSION TAG	UNP E1K247
S	-6	PHE	-	EXPRESSION TAG	UNP E1K247
S	-5	GLU	-	EXPRESSION TAG	UNP E1K247
S	-4	LYS	-	EXPRESSION TAG	UNP E1K247
S	-3	GLY	-	EXPRESSION TAG	UNP E1K247
S	-2	ALA	-	EXPRESSION TAG	UNP E1K247
S	-1	SER	-	EXPRESSION TAG	UNP E1K247
S	0	GLY	-	EXPRESSION TAG	UNP E1K247
S	1	ALA	-	EXPRESSION TAG	UNP E1K247

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



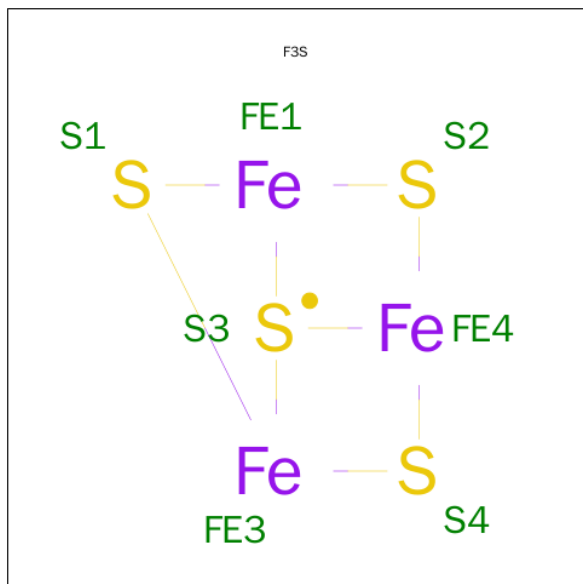
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	Q	1	Total	C	O	0	0
			6	3	3		
3	Q	1	Total	C	O	0	0
			6	3	3		
3	R	1	Total	C	O	0	0
			6	3	3		
3	S	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe_4S_4).



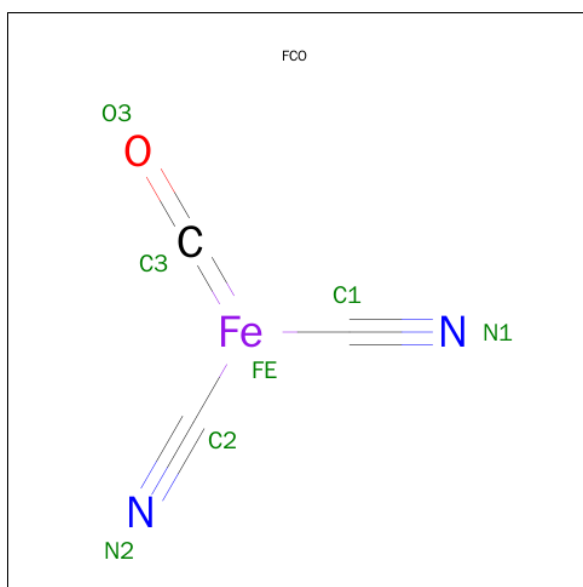
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	Fe	S	0	0
			8	4	4		
4	A	1	Total	Fe	S	0	0
			8	4	4		
4	B	1	Total	Fe	S	0	0
			8	4	4		
4	B	1	Total	Fe	S	0	0
			8	4	4		
4	C	1	Total	Fe	S	0	0
			8	4	4		
4	C	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 5 is FE3-S4 CLUSTER (three-letter code: F3S) (formula: Fe_3S_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	Fe	S	0	0
			7	3	4		
5	B	1	Total	Fe	S	0	0
			7	3	4		
5	C	1	Total	Fe	S	0	0
			7	3	4		

- Molecule 6 is CARBONMONOXIDE-(DICYANO) IRON (three-letter code: FCO) (formula: $\text{C}_3\text{FeN}_2\text{O}$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	Q	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		
6	R	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		
6	S	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		

- Molecule 7 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	R	1	Total	Ni	0	0
			1	1		
7	Q	1	Total	Ni	0	0
			1	1		
7	S	1	Total	Ni	0	0
			1	1		

- Molecule 8 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	R	1	Total	Mg	0	0
			1	1		
8	Q	1	Total	Mg	0	0
			1	1		
8	S	1	Total	Mg	0	0
			1	1		

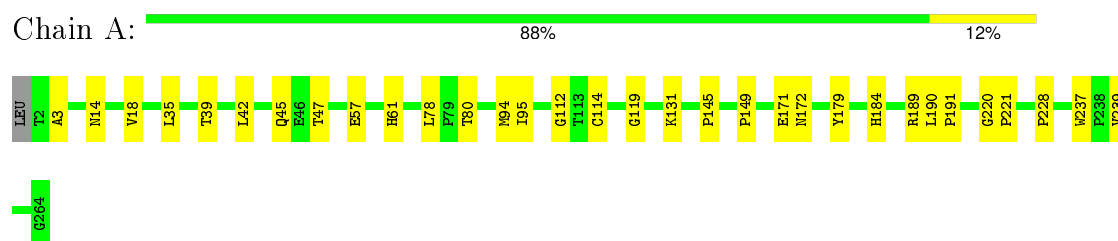
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	95	Total 95	O 95	0	0
9	B	77	Total 77	O 77	0	0
9	C	57	Total 57	O 57	0	0
9	Q	141	Total 141	O 141	0	0
9	R	102	Total 102	O 102	0	0
9	S	96	Total 96	O 96	0	0

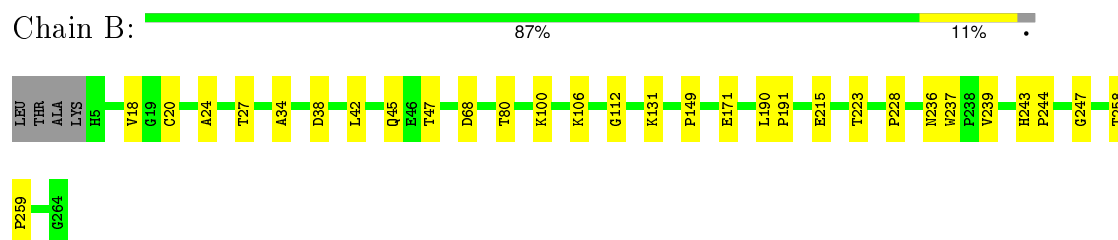
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

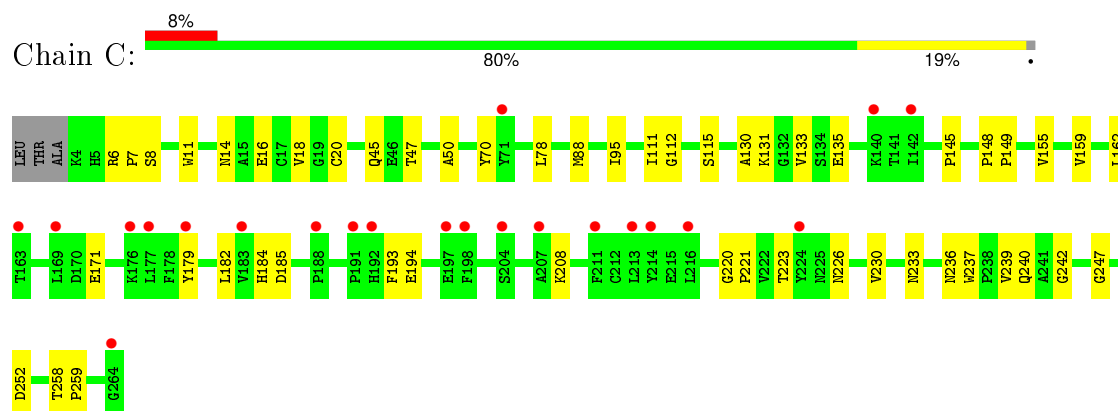
- Molecule 1: HYDROGENASE (NIFE) SMALL SUBUNIT HYDA



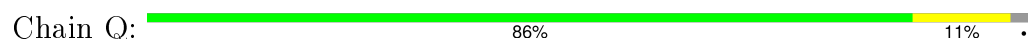
- Molecule 1: HYDROGENASE (NIFE) SMALL SUBUNIT HYDA

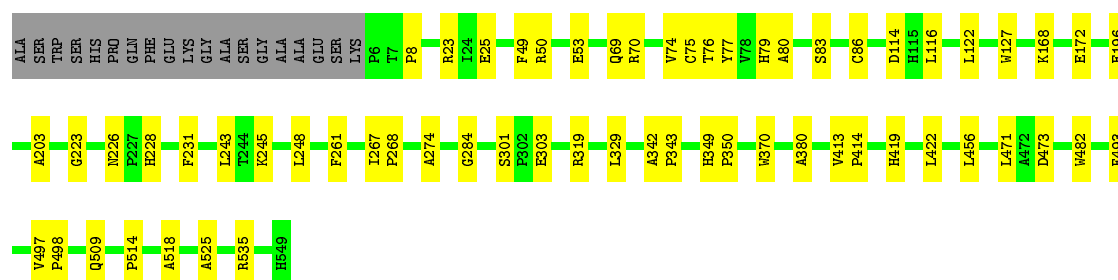


- Molecule 1: HYDROGENASE (NIFE) SMALL SUBUNIT HYDA

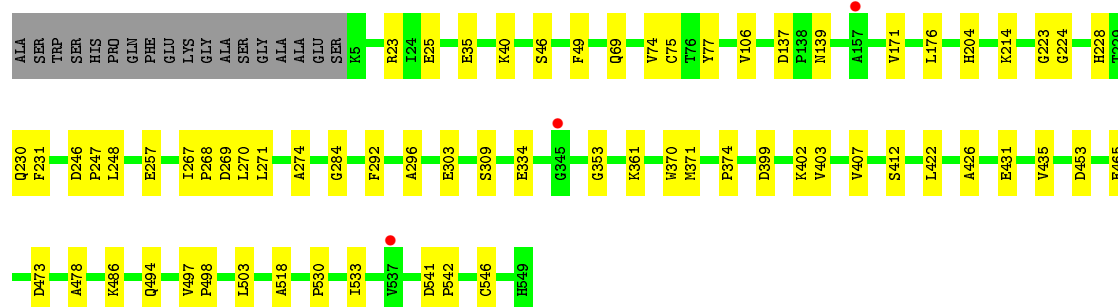
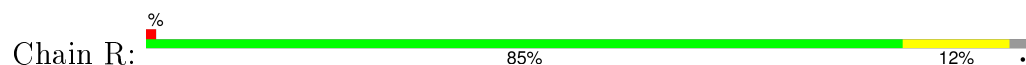


- Molecule 2: NICKEL-DEPENDENT HYDROGENASE LARGE SUBUNIT

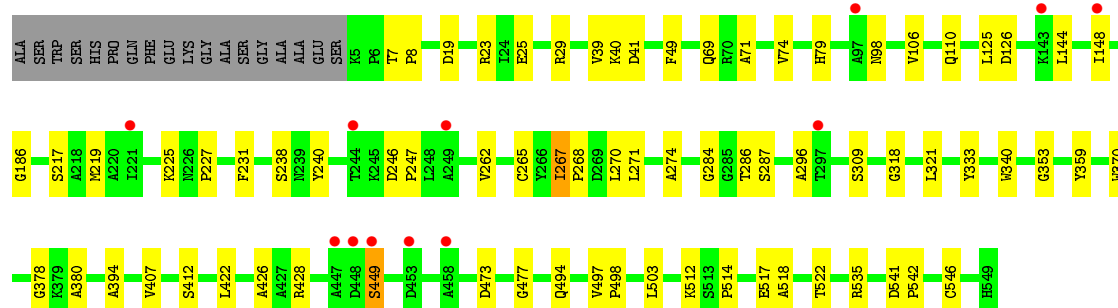
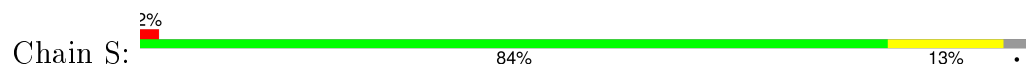




• Molecule 2: NICKEL-DEPENDENT HYDROGENASE LARGE SUBUNIT



• Molecule 2: NICKEL-DEPENDENT HYDROGENASE LARGE SUBUNIT



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	62.20Å 99.59Å 181.84Å 90.00° 92.47° 90.00°	Depositor
Resolution (Å)	25.00 – 2.30 39.06 – 2.30	Depositor EDS
% Data completeness (in resolution range)	93.4 (25.00-2.30) 83.2 (39.06-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.61 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.8.0071	Depositor
R, R_{free}	0.192 , 0.236 0.192 , 0.236	Depositor DCC
R_{free} test set	4581 reflections (5.24%)	DCC
Wilson B-factor (Å ²)	19.4	Xtriage
Anisotropy	0.534	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	(Not available) , (Not available)	EDS
Estimated twinning fraction	0.057 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.25$	Xtriage
Outliers	0 of 92096 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	19144	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: GOL, CSO, NI, SF4, MG, F3S, FCO

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	0.68	0/2040	0.73	1/2776 (0.0%)
1	B	0.61	0/2032	0.70	0/2767
1	C	0.57	0/2030	0.67	0/2765
2	Q	0.66	0/4303	0.75	3/5839 (0.1%)
2	R	0.57	0/4275	0.71	0/5802
2	S	0.59	0/4276	0.71	0/5804
All	All	0.61	0/18956	0.72	4/25753 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	Q	114	ASP	CB-CG-OD1	6.98	124.58	118.30
2	Q	50	ARG	NE-CZ-NH2	-6.72	116.94	120.30
2	Q	70	ARG	NE-CZ-NH2	-5.24	117.68	120.30
1	A	189	ARG	NE-CZ-NH2	-5.06	117.77	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1981	0	1916	17	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1968	0	1903	16	0
1	C	1970	0	1904	34	0
2	Q	4183	0	4159	38	0
2	R	4174	0	4145	40	0
2	S	4174	0	4141	47	0
3	A	6	0	8	0	0
3	Q	12	0	16	1	0
3	R	6	0	8	0	0
3	S	6	0	8	0	0
4	A	16	0	0	0	0
4	B	16	0	0	0	0
4	C	16	0	0	1	0
5	A	7	0	0	0	0
5	B	7	0	0	0	0
5	C	7	0	0	1	0
6	Q	7	0	0	1	0
6	R	7	0	0	1	0
6	S	7	0	0	0	0
7	Q	1	0	0	0	0
7	R	1	0	0	0	0
7	S	1	0	0	0	0
8	Q	1	0	0	0	0
8	R	1	0	0	0	0
8	S	1	0	0	0	0
9	A	95	0	0	1	0
9	B	77	0	0	1	0
9	C	57	0	0	2	0
9	Q	141	0	0	2	0
9	R	102	0	0	0	0
9	S	96	0	0	2	0
All	All	19144	0	18208	183	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 (183) 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:S:186:GLY:HA3	9:S:2045:HOH:O	1.74	0.86
2:R:214:LYS:HE3	2:R:257:GLU:OE2	1.79	0.82
2:R:25:GLU:CG	2:R:74:VAL:HG21	2.17	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:S:267:ILE:HG23	2:S:268:PRO:HD3	1.71	0.71
2:R:25:GLU:HG3	2:R:74:VAL:HG21	1.75	0.69
1:C:252:ASP:HA	9:C:2034:HOH:O	1.92	0.68
2:S:25:GLU:CG	2:S:74:VAL:HG21	2.26	0.65
1:C:50:ALA:HB2	2:S:125:LEU:HD13	1.78	0.64
1:B:215[A]:GLU:HG3	3:Q:1562:GOL:H32	1.78	0.64
2:Q:25:GLU:CG	2:Q:74:VAL:HG21	2.29	0.63
2:S:106:VAL:HG21	2:S:231:PHE:CZ	2.34	0.62
2:S:41:ASP:OD1	2:S:512:LYS:NZ	2.29	0.61
1:C:155:VAL:O	1:C:159:VAL:HG23	2.01	0.61
1:B:47:THR:O	2:R:23:ARG:HA	2.00	0.61
1:C:184:HIS:ND1	1:C:221:PRO:HG3	2.16	0.61
1:A:14:ASN:ND2	1:A:94:MET:HB3	2.15	0.61
1:A:171:GLU:CD	1:A:171:GLU:H	2.04	0.60
1:C:131:LYS:HB3	1:C:135:GLU:HG3	1.82	0.60
1:C:112:GLY:HA2	1:C:149:PRO:HD3	1.84	0.60
2:Q:116[B]:LEU:HD11	2:Q:261:PHE:HE1	1.67	0.60
1:C:237:TRP:CH2	1:C:239:VAL:HB	2.37	0.60
2:Q:116[B]:LEU:HD11	2:Q:261:PHE:CE1	2.37	0.59
2:Q:203:ALA:HB1	9:Q:2058:HOH:O	2.02	0.59
2:Q:413:VAL:HB	2:Q:414:PRO:HD2	1.86	0.57
1:B:42:LEU:HD21	1:B:45:GLN:HG3	1.86	0.56
1:C:148:PRO:HG3	2:S:227:PRO:HB2	1.86	0.56
2:S:497:VAL:HG11	2:S:546:CYS:HB3	1.88	0.56
1:C:47:THR:O	2:S:23:ARG:HA	2.05	0.56
2:S:353:GLY:HA3	2:S:494:GLN:HG3	1.88	0.56
2:S:69:GLN:HA	2:S:79:HIS:HB2	1.87	0.56
2:S:25:GLU:HG2	2:S:74:VAL:HG21	1.87	0.55
2:Q:243:LEU:HD11	2:Q:456:LEU:HD13	1.88	0.55
2:S:49:PHE:HB2	2:S:370:TRP:CD2	2.42	0.55
2:Q:196:GLU:OE1	2:Q:196:GLU:N	2.37	0.54
2:R:465:GLU:HG3	2:R:486:LYS:HA	1.89	0.54
2:S:110:GLN:HB2	2:S:219:MET:HE3	1.89	0.54
2:R:171:VAL:HG13	2:R:176:LEU:HD21	1.88	0.54
1:C:78:LEU:O	1:C:130:ALA:HA	2.08	0.54
2:S:25:GLU:HG2	2:S:74:VAL:CG2	2.38	0.53
1:A:78:LEU:HD22	1:A:95:ILE:HA	1.90	0.53
1:B:258:THR:HA	1:B:259:PRO:C	2.28	0.53
1:A:42:LEU:HD21	1:A:45:GLN:HG3	1.90	0.53
2:R:274:ALA:HA	2:R:422:LEU:HD11	1.91	0.53
2:R:223:GLY:HA2	2:R:231:PHE:CD1	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:R:25:GLU:HG2	2:R:74:VAL:CG2	2.39	0.52
1:B:80:THR:HG21	1:B:131:LYS:HD2	1.92	0.51
2:R:267:ILE:HB	2:R:268:PRO:HD3	1.93	0.51
1:B:236:ASN:HB3	2:R:224:GLY:O	2.10	0.51
2:S:267:ILE:CG2	2:S:268:PRO:HD3	2.41	0.51
2:Q:69:GLN:HB2	2:Q:80:ALA:HB2	1.93	0.51
2:S:19:ASP:HB2	2:S:29:ARG:HG3	1.93	0.51
2:R:399:ASP:O	2:R:403:VAL:HG23	2.11	0.50
1:A:112:GLY:HA2	1:A:149:PRO:HD3	1.92	0.50
2:S:296:ALA:HA	2:S:309:SER:HA	1.91	0.50
1:B:228:PRO:HB3	1:B:237:TRP:CZ2	2.45	0.50
1:A:184:HIS:HB2	1:A:220:GLY:C	2.32	0.50
2:Q:83:SER:O	2:Q:86:CYS:HB3	2.11	0.50
2:Q:380:ALA:HB1	2:Q:514:PRO:HD3	1.94	0.50
2:R:25:GLU:HG2	2:R:74:VAL:HG21	1.92	0.50
2:Q:8:PRO:HB2	2:Q:525:ALA:HB2	1.94	0.50
2:Q:509[A]:GLN:NE2	9:Q:2118:HOH:O	2.45	0.50
2:S:428:ARG:NH2	2:S:541:ASP:HB2	2.26	0.50
1:A:190:LEU:HB3	1:A:191:PRO:HD3	1.94	0.50
1:C:11:TRP:HB3	1:C:45:GLN:HG2	1.94	0.49
2:Q:301:SER:HB2	2:Q:303:GLU:OE1	2.13	0.49
2:S:144:LEU:O	2:S:148:ILE:HG12	2.12	0.49
2:Q:25:GLU:HG3	2:Q:74:VAL:HG21	1.93	0.49
1:B:112:GLY:HA2	1:B:149:PRO:HD3	1.94	0.49
1:A:80:THR:HG21	1:A:131:LYS:HD2	1.95	0.49
1:A:114:CYS:HA	1:A:119:GLY:HA3	1.94	0.49
2:R:204:HIS:ND1	2:R:269:ASP:OD2	2.46	0.49
1:B:24:ALA:O	1:B:27:THR:HG22	2.13	0.49
2:S:39:VAL:CG2	2:S:522:THR:HB	2.44	0.48
1:C:184:HIS:HB2	1:C:220:GLY:C	2.34	0.48
2:S:25:GLU:HG3	2:S:74:VAL:HG21	1.95	0.48
1:C:8:SER:HB3	1:C:70:TYR:HA	1.96	0.48
1:B:34:ALA:O	1:B:38:ASP:HB2	2.13	0.48
1:C:220:GLY:N	1:C:221:PRO:CD	2.76	0.47
1:B:68:ASP:O	1:B:106:LYS:HE3	2.14	0.47
2:Q:342:ALA:HB1	2:Q:343:PRO:HD2	1.95	0.47
2:R:284:GLY:HA2	2:R:518:ALA:O	2.15	0.46
2:R:292:PHE:CD2	2:R:478:ALA:HB1	2.50	0.46
2:R:246:ASP:HB2	2:R:247:PRO:HD3	1.97	0.46
2:S:274:ALA:HA	2:S:422:LEU:HD11	1.97	0.46
2:R:292:PHE:CE2	2:R:478:ALA:HB1	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:S:333:TYR:OH	2:S:378:GLY:HA2	2.16	0.46
1:C:233:ASN:HB3	2:S:217:SER:HA	1.98	0.46
2:Q:274:ALA:CB	2:Q:422:LEU:HD11	2.46	0.46
1:C:193:PHE:HD2	1:C:194:GLU:HG3	1.81	0.46
2:R:214:LYS:HE3	2:R:257:GLU:CD	2.35	0.46
1:B:237:TRP:CH2	1:B:239:VAL:HB	2.51	0.46
2:Q:53:GLU:HG2	2:Q:493:PHE:O	2.16	0.46
2:S:262:VAL:O	2:S:267:ILE:HG22	2.17	0.45
2:S:271:LEU:HD21	2:S:407:VAL:HG22	1.98	0.45
1:B:20:CYS:SG	1:B:112:GLY:HA3	2.57	0.45
1:C:236:ASN:OD1	1:C:240:GLN:HB3	2.17	0.45
2:Q:76:THR:O	2:Q:77:TYR:HB3	2.15	0.45
2:R:46:SER:OG	2:R:371:MET:HG3	2.16	0.45
2:R:77:TYR:CD2	2:R:106:VAL:HG12	2.52	0.45
1:B:223:THR:OG1	1:B:247:GLY:HA2	2.17	0.45
1:C:145:PRO:HD2	1:C:179:TYR:CZ	2.52	0.45
1:C:184:HIS:CE1	1:C:221:PRO:HG3	2.51	0.45
1:C:184:HIS:HA	4:C:1265:SF4:S1	2.57	0.45
1:C:182:LEU:HB2	1:C:185:ASP:OD2	2.17	0.45
1:C:242:GLY:HA3	2:S:238:SER:O	2.16	0.45
1:C:258:THR:HA	1:C:259:PRO:C	2.37	0.45
2:S:71:ALA:HB3	2:S:79:HIS:CD2	2.52	0.44
2:S:284:GLY:HA2	2:S:518:ALA:O	2.16	0.44
2:Q:168:LYS:O	2:Q:172:GLU:HG3	2.17	0.44
2:S:286:THR:HB	9:S:2058:HOH:O	2.17	0.44
1:A:47:THR:O	2:Q:23:ARG:HA	2.17	0.44
2:Q:127:TRP:CZ3	2:Q:535:ARG:HD3	2.52	0.44
2:S:110:GLN:HB2	2:S:219:MET:CE	2.48	0.44
2:S:477:GLY:O	2:S:498:PRO:HG3	2.17	0.44
2:S:246:ASP:HB2	2:S:247:PRO:HD3	1.99	0.44
2:Q:349:HIS:CG	2:Q:350:PRO:HD2	2.53	0.44
1:C:237:TRP:CZ2	1:C:239:VAL:HB	2.53	0.43
1:B:243:HIS:CG	1:B:244:PRO:HD2	2.53	0.43
2:S:98:ASN:OD1	2:S:449:SER:HB3	2.18	0.43
2:R:35:GLU:OE1	2:R:40:LYS:HD3	2.18	0.43
2:Q:319:ARG:HG2	2:Q:419:HIS:CE1	2.54	0.43
2:S:7:THR:HA	2:S:8:PRO:HD3	1.90	0.43
2:R:137:ASP:OD2	2:R:139:ASN:HB2	2.18	0.43
1:A:237:TRP:CH2	1:A:239:VAL:HB	2.54	0.43
2:S:340:TRP:CD1	2:S:359:TYR:HA	2.53	0.43
2:R:541:ASP:N	2:R:542:PRO:HD3	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Q:75:CSO:CB	6:Q:1550:FCO:C2	2.97	0.43
2:R:303:GLU:H	2:R:303:GLU:CD	2.22	0.43
2:Q:497:VAL:CG1	2:Q:498:PRO:HD2	2.49	0.43
1:A:172:ASN:ND2	9:A:2067:HOH:O	2.49	0.42
2:Q:226:ASN:C	2:Q:226:ASN:OD1	2.58	0.42
2:S:265:CYS:O	2:S:268:PRO:HD2	2.20	0.42
2:R:497:VAL:HG13	2:R:498:PRO:HD2	2.01	0.42
2:R:296:ALA:HA	2:R:309:SER:HA	1.99	0.42
2:Q:329:LEU:HD11	2:Q:471:LEU:HD11	2.00	0.42
1:A:184:HIS:HB2	1:A:221:PRO:N	2.35	0.42
1:A:228:PRO:HB3	1:A:237:TRP:CZ2	2.54	0.42
1:C:208:LYS:HG2	2:S:240:TYR:CE1	2.54	0.42
2:Q:69:GLN:HA	2:Q:79:HIS:HB2	2.01	0.42
1:A:57:GLU:HG2	1:A:61:HIS:CE1	2.55	0.42
9:B:2071:HOH:O	2:R:230:GLN:HA	2.18	0.42
2:S:270:LEU:HD21	2:S:426:ALA:HA	2.01	0.42
2:S:380:ALA:HB1	2:S:514:PRO:HD3	2.01	0.42
2:S:541:ASP:N	2:S:542:PRO:HD3	2.34	0.42
2:Q:350:PRO:HB2	2:Q:482:TRP:CG	2.55	0.42
2:Q:267:ILE:HB	2:Q:268:PRO:HD3	2.00	0.42
2:Q:284:GLY:HA2	2:Q:518:ALA:O	2.20	0.42
2:Q:49:PHE:HB2	2:Q:370:TRP:CD2	2.55	0.42
2:R:49:PHE:HB2	2:R:370:TRP:CD2	2.55	0.42
2:Q:25:GLU:HG2	2:Q:74:VAL:CG2	2.50	0.41
2:R:271:LEU:HD21	2:R:407:VAL:HG22	2.02	0.41
1:C:233:ASN:HA	9:C:2010:HOH:O	2.20	0.41
1:C:115:SER:O	1:C:133:VAL:HG23	2.21	0.41
2:Q:349:HIS:CE1	2:Q:350:PRO:HD2	2.56	0.41
2:R:69:GLN:OE1	2:R:228:HIS:HA	2.20	0.41
2:Q:76:THR:OG1	2:Q:228:HIS:N	2.48	0.41
1:C:7:PRO:HG2	1:C:162:LEU:HD13	2.03	0.41
1:C:20:CYS:HB3	1:C:111:ILE:HG12	2.03	0.41
2:R:530:PRO:HB2	2:R:533:ILE:HD12	2.03	0.41
2:Q:25:GLU:HG2	2:Q:74:VAL:HG21	2.02	0.41
2:R:497:VAL:CG1	2:R:498:PRO:HD2	2.51	0.41
1:B:190:LEU:HB3	1:B:191:PRO:HD3	2.02	0.41
2:Q:248:LEU:HD23	2:Q:248:LEU:HA	1.91	0.41
2:S:126:ASP:HB3	2:S:535:ARG:HG3	2.03	0.41
1:C:226:ASN:O	1:C:230:VAL:HG22	2.20	0.41
1:A:35:LEU:HD12	1:A:39:THR:HB	2.03	0.41
2:Q:223:GLY:HA2	2:Q:231:PHE:CD1	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:223:THR:OG1	1:C:247:GLY:HA2	2.21	0.41
2:R:497:VAL:HG11	2:R:546:CYS:HB3	2.04	0.40
1:C:14:ASN:O	1:C:16:GLU:HG3	2.21	0.40
2:S:321:LEU:O	2:S:394:ALA:HB1	2.22	0.40
2:R:75:CSO:CB	6:R:1550:FCO:C2	2.99	0.40
2:R:270:LEU:HD21	2:R:426:ALA:HA	2.03	0.40
2:S:287:SER:HB3	2:S:318:GLY:HA2	2.03	0.40
5:C:1266:F3S:S2	2:S:225:LYS:HE2	2.62	0.40
2:S:39:VAL:HG23	2:S:522:THR:HB	2.04	0.40
1:A:145:PRO:HD2	1:A:179:TYR:CZ	2.56	0.40
2:R:353:GLY:HA3	2:R:494:GLN:HG3	2.03	0.40
1:C:184:HIS:HB2	1:C:221:PRO:HA	2.04	0.40
1:C:6:ARG:HA	1:C:7:PRO:HD3	1.93	0.40
2:R:248:LEU:HA	2:R:248:LEU:HD23	1.96	0.40
2:R:431:GLU:O	2:R:435:VAL:HG23	2.21	0.40
2:R:334:GLU:HA	2:R:374:PRO:HA	2.03	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	262/264 (99%)	252 (96%)	8 (3%)	2 (1%)	24	27
1	B	260/264 (98%)	251 (96%)	8 (3%)	1 (0%)	39	48
1	C	260/264 (98%)	243 (94%)	16 (6%)	1 (0%)	39	48
2	Q	546/563 (97%)	531 (97%)	15 (3%)	0	100	100
2	R	543/563 (96%)	525 (97%)	18 (3%)	0	100	100
2	S	543/563 (96%)	523 (96%)	20 (4%)	0	100	100
All	All	2414/2481 (97%)	2325 (96%)	85 (4%)	4 (0%)	52	64

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	3	ALA
1	A	18	VAL
1	B	18	VAL
1	C	18	VAL

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	209/210 (100%)	209 (100%)	0	100	100
1	B	209/210 (100%)	207 (99%)	2 (1%)	82	91
1	C	208/210 (99%)	205 (99%)	3 (1%)	74	86
2	Q	439/447 (98%)	436 (99%)	3 (1%)	88	95
2	R	435/447 (97%)	429 (99%)	6 (1%)	74	86
2	S	435/447 (97%)	428 (98%)	7 (2%)	70	84
All	All	1935/1971 (98%)	1914 (99%)	21 (1%)	80	90

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	100	LYS
1	B	171	GLU
1	C	88	MET
1	C	95	ILE
1	C	171	GLU
2	Q	122	LEU
2	Q	245	LYS
2	Q	473	ASP
2	R	361	LYS
2	R	402	LYS
2	R	412	SER
2	R	453	ASP
2	R	473	ASP

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Mol	Chain	Res	Type
2	R	503	LEU
2	S	40	LYS
2	S	267	ILE
2	S	412	SER
2	S	449	SER
2	S	473	ASP
2	S	503	LEU
2	S	517	GLU

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

Mol	Chain	Res	Type
1	A	14	ASN
1	A	172	ASN
1	B	14	ASN
1	C	14	ASN
1	C	172	ASN
2	R	123	HIS
2	S	123	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 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	CSO	Q	75	2,7,6	3,6,7	0.45	0	1,6,8	1.46	0
2	CSO	R	75	2,7,6	3,6,7	1.04	0	1,6,8	1.54	0
2	CSO	S	75	2,7,6	3,6,7	0.76	0	1,6,8	1.16	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	CSO	Q	75	2,7,6	-	0/1/5/7	0/0/0/0
2	CSO	R	75	2,7,6	-	0/1/5/7	0/0/0/0
2	CSO	S	75	2,7,6	-	0/1/5/7	0/0/0/0

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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	Q	75	CSO	1	0
2	R	75	CSO	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 23 ligands modelled in this entry, 6 are monoatomic - leaving 17 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	$\# Z > 2$
4	SF4	A	1265	1	0,12,12	0.00	-	0,24,24	0.00	-
5	F3S	A	1266	1	0,9,9	0.00	-	0,15,15	0.00	-
4	SF4	A	1267	1	0,12,12	0.00	-	0,24,24	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	A	271	-	5,5,5	0.48	0	5,5,5	1.49	1 (20%)
4	SF4	B	1265	1	0,12,12	0.00	-	0,24,24	0.00	-
5	F3S	B	1266	1	0,9,9	0.00	-	0,15,15	0.00	-
4	SF4	B	1267	1	0,12,12	0.00	-	0,24,24	0.00	-
4	SF4	C	1265	1	0,12,12	0.00	-	0,24,24	0.00	-
5	F3S	C	1266	1	0,9,9	0.00	-	0,15,15	0.00	-
4	SF4	C	1267	1	0,12,12	0.00	-	0,24,24	0.00	-
6	FCO	Q	1550	2	0,6,6	0.00	-	0,6,6	0.00	-
3	GOL	Q	1561	-	5,5,5	0.32	0	5,5,5	0.55	0
3	GOL	Q	1562	-	5,5,5	0.31	0	5,5,5	0.30	0
6	FCO	R	1550	2	0,6,6	0.00	-	0,6,6	0.00	-
3	GOL	R	1563	-	5,5,5	0.55	0	5,5,5	0.48	0
6	FCO	S	1550	2	0,6,6	0.00	-	0,6,6	0.00	-
3	GOL	S	1561	-	5,5,5	0.40	0	5,5,5	0.63	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
4	SF4	A	1265	1	-	0/0/48/48	0/6/5/5
5	F3S	A	1266	1	-	0/0/24/24	0/0/3/3
4	SF4	A	1267	1	-	0/0/48/48	0/6/5/5
3	GOL	A	271	-	-	0/4/4/4	0/0/0/0
4	SF4	B	1265	1	-	0/0/48/48	0/6/5/5
5	F3S	B	1266	1	-	0/0/24/24	0/0/3/3
4	SF4	B	1267	1	-	0/0/48/48	0/6/5/5
4	SF4	C	1265	1	-	0/0/48/48	0/6/5/5
5	F3S	C	1266	1	-	0/0/24/24	0/0/3/3
4	SF4	C	1267	1	-	0/0/48/48	0/6/5/5
6	FCO	Q	1550	2	-	0/0/6/6	0/0/0/0
3	GOL	Q	1561	-	-	0/4/4/4	0/0/0/0
3	GOL	Q	1562	-	-	0/4/4/4	0/0/0/0
6	FCO	R	1550	2	-	0/0/6/6	0/0/0/0
3	GOL	R	1563	-	-	0/4/4/4	0/0/0/0
6	FCO	S	1550	2	-	0/0/6/6	0/0/0/0
3	GOL	S	1561	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	A	271	GOL	O1-C1-C2	-2.12	99.88	110.18

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	1265	SF4	1	0
5	C	1266	F3S	1	0
6	Q	1550	FCO	1	0
3	Q	1562	GOL	1	0
6	R	1550	FCO	1	0

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	263/264 (99%)	-0.36	0 100 100	10, 20, 34, 66	7 (2%)
1	B	260/264 (98%)	-0.32	0 100 100	14, 22, 35, 61	7 (2%)
1	C	261/264 (98%)	0.48	22 (8%) 14 19	21, 46, 81, 102	5 (1%)
2	Q	543/563 (96%)	-0.43	0 100 100	9, 19, 31, 43	12 (2%)
2	R	544/563 (96%)	-0.22	3 (0%) 90 93	14, 28, 47, 69	12 (2%)
2	S	544/563 (96%)	0.04	12 (2%) 65 73	15, 32, 59, 82	14 (2%)
All	All	2415/2481 (97%)	-0.16	37 (1%) 76 81	9, 25, 56, 102	57 (2%)

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	191	PRO	6.5
2	S	449	SER	5.1
1	C	213	LEU	4.6
1	C	211	PHE	4.0
1	C	169	LEU	4.0
1	C	264	GLY	3.8
1	C	197	GLU	3.5
1	C	176	LYS	3.4
2	S	148	ILE	3.3
2	S	458	ALA	3.3
1	C	140	LYS	3.3
1	C	198	PHE	3.2
1	C	177	LEU	3.0
2	S	448	ASP	2.7
2	S	97	ALA	2.6
1	C	192	HIS	2.6
1	C	142	ILE	2.5
1	C	216	LEU	2.4
1	C	207	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	71	TYR	2.4
1	C	183	VAL	2.4
2	S	143	LYS	2.4
1	C	224	TYR	2.3
2	S	453	ASP	2.3
1	C	214	TYR	2.3
2	S	249	ALA	2.2
2	S	447	ALA	2.2
2	R	537	VAL	2.2
2	R	345	GLY	2.1
1	C	204	SER	2.1
2	S	221	ILE	2.1
2	R	157	ALA	2.1
1	C	163	THR	2.0
2	S	244	THR	2.0
2	S	297	THR	2.0
1	C	188	PRO	2.0
1	C	179	TYR	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	CSO	R	75	7/8	0.98	0.09	-	13,14,15,15	0
2	CSO	S	75	7/8	0.98	0.08	-	19,21,24,25	0
2	CSO	Q	75	7/8	0.96	0.11	-	8,9,11,13	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron

density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	GOL	A	271	6/6	0.87	0.24	8.68	23,26,28,30	0
3	GOL	Q	1562	6/6	0.94	0.17	2.54	20,22,25,25	0
3	GOL	S	1561	6/6	0.90	0.19	0.64	20,24,26,29	0
6	FCO	R	1550	7/7	0.99	0.09	-0.19	14,15,16,16	0
4	SF4	A	1267	8/8	0.99	0.08	-0.76	10,11,11,13	0
8	MG	R	1553	1/1	0.94	0.11	-0.76	17,17,17,17	0
8	MG	Q	1553	1/1	0.94	0.08	-0.86	13,13,13,13	0
4	SF4	A	1265	8/8	0.98	0.09	-1.12	20,21,22,22	0
4	SF4	B	1267	8/8	0.98	0.08	-1.12	14,15,16,17	0
3	GOL	Q	1561	6/6	0.96	0.09	-1.16	16,17,18,19	0
6	FCO	Q	1550	7/7	0.99	0.08	-1.31	11,12,15,18	0
4	SF4	C	1267	8/8	0.98	0.08	-1.60	27,30,34,36	0
4	SF4	B	1265	8/8	0.99	0.08	-1.67	18,19,20,20	0
4	SF4	C	1265	8/8	0.95	0.07	-1.72	66,70,75,76	0
3	GOL	R	1563	6/6	0.97	0.08	-1.78	14,17,18,19	0
6	FCO	S	1550	7/7	0.99	0.08	-1.87	16,17,17,17	0
5	F3S	C	1266	7/7	0.98	0.05	-2.02	42,44,47,49	0
7	NI	R	1551	1/1	0.98	0.06	-2.07	17,17,17,17	0
7	NI	S	1551	1/1	0.98	0.05	-2.14	23,23,23,23	0
7	NI	Q	1551	1/1	0.98	0.06	-2.70	13,13,13,13	0
5	F3S	B	1266	7/7	0.99	0.05	-2.88	16,17,17,18	0
5	F3S	A	1266	7/7	0.99	0.06	-2.89	14,15,17,17	0
8	MG	S	1553	1/1	0.96	0.04	-3.76	19,19,19,19	0

6.5 Other polymers [i](#)

There are no such residues in this entry.