



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 09:24 PM GMT

PDB ID : 5A4M
Title : Mechanism of Hydrogen activation by NiFe-hydrogenases
Authors : Evans, R.M.; Brooke, E.J.; Wehlin, S.A.M.; Nomerotskaia, E.; Sergent, F.;
Carr, S.B.; Philips, S.E.V.; Armstrong, F.A.
Deposited on : 2015-06-10
Resolution : 1.70 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
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) : rb-20026982

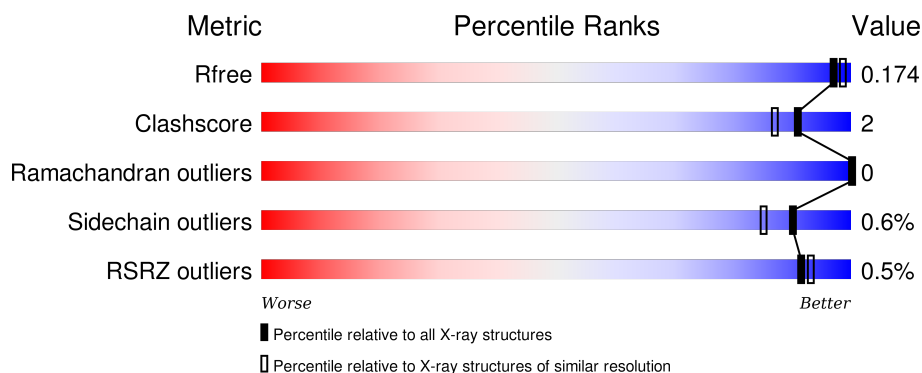
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	3190 (1.70-1.70)
Clashscore	102246	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)
RSRZ outliers	91569	3200 (1.70-1.70)

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 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	L	582	<div> <div>%</div> <div> <div></div> <div>95%</div> <div>5%</div> </div> </div>
1	M	582	<div> <div></div> <div>94%</div> <div>6%</div> </div>
2	S	278	<div> <div>%</div> <div> <div></div> <div>88%</div> <div>6% • 5%</div> </div> </div>
2	T	278	<div> <div></div> <div>85%</div> <div>8% •• 6%</div> </div>

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
8	CL	T	409	-	-	-	X

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 14827 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-1 LARGE CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	581	Total	C	N	O	S	0	8	0
			4600	2919	809	844	28			
1	M	581	Total	C	N	O	S	0	7	0
			4578	2906	798	846	28			

- Molecule 2 is a protein called HYDROGENASE-1 SMALL CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	S	263	Total	C	N	O	S	4	2	0
			2045	1298	353	374	20			
2	T	262	Total	C	N	O	S	0	6	0
			2046	1301	346	378	21			

There are 24 discrepancies between the modelled and reference sequences:

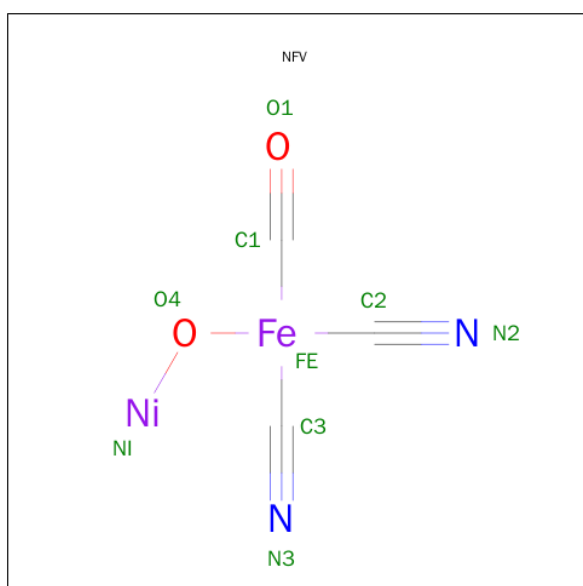
Chain	Residue	Modelled	Actual	Comment	Reference
S	267	GLY	-	EXPRESSION TAG	UNP P69739
S	268	SER	-	EXPRESSION TAG	UNP P69739
S	269	PHE	-	EXPRESSION TAG	UNP P69739
S	270	TYR	-	EXPRESSION TAG	UNP P69739
S	271	SER	-	EXPRESSION TAG	UNP P69739
S	272	ARG	-	EXPRESSION TAG	UNP P69739
S	273	HIS	-	EXPRESSION TAG	UNP P69739
S	274	HIS	-	EXPRESSION TAG	UNP P69739
S	275	HIS	-	EXPRESSION TAG	UNP P69739
S	276	HIS	-	EXPRESSION TAG	UNP P69739
S	277	HIS	-	EXPRESSION TAG	UNP P69739
S	278	HIS	-	EXPRESSION TAG	UNP P69739
T	267	GLY	-	EXPRESSION TAG	UNP P69739
T	268	SER	-	EXPRESSION TAG	UNP P69739
T	269	PHE	-	EXPRESSION TAG	UNP P69739
T	270	TYR	-	EXPRESSION TAG	UNP P69739

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Chain	Residue	Modelled	Actual	Comment	Reference
T	271	SER	-	EXPRESSION TAG	UNP P69739
T	272	ARG	-	EXPRESSION TAG	UNP P69739
T	273	HIS	-	EXPRESSION TAG	UNP P69739
T	274	HIS	-	EXPRESSION TAG	UNP P69739
T	275	HIS	-	EXPRESSION TAG	UNP P69739
T	276	HIS	-	EXPRESSION TAG	UNP P69739
T	277	HIS	-	EXPRESSION TAG	UNP P69739
T	278	HIS	-	EXPRESSION TAG	UNP P69739

- Molecule 3 is NI-FE OXIDIZED ACTIVE CENTER (three-letter code: NFV) (formula: $C_3FeN_2NiO_2$).

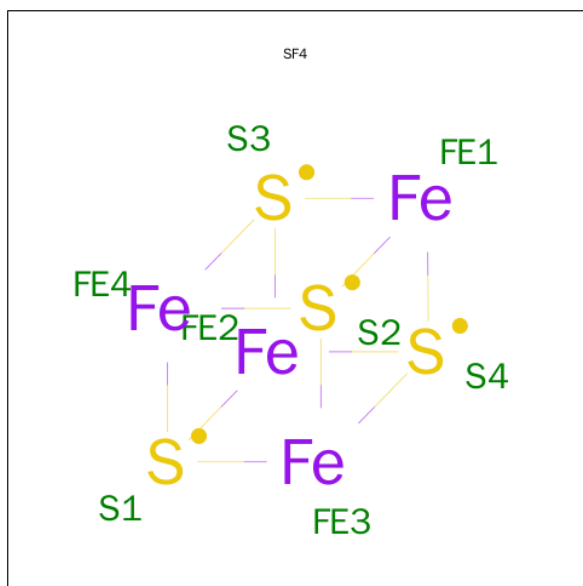


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	L	1	Total	C	Fe	N	Ni	O	
			9	3	1	2	1	2	0
3	M	1	Total	C	Fe	N	Ni	O	
			9	3	1	2	1	2	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

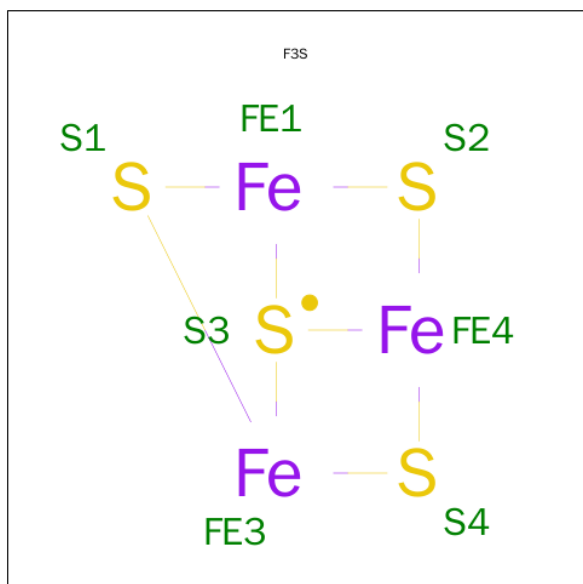
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	1	Total	Mg		
			1	1	0	0
4	M	1	Total	Mg		
			1	1	0	0

- Molecule 5 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe_4S_4).



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

- Molecule 6 is FE3-S4 CLUSTER (three-letter code: F3S) (formula: Fe_3S_4).



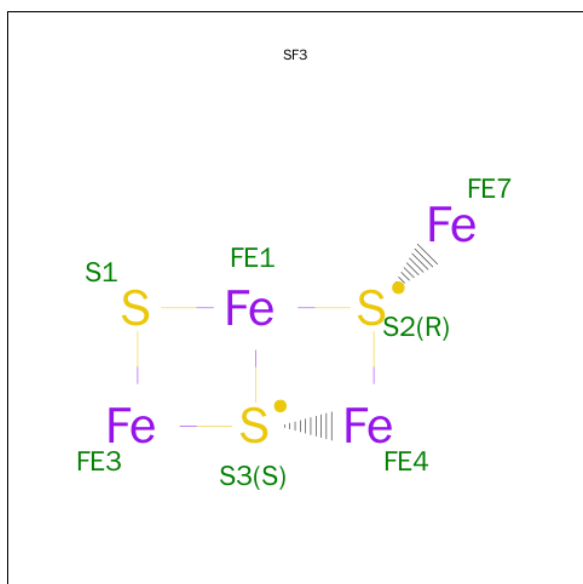
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	S	1	Total	Fe	S	0	0
			7	3	4		

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

- Molecule 7 is FE4-S3 CLUSTER (three-letter code: SF3) (formula: Fe_4S_3).



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

- Molecule 8 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	T	1	Total	Cl	0	0
			1	1		
8	S	1	Total	Cl	0	0
			1	1		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	L	570	Total	O	0	0
			570	570		
9	M	560	Total	O	0	0
			560	560		

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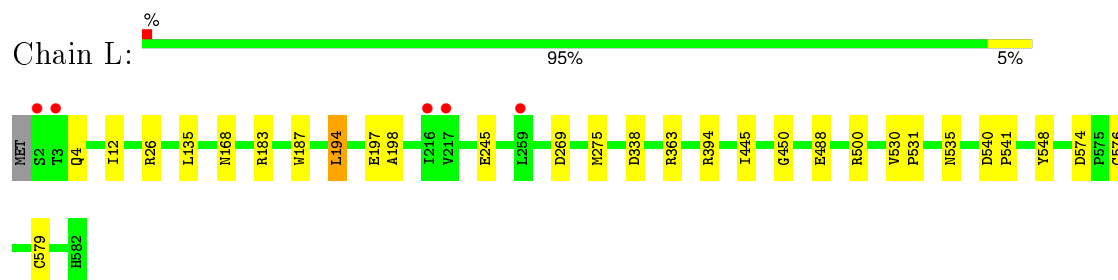
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	S	189	Total	O	0	0
			189	189		
9	T	173	Total	O	0	0
			173	173		

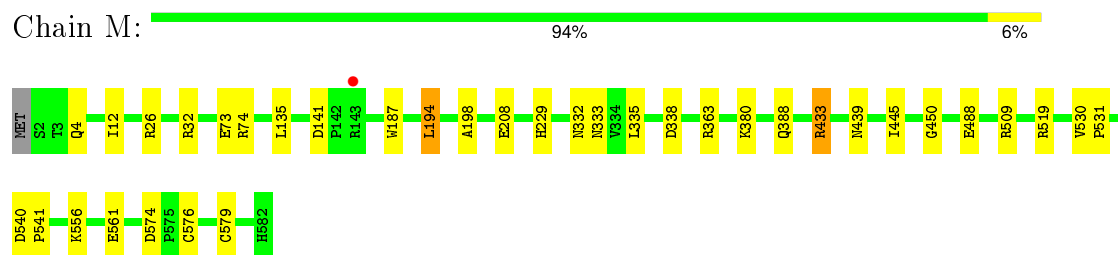
3 Residue-property plots [i](#)

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.

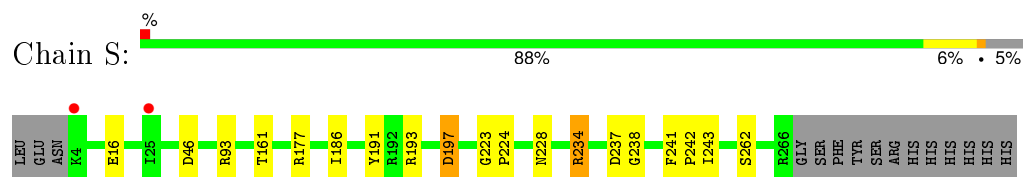
• Molecule 1: HYDROGENASE-1 LARGE CHAIN



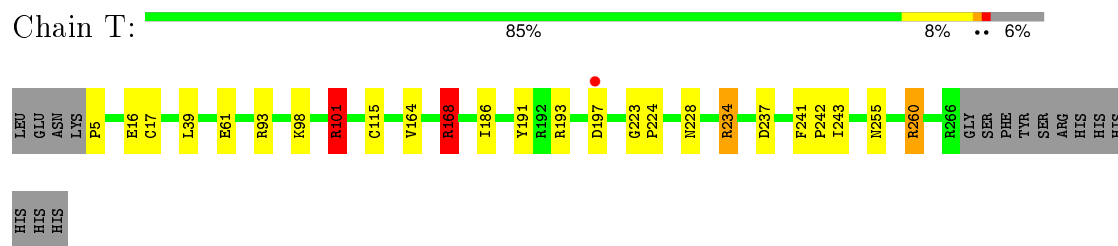
• Molecule 1: HYDROGENASE-1 LARGE CHAIN



• Molecule 2: HYDROGENASE-1 SMALL CHAIN



• Molecule 2: HYDROGENASE-1 SMALL CHAIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	93.67Å 97.16Å 182.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	91.62 – 1.70 46.95 – 1.70	Depositor EDS
% Data completeness (in resolution range)	98.9 (91.62-1.70) 98.9 (46.95-1.70)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.8.0123	Depositor
R, R_{free}	0.134 , 0.162 0.149 , 0.174	Depositor DCC
R_{free} test set	9069 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	17.9	Xtriage
Anisotropy	0.076	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.1	EDS
Estimated twinning fraction	0.018 for k,h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 181398 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	14827	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.16% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: MG, CL, SF4, SF3, NFV, F3S

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	L	0.90	3/4740 (0.1%)	0.94	11/6448 (0.2%)
1	M	0.90	3/4718 (0.1%)	0.93	13/6421 (0.2%)
2	S	0.95	2/2109 (0.1%)	1.00	12/2864 (0.4%)
2	T	0.94	0/2119	1.01	11/2878 (0.4%)
All	All	0.91	8/13686 (0.1%)	0.96	47/18611 (0.3%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	245	GLU	CD-OE2	7.11	1.33	1.25
1	M	73	GLU	CD-OE1	6.45	1.32	1.25
2	S	262	SER	CB-OG	5.70	1.49	1.42
2	S	197	ASP	CB-CG	5.32	1.62	1.51
1	L	197	GLU	CD-OE1	5.29	1.31	1.25
1	M	561	GLU	CD-OE2	5.22	1.31	1.25
1	L	363	ARG	CZ-NH1	-5.11	1.26	1.33
1	M	488	GLU	CD-OE1	5.00	1.31	1.25

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	260	ARG	NE-CZ-NH1	13.41	127.00	120.30
2	T	260	ARG	NE-CZ-NH2	-11.41	114.59	120.30
2	S	234	ARG	NE-CZ-NH1	9.95	125.28	120.30
2	T	168	ARG	NE-CZ-NH1	8.86	124.73	120.30
2	T	193	ARG	NE-CZ-NH2	-8.12	116.24	120.30
2	T	93	ARG	NE-CZ-NH1	8.00	124.30	120.30
1	M	576	CYS	CA-CB-SG	-7.90	99.77	114.00
1	M	519	ARG	NE-CZ-NH2	-7.55	116.52	120.30
1	L	576	CYS	CA-CB-SG	-7.50	100.49	114.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	275[A]	MET	CG-SD-CE	7.41	112.05	100.20
1	L	275[B]	MET	CG-SD-CE	7.41	112.05	100.20
1	L	574	ASP	CB-CG-OD2	-7.29	111.73	118.30
1	M	574	ASP	CB-CG-OD2	-7.25	111.78	118.30
2	T	193	ARG	NE-CZ-NH1	6.96	123.78	120.30
2	S	193	ARG	NE-CZ-NH2	-6.86	116.87	120.30
2	S	197	ASP	CB-CG-OD1	6.79	124.41	118.30
1	M	363	ARG	NE-CZ-NH1	6.77	123.69	120.30
1	L	338	ASP	CB-CG-OD1	6.71	124.34	118.30
1	M	338	ASP	CB-CG-OD1	6.70	124.33	118.30
1	L	363	ARG	NE-CZ-NH2	-6.66	116.97	120.30
1	M	519	ARG	NE-CZ-NH1	6.62	123.61	120.30
2	S	93	ARG	NE-CZ-NH1	6.57	123.59	120.30
2	S	234	ARG	CD-NE-CZ	6.57	132.80	123.60
1	M	574	ASP	CB-CG-OD1	6.39	124.05	118.30
1	M	32	ARG	NE-CZ-NH1	6.37	123.48	120.30
1	L	363	ARG	NE-CZ-NH1	6.32	123.46	120.30
2	S	193	ARG	NE-CZ-NH1	6.23	123.41	120.30
2	S	197	ASP	CB-CG-OD2	-6.21	112.71	118.30
1	L	26	ARG	NE-CZ-NH1	6.09	123.34	120.30
2	S	237	ASP	CB-CG-OD2	-5.83	113.05	118.30
1	M	141	ASP	CB-CG-OD2	5.65	123.39	118.30
2	T	237	ASP	CB-CG-OD2	-5.62	113.24	118.30
1	L	194	LEU	CB-CG-CD2	5.50	120.34	111.00
1	M	26	ARG	NE-CZ-NH1	5.49	123.04	120.30
1	L	394	ARG	NE-CZ-NH1	5.48	123.04	120.30
2	T	193	ARG	CG-CD-NE	-5.40	100.46	111.80
2	T	234	ARG	NE-CZ-NH1	5.38	122.99	120.30
2	S	193	ARG	CG-CD-NE	-5.38	100.51	111.80
1	M	74	ARG	NE-CZ-NH2	-5.31	117.65	120.30
1	L	500	ARG	NE-CZ-NH2	-5.28	117.66	120.30
2	S	234	ARG	NE-CZ-NH2	-5.21	117.69	120.30
1	M	433	ARG	NE-CZ-NH1	5.15	122.87	120.30
2	T	101	ARG	NE-CZ-NH1	5.08	122.84	120.30
2	T	93	ARG	NE-CZ-NH2	-5.08	117.76	120.30
2	S	46	ASP	CB-CG-OD1	5.07	122.87	118.30
1	M	194	LEU	CB-CG-CD2	5.05	119.59	111.00
2	S	177	ARG	NE-CZ-NH1	5.00	122.80	120.30

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	L	4600	0	4497	17	0
1	M	4578	0	4465	18	0
2	S	2045	0	1988	10	0
2	T	2046	0	1986	13	0
3	L	9	0	0	0	0
3	M	9	0	0	1	0
4	L	1	0	0	0	0
4	M	1	0	0	0	0
5	S	8	0	0	0	0
5	T	8	0	0	0	0
6	S	7	0	0	0	0
6	T	7	0	0	0	0
7	S	7	0	0	0	0
7	T	7	0	0	0	0
8	S	1	0	0	0	0
8	T	1	0	0	0	0
9	L	570	0	0	9	0
9	M	560	0	0	6	0
9	S	189	0	0	6	0
9	T	173	0	0	4	0
All	All	14827	0	12936	57	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 2.

All (57) 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:T:5:PRO:N	9:T:2002:HOH:O	2.09	0.86
1:M:208[A]:GLU:OE2	9:M:2192:HOH:O	2.00	0.80
9:L:2511:HOH:O	2:T:168:ARG:NH2	1.95	0.76
2:S:234:ARG:NH2	9:S:2177:HOH:O	2.18	0.76
1:L:579:CYS:CB	9:L:2106:HOH:O	2.37	0.72
2:T:168:ARG:NH1	9:T:2100:HOH:O	2.24	0.71
1:L:183[B]:ARG:HG3	9:L:2245:HOH:O	1.90	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:S:161[A]:THR:HG23	9:S:2118:HOH:O	1.91	0.69
2:S:197:ASP:HB3	9:S:2143:HOH:O	1.95	0.66
2:S:234:ARG:NH1	2:S:238:GLY:O	2.29	0.66
1:L:168:ASN:CG	9:L:2217:HOH:O	2.34	0.65
1:M:332:ASN:HB2	9:M:2367:HOH:O	1.97	0.65
2:T:186:ILE:HD11	2:T:228:ASN:HB3	1.84	0.59
1:M:433:ARG:HD2	9:M:2477:HOH:O	2.04	0.57
1:L:579:CYS:N	9:L:2106:HOH:O	2.31	0.57
1:L:168:ASN:CB	9:L:2217:HOH:O	2.51	0.57
2:S:186:ILE:HD11	2:S:228:ASN:HB3	1.86	0.56
2:T:255:ASN:O	2:T:260:ARG:NH2	2.39	0.55
2:T:61[B]:GLU:OE1	2:T:101:ARG:NH2	2.41	0.54
1:M:530:VAL:CG1	1:M:531:PRO:HD2	2.39	0.53
1:L:269[A]:ASP:OD1	9:L:2319:HOH:O	2.19	0.52
2:S:197:ASP:CB	9:S:2143:HOH:O	2.54	0.51
1:L:530:VAL:CG1	1:L:531:PRO:HD2	2.42	0.49
1:M:540:ASP:HB2	1:M:541:PRO:CD	2.42	0.49
2:T:39[B]:LEU:HG	2:T:164:VAL:HG21	1.95	0.48
1:M:380:LYS:HB3	1:M:388:GLN:HB2	1.96	0.47
2:S:234:ARG:NH1	2:T:234:ARG:HE	2.12	0.46
1:M:530:VAL:HG11	1:M:579:CYS:HB3	1.96	0.46
1:M:530:VAL:HG12	1:M:531:PRO:HD2	1.97	0.46
2:T:241:PHE:CE2	2:T:243:ILE:HB	2.51	0.46
1:L:530:VAL:HG11	1:L:579:CYS:HB3	1.97	0.45
1:L:530:VAL:HG12	1:L:531:PRO:HD2	1.98	0.45
2:S:234:ARG:NE	9:S:2174:HOH:O	2.48	0.45
1:L:445:ILE:O	1:L:450:GLY:HA3	2.17	0.45
1:M:135:LEU:HD22	1:M:187:TRP:CD1	2.52	0.44
1:L:135:LEU:HD22	1:L:187:TRP:CD1	2.53	0.44
1:L:488:GLU:HB2	9:L:2519:HOH:O	2.17	0.43
1:L:535:ASN:HB3	1:L:548:TYR:CE1	2.54	0.43
2:T:115:CYS:HB2	9:T:2008:HOH:O	2.19	0.43
9:S:2145:HOH:O	2:T:197:ASP:HB2	2.19	0.42
1:L:194:LEU:HB3	1:L:198:ALA:HB3	2.00	0.42
1:M:333:ASN:HD21	1:M:335:LEU:HD21	1.84	0.42
2:T:98:LYS:NZ	9:T:2029:HOH:O	2.52	0.42
1:M:445:ILE:O	1:M:450:GLY:HA3	2.19	0.42
1:L:4:GLN:HA	1:L:12:ILE:O	2.19	0.42
1:L:579:CYS:HB2	9:L:2106:HOH:O	2.13	0.42
1:M:439:ASN:ND2	9:M:2480:HOH:O	2.53	0.42
1:L:540:ASP:HB2	1:L:541:PRO:CD	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:333:ASN:ND2	1:M:335:LEU:HD21	2.35	0.41
1:M:4:GLN:HA	1:M:12:ILE:O	2.21	0.41
1:M:509:ARG:HD2	3:M:603:NFV:O4	2.21	0.41
1:M:556:LYS:HE3	9:M:2053:HOH:O	2.21	0.41
1:M:194:LEU:HB3	1:M:198:ALA:HB3	2.02	0.41
2:S:241:PHE:CE2	2:S:243:ILE:HB	2.56	0.40
2:S:223:GLY:N	2:S:224:PRO:CD	2.84	0.40
1:M:229:HIS:NE2	9:M:2265:HOH:O	2.30	0.40
2:T:223:GLY:N	2:T:224:PRO:CD	2.85	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	L	587/582 (101%)	570 (97%)	17 (3%)	0	100	100
1	M	586/582 (101%)	571 (97%)	15 (3%)	0	100	100
2	S	264/278 (95%)	253 (96%)	11 (4%)	0	100	100
2	T	266/278 (96%)	254 (96%)	12 (4%)	0	100	100
All	All	1703/1720 (99%)	1648 (97%)	55 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	L	488/481 (102%)	488 (100%)	0	100	100
1	M	487/481 (101%)	487 (100%)	0	100	100
2	S	218/230 (95%)	215 (99%)	3 (1%)	74	59
2	T	220/230 (96%)	214 (97%)	6 (3%)	52	31
All	All	1413/1422 (99%)	1404 (99%)	9 (1%)	90	85

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	S	16	GLU
2	S	191	TYR
2	S	242	PRO
2	T	16	GLU
2	T	17	CYS
2	T	101	ARG
2	T	168	ARG
2	T	191	TYR
2	T	242	PRO

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

Mol	Chain	Res	Type
1	L	332	ASN
1	L	392	GLN
1	M	333	ASN
1	M	387	GLN
1	M	439	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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
3	NFV	L	601	1	3,8,8	0.91	0	0,10,10	0.00	-
3	NFV	M	603	1	3,8,8	1.48	1 (33%)	0,10,10	0.00	-
5	SF4	S	401	2	0,12,12	0.00	-	0,24,24	0.00	-
6	F3S	S	402	2	0,9,9	0.00	-	0,15,15	0.00	-
7	SF3	S	404	9,2	0,8,8	0.00	-	0,12,12	0.00	-
5	SF4	T	401	2	0,12,12	0.00	-	0,24,24	0.00	-
6	F3S	T	402	2	0,9,9	0.00	-	0,15,15	0.00	-
7	SF3	T	404	9,2	0,8,8	0.00	-	0,12,12	0.00	-

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	NFV	L	601	1	-	0/0/12/12	0/0/0/0
3	NFV	M	603	1	-	0/0/12/12	0/0/0/0
5	SF4	S	401	2	-	0/0/48/48	0/6/5/5
6	F3S	S	402	2	-	0/0/24/24	0/0/3/3
7	SF3	S	404	9,2	-	0/0/17/17	0/2/2/2
5	SF4	T	401	2	-	0/0/48/48	0/6/5/5
6	F3S	T	402	2	-	0/0/24/24	0/0/3/3
7	SF3	T	404	9,2	-	0/0/17/17	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	M	603	NFV	C3-N3	2.18	1.18	1.13

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	M	603	NFV	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	L	581/582 (99%)	-0.46	5 (0%) 85 88	13, 19, 32, 51	3 (0%)
1	M	581/582 (99%)	-0.54	1 (0%) 95 95	13, 19, 30, 53	1 (0%)
2	S	263/278 (94%)	-0.40	2 (0%) 87 90	13, 18, 28, 66	1 (0%)
2	T	262/278 (94%)	-0.34	1 (0%) 93 94	14, 19, 33, 44	1 (0%)
All	All	1687/1720 (98%)	-0.46	9 (0%) 91 93	13, 19, 31, 66	6 (0%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	2	SER	3.2
2	S	4	LYS	2.9
2	T	197	ASP	2.4
1	L	217	VAL	2.4
1	L	216	ILE	2.3
1	L	3	THR	2.2
2	S	25	ILE	2.2
1	L	259[A]	LEU	2.1
1	M	143	ARG	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 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
8	CL	T	409	1/1	0.98	0.12	2.18	32,32,32,32	0
8	CL	S	408	1/1	0.99	0.06	0.44	29,29,29,29	0
4	MG	M	604	1/1	1.00	0.06	-0.65	13,13,13,13	0
5	SF4	S	401	8/8	0.98	0.05	-0.92	15,17,18,20	0
3	NFV	L	601	9/9	0.99	0.06	-0.96	14,15,19,26	0
4	MG	L	602	1/1	0.99	0.06	-1.00	13,13,13,13	0
5	SF4	T	401	8/8	0.98	0.05	-1.10	15,16,17,20	0
3	NFV	M	603	9/9	0.99	0.06	-1.19	14,17,19,26	0
6	F3S	S	402	7/7	1.00	0.03	-1.30	15,15,16,17	0
6	F3S	T	402	7/7	0.99	0.03	-1.61	16,16,17,18	0
7	SF3	S	404	7/7	0.98	0.04	-1.68	19,23,32,34	0
7	SF3	T	404	7/7	0.98	0.04	-1.68	20,24,30,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.