



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 4, 2026 – 10:40 PM EDT

PDB ID : 1UBO / pdb_00001ubo
Title : Three-dimensional Structure of The Carbon Monoxide Complex of [NiFe]hydrogenase From Desulfovibrio vulgaris Miyazaki F
Authors : Ogata, H.; Mizoguchi, Y.; Mizuno, N.; Miki, K.; Adachi, S.; Yasuoka, N.; Yagi, T.; Yamauchi, O.; Hirota, S.; Higuchi, Y.
Deposited on : 2003-04-04
Resolution : 1.35 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

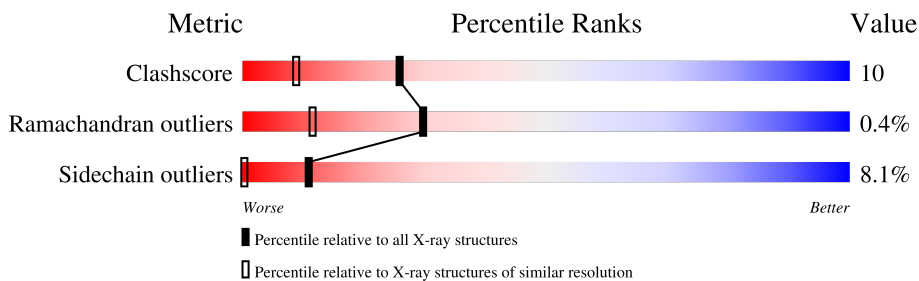
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.35 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))
Clashscore	190562	1232 (1.36-1.36)
Ramachandran outliers	187476	1220 (1.36-1.36)
Sidechain outliers	187428	1220 (1.36-1.36)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	S	267	
2	L	534	

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 7179 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 Periplasmic [NiFe] hydrogenase Small subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	S	267	Total	C	N	O	S	0	0	0
			2019	1282	342	378	17			

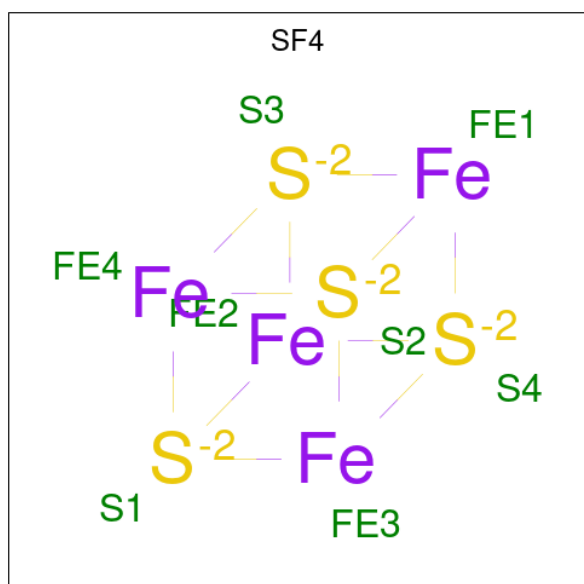
- Molecule 2 is a protein called Periplasmic [NiFe] hydrogenase Large subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	534	Total	C	N	O	S	0	0	0
			4177	2674	725	763	15			

There are 2 discrepancies between the modelled and reference sequences:

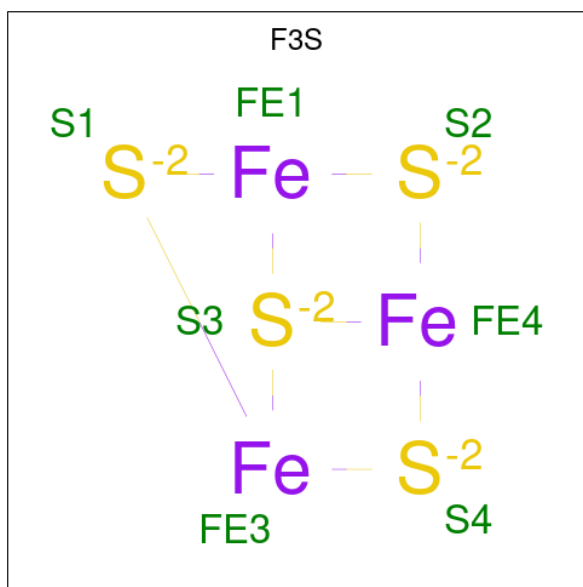
Chain	Residue	Modelled	Actual	Comment	Reference
L	514	LYS	ASN	SEE REMARK 999	UNP P21852
L	515	LEU	VAL	SEE REMARK 999	UNP P21852

- 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		

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

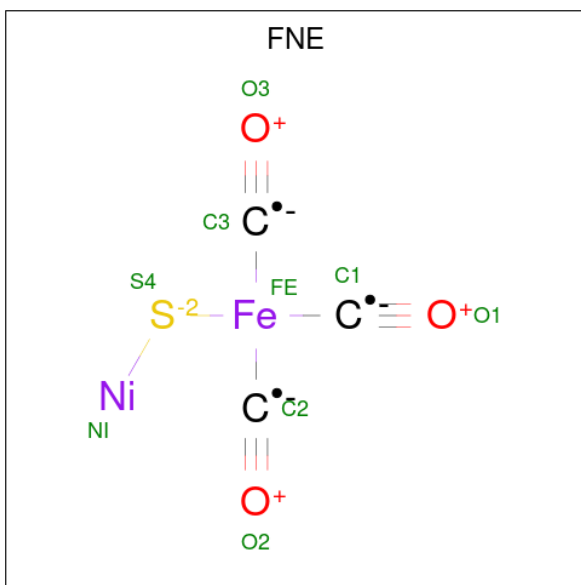


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	S	1	Total	Fe	S	0	0
			7	3	4		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

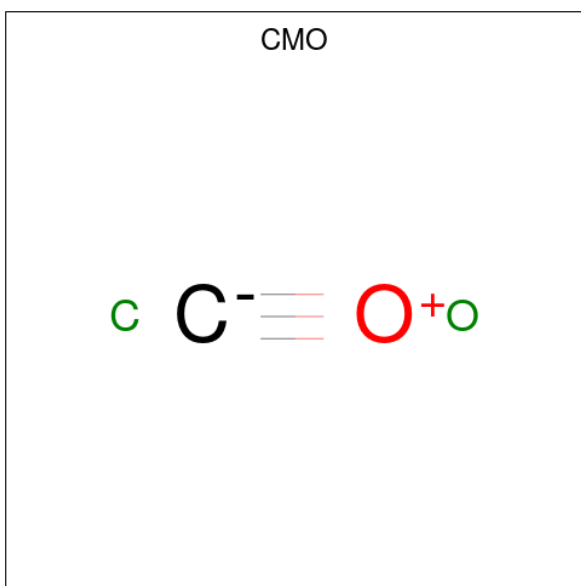
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	L	1	Total	Mg	0	0
			1	1		

- Molecule 6 is (MU-SULPHIDO)-BIS(MU-CYS,S)-[TRICARBONYLIRON-DI-(CYS,S)NICKEL(II)](FE-NI) (CCD ID: FNE) (formula: $\text{C}_3\text{FeNiO}_3\text{S}$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	L	1	Total	C	Fe	Ni	O	0	0
			8	3	1	1	3		

- Molecule 7 is CARBON MONOXIDE (CCD ID: CMO) (formula: CO).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	L	1	Total	C	O	0	0
			2	1	1		

- Molecule 8 is water.

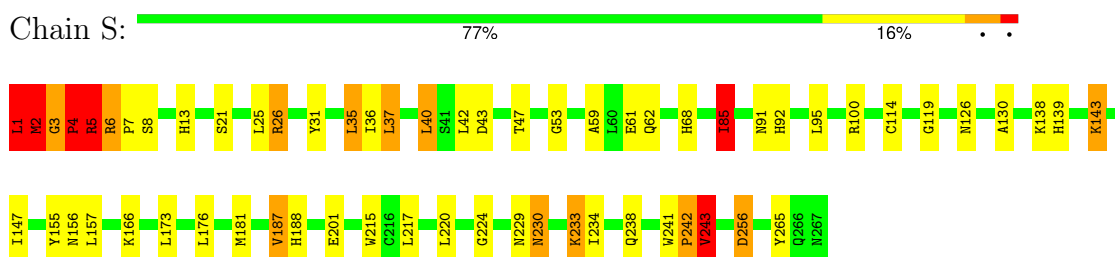
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	S	329	Total 329	O 329	0	0
8	L	620	Total 620	O 620	0	0

3 Residue-property plots

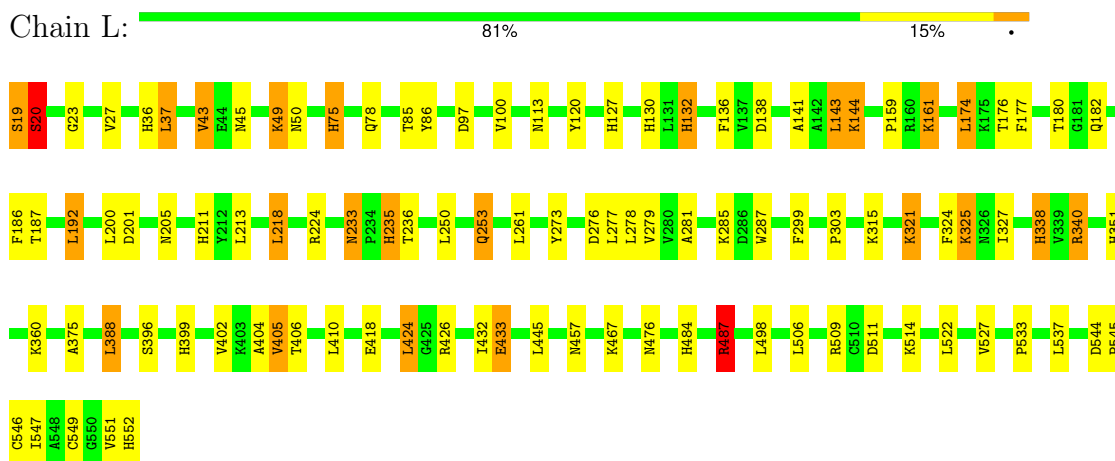
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

- Molecule 1: Periplasmic [NiFe] hydrogenase Small subunit



- Molecule 2: Periplasmic [NiFe] hydrogenase Large subunit



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	98.00Å 125.98Å 66.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.35	Depositor
% Data completeness (in resolution range)	87.4 (20.00-1.35)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS, SHELXL-97	Depositor
R, R_{free}	0.117 , 0.162	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7179	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.80	0/2075	1.60	23/2830 (0.8%)
2	L	0.78	3/4288 (0.1%)	1.43	33/5831 (0.6%)
All	All	0.78	3/6363 (0.0%)	1.48	56/8661 (0.6%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	235	HIS	ND1-CE1	5.47	1.38	1.32
2	L	75	HIS	ND1-CE1	5.46	1.38	1.32
2	L	552	HIS	C-O	5.22	1.33	1.23

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	S	6	ARG	CD-NE-CZ	30.10	166.54	124.40
2	L	487	ARG	CB-CG-CD	9.62	133.43	111.30
2	L	487	ARG	NE-CZ-NH2	9.48	127.73	119.20
1	S	5	ARG	CA-C-O	-9.22	110.78	120.55
1	S	1	LEU	O-C-N	8.94	137.30	123.00
1	S	2	MET	CA-CB-CG	8.71	131.52	114.10
2	L	340	ARG	CG-CD-NE	8.60	130.92	112.00
1	S	2	MET	CA-C-N	8.40	135.05	121.87
1	S	2	MET	C-N-CA	8.40	135.05	121.87
2	L	404	ALA	CA-C-N	-8.27	108.54	121.02
2	L	404	ALA	C-N-CA	-8.27	108.54	121.02
2	L	273	TYR	CA-C-N	-8.03	115.18	120.24
2	L	273	TYR	C-N-CA	-8.03	115.18	120.24
2	L	132	HIS	CB-CG-CD2	-7.81	121.04	131.20
1	S	5	ARG	CD-NE-CZ	7.68	135.16	124.40
2	L	506	LEU	CA-CB-CG	7.65	143.07	116.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	S	4	PRO	CA-C-N	7.62	130.50	120.28
1	S	4	PRO	C-N-CA	7.62	130.50	120.28
2	L	130	HIS	CA-CB-CG	-7.28	106.52	113.80
1	S	242	PRO	CA-C-N	-6.83	109.43	120.64
1	S	242	PRO	C-N-CA	-6.83	109.43	120.64
2	L	75	HIS	CA-CB-CG	-6.68	107.12	113.80
2	L	75	HIS	CB-CG-CD2	-6.65	122.55	131.20
2	L	49	LYS	CA-CB-CG	6.65	127.40	114.10
2	L	287	TRP	CA-CB-CG	-6.55	101.15	113.60
2	L	45	ASN	OD1-CG-ND2	6.45	129.05	122.60
1	S	47	THR	CA-C-N	-6.28	115.17	122.14
1	S	47	THR	C-N-CA	-6.28	115.17	122.14
2	L	551	VAL	CB-CA-C	-6.24	104.37	111.55
1	S	85	ILE	CA-CB-CG1	-6.04	100.13	110.40
2	L	218	LEU	CA-CB-CG	-5.95	95.48	116.30
2	L	351	HIS	CA-CB-CG	-5.88	107.92	113.80
2	L	511	ASP	CA-CB-CG	5.77	118.37	112.60
1	S	26	ARG	CG-CD-NE	-5.68	99.50	112.00
2	L	433	GLU	CA-CB-CG	5.61	125.32	114.10
1	S	2	MET	CB-CG-SD	5.61	129.51	112.70
1	S	233	LYS	CD-CE-NZ	5.53	129.61	111.90
2	L	201	ASP	CA-CB-CG	5.50	118.10	112.60
1	S	243	VAL	CA-CB-CG1	5.49	119.73	110.40
2	L	75	HIS	CB-CG-ND1	5.45	130.88	122.70
2	L	127	HIS	CA-CB-CG	-5.41	108.39	113.80
2	L	324	PHE	CA-CB-CG	-5.39	108.41	113.80
2	L	45	ASN	CA-C-O	-5.38	114.88	121.28
2	L	97	ASP	CA-CB-CG	5.33	117.93	112.60
1	S	100	ARG	CD-NE-CZ	5.32	131.84	124.40
2	L	136	PHE	CA-CB-CG	5.29	119.09	113.80
1	S	256	ASP	N-CA-CB	-5.28	104.28	112.30
1	S	187	VAL	CA-CB-CG1	5.18	119.20	110.40
1	S	92	HIS	CA-CB-CG	-5.14	108.66	113.80
2	L	100	VAL	CA-CB-CG2	-5.13	101.68	110.40
2	L	186	PHE	CA-CB-CG	-5.09	108.70	113.80
1	S	6	ARG	NE-CZ-NH2	5.09	123.78	119.20
2	L	120	TYR	CA-CB-CG	5.09	123.06	113.90
2	L	338	HIS	CA-CB-CG	-5.02	108.78	113.80
2	L	192	LEU	CA-C-N	-5.01	115.14	122.46
2	L	192	LEU	C-N-CA	-5.01	115.14	122.46

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	2019	0	1949	64	0
2	L	4177	0	4129	71	0
3	S	16	0	0	0	0
4	S	7	0	0	0	0
5	L	1	0	0	0	0
6	L	8	0	0	0	0
7	L	2	0	0	0	0
8	L	620	0	0	13	0
8	S	329	0	0	10	0
All	All	7179	0	6078	117	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 10.

All (117) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:2:MET:HA	2:L:182:GLN:HG2	1.51	0.92
1:S:26:ARG:HH21	2:L:233:ASN:HD21	1.18	0.87
1:S:238:GLN:HE21	2:L:224:ARG:HH21	1.24	0.86
1:S:2:MET:HG3	1:S:8:SER:HB2	1.58	0.86
1:S:126:ASN:HD21	1:S:130:ALA:H	1.24	0.83
1:S:230:ASN:HB2	1:S:233:LYS:HE2	1.62	0.80
1:S:2:MET:HG2	2:L:182:GLN:HE21	1.48	0.78
2:L:402:VAL:O	2:L:406:THR:HG23	1.84	0.77
1:S:1:LEU:HD12	1:S:1:LEU:H3	1.50	0.77
2:L:141:ALA:O	2:L:144:LYS:HG3	1.85	0.77
2:L:218:LEU:O	2:L:218:LEU:HG	1.85	0.76
1:S:25:LEU:HD11	8:S:3195:HOH:O	1.85	0.76
2:L:192:LEU:HD13	8:L:3838:HOH:O	1.86	0.75
2:L:176:THR:O	2:L:180:THR:HG23	1.87	0.74
1:S:155:TYR:HB2	8:S:3868:HOH:O	1.87	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:78:GLN:HE21	2:L:86:TYR:H	1.36	0.72
1:S:2:MET:HB3	8:S:3088:HOH:O	1.91	0.70
1:S:2:MET:HG2	2:L:182:GLN:NE2	2.06	0.69
1:S:1:LEU:HG	2:L:187:THR:HG21	1.74	0.69
1:S:1:LEU:HD11	1:S:62:GLN:HG2	1.74	0.69
1:S:256:ASP:HB3	8:S:3293:HOH:O	1.92	0.69
1:S:1:LEU:HD21	1:S:59:ALA:O	1.94	0.68
2:L:37:LEU:HG	2:L:547:ILE:HD11	1.76	0.68
2:L:467:LYS:HB3	8:L:3589:HOH:O	1.94	0.68
2:L:484:HIS:CD2	2:L:498:LEU:HD22	2.32	0.65
2:L:159:PRO:O	2:L:161:LYS:HE2	1.98	0.64
2:L:406:THR:HG22	2:L:432:ILE:HG21	1.80	0.64
1:S:26:ARG:HH21	2:L:233:ASN:ND2	1.95	0.63
1:S:230:ASN:HD22	1:S:230:ASN:H	1.46	0.62
1:S:6:ARG:HH12	2:L:180:THR:HG21	1.64	0.62
1:S:21:SER:O	1:S:25:LEU:HD13	2.00	0.61
1:S:256:ASP:HB3	8:S:3292:HOH:O	2.01	0.61
1:S:13:HIS:HD2	8:S:3020:HOH:O	1.84	0.59
1:S:31:TYR:O	1:S:35:LEU:HD22	2.02	0.59
2:L:325:LYS:HD2	2:L:325:LYS:N	2.17	0.59
2:L:261:LEU:HD11	8:L:3938:HOH:O	2.04	0.58
1:S:233:LYS:HE3	1:S:234:ILE:CD1	2.34	0.58
1:S:1:LEU:HB3	2:L:187:THR:OG1	2.04	0.57
2:L:211:HIS:HE1	8:L:3171:HOH:O	1.87	0.56
2:L:340:ARG:O	2:L:340:ARG:HD3	2.05	0.56
1:S:6:ARG:NH1	2:L:180:THR:HG21	2.20	0.56
1:S:126:ASN:HD21	1:S:130:ALA:N	1.99	0.56
2:L:321:LYS:N	2:L:321:LYS:HD3	2.21	0.56
2:L:36:HIS:HD2	8:L:3047:HOH:O	1.89	0.56
2:L:75:HIS:HD2	8:L:3017:HOH:O	1.89	0.55
2:L:321:LYS:HE2	8:L:3487:HOH:O	2.04	0.55
1:S:1:LEU:HD13	8:S:3088:HOH:O	2.06	0.55
2:L:487:ARG:HG3	8:L:3130:HOH:O	2.07	0.55
1:S:233:LYS:HE3	1:S:234:ILE:HD11	1.89	0.54
1:S:238:GLN:NE2	2:L:224:ARG:HH21	1.99	0.54
2:L:211:HIS:HD2	2:L:276:ASP:OD2	1.91	0.54
2:L:388:LEU:HG	2:L:426:ARG:HA	1.90	0.53
2:L:405:VAL:HG22	2:L:432:ILE:HG23	1.90	0.53
1:S:1:LEU:CG	2:L:187:THR:HG21	2.39	0.53
1:S:1:LEU:HD22	1:S:43:ASP:HB3	1.89	0.53
2:L:527:VAL:HG13	2:L:527:VAL:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:2:MET:HA	2:L:182:GLN:CG	2.33	0.53
2:L:253:GLN:NE2	2:L:253:GLN:H	2.07	0.52
1:S:217:LEU:HD12	1:S:243:VAL:HB	1.91	0.52
1:S:6:ARG:HH12	2:L:180:THR:CB	2.23	0.51
1:S:13:HIS:HE1	1:S:21:SER:OG	1.93	0.51
2:L:299:PHE:H	2:L:476:ASN:ND2	2.08	0.51
1:S:7:PRO:HG2	1:S:40:LEU:HD12	1.93	0.50
1:S:2:MET:CG	1:S:8:SER:HB2	2.37	0.50
2:L:253:GLN:H	2:L:253:GLN:HE21	1.58	0.49
1:S:156:ASN:HD21	1:S:230:ASN:HD21	1.59	0.49
1:S:3:GLY:O	1:S:4:PRO:O	2.30	0.49
2:L:50:ASN:HD21	2:L:509:ARG:NH2	2.09	0.49
2:L:281:ALA:CB	2:L:424:LEU:HD21	2.43	0.48
1:S:143:LYS:HB3	1:S:143:LYS:HE2	1.61	0.48
1:S:265:TYR:OH	2:L:75:HIS:HE1	1.95	0.48
1:S:37:LEU:HD12	2:L:177:PHE:CD2	2.49	0.48
1:S:5:ARG:HB3	1:S:5:ARG:NH1	2.29	0.48
1:S:6:ARG:HH12	2:L:180:THR:CG2	2.26	0.47
2:L:233:ASN:C	2:L:233:ASN:HD22	2.22	0.47
2:L:406:THR:HG22	2:L:432:ILE:HD13	1.96	0.47
1:S:36:ILE:HD12	1:S:42:LEU:HB2	1.96	0.47
1:S:229:ASN:HD22	1:S:230:ASN:H	1.62	0.47
2:L:37:LEU:HG	2:L:547:ILE:CD1	2.43	0.47
2:L:399:HIS:HD2	2:L:402:VAL:H	1.63	0.47
2:L:487:ARG:HB2	8:L:3130:HOH:O	2.14	0.47
1:S:181:MET:HE3	8:S:3656:HOH:O	2.14	0.46
1:S:2:MET:HE1	1:S:68:HIS:HB3	1.96	0.46
2:L:285:LYS:NZ	8:L:3676:HOH:O	2.48	0.46
1:S:85:ILE:HD11	8:S:3638:HOH:O	2.15	0.46
2:L:23:GLY:O	2:L:43:VAL:HG13	2.15	0.46
2:L:78:GLN:HE22	2:L:236:THR:H	1.64	0.45
1:S:139:HIS:H	1:S:139:HIS:CD2	2.34	0.45
2:L:143:LEU:HD13	2:L:174:LEU:HD13	1.99	0.45
1:S:1:LEU:HA	2:L:187:THR:OG1	2.17	0.45
1:S:53:GLY:HA3	8:L:3891:HOH:O	2.16	0.45
2:L:113:ASN:HD21	2:L:303:PRO:HD2	1.80	0.45
2:L:19:SER:HB3	2:L:20:SER:H	1.46	0.44
1:S:241:TRP:CH2	1:S:243:VAL:HG13	2.52	0.44
1:S:3:GLY:HA2	1:S:68:HIS:CD2	2.52	0.44
1:S:114:CYS:HA	1:S:119:GLY:HA3	1.99	0.44
2:L:138:ASP:H	2:L:205:ASN:ND2	2.14	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:340:ARG:HD3	2:L:340:ARG:HH11	1.61	0.43
2:L:514:LYS:NZ	8:L:3897:HOH:O	2.50	0.43
1:S:230:ASN:HD22	1:S:230:ASN:N	2.11	0.43
1:S:188:HIS:HB2	1:S:224:GLY:C	2.44	0.43
1:S:147:ILE:HD13	1:S:157:LEU:HA	1.99	0.43
1:S:201:GLU:HB3	1:S:215:TRP:CE3	2.54	0.42
1:S:241:TRP:CZ3	1:S:243:VAL:HG13	2.55	0.42
2:L:327:ILE:HD11	2:L:396:SER:HB2	2.02	0.42
1:S:230:ASN:H	1:S:230:ASN:ND2	2.14	0.42
2:L:533:PRO:O	2:L:537:LEU:HD13	2.19	0.41
2:L:50:ASN:HD21	2:L:509:ARG:HH22	1.68	0.41
2:L:85:THR:OG1	2:L:235:HIS:HD2	2.03	0.41
2:L:484:HIS:CG	2:L:498:LEU:HD22	2.55	0.41
2:L:27:VAL:HG22	8:L:3968:HOH:O	2.20	0.41
2:L:544:ASP:N	2:L:545:PRO:HD3	2.35	0.41
1:S:2:MET:HE2	1:S:3:GLY:O	2.20	0.41
2:L:546:CYS:SG	2:L:549:CYS:HB2	2.61	0.41
2:L:143:LEU:HD13	2:L:174:LEU:CD1	2.52	0.40
1:S:6:ARG:NH2	8:S:3174:HOH:O	2.47	0.40
2:L:338:HIS:HB2	2:L:375:ALA:HB3	2.04	0.40

There are no symmetry-related clashes.

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	S	265/267 (99%)	258 (97%)	5 (2%)	2 (1%)	16	3
2	L	532/534 (100%)	518 (97%)	13 (2%)	1 (0%)	43	19
All	All	797/801 (100%)	776 (97%)	18 (2%)	3 (0%)	30	11

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	S	4	PRO
2	L	20	SER
1	S	3	GLY

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	S	213/213 (100%)	192 (90%)	21 (10%)	7	0
2	L	438/438 (100%)	406 (93%)	32 (7%)	13	0
All	All	651/651 (100%)	598 (92%)	53 (8%)	11	0

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	S	1	LEU
1	S	2	MET
1	S	4	PRO
1	S	5	ARG
1	S	35	LEU
1	S	37	LEU
1	S	40	LEU
1	S	61	GLU
1	S	85	ILE
1	S	91	ASN
1	S	95	LEU
1	S	138	LYS
1	S	143	LYS
1	S	166	LYS
1	S	173	LEU
1	S	176	LEU
1	S	187	VAL
1	S	220	LEU
1	S	230	ASN
1	S	242	PRO
1	S	243	VAL

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Mol	Chain	Res	Type
2	L	19	SER
2	L	20	SER
2	L	37	LEU
2	L	43	VAL
2	L	49	LYS
2	L	132	HIS
2	L	143	LEU
2	L	144	LYS
2	L	161	LYS
2	L	174	LEU
2	L	200	LEU
2	L	213	LEU
2	L	233	ASN
2	L	250	LEU
2	L	253	GLN
2	L	277	LEU
2	L	278	LEU
2	L	279	VAL
2	L	315	LYS
2	L	321	LYS
2	L	325	LYS
2	L	360	LYS
2	L	388	LEU
2	L	405	VAL
2	L	410	LEU
2	L	418	GLU
2	L	424	LEU
2	L	433	GLU
2	L	445	LEU
2	L	457	ASN
2	L	487	ARG
2	L	522	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	S	13	HIS
1	S	68	HIS
1	S	91	ASN
1	S	126	ASN
1	S	139	HIS
1	S	156	ASN

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Mol	Chain	Res	Type
1	S	190	GLN
1	S	229	ASN
1	S	238	GLN
1	S	266	GLN
2	L	36	HIS
2	L	50	ASN
2	L	75	HIS
2	L	78	GLN
2	L	113	ASN
2	L	132	HIS
2	L	188	ASN
2	L	205	ASN
2	L	211	HIS
2	L	233	ASN
2	L	235	HIS
2	L	253	GLN
2	L	334	GLN
2	L	457	ASN
2	L	468	GLN
2	L	476	ASN
2	L	513	ASN

5.3.3 RNA [i](#)

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 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 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	FNE	L	1004	2,7	3,6,8	1.58	1 (33%)	-		
3	SF4	S	1001	1	0,12,12	-	-	-		
3	SF4	S	1002	1	0,12,12	-	-	-		
4	F3S	S	1003	1	0,9,9	-	-	-		
7	CMO	L	1006	6	0,1,1	-	-	-		

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	F3S	S	1003	1	-	-	0/3/3/3
3	SF4	S	1002	1	-	-	0/6/5/5
3	SF4	S	1001	1	-	-	0/6/5/5

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	L	1004	FNE	O3-C3	-2.20	1.13	1.16

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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.