



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:25 AM UTC

PDB ID : 4KN9 / pdb_00004kn9
Title : High-resolution structure of H₂-activated anaerobically purified *Dm. baculatum* [NiFeSe]-hydrogenase after crystallization under air
Authors : Volbeda, A.; Cavazza, C.; Fontecilla-Camps, J.C.
Deposited on : 2013-05-09
Resolution : 1.40 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

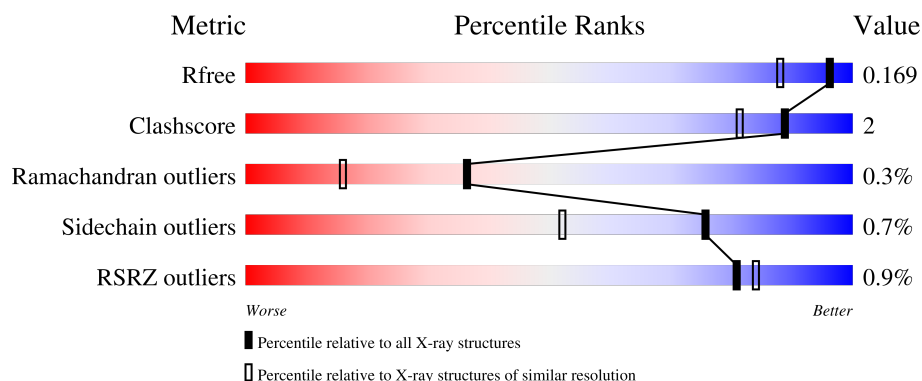
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.40 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	2563 (1.40-1.40)
Clashscore	190562	2660 (1.40-1.40)
Ramachandran outliers	187476	2611 (1.40-1.40)
Sidechain outliers	187428	2610 (1.40-1.40)
RSRZ outliers	180081	2561 (1.40-1.40)

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	283	2% 93% 6%
1	T	283	2% 94%
2	L	499	93%
2	M	499	93% 5%

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 14144 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.

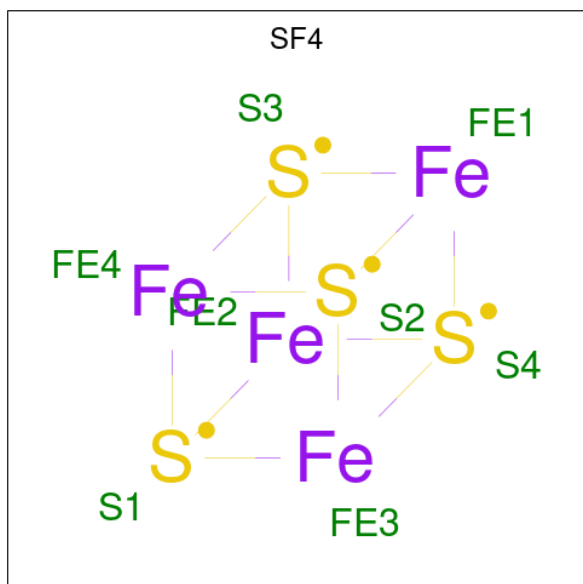
- Molecule 1 is a protein called Periplasmic [NiFeSe] hydrogenase small subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	S	279	Total	C	N	O	S	0	17	0
			2219	1415	364	422	18			
1	T	279	Total	C	N	O	S	0	11	0
			2184	1392	364	410	18			

- Molecule 2 is a protein called Nickel-dependent hydrogenase large subunit.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	L	489	Total	C	N	O	S	Se	0	24	0
			3925	2514	668	720	20	3			
2	M	489	Total	C	N	O	S	Se	0	27	0
			3935	2518	669	727	18	3			

- 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	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		
3	T	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	S	2	Total	Ca	0	0
			2	2		
4	L	1	Total	Ca	0	0
			1	1		
4	T	3	Total	Ca	0	0
			3	3		
4	M	1	Total	Ca	0	0
			1	1		

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

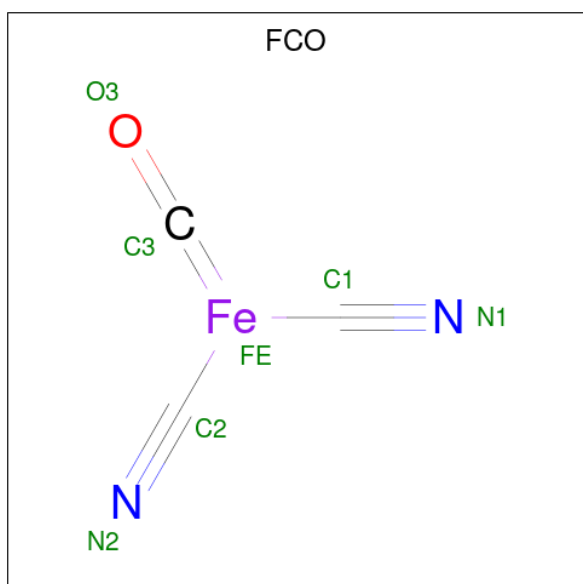
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	S	1	Total	Cl	0	0
			1	1		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	S	1	Total	C	O	0	0
			6	3	3		
6	S	1	Total	C	O	0	0
			6	3	3		
6	M	1	Total	C	O	0	0
			6	3	3		
6	M	1	Total	C	O	0	0
			6	3	3		

- Molecule 7 is CARBONMONOXIDE-(DICYANO) IRON (CCD ID: FCO) (formula: C_3FeN_2O).

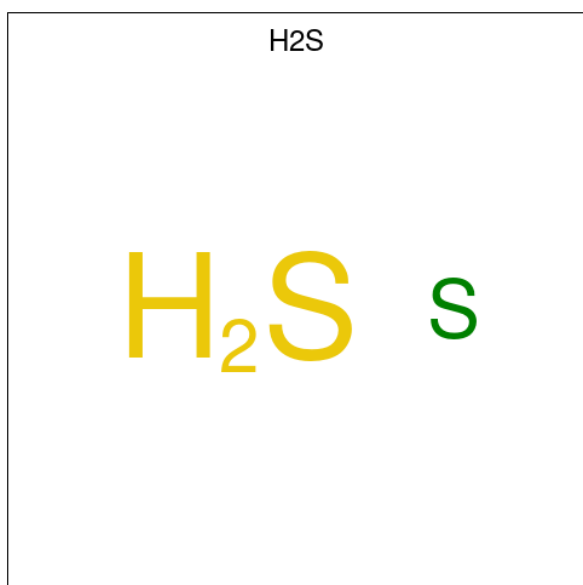


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	L	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		
7	M	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		

- Molecule 8 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	L	1	Total	Ni	0	1
			3	3		
8	M	1	Total	Ni	0	1
			3	3		

- Molecule 9 is HYDROSULFURIC ACID (CCD ID: H2S) (formula: H₂S).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	L	1	Total	S	0	0
			1	1		
9	M	1	Total	S	0	0
			1	1		

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	S	366	Total	O	0	5
			367	367		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	L	577	Total 578	O 578	0	2
10	T	347	Total 349	O 349	0	2
10	M	485	Total 485	O 485	0	3

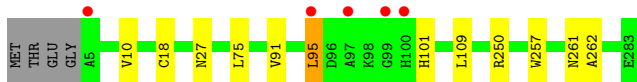
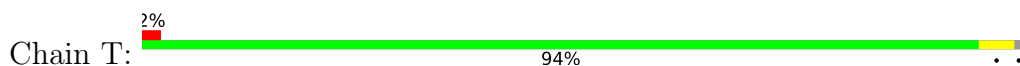
3 Residue-property plots

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 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: Periplasmic [NiFeSe] hydrogenase small subunit



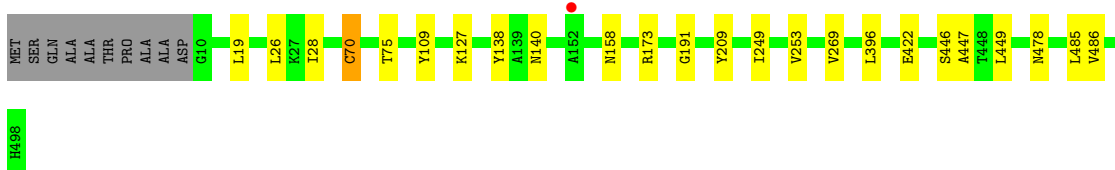
- Molecule 1: Periplasmic [NiFeSe] hydrogenase small subunit



- 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 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	106.26Å 108.73Å 136.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 1.40 25.00 – 1.40	Depositor EDS
% Data completeness (in resolution range)	91.5 (25.00-1.40) 91.5 (25.00-1.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.62 (at 1.40Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.127 , 0.168 0.128 , 0.169	Depositor DCC
R_{free} test set	14304 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å ²)	11.3	Xtriage
Anisotropy	0.360	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.015 for k,h,-l	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	14144	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.58% 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: FCO, H2S, CL, SF4, GOL, NI, CA, SEC

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	0.67	0/2341	0.75	0/3177
1	T	0.60	0/2288	0.75	0/3108
2	L	0.64	0/4090	0.75	0/5534
2	M	0.60	0/4112	0.74	0/5562
All	All	0.63	0/12831	0.75	0/17381

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [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	S	2219	0	2131	10	0
1	T	2184	0	2111	7	0
2	L	3925	0	3929	16	0
2	M	3935	0	3930	16	0
3	S	24	0	0	0	0
3	T	24	0	0	0	0
4	L	1	0	0	0	0
4	M	1	0	0	0	0
4	S	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	T	3	0	0	0	0
5	S	1	0	0	0	0
6	M	12	0	16	0	0
6	S	12	0	16	0	0
7	L	7	0	0	1	0
7	M	7	0	0	0	0
8	L	3	0	0	0	0
8	M	3	0	0	0	0
9	L	1	0	0	0	0
9	M	1	0	0	0	0
10	L	578	0	0	1	0
10	M	485	0	0	0	0
10	S	367	0	0	1	0
10	T	349	0	0	0	0
All	All	14144	0	12133	44	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 (44) 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:L:127:LYS:H	2:L:158:ASN:HD21	1.19	0.91
2:M:127:LYS:H	2:M:158:ASN:HD21	1.13	0.91
1:S:75[B]:LEU:HD21	1:S:175[B]:LEU:HD11	1.65	0.79
2:M:19[B]:LEU:HD23	2:M:26:LEU:HD23	1.70	0.73
2:M:19[B]:LEU:CD2	2:M:486:VAL:HG11	2.18	0.73
1:T:18:CYS:HB2	2:M:70[D]:CYS:HA	1.77	0.67
1:T:250[A]:ARG:HD2	1:T:257:TRP:CD1	2.36	0.60
1:S:82:THR:HG21	1:S:143[A]:LYS:HD2	1.84	0.59
2:M:127:LYS:N	2:M:158:ASN:HD21	1.95	0.58
1:S:118:LEU:HD21	1:S:182:VAL:HG21	1.86	0.57
1:S:18:CYS:HB2	2:L:70[D]:CYS:HA	1.87	0.56
2:M:19[A]:LEU:HD13	2:M:28:ILE:HG13	1.87	0.55
1:T:91:VAL:HG21	1:T:109:LEU:HD22	1.89	0.54
2:M:253:VAL:HB	2:M:422:GLU:HB2	1.89	0.54
2:M:19[B]:LEU:HD21	2:M:486:VAL:HG11	1.90	0.53
2:M:138:TYR:HB3	2:M:478:ASN:HD22	1.76	0.51
1:S:75[B]:LEU:HD11	1:S:175[B]:LEU:HD21	1.93	0.50
2:L:138:TYR:HB3	2:L:478:ASN:HD22	1.77	0.50
2:M:485:LEU:C	2:M:485:LEU:HD13	2.37	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:447[A]:ALA:HB3	7:L:501:FCO:N1	2.27	0.50
2:L:127:LYS:N	2:L:158:ASN:HD21	2.00	0.49
2:M:140:ASN:H	2:M:478:ASN:ND2	2.12	0.47
1:S:250:ARG:HD2	1:S:257:TRP:CD1	2.50	0.47
2:L:127:LYS:H	2:L:158:ASN:ND2	2.01	0.47
2:L:446[A]:SER:O	2:L:449:LEU:HB3	2.15	0.46
2:L:253:VAL:HB	2:L:422:GLU:HB2	1.98	0.46
2:M:249[B]:ILE:HG22	2:M:269:VAL:HG22	1.98	0.45
2:L:140:ASN:H	2:L:478:ASN:ND2	2.13	0.45
2:L:242:GLY:HA3	2:L:485:LEU:HD23	1.98	0.45
1:S:277:SER:HA	1:S:278:PRO:C	2.41	0.44
2:M:446[A]:SER:O	2:M:449:LEU:HB3	2.18	0.44
2:L:485:LEU:C	2:L:485:LEU:HD13	2.43	0.44
2:L:62:SER:O	2:L:66:VAL:HG22	2.18	0.43
1:S:261:ASN:HD21	2:L:191:GLY:CA	2.31	0.43
1:S:276:LYS:NZ	10:S:614:HOH:O	2.51	0.43
1:T:10:VAL:HG13	1:T:75[B]:LEU:HD23	2.01	0.43
2:L:23:GLU:HB3	2:L:492[D]:SEC:SE	2.70	0.42
1:S:143[B]:LYS:HE2	1:S:143[B]:LYS:HB2	1.84	0.42
1:T:27:ASN:HA	2:M:173:ARG:HG3	2.01	0.41
2:M:209:TYR:CD1	2:M:396:LEU:HD21	2.55	0.41
2:L:344:GLU:HA	2:L:363:ASN:HD22	1.84	0.41
1:T:95:LEU:HG	1:T:101:HIS:CE1	2.55	0.41
2:L:425:ARG:HD2	10:L:607[N]:HOH:O	2.21	0.40
1:T:261:ASN:HD21	2:M:191:GLY:CA	2.35	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	293/283 (104%)	286 (98%)	5 (2%)	2 (1%)	18 4

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	T	288/283 (102%)	280 (97%)	7 (2%)	1 (0%)	36	16
2	L	509/499 (102%)	494 (97%)	13 (3%)	2 (0%)	30	10
2	M	511/499 (102%)	496 (97%)	13 (2%)	2 (0%)	30	10
All	All	1601/1564 (102%)	1556 (97%)	38 (2%)	7 (0%)	36	10

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	S	96	ASP
2	L	447[A]	ALA
2	L	447[B]	ALA
2	M	447[A]	ALA
2	M	447[B]	ALA
1	S	262	ALA
1	T	262	ALA

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	247/233 (106%)	246 (100%)	1 (0%)	84	67
1	T	241/233 (103%)	240 (100%)	1 (0%)	84	67
2	L	426/410 (104%)	421 (99%)	5 (1%)	63	34
2	M	428/410 (104%)	424 (99%)	4 (1%)	70	44
All	All	1342/1286 (104%)	1331 (99%)	11 (1%)	76	48

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	S	207	ILE
2	L	70[D]	CYS
2	L	70[N]	CYS
2	L	75	THR

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Mol	Chain	Res	Type
2	L	109	TYR
2	L	276	GLU
1	T	95	LEU
2	M	70[D]	CYS
2	M	70[N]	CYS
2	M	75	THR
2	M	109	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	S	16	GLN
1	S	101	HIS
1	S	261	ASN
2	L	52	GLN
2	L	140	ASN
2	L	158	ASN
2	L	188	HIS
2	L	363	ASN
2	L	478	ASN
1	T	101	HIS
1	T	188	HIS
1	T	261	ASN
2	M	102	ASN
2	M	140	ASN
2	M	158	ASN
2	M	188	HIS
2	M	400	GLN
2	M	478	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 28 ligands modelled in this entry, 14 are monoatomic and 2 are modelled with single atom - leaving 12 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	$\# Z > 2$
7	FCO	M	501	10,2	0,6,6	-	-	-		
3	SF4	S	303	1	0,12,12	-	-	-		
6	GOL	M	505	-	5,5,5	0.50	0	5,5,5	0.67	0
3	SF4	S	301	1	0,12,12	-	-	-		
7	FCO	L	501	10,2	0,6,6	-	-	-		
3	SF4	S	302	1	0,12,12	-	-	-		
6	GOL	M	506	-	5,5,5	0.40	0	5,5,5	0.20	0
3	SF4	T	301	1	0,12,12	-	-	-		
6	GOL	S	307	-	5,5,5	0.33	0	5,5,5	0.42	0
3	SF4	T	302	1	0,12,12	-	-	-		
6	GOL	S	306	-	5,5,5	0.28	0	5,5,5	0.53	0
3	SF4	T	303	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
6	GOL	M	505	-	-	1/4/4/4	-
3	SF4	S	303	1	-	-	0/6/5/5
3	SF4	S	301	1	-	-	0/6/5/5
3	SF4	S	302	1	-	-	0/6/5/5
6	GOL	M	506	-	-	3/4/4/4	-
3	SF4	T	301	1	-	-	0/6/5/5
6	GOL	S	307	-	-	0/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SF4	T	302	1	-	-	0/6/5/5
6	GOL	S	306	-	-	0/4/4/4	-
3	SF4	T	303	1	-	-	0/6/5/5

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	M	506	GOL	C1-C2-C3-O3
6	M	506	GOL	O2-C2-C3-O3
6	M	505	GOL	O1-C1-C2-O2
6	M	506	GOL	O1-C1-C2-C3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	L	501	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 [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	S	279/283 (98%)	-0.80	7 (2%) 58 61	6, 12, 23, 38	29 (10%)
1	T	279/283 (98%)	-0.60	5 (1%) 67 71	9, 16, 28, 51	27 (9%)
2	L	488/499 (97%)	-0.94	1 (0%) 91 93	5, 13, 21, 43	31 (6%)
2	M	488/499 (97%)	-0.61	1 (0%) 91 93	8, 18, 30, 46	33 (6%)
All	All	1534/1564 (98%)	-0.75	14 (0%) 81 84	5, 15, 27, 51	120 (7%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	S	95	LEU	6.0
1	S	100	HIS	4.4
1	T	97	ALA	4.4
1	S	97	ALA	4.0
1	S	5	ALA	3.7
1	S	99	GLY	3.4
1	S	96	ASP	3.4
1	T	95	LEU	3.4
1	T	99	GLY	3.3
1	S	98	LYS	2.7
1	T	100	HIS	2.7
2	L	10	GLY	2.6
1	T	5	ALA	2.3
2	M	152	ALA	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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	GOL	M	506	6/6	0.88	0.12	64,65,66,67	0
4	CA	T	306	1/1	0.93	0.15	82,82,82,82	0
6	GOL	S	307	6/6	0.94	0.07	29,30,31,34	0
6	GOL	M	505	6/6	0.97	0.06	19,24,30,34	0
6	GOL	S	306	6/6	0.97	0.05	23,24,25,30	0
8	NI	M	502[D]	1/1	0.97	0.09	19,19,19,19	1
8	NI	M	502[N]	1/1	0.97	0.09	20,20,20,20	1
8	NI	M	502[Q]	1/1	0.97	0.09	16,16,16,16	1
8	NI	L	502[N]	1/1	0.99	0.07	16,16,16,16	1
8	NI	L	502[Q]	1/1	0.99	0.07	10,10,10,10	1
4	CA	T	305	1/1	0.99	0.11	29,29,29,29	1
4	CA	S	308	1/1	0.99	0.08	33,33,33,33	1
8	NI	L	502[D]	1/1	0.99	0.07	17,17,17,17	1
5	CL	S	305	1/1	1.00	0.02	17,17,17,17	1
3	SF4	T	301	8/8	1.00	0.01	16,18,19,20	0
3	SF4	T	302	8/8	1.00	0.01	12,13,14,14	0
3	SF4	T	303	8/8	1.00	0.01	11,12,12,12	0
4	CA	S	304	1/1	1.00	0.01	12,12,12,12	0
7	FCO	L	501	7/7	1.00	0.02	9,10,12,14	0
7	FCO	M	501	7/7	1.00	0.02	11,13,14,17	0
3	SF4	S	301	8/8	1.00	0.01	11,12,13,13	0
4	CA	L	503	1/1	1.00	0.01	13,13,13,13	1
4	CA	T	304	1/1	1.00	0.03	20,20,20,20	1
3	SF4	S	302	8/8	1.00	0.01	9,9,10,10	0
3	SF4	S	303	8/8	1.00	0.01	8,8,9,9	0
4	CA	M	503	1/1	1.00	0.03	14,14,14,14	1
9	H2S	L	504	1/1	1.00	0.02	9,9,9,9	0
9	H2S	M	504	1/1	1.00	0.01	14,14,14,14	0

6.5 Other polymers ⓘ

There are no such residues in this entry.