



wwPDB X-ray Structure Validation Summary Report i

Feb 28, 2014 – 02:21 PM GMT

PDB ID : 1YQW
Title : Structure of the Oxidized Unready Form of Ni-Fe Hydrogenase
Authors : Volbeda, A.
Deposited on : 2005-02-02
Resolution : 1.83 Å(reported)

This is a wwPDB validation summary 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/ValidationPDFNotes.html>

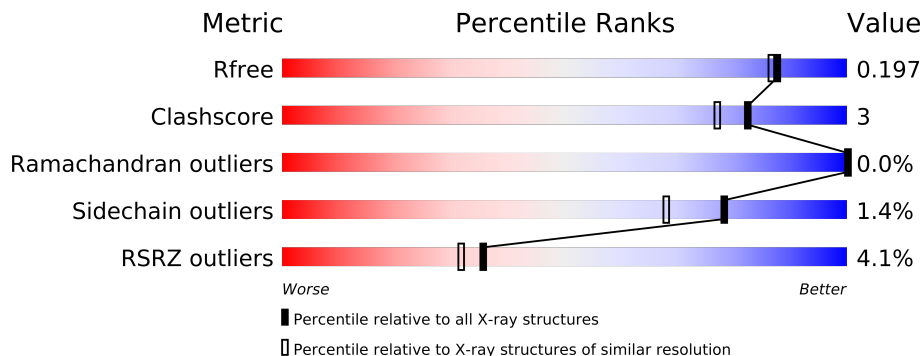
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	FAILED
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.83 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}	66092	1857 (1.86-1.82)
Clashscore	79885	2149 (1.86-1.82)
Ramachandran outliers	78287	2124 (1.86-1.82)
Sidechain outliers	78261	2125 (1.86-1.82)
RSRZ outliers	66119	1857 (1.86-1.82)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	264	
1	B	264	
1	C	264	
2	Q	549	
2	R	549	
2	S	549	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
11	GOL	A	268	-	X
11	GOL	A	269	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
11	GOL	Q	559	-	X
11	GOL	Q	560	-	X
11	GOL	R	1304	-	X
11	GOL	R	1305	-	X
11	GOL	R	1306	-	X
11	GOL	S	2303	-	X
5	BCT	Q	554	-	X
6	MG	R	1308	-	X

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 20497 atoms, of which 0 are hydrogen and 0 are deuterium.

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	A	262	Total	C	N	O	S	0	0	0
			1973	1256	330	372	15			
1	B	262	Total	C	N	O	S	0	2	0
			1980	1260	330	375	15			
1	C	260	Total	C	N	O	S	0	0	0
			1960	1248	327	370	15			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	Q	544	Total	C	N	O	S	0	8	0
			4187	2669	728	767	23			
2	R	545	Total	C	N	O	S	0	3	0
			4173	2660	725	765	23			
2	S	544	Total	C	N	O	S	0	2	0
			4165	2652	725	765	23			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Q	198	ASN	ASP	CONFLICT	UNP P18188
Q	303	SER	GLU	CONFLICT	UNP P18188
Q	499	ALA	SER	ENGINEERED MUTATION	UNP P18188
R	198	ASN	ASP	CONFLICT	UNP P18188
R	303	SER	GLU	CONFLICT	UNP P18188
R	499	ALA	SER	ENGINEERED MUTATION	UNP P18188
S	198	ASN	ASP	CONFLICT	UNP P18188
S	303	SER	GLU	CONFLICT	UNP P18188
S	499	ALA	SER	ENGINEERED MUTATION	UNP P18188

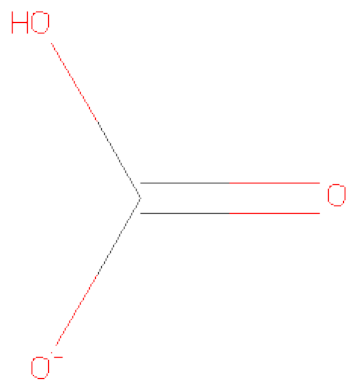
- Molecule 3 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

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

- Molecule 4 is FE (II) ION (three-letter code: FE2) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	R	1	Total Fe 1 1	0	0
4	Q	1	Total Fe 1 1	0	0
4	S	1	Total Fe 1 1	0	0

- Molecule 5 is BICARBONATE ION (three-letter code: BCT) (formula: CHO_3).

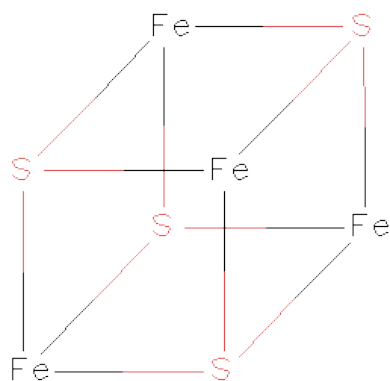


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	Q	1	Total C O 4 1 3	0	0

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

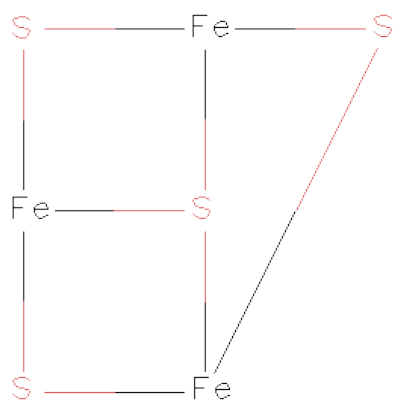
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	R	1	Total Mg 1 1	0	0
6	Q	2	Total Mg 2 2	0	0

- Molecule 7 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe_4S_4).



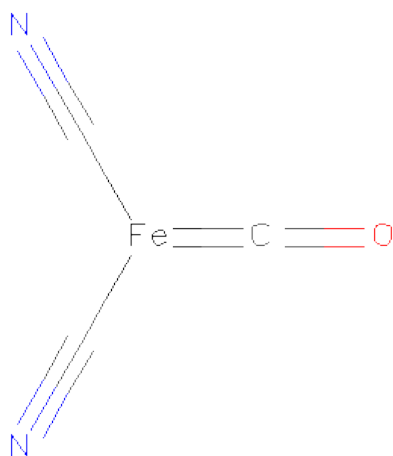
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total Fe S 8 4 4	0	0
7	A	1	Total Fe S 8 4 4	0	0
7	B	1	Total Fe S 8 4 4	0	0
7	B	1	Total Fe S 8 4 4	0	0
7	C	1	Total Fe S 8 4 4	0	0
7	C	1	Total Fe S 8 4 4	0	0

- Molecule 8 is FE3-S4 CLUSTER (three-letter code: F3S) (formula: Fe_3S_4).



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

- Molecule 9 is CARBONMONOXIDE-(DICYANO)IRON (three-letter code: FCO) (formula: $\text{C}_3\text{FeN}_2\text{O}$).



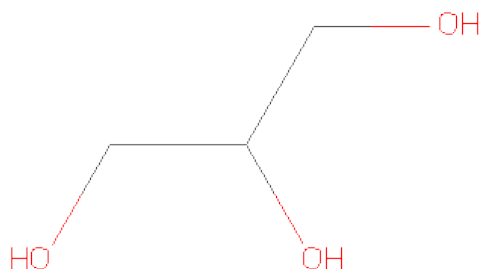
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	Q	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		
9	R	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		
9	S	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		

- Molecule 10 is PEROXIDE ION (three-letter code: PER) (formula: O₂).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	Q	1	Total	O	0	0
			2	2		
10	R	1	Total	O	0	0
			2	2		
10	S	1	Total	O	0	0
			2	2		

- Molecule 11 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

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

- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	275	Total	O	0	0
			275	275		
12	B	194	Total	O	0	0
			194	194		
12	C	185	Total	O	0	0
			185	185		
12	Q	482	Total	O	0	0
			482	482		
12	R	419	Total	O	0	0
			419	419		
12	S	293	Total	O	0	0
			293	293		

3 Residue-property plots

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.

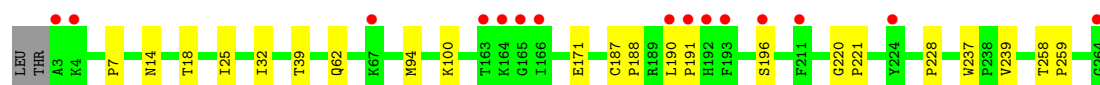
- Molecule 1: Periplasmic [NiFe] hydrogenase small subunit

Chain A: 



- Molecule 1: Periplasmic [NiFe] hydrogenase small subunit

Chain B: 



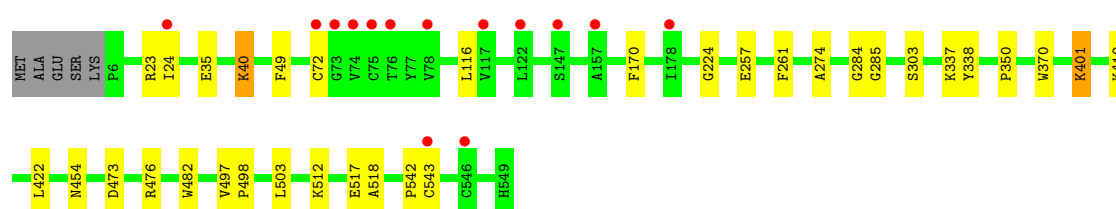
- Molecule 1: Periplasmic [NiFe] hydrogenase small subunit

Chain C: 



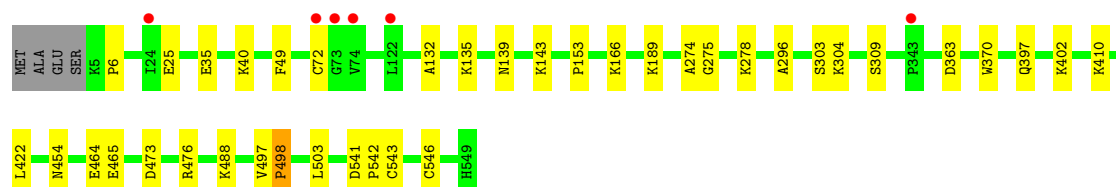
- Molecule 2: Periplasmic [NiFe] hydrogenase large subunit

Chain Q: 



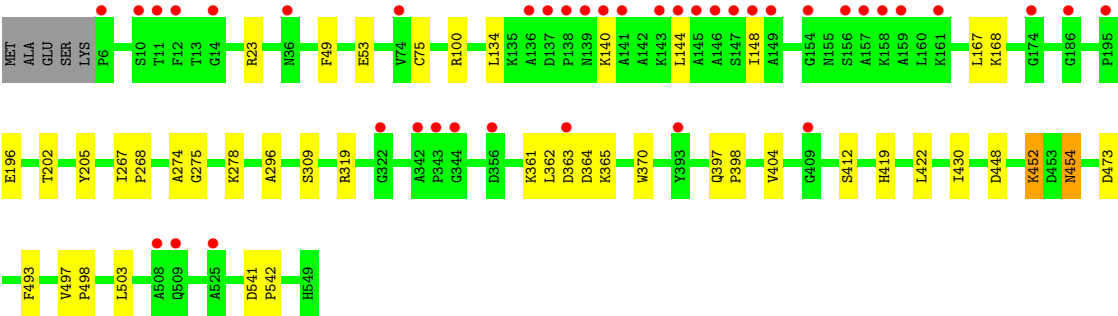
- Molecule 2: Periplasmic [NiFe] hydrogenase large subunit

Chain R: 



● Molecule 2: Periplasmic [NiFe] hydrogenase large subunit

Chain S: 



4 Data and refinement statistics

Xtriage (Phenix) failed to run properly - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.41Å 99.70Å 182.83Å 90.00° 91.42° 90.00°	Depositor
Resolution (Å)	20.00 – 1.83 46.98 – 1.78	Depositor EDS
% Data completeness (in resolution range)	83.0 (20.00-1.83) 81.0 (46.98-1.78)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.146 , 0.185 0.166 , 0.197	Depositor DCC
R_{free} test set	8482 reflections (5.03%)	DCC
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 35.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	20497	wwPDB-VP
Average B, all atoms (Å ²)	11.0	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: GOL, MG, NI, SF4, PER, F3S, FE2, BCT, 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	0.50	0/2027	0.60	0/2759
1	B	0.42	0/2043	0.55	0/2781
1	C	0.41	0/2014	0.53	0/2743
2	Q	0.48	0/4330	0.60	0/5876
2	R	0.45	0/4292	0.58	0/5827
2	S	0.40	0/4280	0.55	0/5810
All	All	0.44	0/18986	0.57	0/25796

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1973	0	1911	16	0
1	B	1980	0	1918	11	0
1	C	1960	0	1897	8	0
2	Q	4187	0	4180	25	0
2	R	4173	0	4154	31	0
2	S	4165	0	4142	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	Q	1	0	0	0	0
3	R	1	0	0	0	0
3	S	1	0	0	0	0
4	Q	1	0	0	0	0
4	R	1	0	0	0	0
4	S	1	0	0	0	0
5	Q	4	0	0	0	0
6	Q	2	0	0	0	0
6	R	1	0	0	0	0
7	A	16	0	0	0	0
7	B	16	0	0	0	0
7	C	16	0	0	0	0
8	A	7	0	0	0	0
8	B	7	0	0	0	0
8	C	7	0	0	0	0
9	Q	7	0	0	1	0
9	R	7	0	0	0	0
9	S	7	0	0	0	0
10	Q	2	0	0	1	0
10	R	2	0	0	1	0
10	S	2	0	0	0	0
11	A	18	0	24	2	0
11	Q	24	0	32	3	0
11	R	42	0	56	5	0
11	S	18	0	24	0	0
12	A	275	0	0	6	0
12	B	194	0	0	1	0
12	C	185	0	0	0	0
12	Q	482	0	0	3	0
12	R	419	0	0	9	0
12	S	293	0	0	2	0
All	All	20497	0	18338	113	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 3.

The worst 5 of 113 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:Q:543[B]:CYS:SG	10:Q:552:PER:O2	2.11	1.08
2:R:410:LYS:HD3	11:R:1302:GOL:H11	1.59	0.84
2:S:134:LEU:HD21	2:S:167:LEU:HD23	1.68	0.74
12:Q:1038:HOH:O	2:R:6:PRO:HB3	1.88	0.71

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:61:HIS:HE1	12:A:469:HOH:O	1.73	0.71

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	A	260/264 (98%)	253 (97%)	7 (3%)	0	100	100
1	B	262/264 (99%)	255 (97%)	7 (3%)	0	100	100
1	C	258/264 (98%)	252 (98%)	6 (2%)	0	100	100
2	Q	550/549 (100%)	539 (98%)	11 (2%)	0	100	100
2	R	546/549 (100%)	529 (97%)	17 (3%)	0	100	100
2	S	544/549 (99%)	525 (96%)	18 (3%)	1 (0%)	56	38
All	All	2420/2439 (99%)	2353 (97%)	66 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	S	364	ASP

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	208/210 (99%)	207 (100%)	1 (0%)	94	91
1	B	210/210 (100%)	207 (99%)	3 (1%)	78	68

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	207/210 (99%)	205 (99%)	2 (1%)	85	80
2	Q	442/438 (101%)	436 (99%)	6 (1%)	78	68
2	R	437/438 (100%)	430 (98%)	7 (2%)	75	62
2	S	436/438 (100%)	429 (98%)	7 (2%)	75	62
All	All	1940/1944 (100%)	1914 (99%)	26 (1%)	78	71

5 of 26 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	R	397	GLN
2	R	473	ASP
2	S	473	ASP
2	R	402	LYS
2	R	454	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 15 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	14	ASN
1	B	62	GLN
2	R	509	GLN
2	Q	509	GLN
2	R	454	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 42 ligands modelled in this entry, 9 are monoatomic - leaving 33 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	SF4	A	265	1	12,12,12	4.62	8 (66%)	0,24,24	0.00	-
8	F3S	A	266	1	3,9,9	14.22	3 (100%)	0,15,15	0.00	-
7	SF4	A	267	1	12,12,12	5.06	7 (58%)	0,24,24	0.00	-
11	GOL	A	268	-	5,5,5	0.44	0	5,5,5	0.72	0
11	GOL	A	269	-	5,5,5	0.20	0	5,5,5	0.73	0
11	GOL	A	270	-	5,5,5	0.33	0	5,5,5	1.01	0
7	SF4	B	265	1	12,12,12	3.77	11 (91%)	0,24,24	0.00	-
8	F3S	B	266	1	3,9,9	16.19	3 (100%)	0,15,15	0.00	-
7	SF4	B	267	1	12,12,12	4.84	10 (83%)	0,24,24	0.00	-
7	SF4	C	265	1	12,12,12	4.05	10 (83%)	0,24,24	0.00	-
8	F3S	C	266	1	3,9,9	14.37	3 (100%)	0,15,15	0.00	-
7	SF4	C	267	1	12,12,12	4.52	12 (100%)	0,24,24	0.00	-
9	FCO	Q	550	10,2	2,6,6	0.12	0	0,6,6	0.00	-
10	PER	Q	552	9,3	1,1,1	0.93	0	0,0,0	0.00	-
5	BCT	Q	554	-	0,3,3	0.00	-	0,3,3	0.00	-
11	GOL	Q	557	-	5,5,5	0.34	0	5,5,5	0.14	0
11	GOL	Q	558	-	5,5,5	0.44	0	5,5,5	0.27	0
11	GOL	Q	559	-	5,5,5	0.36	0	5,5,5	0.22	0
11	GOL	Q	560	-	5,5,5	0.40	0	5,5,5	0.45	0
11	GOL	R	1301	-	5,5,5	0.35	0	5,5,5	0.44	0
11	GOL	R	1302	-	5,5,5	0.37	0	5,5,5	0.17	0
11	GOL	R	1303	-	5,5,5	0.22	0	5,5,5	0.46	0
11	GOL	R	1304	-	5,5,5	0.33	0	5,5,5	0.27	0
11	GOL	R	1305	-	5,5,5	0.39	0	5,5,5	0.25	0
11	GOL	R	1306	-	5,5,5	0.35	0	5,5,5	0.29	0
11	GOL	R	1307	-	5,5,5	0.24	0	5,5,5	0.39	0
9	FCO	R	550	10,2	2,6,6	0.09	0	0,6,6	0.00	-
10	PER	R	552	9,3	1,1,1	0.97	0	0,0,0	0.00	-
11	GOL	S	2301	-	5,5,5	0.40	0	5,5,5	0.08	0
11	GOL	S	2302	-	5,5,5	0.40	0	5,5,5	0.21	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
11	GOL	S	2303	-	5,5,5	0.35	0	5,5,5	0.26	0
9	FCO	S	550	10,2	2,6,6	0.04	0	0,6,6	0.00	-
10	PER	S	552	9,3	1,1,1	1.02	0	0,0,0	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
7	SF4	A	265	1	-	0/0/48/48	0/0/5/5
8	F3S	A	266	1	-	0/0/24/24	0/0/3/3
7	SF4	A	267	1	-	0/0/48/48	0/0/5/5
11	GOL	A	268	-	-	0/4/4/4	0/0/0/0
11	GOL	A	269	-	-	0/4/4/4	0/0/0/0
11	GOL	A	270	-	-	0/4/4/4	0/0/0/0
7	SF4	B	265	1	-	0/0/48/48	0/0/5/5
8	F3S	B	266	1	-	0/0/24/24	0/0/3/3
7	SF4	B	267	1	-	0/0/48/48	0/0/5/5
7	SF4	C	265	1	-	0/0/48/48	0/0/5/5
8	F3S	C	266	1	-	0/0/24/24	0/0/3/3
7	SF4	C	267	1	-	0/0/48/48	0/0/5/5
9	FCO	Q	550	10,2	-	0/0/6/6	0/0/0/0
10	PER	Q	552	9,3	-	0/0/0/0	0/0/0/0
5	BCT	Q	554	-	-	0/0/0/0	0/0/0/0
11	GOL	Q	557	-	-	0/4/4/4	0/0/0/0
11	GOL	Q	558	-	-	0/4/4/4	0/0/0/0
11	GOL	Q	559	-	-	0/4/4/4	0/0/0/0
11	GOL	Q	560	-	-	0/4/4/4	0/0/0/0
11	GOL	R	1301	-	-	0/4/4/4	0/0/0/0
11	GOL	R	1302	-	-	0/4/4/4	0/0/0/0
11	GOL	R	1303	-	-	0/4/4/4	0/0/0/0
11	GOL	R	1304	-	-	0/4/4/4	0/0/0/0
11	GOL	R	1305	-	-	0/4/4/4	0/0/0/0
11	GOL	R	1306	-	-	0/4/4/4	0/0/0/0
11	GOL	R	1307	-	-	0/4/4/4	0/0/0/0
9	FCO	R	550	10,2	-	0/0/6/6	0/0/0/0
10	PER	R	552	9,3	-	0/0/0/0	0/0/0/0
11	GOL	S	2301	-	-	0/4/4/4	0/0/0/0
11	GOL	S	2302	-	-	0/4/4/4	0/0/0/0
11	GOL	S	2303	-	-	0/4/4/4	0/0/0/0
9	FCO	S	550	10,2	-	0/0/6/6	0/0/0/0
10	PER	S	552	9,3	-	0/0/0/0	0/0/0/0

The worst 5 of 67 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	266	F3S	S3-FE3	-18.13	2.21	2.33
8	C	266	F3S	S3-FE3	-15.92	2.22	2.33
8	A	266	F3S	S3-FE3	-15.63	2.22	2.33
8	B	266	F3S	S3-FE1	-15.45	2.22	2.33
8	A	266	F3S	S3-FE4	-15.36	2.22	2.33

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	262/264 (99%)	0.13	13 (4%)	28 24	4, 9, 16, 32	8 (3%)
1	B	262/264 (99%)	0.15	15 (5%)	23 19	5, 10, 17, 29	9 (3%)
1	C	260/264 (98%)	0.16	11 (4%)	35 31	6, 11, 20, 28	6 (2%)
2	Q	544/549 (99%)	0.02	14 (2%)	53 49	4, 9, 15, 22	12 (2%)
2	R	545/549 (99%)	-0.12	6 (1%)	77 76	5, 9, 16, 40	13 (2%)
2	S	544/549 (99%)	0.42	40 (7%)	14 13	6, 11, 18, 28	15 (2%)
All	All	2417/2439 (99%)	0.12	99 (4%)	35 32	4, 10, 17, 40	63 (2%)

The worst 5 of 99 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	3	ALA	5.8
2	S	145	ALA	4.7
1	C	67	LYS	4.7
1	A	3	ALA	4.7
1	C	162	LEU	4.3

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates 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. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
11	GOL	S	2303	6/6	0.23	9.92	34,35,36,38	6
11	GOL	R	1304	6/6	0.20	7.80	28,31,33,36	0
11	GOL	Q	559	6/6	0.18	6.90	29,31,33,35	0
5	BCT	Q	554	4/4	0.15	4.12	24,29,31,32	0
11	GOL	R	1305	6/6	0.12	3.65	21,23,25,26	0
6	MG	R	1308	1/1	0.14	3.21	30,30,30,30	0
11	GOL	A	269	6/6	0.13	3.03	21,27,28,36	0
11	GOL	A	268	6/6	0.13	2.91	16,20,24,28	0
11	GOL	Q	560	6/6	0.18	2.85	25,38,41,47	0
11	GOL	R	1306	6/6	0.17	2.07	30,37,39,41	0
11	GOL	A	270	6/6	0.24	1.77	17,29,31,34	0
11	GOL	R	1307	6/6	0.20	1.50	28,31,32,36	0
11	GOL	R	1302	6/6	0.16	1.40	27,28,30,33	0
11	GOL	R	1303	6/6	0.11	0.83	9,10,11,13	0
11	GOL	S	2302	6/6	0.14	0.69	32,35,35,35	6
6	MG	Q	555	1/1	0.09	0.10	29,29,29,29	0
11	GOL	R	1301	6/6	0.09	-0.16	7,8,9,10	0
11	GOL	S	2301	6/6	0.07	-0.61	7,10,11,12	0
11	GOL	Q	558	6/6	0.08	-0.66	8,10,11,12	0
8	F3S	A	266	7/7	0.11	-0.66	9,9,10,10	0
8	F3S	C	266	7/7	0.08	-0.86	10,10,11,12	0
6	MG	Q	556	1/1	0.08	-0.99	35,35,35,35	0
4	FE2	R	553	1/1	0.06	-1.07	9,9,9,9	1
7	SF4	A	267	8/8	0.16	-1.09	8,9,9,9	0
9	FCO	Q	550	7/7	0.16	-1.12	7,8,9,10	0
9	FCO	R	550	7/7	0.13	-1.14	5,7,9,11	0
7	SF4	A	265	8/8	0.07	-1.18	7,8,8,10	0
7	SF4	B	267	8/8	0.08	-1.20	9,9,11,12	0
10	PER	S	552	2/2	0.11	-1.34	12,12,12,12	1
7	SF4	C	267	8/8	0.09	-1.39	9,10,10,10	0
4	FE2	Q	553	1/1	0.08	-1.42	9,9,9,9	1
7	SF4	C	265	8/8	0.06	-1.42	8,9,10,11	0
11	GOL	Q	557	6/6	0.05	-1.53	8,9,11,12	0
7	SF4	B	265	8/8	0.06	-1.54	8,10,11,11	0
9	FCO	S	550	7/7	0.10	-1.63	8,11,13,13	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
10	PER	R	552	2/2	0.14	-1.65	14,14,14,16	1
8	F3S	B	266	7/7	0.04	-2.04	9,9,11,12	0
3	NI	S	551	1/1	0.08	-2.38	15,15,15,15	0
4	FE2	S	553	1/1	0.02	-2.97	9,9,9,9	1
3	NI	R	551	1/1	0.10	-3.26	14,14,14,14	0
3	NI	Q	551	1/1	0.13	-3.43	14,14,14,14	0
10	PER	Q	552	2/2	0.12	-3.88	14,14,14,15	1

6.5 Other polymers ⓘ

There are no such residues in this entry.