Full wwPDB X-ray Structure Validation Report ⓘ

Aug 26, 2026 – 02:58 am BST

PDB ID	:	3ZE7 / pdb_00003ze7
Title	:	3D structure of the Ni-Fe-Se hydrogenase from D. vulgaris Hildenborough in the reduced state at 1.95 Angstroms
Authors	:	Marques, M.C.; Coelho, R.; Pereira, I.A.C.; Matias, P.M.
Deposited on	:	2012-12-03
Resolution	:	1.95 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:

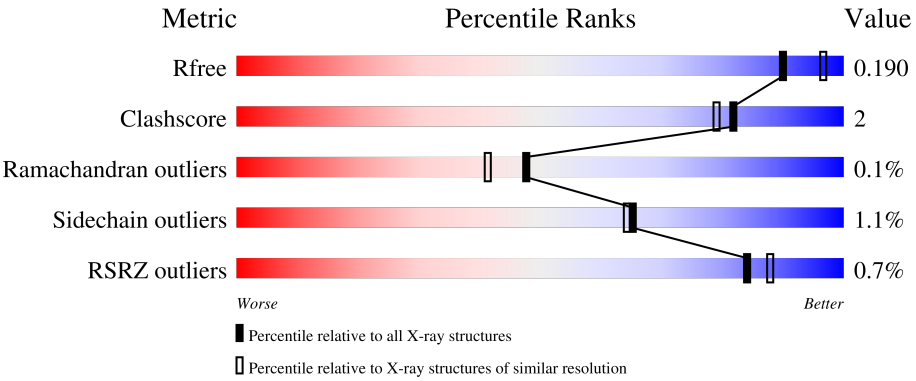
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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 <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	283	91% 6% .
2	B	484	% 92% 7% .

2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 6625 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.

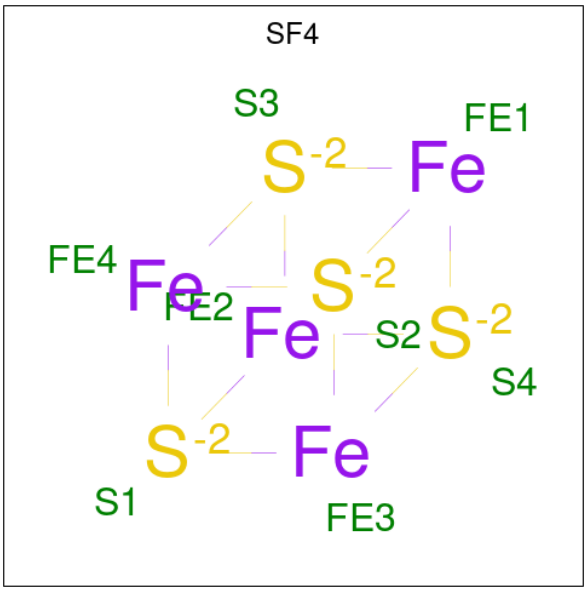
- Molecule 1 is a protein called PERIPLASMIC [NIFESE] HYDROGENASE, SMALL SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	277	2112	1350	352	390	20	0	5	0

- Molecule 2 is a protein called PERIPLASMIC [NIFESE] HYDROGENASE, LARGE SUBUNIT, SELENOCYSTEINE-CONTAINING.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
2	B	481	3745	2391	653	683	17	1	0	2	0

- Molecule 3 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4).



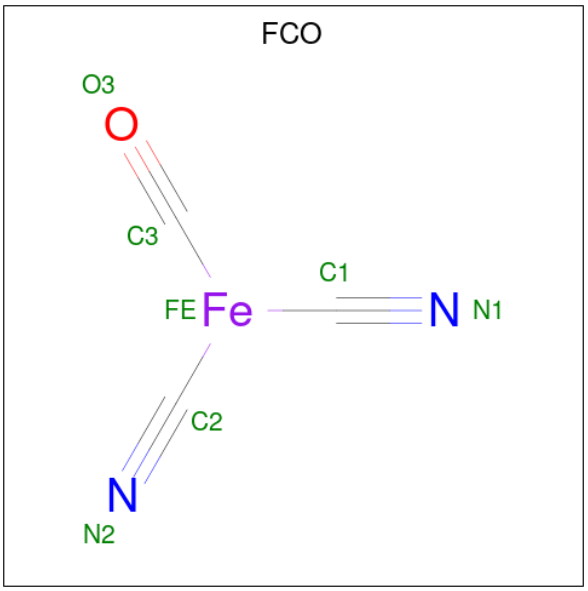
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	Fe	S		
3	A	1	8	4	4	0	0

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

- Molecule 4 is CARBONMONOXIDE-(DICYANO) IRON (CCD ID: FCO) (formula: C_3FeN_2O).



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

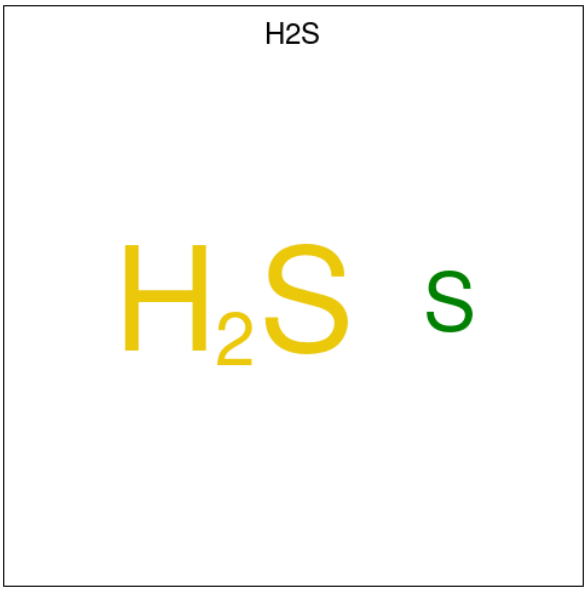
- Molecule 5 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Ni 1 1	0	0

- Molecule 6 is FE (II) ION (CCD ID: FE2) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total Fe 1 1	0	0

- Molecule 7 is HYDROSULFURIC ACID (CCD ID: H2S) (formula: H_2S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	B	1	Total S 1 1	0	0

- Molecule 8 is water.

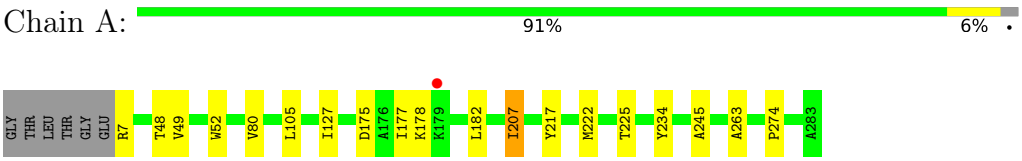
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	280	Total O 280 280	0	0
8	B	454	Total O 454 454	0	0

3 Residue-property plots (i)

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.

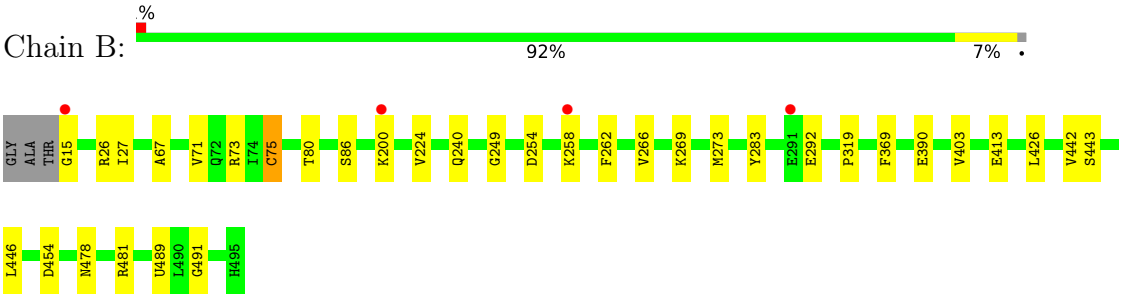
• Molecule 1: PERIPLASMIC [NIFESE] HYDROGENASE, SMALL SUBUNIT

Chain A:



• Molecule 2: PERIPLASMIC [NIFESE] HYDROGENASE, LARGE SUBUNIT, SELENOCYSTEINE-CONTAINING

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	72.30Å 96.99Å 103.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.78 – 1.95 45.78 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.0 (45.78-1.95) 99.2 (45.78-1.95)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.31 (at 1.95Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.153 , 0.190 0.152 , 0.190	Depositor DCC
R_{free} test set	2686 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	12.9	Xtriage
Anisotropy	0.639	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 51.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	6625	wwPDB-VP
Average B, all atoms (Å^2)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.16% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: FE2, SF4, H2S, NI, SEC, FCO, OCS

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	0.49	0/2184	0.77	1/2968 (0.0%)
2	B	0.51	0/3821	0.77	1/5161 (0.0%)
All	All	0.50	0/6005	0.77	2/8129 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	27	ILE	N-CA-C	-5.49	101.84	108.53
1	A	49	VAL	N-CA-C	5.09	118.61	113.47

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	A	2112	0	2060	11	0
2	B	3745	0	3762	20	0
3	A	24	0	0	0	0
4	B	7	0	0	0	0
5	B	1	0	0	0	0
6	B	1	0	0	0	0
7	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	A	280	0	0	2	0
8	B	454	0	0	5	0
All	All	6625	0	5822	29	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 (29) 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:B:390:GLU:OE2	8:B:2397:HOH:O	2.02	0.77
1:A:175:ASP:HA	1:A:178:LYS:HE3	1.76	0.66
2:B:15:GLY:O	8:B:2004:HOH:O	2.14	0.64
2:B:319:PRO:HG3	2:B:446:LEU:HG	1.85	0.58
2:B:75:OCS:OD2	2:B:489:SEC:SE	2.73	0.56
1:A:207:ILE:HD11	1:A:245:ALA:HB3	1.90	0.52
1:A:7:ARG:HB2	8:A:2001:HOH:O	2.09	0.50
2:B:283:TYR:HB3	2:B:454:ASP:OD1	2.14	0.48
8:A:2012:HOH:O	2:B:491:GLY:HA3	2.13	0.47
1:A:127:ILE:HD12	2:B:73:ARG:HG2	1.97	0.47
2:B:200:LYS:HG2	8:B:2232:HOH:O	2.15	0.46
2:B:71:VAL:HG21	2:B:86:SER:HB2	1.97	0.45
2:B:67:ALA:O	2:B:71:VAL:HG22	2.16	0.45
2:B:443:SER:O	2:B:446:LEU:HB3	2.17	0.45
1:A:48:THR:O	2:B:26:ARG:HA	2.17	0.45
1:A:80:VAL:HG21	1:A:105:LEU:HD12	1.99	0.45
2:B:240:GLN:HA	2:B:369:PHE:O	2.17	0.45
1:A:175:ASP:O	1:A:178:LYS:HG2	2.18	0.44
1:A:127:ILE:HG12	1:A:274:PRO:HG3	2.00	0.44
2:B:478:ASN:HA	2:B:481:ARG:HB2	2.00	0.43
2:B:413:GLU:OE2	8:B:2334:HOH:O	2.22	0.42
1:A:177:ILE:HG23	1:A:182:LEU:HA	2.01	0.42
2:B:249:GLY:HA2	2:B:262:PHE:O	2.19	0.42
2:B:269[A]:LYS:HD2	8:B:2301:HOH:O	2.19	0.42
2:B:426:LEU:HG	2:B:442:VAL:HB	2.02	0.42
2:B:266:VAL:HB	2:B:273:MET:HG2	2.02	0.41
2:B:254:ASP:OD2	2:B:258:LYS:HG3	2.20	0.41
1:A:217:TYR:HA	1:A:222[B]:MET:HE1	2.01	0.41
1:A:222[A]:MET:SD	1:A:234:TYR:HB3	2.61	0.41

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	280/283 (99%)	271 (97%)	8 (3%)	1 (0%)	30 21
2	B	479/484 (99%)	466 (97%)	13 (3%)	0	100 100
All	All	759/767 (99%)	737 (97%)	21 (3%)	1 (0%)	48 41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	263	ALA

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	A	227/226 (100%)	224 (99%)	3 (1%)	61 58
2	B	392/391 (100%)	388 (99%)	4 (1%)	68 67
All	All	619/617 (100%)	612 (99%)	7 (1%)	65 64

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	TRP
1	A	207	ILE
1	A	225	THR
2	B	80	THR
2	B	224	VAL

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Mol	Chain	Res	Type
2	B	292	GLU
2	B	403	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	GLN
1	A	96	HIS
1	A	111	ASN
2	B	255	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)

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	OCS	B	75	2,5	3,7,9	3.30	1 (33%)	1,8,13	1.18	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	OCS	B	75	2,5	-	0/2/6/9	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	75	OCS	O-C	5.38	1.41	1.19

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
2	B	75	OCS	1	0

5.5 Carbohydrates (i)

There are no oligosaccharides in this entry.

5.6 Ligand geometry (i)

Of 7 ligands modelled in this entry, 2 are monoatomic and 1 is modelled with single atom - leaving 4 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	SF4	A	284	1	0,12,12	-	-	-		
3	SF4	A	286	1	0,12,12	-	-	-		
4	FCO	B	500	2	1,6,6	2.21	1 (100%)	-		
3	SF4	A	285	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
3	SF4	A	284	1	-	-	0/6/5/5
3	SF4	A	286	1	-	-	0/6/5/5
3	SF4	A	285	1	-	-	0/6/5/5

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	500	FCO	O3-C3	-2.21	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

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, 95^{th} 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	277/283 (97%)	-0.61	1 (0%)	88	91	7, 13, 27, 46	11 (3%)
2	B	479/484 (98%)	-0.69	4 (0%)	82	86	7, 11, 23, 44	13 (2%)
All	All	756/767 (98%)	-0.66	5 (0%)	84	88	7, 12, 25, 46	24 (3%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	291	GLU	3.1
2	B	15	GLY	3.0
2	B	258	LYS	2.8
2	B	200	LYS	2.7
1	A	179	LYS	2.5

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

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, 95^{th} 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
2	OCS	B	75	8/10	0.98	0.05	7,9,12,13	0

6.3 Carbohydrates

There are no oligosaccharides 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. The B-factors column lists the minimum, median, 95^{th} 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
7	H2S	B	503	1/1	0.99	0.02	8,8,8,8	0
3	SF4	A	285	8/8	1.00	0.01	11,11,12,12	0
3	SF4	A	286	8/8	1.00	0.02	10,10,10,11	0
4	FCO	B	500	7/7	1.00	0.03	10,10,12,15	0
5	NI	B	501	1/1	1.00	0.01	12,12,12,12	0
6	FE2	B	502	1/1	1.00	0.01	11,11,11,11	0
3	SF4	A	284	8/8	1.00	0.02	13,13,14,14	0

6.5 Other polymers (i)

There are no such residues in this entry.