Full wwPDB X-ray Structure Validation Report

Jan 31, 2016 – 10:35 PM GMT

PDB ID	:	1UBJ
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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

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

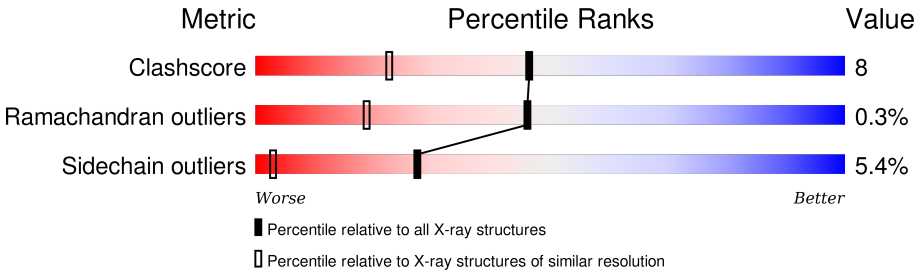
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

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 1.35 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	102246	2337 (1.40-1.32)
Ramachandran outliers	100387	2280 (1.40-1.32)
Sidechain outliers	100360	2279 (1.40-1.32)

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 <=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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	S	267	81% 14%
2	L	534	83% 14%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 7163 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			

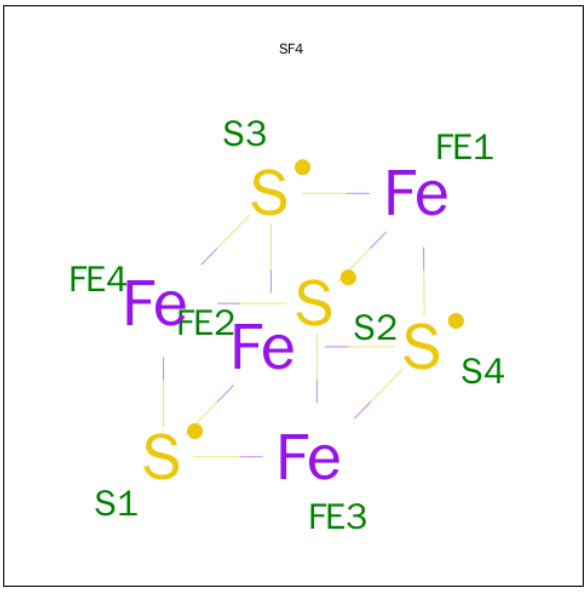
There are 2 discrepancies between the modelled and reference sequences:

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

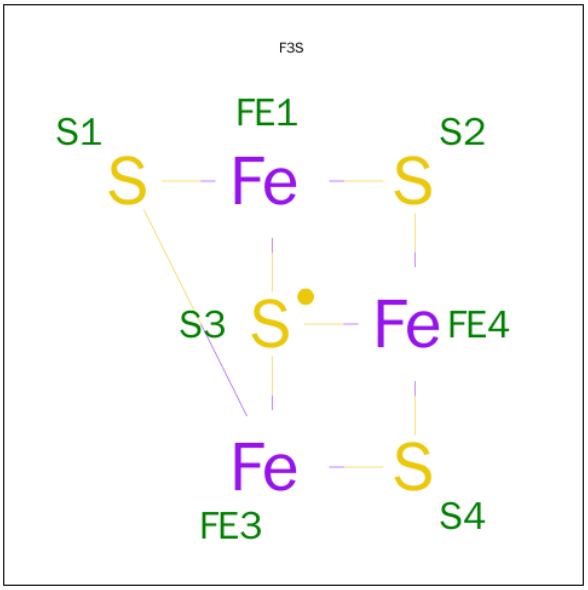
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	L	1	Total	Mg	0	0
			1	1		

- Molecule 4 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe_4S_4).



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

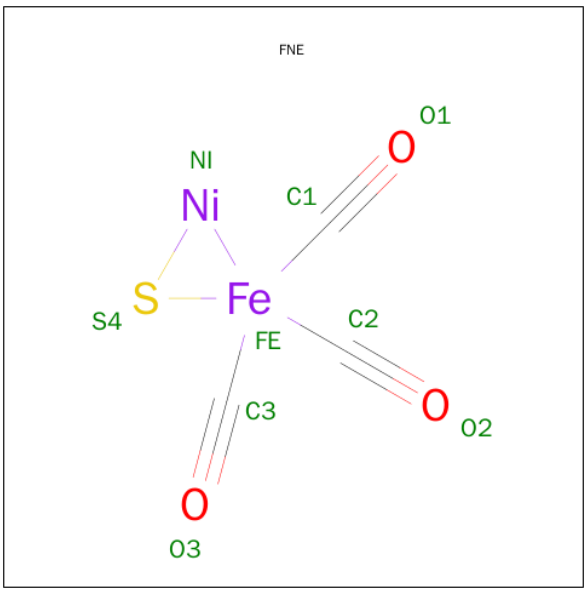
- Molecule 5 is FE3-S4 CLUSTER (three-letter code: F3S) (formula: Fe_3S_4).



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

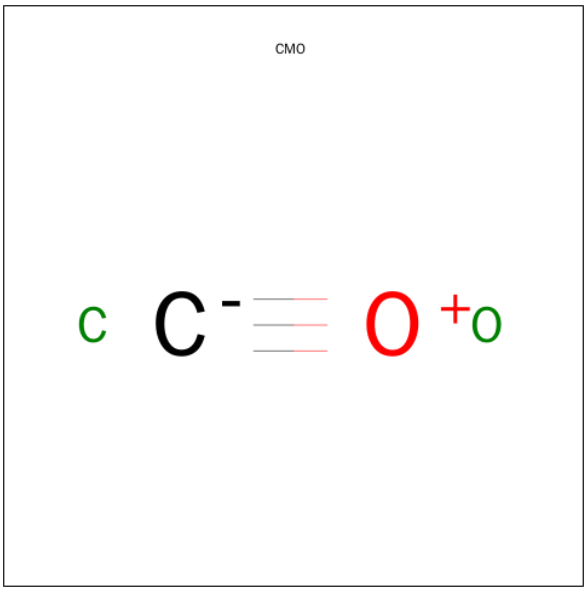
- Molecule 6 is (MU-SULPHIDO)-BIS(MU-CYS,S)-[TRICARBONYLIRON-DI-(CYS,S)NIC

KEL(II)](FE-NI) (three-letter code: FNE) (formula: C_3FeNiO_3S).



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 (three-letter code: 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	L	607	Total O 607 607	0	0
8	S	326	Total O 326 326	0	0

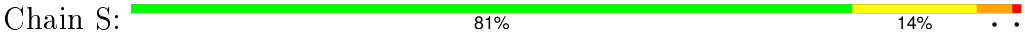
3 Residue-property plots (i)

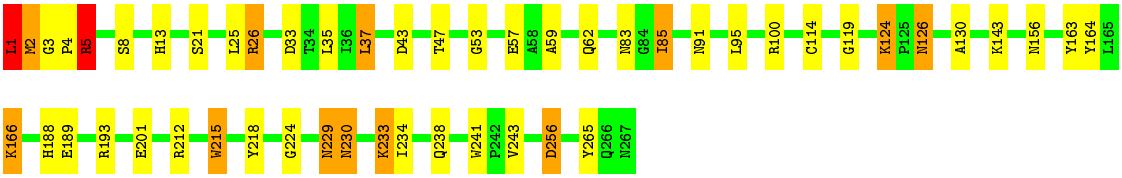
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

Note EDS was not executed.

- Molecule 1: Periplasmic [NiFe] hydrogenase small subunit

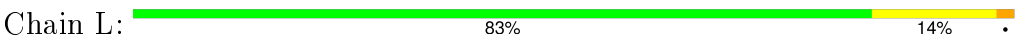
Chain S:

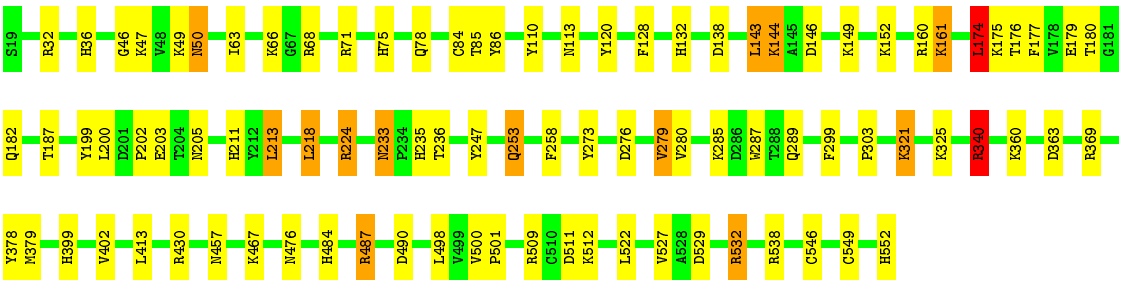




- Molecule 2: Periplasmic [NiFe] hydrogenase large subunit

Chain L:





4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	97.74Å 125.86Å 66.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.35	Depositor
% Data completeness (in resolution range)	71.5 (20.00-1.35)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.121 , 0.183	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7163	wwPDB-VP
Average B, all atoms (Å^2)	21.0	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: CMO, MG, F3S, FNE, SF4

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	0.62	0/2075	1.26	21/2830 (0.7%)
2	L	0.60	1/4288 (0.0%)	1.34	33/5831 (0.6%)
All	All	0.60	1/6363 (0.0%)	1.32	54/8661 (0.6%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	552	HIS	C-O	5.44	1.33	1.23

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	L	532	ARG	CD-NE-CZ	30.12	165.76	123.60
2	L	532	ARG	NE-CZ-NH2	-15.06	112.77	120.30
2	L	273	TYR	CB-CG-CD1	12.45	128.47	121.00
2	L	532	ARG	NE-CZ-NH1	12.13	126.37	120.30
2	L	160	ARG	NE-CZ-NH2	-9.37	115.62	120.30
2	L	71	ARG	NE-CZ-NH2	-9.21	115.69	120.30
2	L	224	ARG	NE-CZ-NH2	8.53	124.56	120.30
2	L	340	ARG	NE-CZ-NH1	8.32	124.46	120.30
1	S	212	ARG	NE-CZ-NH2	7.81	124.20	120.30
2	L	340	ARG	NE-CZ-NH2	-7.62	116.49	120.30
2	L	174	LEU	CA-CB-CG	7.55	132.66	115.30
2	L	128	PHE	CB-CG-CD1	-7.37	115.64	120.80
2	L	199	TYR	CB-CG-CD2	-7.29	116.63	121.00
2	L	273	TYR	CB-CG-CD2	-7.25	116.65	121.00
1	S	256	ASP	CB-CG-OD1	7.23	124.81	118.30
1	S	218	TYR	CB-CG-CD1	7.21	125.32	121.00
1	S	163	TYR	CB-CG-CD2	-7.10	116.74	121.00
1	S	163	TYR	CB-CG-CD1	7.08	125.25	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	L	218	LEU	CA-CB-CG	-6.95	99.31	115.30
1	S	212	ARG	NE-CZ-NH1	-6.74	116.93	120.30
1	S	1	LEU	O-C-N	6.68	133.39	122.70
2	L	538	ARG	NE-CZ-NH1	6.56	123.58	120.30
2	L	369	ARG	NE-CZ-NH2	-6.54	117.03	120.30
2	L	146	ASP	CB-CG-OD2	-6.30	112.63	118.30
2	L	110	TYR	CB-CG-CD1	6.23	124.74	121.00
1	S	85	ILE	CA-CB-CG2	6.12	123.13	110.90
1	S	218	TYR	CB-CG-CD2	-6.08	117.35	121.00
2	L	538	ARG	NE-CZ-NH2	-6.01	117.29	120.30
2	L	363	ASP	CB-CG-OD2	5.92	123.63	118.30
1	S	33	ASP	CB-CG-OD1	5.88	123.59	118.30
1	S	164	TYR	CB-CG-CD1	-5.88	117.47	121.00
2	L	287	TRP	CA-CB-CG	-5.79	102.69	113.70
1	S	193	ARG	NE-CZ-NH1	5.73	123.16	120.30
2	L	430	ARG	NE-CZ-NH2	-5.62	117.49	120.30
2	L	146	ASP	CB-CG-OD1	5.62	123.36	118.30
2	L	120	TYR	CA-CB-CG	5.60	124.04	113.40
1	S	5	ARG	CA-C-N	5.60	129.52	117.20
1	S	233	LYS	CD-CE-NZ	5.60	124.57	111.70
2	L	247	TYR	CB-CG-CD1	5.58	124.35	121.00
2	L	160	ARG	NH1-CZ-NH2	5.57	125.52	119.40
2	L	273	TYR	CG-CD1-CE1	5.53	125.73	121.30
1	S	215	TRP	CH2-CZ2-CE2	-5.45	111.95	117.40
1	S	83	ASN	CA-CB-CG	-5.41	101.50	113.40
1	S	26	ARG	CG-CD-NE	-5.41	100.44	111.80
1	S	189	GLU	OE1-CD-OE2	-5.35	116.88	123.30
1	S	5	ARG	CA-C-O	-5.33	108.91	120.10
2	L	487	ARG	NE-CZ-NH1	5.29	122.95	120.30
2	L	213	LEU	CA-CB-CG	5.24	127.35	115.30
2	L	247	TYR	CB-CG-CD2	-5.11	117.93	121.00
1	S	100	ARG	NE-CZ-NH1	5.09	122.85	120.30
1	S	57	GLU	CA-CB-CG	5.09	124.60	113.40
2	L	378	TYR	CB-CG-CD2	-5.06	117.97	121.00
2	L	511	ASP	CB-CG-OD2	5.04	122.83	118.30
2	L	258	PHE	CB-CG-CD1	5.03	124.32	120.80

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	2019	0	1949	48	0
2	L	4177	0	4129	61	0
3	L	1	0	0	0	0
4	S	16	0	0	0	0
5	S	7	0	0	0	0
6	L	8	0	0	1	0
7	L	2	0	0	0	0
8	L	607	0	0	16	8
8	S	326	0	0	2	4
All	All	7163	0	6078	94	8

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 8.

All (94) 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:230:ASN:HB2	1:S:233:LYS:HE2	1.53	0.89
1:S:26:ARG:HH21	2:L:233:ASN:HD21	1.22	0.85
1:S:1:LEU:HD11	1:S:62:GLN:HG2	1.58	0.84
1:S:2:MET:HG2	2:L:182:GLN:HE21	1.47	0.80
1:S:126:ASN:HD21	1:S:130:ALA:H	1.26	0.79
2:L:467:LYS:HB3	8:L:3930:HOH:O	1.83	0.78
1:S:238:GLN:HE21	2:L:224:ARG:HH21	1.32	0.77
2:L:218:LEU:HG	2:L:218:LEU:O	1.84	0.76
1:S:2:MET:HG3	1:S:8:SER:HB2	1.70	0.74
2:L:176:THR:O	2:L:180:THR:HG23	1.89	0.73
2:L:161:LYS:HD3	8:L:3446:HOH:O	1.88	0.72
1:S:2:MET:HG2	2:L:182:GLN:NE2	2.03	0.72
2:L:78:GLN:HE21	2:L:86:TYR:H	1.37	0.72
1:S:2:MET:HB2	1:S:43:ASP:OD1	1.91	0.69
1:S:256:ASP:HB3	8:S:3427:HOH:O	1.96	0.66
2:L:484:HIS:CD2	2:L:498:LEU:HD22	2.32	0.64
2:L:175:LYS:O	2:L:179:GLU:HG3	1.98	0.64
1:S:1:LEU:HD21	1:S:59:ALA:O	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:230:ASN:HD22	1:S:230:ASN:H	1.44	0.64
2:L:289:GLN:HG3	8:L:3392:HOH:O	2.02	0.60
2:L:253:GLN:HE21	2:L:253:GLN:H	1.49	0.60
2:L:490:ASP:HA	8:L:3930:HOH:O	2.01	0.59
2:L:143:LEU:HD13	2:L:174:LEU:HD13	1.83	0.59
2:L:211:HIS:HE1	8:L:3257:HOH:O	1.86	0.59
1:S:13:HIS:HD2	8:S:3055:HOH:O	1.84	0.59
2:L:529:ASP:OD1	2:L:532:ARG:HD3	2.01	0.58
2:L:149:LYS:HE2	2:L:203:GLU:OE2	2.04	0.57
1:S:53:GLY:HA3	8:L:3859:HOH:O	2.03	0.56
1:S:233:LYS:HE3	1:S:234:ILE:CD1	2.36	0.56
2:L:75:HIS:HD2	8:L:3039:HOH:O	1.89	0.55
1:S:1:LEU:HD22	1:S:43:ASP:HB3	1.89	0.54
1:S:1:LEU:HG	2:L:187:THR:HG21	1.90	0.54
1:S:126:ASN:HD21	1:S:130:ALA:N	2.03	0.53
2:L:321:LYS:HD3	2:L:321:LYS:N	2.23	0.53
2:L:50:ASN:HD21	2:L:509:ARG:NH2	2.06	0.53
1:S:37:LEU:HD12	2:L:177:PHE:CD2	2.44	0.53
1:S:156:ASN:HD21	1:S:230:ASN:HD21	1.56	0.52
2:L:36:HIS:HD2	8:L:3157:HOH:O	1.93	0.52
2:L:253:GLN:NE2	2:L:253:GLN:H	2.08	0.52
2:L:66:LYS:HE3	8:L:3952:HOH:O	2.10	0.51
2:L:299:PHE:H	2:L:476:ASN:ND2	2.08	0.51
1:S:233:LYS:HE3	1:S:234:ILE:HD11	1.93	0.51
2:L:211:HIS:HD2	2:L:276:ASP:OD2	1.94	0.51
1:S:2:MET:CG	1:S:8:SER:HB2	2.38	0.51
1:S:1:LEU:HD12	1:S:1:LEU:H3	1.75	0.50
2:L:512:LYS:HE2	8:L:3471:HOH:O	2.11	0.50
1:S:230:ASN:ND2	1:S:230:ASN:H	2.09	0.50
1:S:2:MET:HA	2:L:182:GLN:HG2	1.93	0.49
1:S:2:MET:C	2:L:182:GLN:HG2	2.33	0.49
1:S:26:ARG:HH21	2:L:233:ASN:ND2	2.01	0.48
2:L:340:ARG:HD2	8:L:3520:HOH:O	2.14	0.48
2:L:399:HIS:HD2	2:L:402:VAL:H	1.61	0.47
2:L:487:ARG:HG3	8:L:3230:HOH:O	2.13	0.47
1:S:230:ASN:N	1:S:230:ASN:HD22	2.07	0.47
1:S:13:HIS:HE1	1:S:21:SER:OG	1.97	0.47
1:S:241:TRP:CH2	1:S:243:VAL:HB	2.50	0.47
1:S:201:GLU:HB3	1:S:215:TRP:CE3	2.49	0.47
1:S:5:ARG:HB2	8:L:3759:HOH:O	2.14	0.46
1:S:2:MET:SD	1:S:8:SER:HB2	2.55	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:238:GLN:NE2	2:L:224:ARG:HH21	2.05	0.46
1:S:265:TYR:OH	2:L:75:HIS:HE1	1.99	0.46
1:S:124:LYS:HG2	1:S:124:LYS:HZ2	1.45	0.46
1:S:233:LYS:HE3	1:S:234:ILE:HG12	1.97	0.46
2:L:399:HIS:CD2	2:L:402:VAL:H	2.35	0.45
2:L:85:THR:OG1	2:L:235:HIS:HD2	1.99	0.45
2:L:113:ASN:HD21	2:L:303:PRO:HD2	1.81	0.45
1:S:1:LEU:HB3	2:L:187:THR:OG1	2.18	0.44
2:L:50:ASN:HD21	2:L:509:ARG:HH22	1.66	0.44
2:L:500:VAL:CG1	2:L:501:PRO:HD2	2.48	0.44
2:L:144:LYS:HG2	2:L:202:PRO:HB3	2.00	0.44
1:S:2:MET:HA	2:L:182:GLN:HE21	1.83	0.44
1:S:1:LEU:HD11	1:S:62:GLN:CG	2.40	0.43
1:S:229:ASN:HD22	1:S:230:ASN:H	1.66	0.43
1:S:166:LYS:HB2	1:S:166:LYS:HE2	1.57	0.43
2:L:138:ASP:H	2:L:205:ASN:ND2	2.15	0.43
2:L:546:CYS:SG	2:L:549:CYS:HB2	2.59	0.43
2:L:78:GLN:HE22	2:L:236:THR:H	1.65	0.43
2:L:63:ILE:HA	2:L:66:LYS:HE2	2.01	0.43
1:S:114:CYS:HA	1:S:119:GLY:HA3	2.00	0.43
2:L:47:LYS:NZ	8:L:3716:HOH:O	2.51	0.42
2:L:285:LYS:NZ	2:L:413:LEU:O	2.49	0.42
2:L:68:ARG:HG2	8:L:3738:HOH:O	2.18	0.42
2:L:143:LEU:HD12	2:L:143:LEU:HA	1.71	0.42
1:S:2:MET:CG	1:S:3:GLY:H	2.33	0.42
2:L:484:HIS:CG	2:L:498:LEU:HD22	2.54	0.42
2:L:299:PHE:H	2:L:476:ASN:HD22	1.68	0.42
2:L:279:VAL:HG12	2:L:280:VAL:N	2.35	0.42
1:S:47:THR:O	2:L:32:ARG:HA	2.19	0.41
1:S:5:ARG:HA	1:S:5:ARG:CZ	2.50	0.41
2:L:46:GLY:C	2:L:47:LYS:HD2	2.40	0.41
1:S:2:MET:CA	2:L:182:GLN:HG2	2.50	0.41
2:L:84:CYS:CB	6:L:1004:FNE:C2	2.99	0.41
2:L:379:MET:HE1	8:L:3911:HOH:O	2.22	0.40
1:S:188:HIS:HB2	1:S:224:GLY:C	2.42	0.40

All (8) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:L:3457:HOH:O	8:L:3715:HOH:O[3_647]	1.90	0.30

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:L:3582:HOH:O	8:L:3700:HOH:O[2_664]	1.97	0.23
8:L:3270:HOH:O	8:L:3670:HOH:O[3_657]	2.01	0.19
8:S:3184:HOH:O	8:L:3740:HOH:O[2_664]	2.02	0.18
8:L:3670:HOH:O	8:L:3715:HOH:O[3_647]	2.06	0.14
8:S:3496:HOH:O	8:L:3419:HOH:O[2_664]	2.12	0.08
8:S:3483:HOH:O	8:L:3482:HOH:O[2_664]	2.13	0.07
8:S:3416:HOH:O	8:L:3779:HOH:O[2_664]	2.18	0.02

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	S	265/267 (99%)	256 (97%)	7 (3%)	2 (1%)	24 5
2	L	532/534 (100%)	519 (98%)	13 (2%)	0	100 100
All	All	797/801 (100%)	775 (97%)	20 (2%)	2 (0%)	46 19

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	S	5	ARG
1	S	4	PRO

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	S	213/213 (100%)	198 (93%)	15 (7%)	19 1
2	L	438/438 (100%)	418 (95%)	20 (5%)	33 4
All	All	651/651 (100%)	616 (95%)	35 (5%)	27 3

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	S	1	LEU
1	S	2	MET
1	S	5	ARG
1	S	25	LEU
1	S	35	LEU
1	S	37	LEU
1	S	85	ILE
1	S	91	ASN
1	S	95	LEU
1	S	124	LYS
1	S	126	ASN
1	S	143	LYS
1	S	166	LYS
1	S	229	ASN
1	S	230	ASN
2	L	49	LYS
2	L	50	ASN
2	L	132	HIS
2	L	143	LEU
2	L	144	LYS
2	L	152	LYS
2	L	161	LYS
2	L	174	LEU
2	L	200	LEU
2	L	213	LEU
2	L	233	ASN
2	L	253	GLN
2	L	279	VAL
2	L	321	LYS
2	L	325	LYS
2	L	340	ARG
2	L	360	LYS
2	L	457	ASN
2	L	522	LEU
2	L	527	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	S	13	HIS
1	S	91	ASN
1	S	126	ASN
1	S	139	HIS
1	S	190	GLN
1	S	229	ASN
1	S	230	ASN
1	S	238	GLN
1	S	266	GLN
1	S	267	ASN
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	310	ASN
2	L	334	GLN
2	L	390	GLN
2	L	399	HIS
2	L	446	GLN
2	L	451	ASN
2	L	457	ASN
2	L	476	ASN
2	L	513	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 carbohydrates 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	3,7,9	1.13	0	0,9,15	0.00	-
7	CMO	L	1006	6	0,1,1	0.00	-	0,0,0	0.00	-
4	SF4	S	1001	1	0,12,12	0.00	-	0,24,24	0.00	-
4	SF4	S	1002	1	0,12,12	0.00	-	0,24,24	0.00	-
5	F3S	S	1003	1	0,9,9	0.00	-	0,15,15	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
6	FNE	L	1004	2,7	-	0/0/9/18	0/0/0/1
7	CMO	L	1006	6	-	0/0/0/0	0/0/0/0
4	SF4	S	1001	1	-	0/0/48/48	2/6/5/5
4	SF4	S	1002	1	-	0/0/48/48	2/6/5/5
5	F3S	S	1003	1	-	0/0/24/24	0/0/3/3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

All (4) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	S	1001	SF4	FE1-FE2-S3-S4
4	S	1001	SF4	FE3-FE4-S1-S2
4	S	1002	SF4	FE3-FE4-S1-S2
4	S	1002	SF4	FE1-FE2-S3-S4

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	L	1004	FNE	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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.