



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 13, 2018 – 02:03 pm GMT

PDB ID : 5VQ4
Title : Nitrogenase Av1 at pH 5
Authors : Morrison, C.N.; Spatzal, T.; Rees, D.C.
Deposited on : 2017-05-07
Resolution : 2.30 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : 1.13
EDS : trunk31020
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk31020

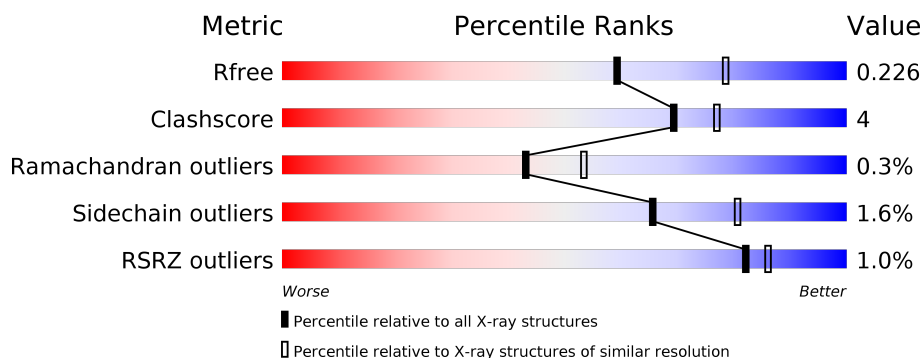
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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.



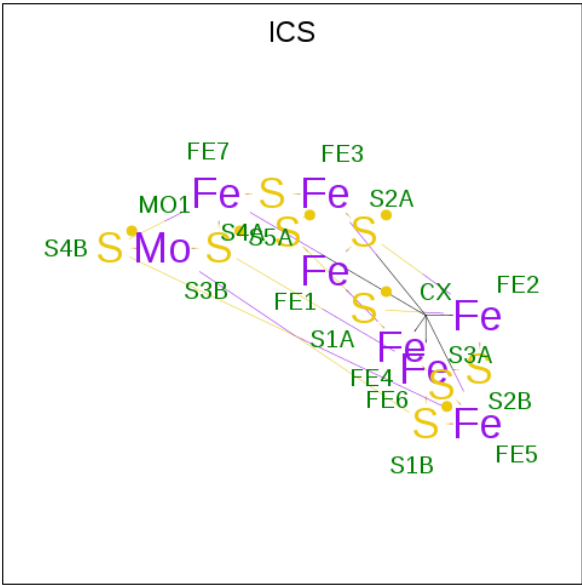
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	111664	4477 (2.30-2.30)
Clashscore	122126	5072 (2.30-2.30)
Ramachandran outliers	120053	5022 (2.30-2.30)
Sidechain outliers	120020	5021 (2.30-2.30)
RSRZ outliers	108989	4374 (2.30-2.30)

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 $\leq 5\%$. 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	A	492	<div> <div>2%</div> <div> <div></div> <div>85%</div> <div>11%</div> <div>••</div> </div> </div>
1	C	492	<div> <div>%</div> <div> <div></div> <div>85%</div> <div>11%</div> <div>••</div> </div> </div>
2	B	523	<div> <div></div> <div> <div>92%</div> <div>8%</div> </div> </div>
2	D	523	<div> <div></div> <div> <div>91%</div> <div>8%</div> </div> </div>

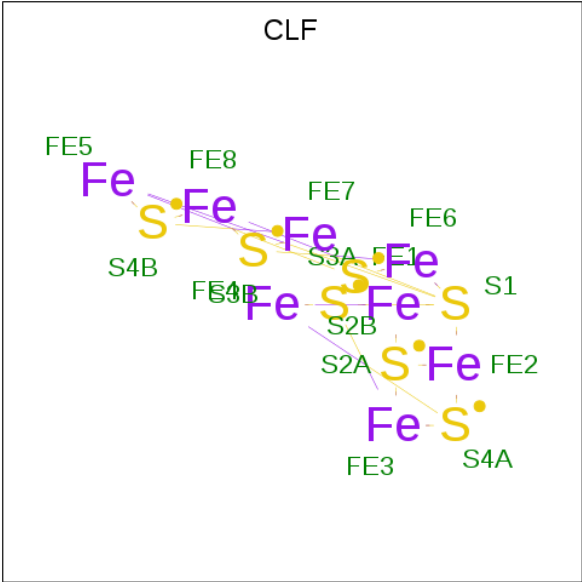
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			14	7	7		
3	C	1	Total	C	O	0	0
			14	7	7		

- Molecule 4 is iron-sulfur-molybdenum cluster with interstitial carbon (three-letter code: ICS) (formula: CFe₇MoS₉).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	Fe	Mo	S	0	0
			18	1	7	1	9		
4	C	1	Total	C	Fe	Mo	S	0	0
			18	1	7	1	9		

- Molecule 5 is FE(8)-S(7) CLUSTER (three-letter code: CLF) (formula: Fe₈S₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	Fe	S	0	1
			15	8	7		
5	C	1	Total	Fe	S	0	1
			15	8	7		

- Molecule 6 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	2	Total	Fe	0	0
			2	2		

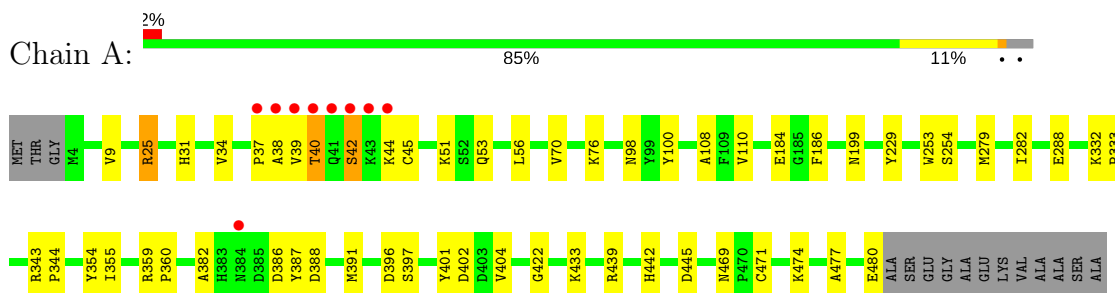
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	74	Total	O	0	0
			74	74		
7	B	114	Total	O	0	0
			114	114		
7	C	76	Total	O	0	0
			76	76		
7	D	98	Total	O	0	0
			98	98		

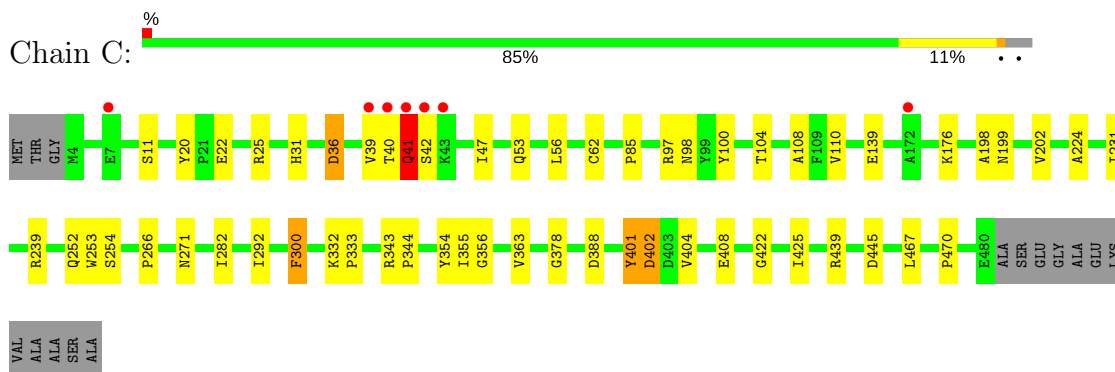
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

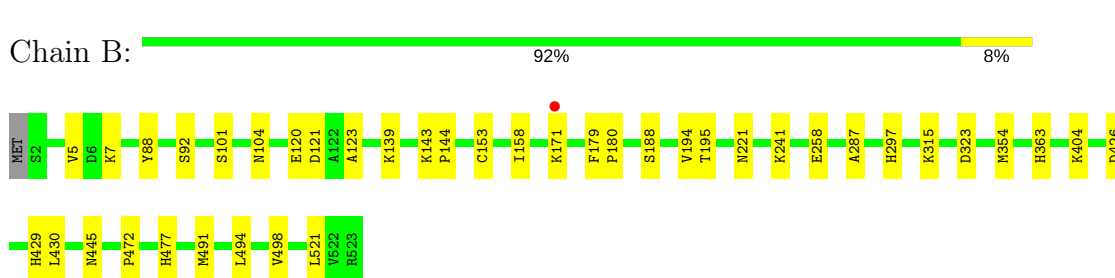
- Molecule 1: Nitrogenase molybdenum-iron protein alpha chain



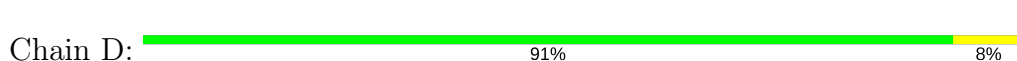
- Molecule 1: Nitrogenase molybdenum-iron protein alpha chain

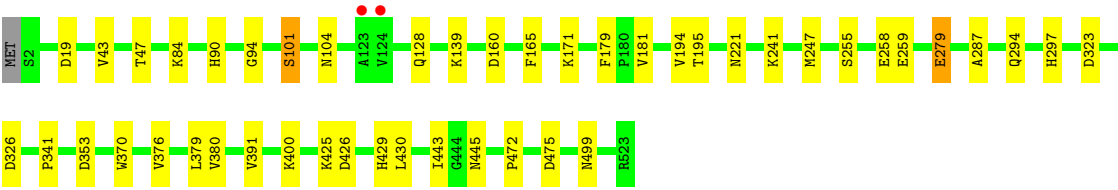


- Molecule 2: Nitrogenase molybdenum-iron protein beta chain



- Molecule 2: Nitrogenase molybdenum-iron protein beta chain





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.31Å 128.86Å 108.36Å 90.00° 110.89° 90.00°	Depositor
Resolution (Å)	39.54 – 2.30 39.54 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.7 (39.54-2.30) 98.7 (39.54-2.30)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.04 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.8.0151	Depositor
R, R_{free}	0.168 , 0.222 0.176 , 0.226	Depositor DCC
R_{free} test set	4320 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å ²)	21.1	Xtriage
Anisotropy	0.484	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 27.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.018 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	16419	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.80% 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: ICS, CLF, HCA, FE

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	0.58	0/3878	0.70	0/5229
1	C	0.59	0/3878	0.70	1/5229 (0.0%)
2	B	0.60	0/4289	0.69	2/5799 (0.0%)
2	D	0.59	0/4310	0.70	1/5827 (0.0%)
All	All	0.59	0/16355	0.70	4/22084 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	171	LYS	CD-CE-NZ	6.88	127.53	111.70
2	B	323	ASP	CB-CG-OD1	5.88	123.60	118.30
2	D	19	ASP	CB-CG-OD1	5.77	123.49	118.30
1	C	402	ASP	CB-CG-OD2	-5.51	113.34	118.30

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	A	3790	0	3729	46	0
1	C	3790	0	3729	34	0
2	B	4180	0	4094	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	4201	0	4107	29	0
3	A	14	0	6	1	0
3	C	14	0	6	1	0
4	A	18	0	0	2	0
4	C	18	0	0	1	0
5	A	15	0	0	0	0
5	C	15	0	0	1	0
6	B	2	0	0	0	0
7	A	74	0	0	1	0
7	B	114	0	0	1	0
7	C	76	0	0	1	0
7	D	98	0	0	1	0
All	All	16419	0	15671	128	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 4.

All (128) 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:A:39:VAL:CG1	1:A:44:LYS:CE	1.75	1.63
1:A:39:VAL:HG11	1:A:44:LYS:CD	1.18	1.61
1:A:39:VAL:CG1	1:A:44:LYS:HD2	1.41	1.37
1:A:39:VAL:CG1	1:A:44:LYS:CD	1.83	1.32
1:A:39:VAL:HG13	1:A:44:LYS:CE	1.42	1.27
1:A:39:VAL:HG12	1:A:44:LYS:HE2	1.23	1.11
1:A:39:VAL:HG12	1:A:44:LYS:CE	1.78	1.05
1:A:25:ARG:HH11	1:A:25:ARG:HG2	0.88	1.00
1:A:39:VAL:CG1	1:A:44:LYS:HE2	1.73	0.98
1:A:37:PRO:HD2	1:A:396:ASP:OD2	1.64	0.97
1:A:25:ARG:NH1	1:A:25:ARG:HG2	1.66	0.97
1:A:25:ARG:HH11	1:A:25:ARG:CG	1.78	0.95
1:A:39:VAL:HG13	1:A:44:LYS:HE3	0.90	0.90
2:B:477:HIS:H	2:D:499:ASN:HD21	1.20	0.89
2:B:426:ASP:H	2:B:429:HIS:HD2	1.29	0.81
1:C:36:ASP:HB3	1:C:39:VAL:HG12	1.62	0.79
2:D:279[A]:GLU:H	2:D:279[A]:GLU:CD	1.84	0.77
1:A:39:VAL:CG1	1:A:44:LYS:HE3	1.69	0.75
1:C:36:ASP:CB	1:C:39:VAL:HG12	2.16	0.75
1:A:42:SER:HB3	1:A:388:ASP:OD1	1.86	0.74
2:D:426:ASP:H	2:D:429:HIS:HD2	1.35	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:ARG:NH1	1:A:25:ARG:CG	2.45	0.71
2:D:279[A]:GLU:N	2:D:279[A]:GLU:OE1	2.28	0.66
1:A:39:VAL:HG13	1:A:44:LYS:CD	1.87	0.65
1:A:288:GLU:OE2	7:A:601:HOH:O	2.15	0.64
2:B:363:HIS:HD2	7:B:813:HOH:O	1.81	0.64
2:B:241:LYS:CE	2:B:258:GLU:OE2	2.49	0.61
1:A:354:TYR:CZ	1:A:404:VAL:HG12	2.37	0.60
1:A:51:LYS:NZ	2:B:121:ASP:OD2	2.35	0.60
1:C:53:GLN:HB2	1:C:56:LEU:HD12	1.84	0.59
1:C:97:ARG:O	1:C:231:ILE:HA	2.03	0.58
2:D:43:VAL:O	2:D:47:THR:HG23	2.04	0.57
2:B:426:ASP:H	2:B:429:HIS:CD2	2.17	0.57
2:D:294:GLN:NE2	7:D:702:HOH:O	2.39	0.55
1:C:40:THR:O	1:C:41:GLN:O	2.25	0.55
1:C:100:TYR:CE1	1:C:110:VAL:HB	2.42	0.55
1:C:22:GLU:OE2	1:C:25:ARG:NH2	2.40	0.54
1:A:253:TRP:HB2	1:A:254:SER:HA	1.89	0.54
1:C:239:ARG:HH11	1:C:252:GLN:HE21	1.55	0.54
1:C:224:ALA:HB3	1:C:271:ASN:HD22	1.73	0.54
1:A:53:GLN:HB2	1:A:56:LEU:HD12	1.90	0.54
1:C:39:VAL:O	1:C:39:VAL:HG13	2.08	0.53
1:C:42:SER:N	1:C:388:ASP:OD1	2.38	0.53
2:D:279[A]:GLU:N	2:D:279[A]:GLU:CD	2.60	0.53
2:D:426:ASP:H	2:D:429:HIS:CD2	2.20	0.53
1:C:199:ASN:HD22	1:C:282:ILE:H	1.57	0.52
2:B:194:VAL:HB	2:B:297:HIS:CG	2.45	0.52
1:A:469:ASN:HD22	1:A:471:CYS:H	1.57	0.51
2:B:123:ALA:HA	2:B:158:ILE:HD11	1.93	0.50
1:A:100:TYR:CE1	1:A:110:VAL:HB	2.46	0.50
1:A:9:VAL:HG11	1:A:34:VAL:HG22	1.94	0.50
1:C:354:TYR:CZ	1:C:404:VAL:HG12	2.47	0.49
2:D:425:LYS:HB3	2:D:429:HIS:CD2	2.47	0.49
2:D:128:GLN:HE22	2:D:165:PHE:HA	1.78	0.49
2:D:90:HIS:NE2	2:D:160:ASP:OD2	2.46	0.49
2:D:445:ASN:HB2	2:D:472:PRO:O	2.14	0.48
2:B:354:MET:SD	2:B:491:MET:HG2	2.52	0.48
1:C:31:HIS:HD2	1:C:402:ASP:OD2	1.96	0.48
2:D:194:VAL:HB	2:D:297:HIS:CG	2.49	0.48
1:A:31:HIS:HD2	1:A:402:ASP:OD2	1.97	0.48
2:B:221:ASN:OD1	2:B:287:ALA:HA	2.13	0.48
2:D:247:MET:HG2	2:D:341:PRO:HD3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:THR:OG1	1:A:40:THR:O	2.31	0.48
1:A:387:TYR:O	1:A:391:MET:HG3	2.14	0.47
2:B:445:ASN:HB2	2:B:472:PRO:O	2.15	0.47
1:C:422:GLY:HA2	1:C:439:ARG:O	2.15	0.47
1:A:37:PRO:HD2	1:A:396:ASP:CG	2.32	0.47
1:C:354:TYR:CE2	1:C:404:VAL:HG12	2.49	0.46
1:C:425:ILE:HG22	7:C:602:HOH:O	2.14	0.46
1:C:139:GLU:OE2	1:C:176:LYS:HD2	2.15	0.46
2:D:376:VAL:O	2:D:380:VAL:HG23	2.15	0.46
2:D:323:ASP:O	2:D:326:ASP:HB2	2.15	0.46
2:B:241:LYS:NZ	2:B:258:GLU:OE2	2.46	0.46
1:C:36:ASP:HB2	1:C:39:VAL:CG1	2.46	0.46
1:A:442:HIS:CG	3:A:501:HCA:H52	2.51	0.46
1:C:239:ARG:HH11	1:C:252:GLN:NE2	2.12	0.46
2:B:120:GLU:O	2:B:123:ALA:HB3	2.16	0.46
3:C:501:HCA:O1	3:C:501:HCA:O7	2.34	0.45
1:C:36:ASP:HB2	1:C:39:VAL:HG12	1.97	0.45
1:C:253:TRP:HB2	1:C:254:SER:HA	1.99	0.45
2:D:101:SER:HA	2:D:104:ASN:HD22	1.81	0.45
2:B:521:LEU:O	2:D:475:ASP:HB3	2.17	0.45
1:A:9:VAL:CG1	1:A:34:VAL:HG22	2.46	0.44
2:D:379:LEU:HD21	2:D:443:ILE:HG21	1.99	0.44
1:A:199:ASN:HD22	1:A:282:ILE:H	1.64	0.44
2:B:101:SER:HA	2:B:104:ASN:HD22	1.83	0.44
1:C:343:ARG:HB3	1:C:344:PRO:HD3	1.98	0.44
2:B:143:LYS:N	2:B:144:PRO:CD	2.81	0.44
2:D:241:LYS:CE	2:D:258:GLU:OE2	2.65	0.44
2:B:5:VAL:HG22	2:B:5:VAL:O	2.18	0.43
1:C:62:CYS:HB3	2:D:94:GLY:HA3	2.00	0.43
1:A:359:ARG:N	1:A:360:PRO:CD	2.81	0.43
1:C:85:PRO:HB2	5:C:503[B]:CLF:S2B	2.58	0.43
1:A:199:ASN:HD21	1:A:279:MET:HA	1.83	0.43
1:A:477:ALA:HB3	1:A:480:GLU:HG3	2.00	0.43
1:C:266:PRO:O	1:C:292:ILE:HD11	2.18	0.43
1:A:76:LYS:O	1:A:108:ALA:HA	2.18	0.43
1:C:47:ILE:HD13	1:C:47:ILE:HA	1.80	0.42
1:A:39:VAL:HG11	1:A:44:LYS:HD2	0.46	0.42
2:B:139:LYS:HE2	2:B:179:PHE:CE1	2.54	0.42
2:D:221:ASN:OD1	2:D:287:ALA:HA	2.19	0.42
2:D:181:VAL:O	2:D:181:VAL:HG12	2.20	0.42
1:C:356:GLY:HA3	4:C:502:ICS:S1B	2.60	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:104:THR:HA	1:C:108:ALA:O	2.20	0.42
1:C:300:PHE:CZ	1:C:363:VAL:HG12	2.55	0.42
2:D:370:TRP:CE3	2:D:430:LEU:CD2	3.02	0.42
2:B:7:LYS:HA	2:B:7:LYS:HD2	1.79	0.41
2:B:153:CYS:HB3	2:B:188:SER:OG	2.21	0.41
1:C:332:LYS:HB3	1:C:333:PRO:CD	2.51	0.41
1:C:198:ALA:O	1:C:202:VAL:HG23	2.20	0.41
1:A:332:LYS:HB3	1:A:333:PRO:CD	2.51	0.41
2:B:179:PHE:HA	2:B:180:PRO:HD3	1.89	0.41
1:A:433:LYS:NZ	2:D:353:ASP:OD2	2.45	0.41
1:A:70:VAL:HG21	4:A:502:ICS:S2B	2.60	0.41
2:B:494:LEU:O	2:B:498:VAL:HG23	2.21	0.41
1:A:422:GLY:HA2	1:A:439:ARG:O	2.21	0.40
2:B:123:ALA:HA	2:B:158:ILE:CD1	2.50	0.40
1:C:20:TYR:OH	1:C:408:GLU:OE2	2.36	0.40
2:D:171:LYS:HB3	2:D:171:LYS:HE2	1.93	0.40
2:D:258:GLU:HG3	2:D:259:GLU:N	2.34	0.40
1:A:229:TYR:CE2	4:A:502:ICS:S2A	3.14	0.40
2:D:426:ASP:N	2:D:429:HIS:HD2	2.12	0.40
1:A:343:ARG:N	1:A:344:PRO:CD	2.84	0.40
1:A:474:LYS:HA	1:A:474:LYS:HD3	1.91	0.40
1:C:378:GLY:HA3	1:C:401:TYR:O	2.21	0.40
1:A:184:GLU:HB3	1:A:186:PHE:CE2	2.56	0.40
1:A:382:ALA:HB1	1:A:386:ASP:HB2	2.04	0.40
2:D:139:LYS:HE2	2:D:179:PHE:CD1	2.57	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	A	475/492 (96%)	451 (95%)	22 (5%)	2 (0%)	36 45

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	475/492 (96%)	450 (95%)	23 (5%)	2 (0%)	36	45
2	B	521/523 (100%)	505 (97%)	16 (3%)	0	100	100
2	D	524/523 (100%)	508 (97%)	14 (3%)	2 (0%)	36	45
All	All	1995/2030 (98%)	1914 (96%)	75 (4%)	6 (0%)	43	53

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	41	GLN
1	A	38	ALA
2	D	84	LYS
2	D	255	SER
1	C	355	ILE
1	A	355	ILE

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	407/415 (98%)	399 (98%)	8 (2%)	58	75
1	C	407/415 (98%)	398 (98%)	9 (2%)	55	72
2	B	455/455 (100%)	449 (99%)	6 (1%)	71	84
2	D	458/455 (101%)	452 (99%)	6 (1%)	71	84
All	All	1727/1740 (99%)	1698 (98%)	29 (2%)	65	79

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	ARG
1	A	40	THR
1	A	42	SER
1	A	45	CYS
1	A	98	ASN

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Mol	Chain	Res	Type
1	A	397	SER
1	A	401	TYR
1	A	445	ASP
2	B	88	TYR
2	B	92	SER
2	B	195	THR
2	B	315	LYS
2	B	404	LYS
2	B	430	LEU
1	C	11	SER
1	C	36	ASP
1	C	41	GLN
1	C	98	ASN
1	C	300	PHE
1	C	401	TYR
1	C	445	ASP
1	C	467	LEU
1	C	470	PRO
2	D	101	SER
2	D	195	THR
2	D	279[A]	GLU
2	D	279[B]	GLU
2	D	391	VAL
2	D	400	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	HIS
1	A	199	ASN
1	A	271	ASN
1	A	384	ASN
1	A	468	ASN
1	A	469	ASN
2	B	45	GLN
2	B	104	ASN
2	B	130	ASN
2	B	225	ASN
2	B	286	ASN
2	B	294	GLN
2	B	363	HIS
2	B	418	ASN

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Mol	Chain	Res	Type
2	B	429	HIS
2	B	518	ASN
1	C	31	HIS
1	C	199	ASN
1	C	252	GLN
1	C	271	ASN
1	C	468	ASN
2	D	104	ASN
2	D	130	ASN
2	D	225	ASN
2	D	286	ASN
2	D	294	GLN
2	D	418	ASN
2	D	429	HIS
2	D	499	ASN
2	D	518	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 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	HCA	A	501	-	4,13,13	0.69	0	3,18,18	2.04	1 (33%)
4	ICS	A	502	1	6,30,30	1.96	3 (50%)	0,78,78	0.00	-
5	CLF	A	503[B]	1,2	0,24,24	0.00	-	0,57,57	0.00	-
3	HCA	C	501	-	4,13,13	1.62	1 (25%)	3,18,18	2.65	2 (66%)
4	ICS	C	502	1	6,30,30	1.97	2 (33%)	0,78,78	0.00	-
5	CLF	C	503[B]	1,2	0,24,24	0.00	-	0,57,57	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
3	HCA	A	501	-	-	0/7/17/17	0/0/0/0
4	ICS	A	502	1	-	0/0/204/204	0/0/13/13
5	CLF	A	503[B]	1,2	-	0/0/132/132	0/12/10/10
3	HCA	C	501	-	-	0/7/17/17	0/0/0/0
4	ICS	C	502	1	-	0/0/204/204	0/0/13/13
5	CLF	C	503[B]	1,2	-	0/0/132/132	0/12/10/10

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	502	ICS	S5A-FE7	-2.85	2.18	2.24
4	A	502	ICS	S2B-FE6	-2.67	2.18	2.24
4	A	502	ICS	S5A-FE7	-2.66	2.18	2.24
4	C	502	ICS	S3A-FE4	-2.54	2.18	2.24
4	A	502	ICS	S2B-FE2	-2.10	2.19	2.24
3	C	501	HCA	O7-C3	3.06	1.47	1.43

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	501	HCA	C3-C2-C1	-3.52	109.67	114.95
3	A	501	HCA	C3-C2-C1	-2.79	110.77	114.95
3	C	501	HCA	C4-C3-C7	-2.40	107.22	111.46

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	501	HCA	1	0
4	A	502	ICS	2	0
3	C	501	HCA	1	0
4	C	502	ICS	1	0
5	C	503[B]	CLF	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

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	477/492 (96%)	-0.32	9 (1%)	66	73	13, 24, 51, 133	13 (2%)
1	C	477/492 (96%)	-0.37	7 (1%)	73	78	12, 23, 49, 107	4 (0%)
2	B	522/523 (99%)	-0.56	1 (0%)	94	96	11, 20, 36, 58	8 (1%)
2	D	522/523 (99%)	-0.53	2 (0%)	92	95	11, 21, 38, 49	3 (0%)
All	All	1998/2030 (98%)	-0.45	19 (0%)	82	86	11, 22, 43, 133	28 (1%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	42	SER	6.6
1	A	39	VAL	6.3
1	C	39	VAL	6.0
1	A	44	LYS	5.5
1	A	37	PRO	5.0
1	C	42	SER	4.7
1	C	40	THR	4.2
1	C	43	LYS	3.3
1	A	40	THR	3.1
1	C	41	GLN	3.0
1	A	43	LYS	2.9
1	A	38	ALA	2.7
2	D	124	VAL	2.7
1	A	41	GLN	2.6
2	D	123	ALA	2.3
1	A	384	ASN	2.2
1	C	7	GLU	2.1
1	C	172	ALA	2.0
2	B	171	LYS	2.0

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 [i](#)

There are no carbohydrates 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, 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(\AA^2)	Q<0.9
3	HCA	C	501	14/14	0.96	0.14	13,14,21,22	0
6	FE	B	602	1/1	0.98	0.07	19,19,19,19	1
5	CLF	A	503[B]	15/15	0.99	0.06	15,17,22,24	0
4	ICS	C	502	18/18	0.99	0.06	14,15,16,16	0
4	ICS	A	502	18/18	0.99	0.08	14,17,20,20	0
5	CLF	C	503[B]	15/15	0.99	0.05	14,18,24,26	0
3	HCA	A	501	14/14	0.99	0.15	15,19,22,22	0
6	FE	B	601	1/1	0.99	0.10	19,19,19,19	1

6.5 Other polymers [i](#)

There are no such residues in this entry.