



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:53 AM GMT

PDB ID : 3K1A
Title : Insights into substrate binding at FeMo-cofactor in nitrogenase from the structure of an alpha-70Ile MoFe protein variant
Authors : Peters, J.W.; Sarma, R.; Barney, B.M.; Keable, S.; Seefeldt, L.C.; Dean, D.R.
Deposited on : 2009-09-26
Resolution : 2.23 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 ⓘ 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

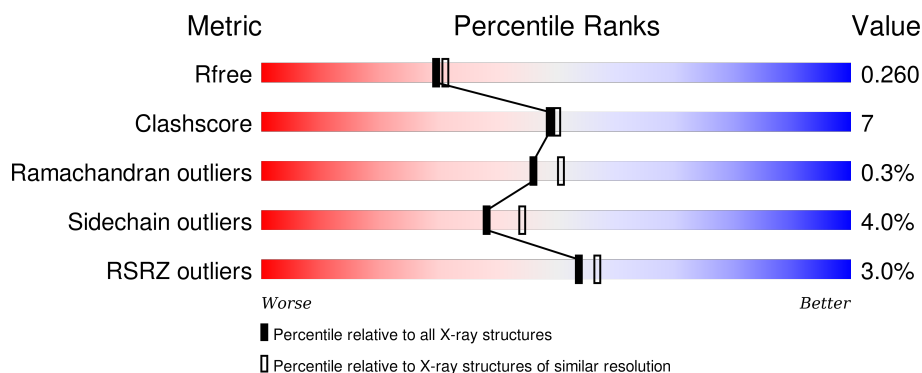
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION




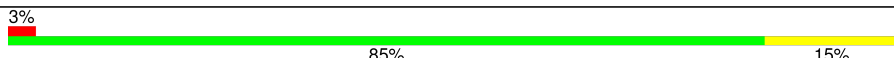
The reported resolution of this entry is 2.23 Å.

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}	91344	1611 (2.26-2.22)
Clashscore	102246	1764 (2.26-2.22)
Ramachandran outliers	100387	1724 (2.26-2.22)
Sidechain outliers	100360	1724 (2.26-2.22)
RSRZ outliers	91569	1616 (2.26-2.22)

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	491	
1	C	491	
2	B	522	
2	D	522	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	CLF	B	6498	-	-	-	X
5	CLF	D	7498	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 16704 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 Nitrogenase molybdenum-iron protein alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	467	Total	C	N	O	S	0	0	0
			3716	2365	633	694	24			
1	C	468	Total	C	N	O	S	0	0	0
			3714	2365	631	694	24			

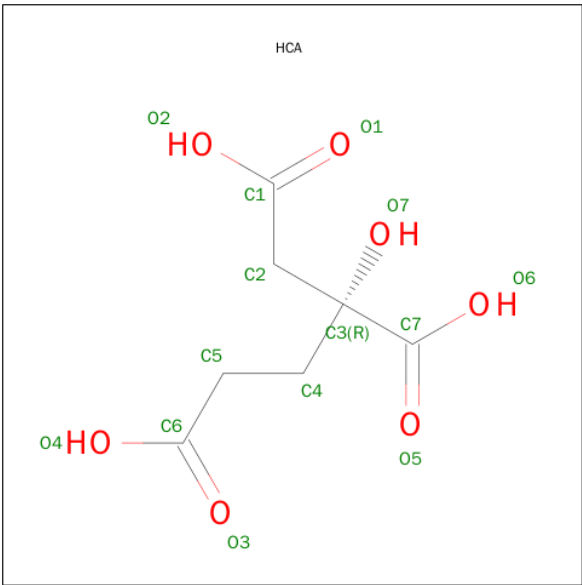
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	70	ILE	VAL	ENGINEERED	UNP P07328
C	70	ILE	VAL	ENGINEERED	UNP P07328

- Molecule 2 is a protein called Nitrogenase molybdenum-iron protein beta chain.

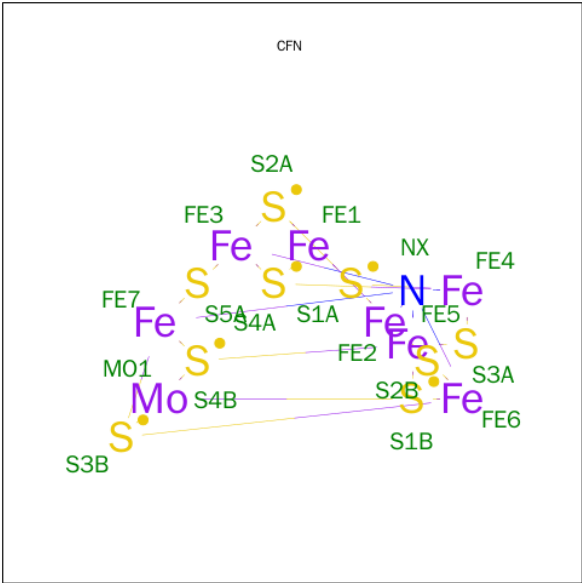
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	522	Total	C	N	O	S	0	0	0
			4174	2666	705	775	28			
2	D	522	Total	C	N	O	S	0	0	0
			4174	2666	705	775	28			

- Molecule 3 is 3-HYDROXY-3-CARBOXY-ADIPIC ACID (three-letter code: HCA) (formula: C₇H₁₀O₇).



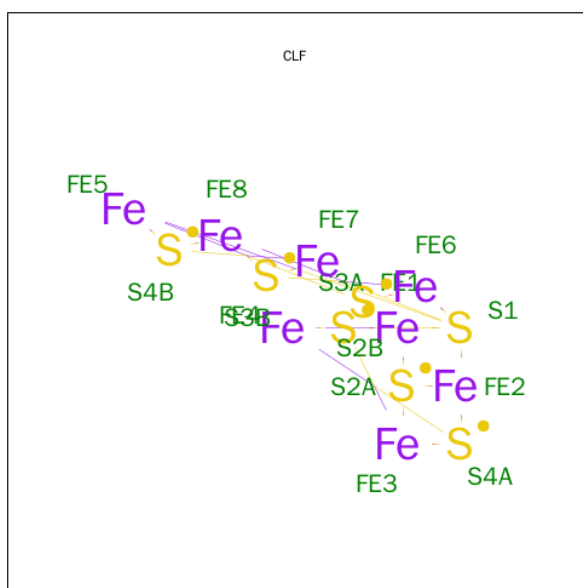
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			14	7	7		
3	C	1	Total	C	O	0	0
			14	7	7		

- Molecule 4 is FE(7)-MO-S(9)-N CLUSTER (three-letter code: CFN) (formula: Fe₇MoNS₉).



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

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	Fe	S	0	0
			15	8	7		
5	D	1	Total	Fe	S	0	0
			15	8	7		

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Ca	0	0
			1	1		
6	D	1	Total	Ca	0	0
			1	1		

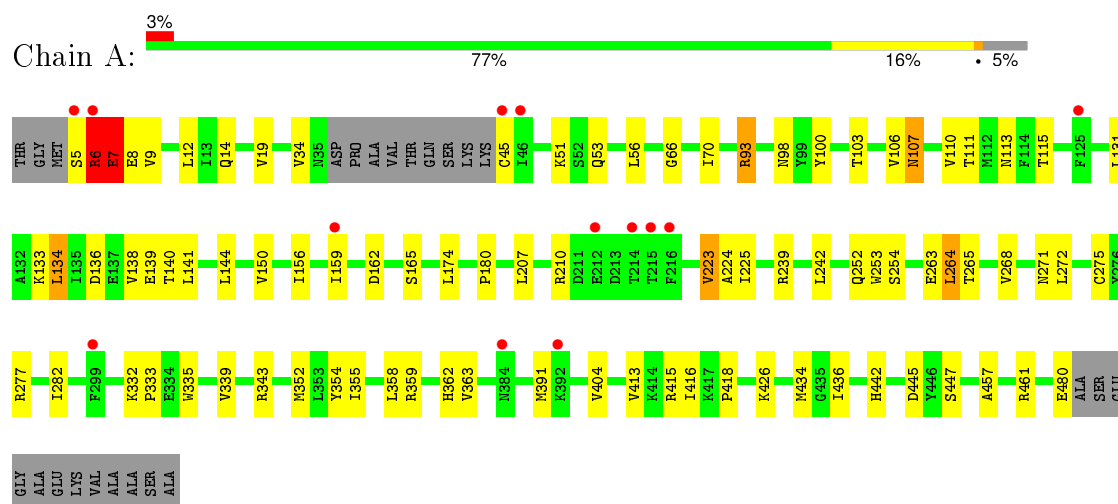
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	166	Total	O	0	0
			166	166		
7	B	255	Total	O	0	0
			255	255		
7	C	175	Total	O	0	0
			175	175		
7	D	234	Total	O	0	0
			234	234		

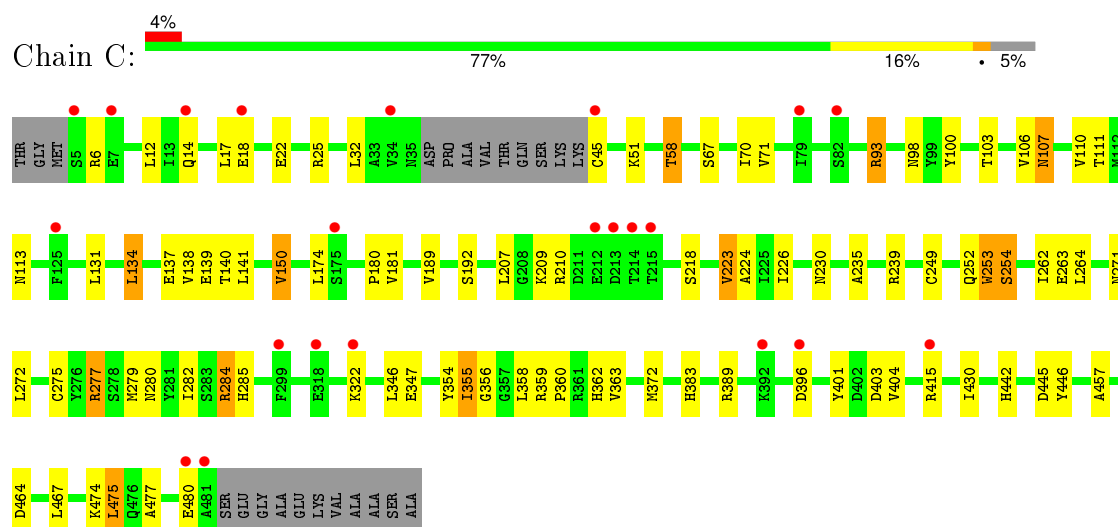
3 Residue-property plots [i](#)

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.

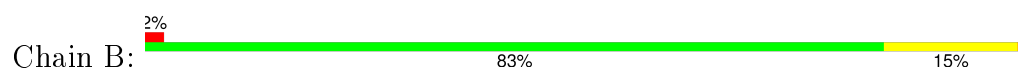
- Molecule 1: Nitrogenase molybdenum-iron protein alpha chain

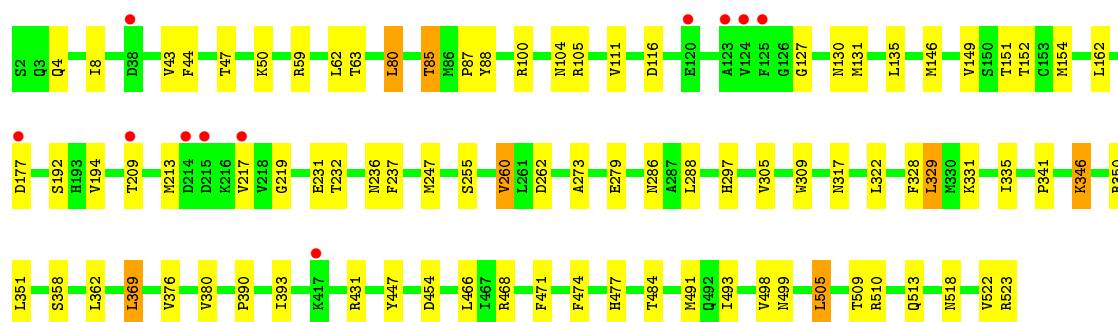


- Molecule 1: Nitrogenase molybdenum-iron protein alpha chain

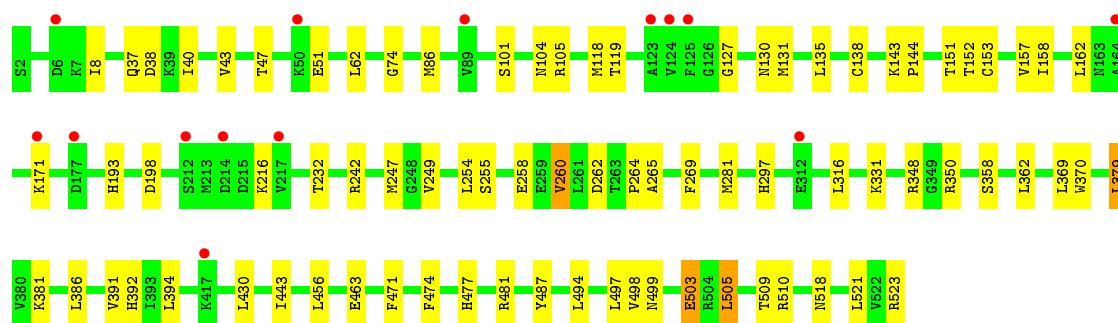
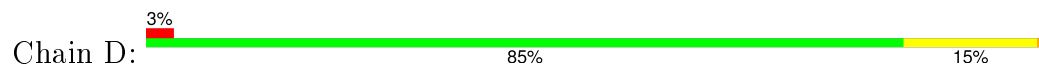


- Molecule 2: Nitrogenase molybdenum-iron protein beta chain





- 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	77.02Å 129.46Å 107.09Å 90.00° 109.01° 90.00°	Depositor
Resolution (Å)	50.00 – 2.23 27.15 – 2.23	Depositor EDS
% Data completeness (in resolution range)	98.5 (50.00-2.23) 98.5 (27.15-2.23)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.40 (at 2.22Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.204 , 0.254 0.209 , 0.260	Depositor DCC
R_{free} test set	4776 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	25.4	Xtriage
Anisotropy	0.141	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 30.4	EDS
Estimated twinning fraction	0.018 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 95443 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	16704	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.65% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: CFN, CLF, HCA, CA

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	0.33	0/3802	0.55	0/5125
1	C	0.34	0/3800	0.56	0/5124
2	B	0.33	0/4280	0.53	0/5786
2	D	0.33	0/4280	0.52	0/5786
All	All	0.33	0/16162	0.54	0/21821

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	C	0	1
All	All	0	3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	5	SER	Peptide
1	A	6	ARG	Peptide
1	C	253	TRP	Peptide

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	3716	0	3650	59	0
1	C	3714	0	3641	60	0
2	B	4174	0	4089	73	0
2	D	4174	0	4089	59	0
3	A	14	0	6	0	0
3	C	14	0	6	0	0
4	A	18	0	0	2	0
4	C	18	0	0	3	0
5	B	15	0	0	1	0
5	D	15	0	0	2	0
6	B	1	0	0	0	0
6	D	1	0	0	0	0
7	A	166	0	0	4	0
7	B	255	0	0	11	0
7	C	175	0	0	6	0
7	D	234	0	0	3	0
All	All	16704	0	15481	225	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 7.

All (225) 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:447:TYR:CE1	7:B:779:HOH:O	1.82	1.33
7:B:779:HOH:O	2:D:521:LEU:HD21	1.14	1.25
7:B:779:HOH:O	2:D:521:LEU:CD2	1.81	0.96
2:D:499:ASN:O	2:D:503:GLU:HG3	1.73	0.88
1:C:239:ARG:HE	1:C:252:GLN:HE21	1.19	0.86
1:A:70:ILE:HD13	4:A:6496:CFN:S2B	2.17	0.85
2:D:131:MET:HE2	2:D:135:LEU:HD11	1.58	0.85
1:C:70:ILE:HD13	4:C:7496:CFN:S2B	2.16	0.84
2:D:499:ASN:O	2:D:503:GLU:CG	2.27	0.82
1:C:457:ALA:HB1	2:D:8:ILE:HD12	1.61	0.82
2:B:513:GLN:HE22	2:D:37:GLN:HE22	1.30	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:499:ASN:HD21	2:D:477:HIS:H	1.29	0.79
1:C:415:ARG:HD2	7:C:692:HOH:O	1.82	0.79
2:B:477:HIS:H	2:D:499:ASN:HD21	1.30	0.78
2:B:131:MET:HE2	2:B:135:LEU:HD11	1.66	0.78
2:B:4:GLN:HG3	7:B:600:HOH:O	1.84	0.77
2:B:209:THR:HG21	2:B:309:TRP:NE1	2.00	0.77
1:A:239:ARG:HE	1:A:252:GLN:HE21	1.31	0.76
1:A:6:ARG:HB3	1:A:7:GLU:HB2	1.67	0.75
2:B:80:LEU:HD13	2:B:87:PRO:HG3	1.67	0.75
2:B:209:THR:HG21	2:B:309:TRP:HE1	1.52	0.74
2:D:151:THR:HG23	2:D:162:LEU:HD11	1.69	0.74
1:A:354:TYR:CZ	1:A:404:VAL:HG12	2.28	0.69
2:B:151:THR:HG23	2:B:162:LEU:HD11	1.75	0.68
1:A:144:LEU:HD13	2:B:43:VAL:HG21	1.77	0.67
1:A:70:ILE:CD1	4:A:6496:CFN:S2B	2.82	0.67
1:C:239:ARG:HE	1:C:252:GLN:NE2	1.92	0.66
1:C:93:ARG:HG3	1:C:113:ASN:HB2	1.78	0.66
1:A:156:ILE:O	1:A:159:ILE:HG22	1.96	0.64
1:A:45:CYS:N	7:A:828:HOH:O	2.30	0.64
2:B:100:ARG:HD2	2:B:111:VAL:O	1.97	0.64
1:A:457:ALA:HB1	2:B:8:ILE:HD12	1.80	0.64
2:D:509:THR:HG21	2:D:518:ASN:HD22	1.63	0.63
1:A:352:MET:HE1	1:A:416:ILE:HB	1.79	0.62
1:C:51:LYS:HD3	1:C:189:VAL:HG12	1.81	0.62
1:C:354:TYR:CZ	1:C:404:VAL:HG12	2.34	0.62
2:B:43:VAL:O	2:B:47:THR:HG23	1.99	0.61
2:B:484:THR:HG22	2:B:493:ILE:HD11	1.83	0.61
2:D:499:ASN:O	2:D:503:GLU:HG2	2.01	0.61
2:D:379:LEU:HD11	2:D:443:ILE:HG21	1.84	0.59
2:B:509:THR:HG21	2:B:518:ASN:HD22	1.66	0.59
2:B:232:THR:HG21	2:B:471:PHE:CD1	2.37	0.59
1:A:66:GLY:HA2	1:A:70:ILE:HD12	1.85	0.59
2:D:316:LEU:HD11	2:D:331:LYS:HG2	1.85	0.59
1:C:70:ILE:CD1	4:C:7496:CFN:S2B	2.91	0.59
1:C:275:CYS:HA	1:C:358:LEU:HD22	1.85	0.59
1:A:103:THR:H	1:A:107:ASN:HD21	1.50	0.58
1:C:224:ALA:HB3	1:C:271:ASN:HD22	1.66	0.58
5:D:7498:CLF:S4B	5:D:7498:CLF:S3A	3.01	0.58
1:C:210:ARG:HD3	1:C:263:GLU:HB3	1.84	0.58
1:C:106:VAL:HG21	2:D:40:ILE:HG23	1.85	0.58
2:B:505:LEU:HD13	2:B:523:ARG:CZ	2.35	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:207:LEU:HD22	1:C:282:ILE:HD11	1.87	0.57
1:C:230:ASN:HD22	1:C:235:ALA:H	1.52	0.56
1:C:51:LYS:HD2	2:D:119:THR:HG21	1.89	0.55
1:A:224:ALA:HB3	1:A:271:ASN:HD22	1.71	0.55
2:D:151:THR:CG2	2:D:162:LEU:HD11	2.37	0.55
2:B:447:TYR:CD1	7:B:779:HOH:O	2.29	0.55
1:A:277:ARG:HD3	7:A:582:HOH:O	2.06	0.55
1:A:93:ARG:HD2	1:A:111:THR:O	2.06	0.55
1:A:275:CYS:HA	1:A:358:LEU:HD22	1.88	0.55
1:C:140:THR:HG23	1:C:141:LEU:HG	1.89	0.54
1:C:137:GLU:HA	1:C:140:THR:HG22	1.90	0.54
1:A:150:VAL:HG13	1:A:180:PRO:HA	1.89	0.54
1:A:134:LEU:HD13	2:B:62:LEU:HD13	1.89	0.54
1:C:253:TRP:CZ2	1:C:262:ILE:HG23	2.42	0.53
1:C:139:GLU:HG3	1:C:174:LEU:HD13	1.89	0.53
1:C:134:LEU:HD13	2:D:62:LEU:HD13	1.89	0.53
2:D:118:MET:HE1	2:D:158:ILE:HD11	1.89	0.53
2:D:232:THR:HG21	2:D:471:PHE:CD1	2.43	0.53
1:C:359:ARG:O	1:C:363:VAL:HG22	2.08	0.53
2:D:394:LEU:HD13	2:D:430:LEU:HB2	1.91	0.53
1:C:346:LEU:HD21	1:C:464:ASP:HA	1.92	0.52
1:A:207:LEU:HD22	1:A:282:ILE:HD11	1.91	0.52
2:D:260:VAL:HG22	7:D:539:HOH:O	2.10	0.52
2:D:43:VAL:O	2:D:47:THR:HG23	2.09	0.51
1:A:6:ARG:HG2	1:A:7:GLU:H	1.75	0.51
1:C:67:SER:OG	1:C:181:VAL:HG11	2.10	0.51
1:C:106:VAL:HG23	1:C:107:ASN:N	2.25	0.51
1:A:100:TYR:CE1	1:A:110:VAL:HB	2.45	0.51
1:C:106:VAL:CG2	1:C:107:ASN:N	2.74	0.50
2:B:358:SER:HB3	2:B:498:VAL:HG21	1.93	0.50
2:B:362:LEU:HG	2:B:498:VAL:HG22	1.93	0.50
2:B:260:VAL:HG22	7:B:532:HOH:O	2.11	0.50
2:B:351:LEU:HD13	2:B:491:MET:HG3	1.94	0.50
1:A:352:MET:HE3	1:A:418:PRO:HG3	1.93	0.50
1:A:93:ARG:HD3	1:A:113:ASN:HB2	1.93	0.50
1:C:17:LEU:HD21	1:C:32:LEU:CD1	2.42	0.50
1:A:141:LEU:CD1	2:B:59:ARG:HD2	2.41	0.50
1:C:22:GLU:CD	1:C:25:ARG:HH12	2.15	0.49
2:B:105:ARG:HB3	2:B:474:PHE:CD1	2.47	0.49
2:B:127:GLY:HA2	2:B:130:ASN:HD22	1.77	0.49
2:B:131:MET:HE1	2:B:149:VAL:HG21	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:ARG:NH2	1:A:264:LEU:HD11	2.27	0.49
2:D:362:LEU:HG	2:D:498:VAL:HG22	1.94	0.49
2:B:493:ILE:HD13	7:B:616:HOH:O	2.12	0.49
1:A:359:ARG:O	1:A:363:VAL:HG22	2.11	0.49
2:D:131:MET:HE2	2:D:135:LEU:CD1	2.38	0.48
2:D:358:SER:HB3	2:D:498:VAL:HG21	1.93	0.48
2:B:322:LEU:HD21	1:C:474:LYS:HB3	1.96	0.48
1:C:477:ALA:HB3	1:C:480:GLU:HB2	1.95	0.48
1:A:103:THR:H	1:A:107:ASN:ND2	2.10	0.48
1:C:192:SER:OG	1:C:383:HIS:HE1	1.96	0.48
2:D:105:ARG:HB3	2:D:474:PHE:CD1	2.48	0.48
1:C:93:ARG:HD3	1:C:111:THR:O	2.14	0.48
2:D:348:ARG:HG3	2:D:487:TYR:CE1	2.49	0.48
2:B:85:THR:HG21	2:B:146:MET:HE1	1.94	0.48
1:A:139:GLU:HG3	1:A:174:LEU:HD13	1.96	0.47
1:A:115:THR:HG23	2:B:63:THR:HB	1.96	0.47
1:A:223:VAL:HG22	1:A:272:LEU:HG	1.96	0.47
1:A:335:TRP:O	1:A:339:VAL:HG12	2.15	0.47
1:A:224:ALA:HB3	1:A:271:ASN:ND2	2.30	0.47
1:A:6:ARG:O	1:A:8:GLU:N	2.48	0.47
1:A:339:VAL:HG23	1:A:343:ARG:HB2	1.97	0.47
2:B:192:SER:OG	2:B:194:VAL:HG22	2.15	0.47
1:C:134:LEU:HD12	1:C:134:LEU:C	2.35	0.46
1:A:210:ARG:NH2	1:A:264:LEU:CD1	2.78	0.46
1:A:447:SER:HA	7:A:859:HOH:O	2.14	0.46
1:A:352:MET:CE	1:A:418:PRO:HD3	2.45	0.46
1:A:207:LEU:CD2	1:A:282:ILE:HD11	2.46	0.46
2:D:86:MET:HG2	2:D:138:CYS:SG	2.56	0.46
2:B:447:TYR:HE1	7:B:779:HOH:O	1.52	0.46
2:B:262:ASP:O	2:D:350:ARG:HD3	2.16	0.46
2:B:131:MET:HE2	2:B:135:LEU:CD1	2.41	0.46
1:A:150:VAL:CG1	1:A:180:PRO:HA	2.45	0.46
1:C:442:HIS:HB2	7:C:500:HOH:O	2.16	0.46
1:C:209:LYS:HE2	7:C:774:HOH:O	2.15	0.46
2:B:317:ASN:ND2	7:B:726:HOH:O	2.41	0.46
1:A:6:ARG:O	1:A:9:VAL:N	2.25	0.46
2:B:151:THR:CG2	2:B:162:LEU:HD11	2.45	0.46
1:C:150:VAL:HG13	1:C:180:PRO:HA	1.98	0.46
1:C:280:ASN:O	1:C:284:ARG:HD3	2.16	0.46
2:B:305:VAL:O	2:B:309:TRP:HB2	2.16	0.45
1:A:210:ARG:HG3	1:A:263:GLU:HB3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:247:MET:CB	2:D:249:VAL:HG23	2.45	0.45
2:D:505:LEU:HD13	2:D:523:ARG:CZ	2.46	0.45
2:D:456:LEU:HD13	2:D:463:GLU:OE1	2.17	0.45
2:B:431:ARG:HD2	2:B:454:ASP:OD2	2.17	0.45
2:B:85:THR:HG21	2:B:146:MET:CE	2.46	0.45
2:D:523:ARG:HD3	7:D:684:HOH:O	2.16	0.45
1:C:475:LEU:HD13	2:D:265:ALA:O	2.16	0.45
1:A:354:TYR:CE1	1:A:404:VAL:HG12	2.51	0.45
1:C:22:GLU:CD	1:C:25:ARG:NH1	2.70	0.45
1:C:71:VAL:O	1:C:254:SER:OG	2.30	0.45
1:C:430:ILE:HG23	2:D:269:PHE:CG	2.51	0.45
1:A:134:LEU:HD12	1:A:134:LEU:C	2.36	0.45
1:A:144:LEU:CD1	2:B:43:VAL:HG21	2.45	0.45
2:B:390:PRO:HB2	2:B:393:ILE:HD11	1.97	0.45
1:C:223:VAL:HG12	1:C:249:CYS:HA	1.99	0.45
2:B:346:LYS:CB	2:B:346:LYS:HZ2	2.29	0.45
1:C:17:LEU:O	1:C:25:ARG:HG3	2.18	0.44
2:B:152:THR:HB	5:B:6498:CLF:S3B	2.57	0.44
2:D:443:ILE:HD11	2:D:497:LEU:HD21	1.99	0.44
1:C:106:VAL:CG2	2:D:40:ILE:HG23	2.47	0.44
2:D:316:LEU:HD11	2:D:331:LYS:CG	2.48	0.44
1:C:103:THR:HB	1:C:106:VAL:HG22	1.99	0.44
2:B:231:GLU:HB3	2:B:237:PHE:CZ	2.53	0.44
2:B:247:MET:HG2	2:B:341:PRO:HD3	1.98	0.44
2:B:209:THR:HG23	2:B:213:MET:SD	2.58	0.44
2:B:491:MET:HE1	2:D:481:ARG:NH1	2.32	0.44
2:B:194:VAL:HB	2:B:297:HIS:CG	2.53	0.44
2:D:509:THR:HG21	2:D:518:ASN:ND2	2.32	0.44
2:D:391:VAL:HG12	2:D:392:HIS:CE1	2.53	0.43
2:B:346:LYS:CA	2:B:346:LYS:HZ2	2.31	0.43
1:A:426:LYS:HA	2:B:104:ASN:ND2	2.33	0.43
1:A:6:ARG:HG2	1:A:7:GLU:N	2.33	0.43
2:D:254:LEU:HD22	2:D:281:MET:SD	2.59	0.43
1:A:9:VAL:CG1	1:A:34:VAL:HG22	2.49	0.43
2:D:242:ARG:NH1	7:D:731:HOH:O	2.51	0.43
1:C:285:HIS:HE1	7:C:558:HOH:O	2.02	0.43
1:C:230:ASN:HD22	1:C:235:ALA:N	2.15	0.43
2:B:219:GLY:CA	2:B:288:LEU:HD23	2.49	0.43
2:B:466:LEU:HD23	2:B:468:ARG:NH1	2.33	0.42
1:A:12:LEU:HD13	1:A:415:ARG:HG2	2.00	0.42
1:A:434:MET:HB2	1:A:436:ILE:HD12	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:198:ASP:HB2	2:D:297:HIS:O	2.20	0.42
1:C:356:GLY:HA3	4:C:7496:CFN:S1B	2.60	0.42
1:C:224:ALA:HB3	1:C:271:ASN:ND2	2.33	0.42
1:C:359:ARG:N	1:C:360:PRO:CD	2.82	0.42
2:B:369:LEU:C	2:B:369:LEU:HD22	2.39	0.42
2:B:509:THR:HG21	2:B:518:ASN:ND2	2.34	0.42
2:D:153:CYS:O	2:D:157:VAL:HG23	2.19	0.42
1:C:6:ARG:CB	7:C:524:HOH:O	2.68	0.42
2:D:101:SER:HA	2:D:104:ASN:HD22	1.84	0.42
2:D:386:LEU:HD21	2:D:494:LEU:HD23	2.00	0.42
1:A:352:MET:HE1	1:A:413:VAL:HA	2.01	0.42
2:B:260:VAL:HG13	2:B:273:ALA:H	1.85	0.42
1:C:277:ARG:HD3	7:C:548:HOH:O	2.20	0.42
2:D:370:TRP:CE3	2:D:430:LEU:HD21	2.55	0.42
1:A:53:GLN:HB2	1:A:56:LEU:HD12	2.02	0.42
2:B:88:TYR:OH	2:B:116:ASP:HB3	2.20	0.42
1:C:58:THR:CG2	1:C:403:ASP:OD1	2.68	0.42
1:A:134:LEU:O	1:A:138:VAL:HG23	2.19	0.41
2:B:219:GLY:HA2	2:B:288:LEU:HD23	2.01	0.41
2:B:328:PHE:HD2	2:B:329:LEU:HD13	1.85	0.41
2:B:376:VAL:O	2:B:380:VAL:HG23	2.21	0.41
1:A:253:TRP:HA	1:A:254:SER:HA	1.84	0.41
2:B:331:LYS:O	2:B:335:ILE:HG12	2.20	0.41
2:B:177:ASP:HB3	7:B:564:HOH:O	2.19	0.41
2:B:510:ARG:HG2	2:D:456:LEU:HD23	2.03	0.41
1:C:100:TYR:CE1	1:C:110:VAL:HB	2.56	0.41
2:B:350:ARG:HD3	2:D:262:ASP:O	2.20	0.41
2:B:369:LEU:CD2	2:B:376:VAL:HG13	2.51	0.41
1:A:136:ASP:O	1:A:140:THR:HG23	2.20	0.41
1:C:134:LEU:O	1:C:138:VAL:HG23	2.21	0.41
2:B:231:GLU:CD	2:B:236:ASN:HD22	2.24	0.41
7:B:552:HOH:O	2:D:510:ARG:HD3	2.20	0.41
1:A:332:LYS:N	1:A:333:PRO:HD2	2.36	0.41
2:B:522:VAL:HG12	1:C:446:TYR:CE1	2.56	0.41
1:C:226:ILE:HG22	1:C:279:MET:HB3	2.02	0.41
1:C:138:VAL:HG23	2:D:62:LEU:HD22	2.03	0.41
1:A:265:THR:O	1:A:268:VAL:HG22	2.20	0.41
2:B:217:VAL:H	2:B:286:ASN:ND2	2.19	0.40
2:B:513:GLN:NE2	2:D:37:GLN:HE22	2.07	0.40
2:B:346:LYS:HE3	2:D:264:PRO:HG3	2.02	0.40
2:B:369:LEU:O	2:B:369:LEU:HD22	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:152:THR:HB	5:D:7498:CLF:S3B	2.60	0.40
1:C:223:VAL:HG22	1:C:272:LEU:HG	2.02	0.40
2:D:143:LYS:N	2:D:144:PRO:HD3	2.35	0.40
1:A:442:HIS:HB2	7:A:516:HOH:O	2.22	0.40
2:D:127:GLY:HA2	2:D:130:ASN:HD22	1.86	0.40
1:C:346:LEU:HD13	1:C:372:MET:CE	2.51	0.40
1:A:225:ILE:HD11	1:A:242:LEU:HD12	2.03	0.40
2:D:74:GLY:HA3	2:D:193:HIS:O	2.21	0.40
1:A:106:VAL:HG21	2:B:44:PHE:HB2	2.03	0.40

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	463/491 (94%)	441 (95%)	20 (4%)	2 (0%)	39	42
1	C	464/491 (94%)	440 (95%)	23 (5%)	1 (0%)	52	60
2	B	520/522 (100%)	509 (98%)	10 (2%)	1 (0%)	52	60
2	D	520/522 (100%)	508 (98%)	11 (2%)	1 (0%)	52	60
All	All	1967/2026 (97%)	1898 (96%)	64 (3%)	5 (0%)	46	51

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	6	ARG
1	A	7	GLU
2	B	255	SER
2	D	255	SER
1	C	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	398/414 (96%)	378 (95%)	20 (5%)	30	31
1	C	396/414 (96%)	369 (93%)	27 (7%)	20	17
2	B	454/454 (100%)	444 (98%)	10 (2%)	60	70
2	D	454/454 (100%)	443 (98%)	11 (2%)	57	67
All	All	1702/1736 (98%)	1634 (96%)	68 (4%)	38	44

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	GLU
1	A	14	GLN
1	A	19	VAL
1	A	51	LYS
1	A	93	ARG
1	A	98	ASN
1	A	107	ASN
1	A	131	LEU
1	A	133	LYS
1	A	134	LEU
1	A	162	ASP
1	A	165	SER
1	A	223	VAL
1	A	264	LEU
1	A	355	ILE
1	A	362	HIS
1	A	391	MET
1	A	445	ASP
1	A	461	ARG
1	A	480	GLU
2	B	50	LYS
2	B	80	LEU
2	B	85	THR
2	B	154	MET

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Mol	Chain	Res	Type
2	B	260	VAL
2	B	279	GLU
2	B	329	LEU
2	B	346	LYS
2	B	369	LEU
2	B	505	LEU
1	C	12	LEU
1	C	14	GLN
1	C	18	GLU
1	C	45	CYS
1	C	58	THR
1	C	93	ARG
1	C	98	ASN
1	C	107	ASN
1	C	131	LEU
1	C	134	LEU
1	C	150	VAL
1	C	218	SER
1	C	223	VAL
1	C	254	SER
1	C	264	LEU
1	C	277	ARG
1	C	284	ARG
1	C	322	LYS
1	C	347	GLU
1	C	355	ILE
1	C	362	HIS
1	C	389	ARG
1	C	396	ASP
1	C	401	TYR
1	C	445	ASP
1	C	467	LEU
1	C	475	LEU
2	D	38	ASP
2	D	51	GLU
2	D	171	LYS
2	D	216	LYS
2	D	258	GLU
2	D	260	VAL
2	D	369	LEU
2	D	379	LEU
2	D	381	LYS

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Mol	Chain	Res	Type
2	D	503	GLU
2	D	505	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	GLN
1	A	49	ASN
1	A	53	GLN
1	A	107	ASN
1	A	196	HIS
1	A	252	GLN
1	A	271	ASN
1	A	285	HIS
1	A	362	HIS
1	A	383	HIS
2	B	37	GLN
2	B	104	ASN
2	B	128	GLN
2	B	130	ASN
2	B	168	ASN
2	B	268	GLN
2	B	286	ASN
2	B	317	ASN
2	B	457	HIS
2	B	499	ASN
2	B	518	ASN
2	B	519	HIS
1	C	107	ASN
1	C	230	ASN
1	C	252	GLN
1	C	271	ASN
1	C	285	HIS
1	C	362	HIS
1	C	383	HIS
2	D	37	GLN
2	D	58	GLN
2	D	104	ASN
2	D	128	GLN
2	D	129	GLN
2	D	130	ASN
2	D	167	ASN

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Mol	Chain	Res	Type
2	D	168	ASN
2	D	268	GLN
2	D	286	ASN
2	D	499	ASN
2	D	513	GLN
2	D	518	ASN
2	D	519	HIS

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	494	-	4,13,13	0.83	0	3,18,18	0.57	0
4	CFN	A	6496	1	6,30,30	0.29	0	0,78,78	0.00	-
5	CLF	B	6498	1,2	0,24,24	0.00	-	0,57,57	0.00	-
3	HCA	C	494	-	4,13,13	0.94	0	3,18,18	1.19	0
4	CFN	C	7496	1	6,30,30	0.24	0	0,78,78	0.00	-
5	CLF	D	7498	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	494	-	-	0/7/17/17	0/0/0/0
4	CFN	A	6496	1	-	0/0/204/204	0/0/13/13
5	CLF	B	6498	1,2	-	0/0/132/132	0/12/10/10
3	HCA	C	494	-	-	0/7/17/17	0/0/0/0
4	CFN	C	7496	1	-	0/0/204/204	0/0/13/13
5	CLF	D	7498	1,2	-	0/0/132/132	0/12/10/10

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	6496	CFN	2	0
5	B	6498	CLF	1	0
4	C	7496	CFN	3	0
5	D	7498	CLF	2	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	467/491 (95%)	0.27	13 (2%) 56 59	25, 37, 53, 73	0
1	C	468/491 (95%)	0.37	22 (4%) 35 38	27, 37, 55, 73	0
2	B	522/522 (100%)	0.04	11 (2%) 67 69	27, 35, 49, 63	0
2	D	522/522 (100%)	0.12	14 (2%) 58 60	27, 37, 50, 67	0
All	All	1979/2026 (97%)	0.19	60 (3%) 54 57	25, 36, 52, 73	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	125	PHE	5.5
1	A	212	GLU	5.0
2	B	214	ASP	5.0
1	C	212	GLU	4.7
1	C	481	ALA	4.6
2	D	123	ALA	4.3
1	A	6	ARG	4.2
1	C	215	THR	3.9
1	C	396	ASP	3.8
1	A	125	PHE	3.7
1	C	415	ARG	3.6
1	C	392	LYS	3.5
2	B	38	ASP	3.4
1	A	216	PHE	3.3
2	D	217	VAL	3.2
1	A	159	ILE	3.2
2	B	123	ALA	3.2
2	D	214	ASP	3.2
2	D	124	VAL	3.1
1	A	215	THR	3.1
1	C	5	SER	3.1

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Mol	Chain	Res	Type	RSRZ
2	B	125	PHE	3.0
1	A	45	CYS	3.0
2	B	177	ASP	3.0
1	C	14	GLN	3.0
1	C	480	GLU	3.0
2	B	124	VAL	3.0
1	A	384	ASN	2.9
1	C	79	ILE	2.9
1	A	392	LYS	2.9
2	B	215	ASP	2.8
1	C	213	ASP	2.8
1	A	299	PHE	2.8
1	C	7	GLU	2.7
1	C	214	THR	2.7
1	C	175	SER	2.6
2	D	212	SER	2.6
2	D	177	ASP	2.6
1	C	18	GLU	2.5
1	A	214	THR	2.5
1	C	125	PHE	2.5
1	C	299	PHE	2.4
2	D	50	LYS	2.4
1	C	45	CYS	2.4
2	B	417	LYS	2.4
1	C	34	VAL	2.3
2	B	209	THR	2.3
1	C	318	GLU	2.3
2	B	217	VAL	2.3
1	C	82	SER	2.2
1	A	5	SER	2.2
2	D	164	ALA	2.2
2	D	171	LYS	2.1
2	D	417	LYS	2.1
2	D	312	GLU	2.1
2	D	89	VAL	2.1
2	D	6	ASP	2.1
1	C	322	LYS	2.0
1	A	46	ILE	2.0
2	B	120	GLU	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. 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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
5	CLF	B	6498	15/15	0.62	0.23	5.08	21,31,40,46	2
5	CLF	D	7498	15/15	0.64	0.22	3.71	20,28,44,54	0
3	HCA	A	494	14/14	0.96	0.13	-0.56	27,30,34,37	0
3	HCA	C	494	14/14	0.97	0.12	-0.84	30,33,36,39	0
4	CFN	C	7496	18/18	0.94	0.09	-1.83	18,25,32,38	0
4	CFN	A	6496	18/18	0.94	0.09	-2.16	22,27,33,42	0
6	CA	D	524	1/1	0.87	0.08	-2.44	61,61,61,61	0
6	CA	B	524	1/1	0.85	0.07	-2.76	57,57,57,57	0

6.5 Other polymers [i](#)

There are no such residues in this entry.