



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 09:03 PM UTC

PDB ID : 3VMG / pdb_00003vmg
Title : Reduced carbazole-bound complex between oxygenase and ferredoxin in carbazole 1,9a-dioxygenase
Authors : Ashikawa, Y.; Nojiri, H.
Deposited on : 2011-12-12
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:

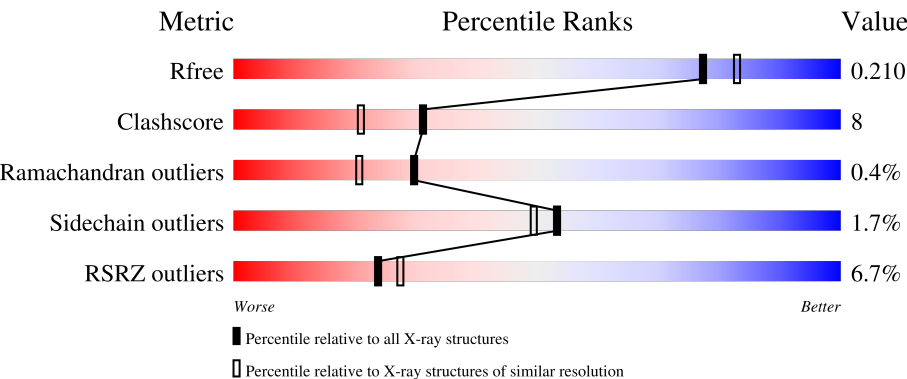
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

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 $\leq 5\%$. 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	392	5% 82% 16% ..
1	B	392	7% 78% 19% ..
1	C	392	4% 79% 20% ..
2	D	115	15% 76% 15% 10%
2	E	115	9% 71% 21% 7%

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Mol	Chain	Length	Quality of chain
2	F	115	

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	9CA	C	2002	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 12864 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 Terminal oxygenase component of carbazole.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	389	Total	C	N	O	S	0	0	0
			3136	2004	535	583	14			
1	B	389	Total	C	N	O	S	0	0	0
			3136	2004	535	583	14			
1	C	389	Total	C	N	O	S	0	0	0
			3136	2004	535	583	14			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	385	LEU	-	expression tag	UNP Q84II6
A	386	GLU	-	expression tag	UNP Q84II6
A	387	HIS	-	expression tag	UNP Q84II6
A	388	HIS	-	expression tag	UNP Q84II6
A	389	HIS	-	expression tag	UNP Q84II6
A	390	HIS	-	expression tag	UNP Q84II6
A	391	HIS	-	expression tag	UNP Q84II6
A	392	HIS	-	expression tag	UNP Q84II6
B	385	LEU	-	expression tag	UNP Q84II6
B	386	GLU	-	expression tag	UNP Q84II6
B	387	HIS	-	expression tag	UNP Q84II6
B	388	HIS	-	expression tag	UNP Q84II6
B	389	HIS	-	expression tag	UNP Q84II6
B	390	HIS	-	expression tag	UNP Q84II6
B	391	HIS	-	expression tag	UNP Q84II6
B	392	HIS	-	expression tag	UNP Q84II6
C	385	LEU	-	expression tag	UNP Q84II6
C	386	GLU	-	expression tag	UNP Q84II6
C	387	HIS	-	expression tag	UNP Q84II6
C	388	HIS	-	expression tag	UNP Q84II6
C	389	HIS	-	expression tag	UNP Q84II6
C	390	HIS	-	expression tag	UNP Q84II6
C	391	HIS	-	expression tag	UNP Q84II6

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Chain	Residue	Modelled	Actual	Comment	Reference
C	392	HIS	-	expression tag	UNP Q84II6

- Molecule 2 is a protein called Ferredoxin component of carbazole.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	104	Total	C	N	O	S	0	0	0
			768	483	129	149	7			
2	E	107	Total	C	N	O	S	0	0	0
			794	499	133	155	7			
2	F	104	Total	C	N	O	S	0	0	0
			768	483	129	149	7			

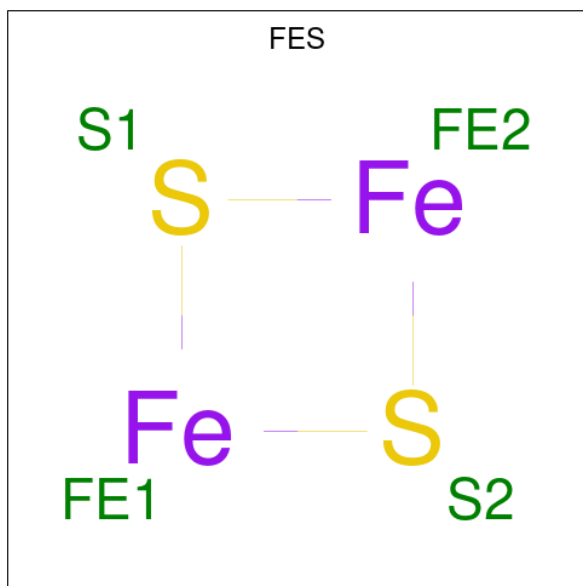
There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	108	LEU	-	expression tag	UNP Q8GI16
D	109	GLU	-	expression tag	UNP Q8GI16
D	110	HIS	-	expression tag	UNP Q8GI16
D	111	HIS	-	expression tag	UNP Q8GI16
D	112	HIS	-	expression tag	UNP Q8GI16
D	113	HIS	-	expression tag	UNP Q8GI16
D	114	HIS	-	expression tag	UNP Q8GI16
D	115	HIS	-	expression tag	UNP Q8GI16
E	108	LEU	-	expression tag	UNP Q8GI16
E	109	GLU	-	expression tag	UNP Q8GI16
E	110	HIS	-	expression tag	UNP Q8GI16
E	111	HIS	-	expression tag	UNP Q8GI16
E	112	HIS	-	expression tag	UNP Q8GI16
E	113	HIS	-	expression tag	UNP Q8GI16
E	114	HIS	-	expression tag	UNP Q8GI16
E	115	HIS	-	expression tag	UNP Q8GI16
F	108	LEU	-	expression tag	UNP Q8GI16
F	109	GLU	-	expression tag	UNP Q8GI16
F	110	HIS	-	expression tag	UNP Q8GI16
F	111	HIS	-	expression tag	UNP Q8GI16
F	112	HIS	-	expression tag	UNP Q8GI16
F	113	HIS	-	expression tag	UNP Q8GI16
F	114	HIS	-	expression tag	UNP Q8GI16
F	115	HIS	-	expression tag	UNP Q8GI16

- Molecule 3 is FE (II) ION (CCD ID: FE2) (formula: Fe).

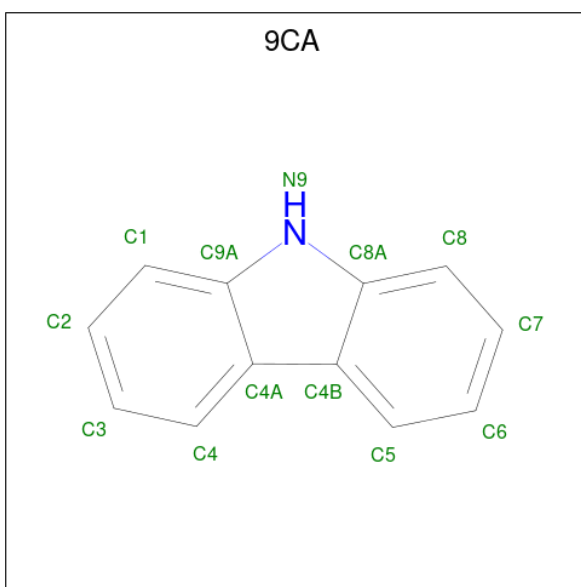
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Fe 1 1	0	0
3	B	1	Total Fe 1 1	0	0
3	C	1	Total Fe 1 1	0	0

- Molecule 4 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Fe S 4 2 2	0	0
4	B	1	Total Fe S 4 2 2	0	0
4	C	1	Total Fe S 4 2 2	0	0
4	D	1	Total Fe S 4 2 2	0	0
4	E	1	Total Fe S 4 2 2	0	0
4	F	1	Total Fe S 4 2 2	0	0

- Molecule 5 is 9H-CARBAZOLE (CCD ID: 9CA) (formula: $\text{C}_{12}\text{H}_9\text{N}$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	N	0	0
			13	12	1		
5	C	1	Total	C	N	0	0
			13	12	1		

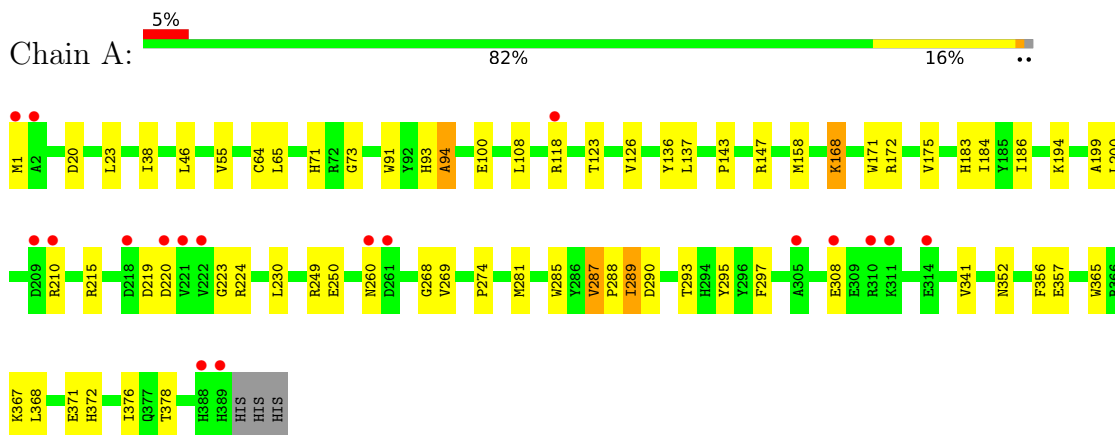
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	339	Total	O	0	0
			339	339		
6	B	291	Total	O	0	0
			291	291		
6	C	299	Total	O	0	0
			299	299		
6	D	39	Total	O	0	0
			39	39		
6	E	45	Total	O	0	0
			45	45		
6	F	60	Total	O	0	0
			60	60		

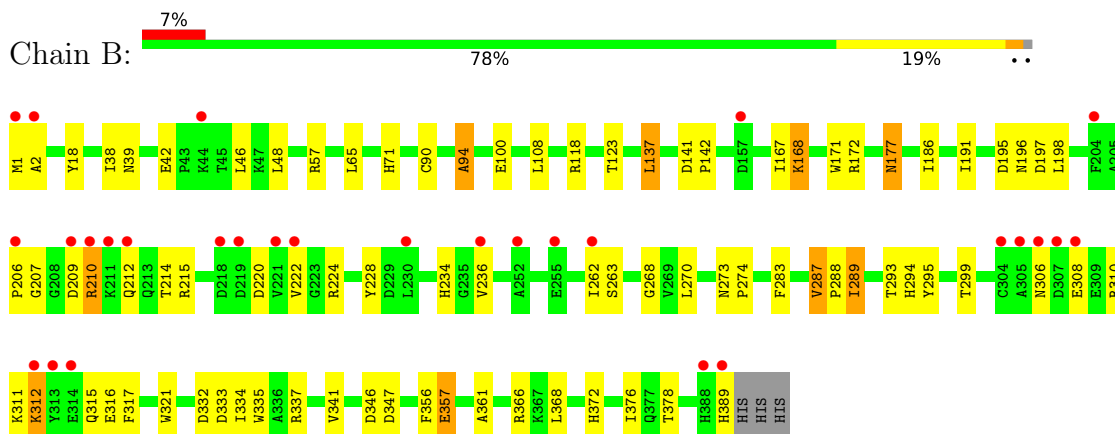
3 Residue-property plots

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.

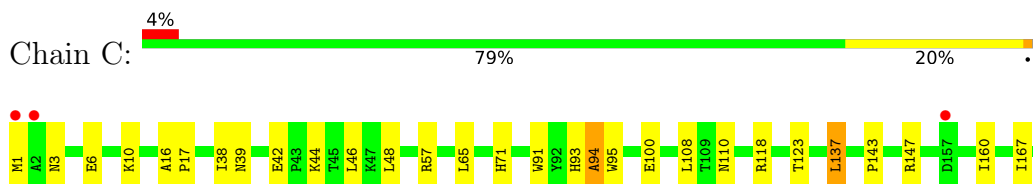
- Molecule 1: Terminal oxygenase component of carbazole

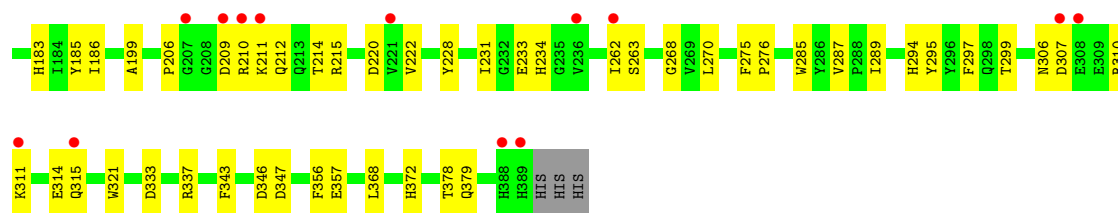


- Molecule 1: Terminal oxygenase component of carbazole

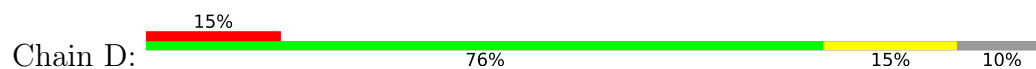


- Molecule 1: Terminal oxygenase component of carbazole





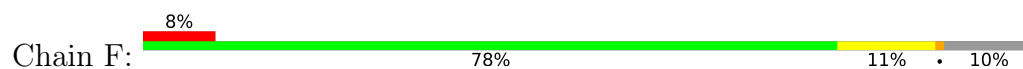
- Molecule 2: Ferredoxin component of carbazole



- Molecule 2: Ferredoxin component of carbazole



- Molecule 2: Ferredoxin component of carbazole



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	98.16Å 89.84Å 105.21Å 90.00° 104.40° 90.00°	Depositor
Resolution (Å)	40.62 – 1.95 40.62 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.7 (40.62-1.95) 99.8 (40.62-1.95)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.96 (at 1.95Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.198 , 0.219 0.191 , 0.210	Depositor DCC
R_{free} test set	6437 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	26.0	Xtriage
Anisotropy	0.033	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 38.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12864	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.79% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FES, FE2, 9CA

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	0.38	0/3221	0.89	10/4372 (0.2%)
1	B	0.35	0/3221	0.88	13/4372 (0.3%)
1	C	0.35	0/3221	0.86	9/4372 (0.2%)
2	D	0.32	0/784	0.84	1/1066 (0.1%)
2	E	0.33	0/810	0.83	1/1101 (0.1%)
2	F	0.33	0/784	0.83	2/1066 (0.2%)
All	All	0.35	0/12041	0.87	36/16349 (0.2%)

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	1

There are no bond length outliers.

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	3	ASN	N-CA-C	-7.65	104.09	113.50
1	A	287	VAL	N-CA-C	7.46	115.30	107.76
1	B	142	PRO	N-CA-C	7.01	118.08	110.58
1	B	287	VAL	N-CA-C	6.94	114.77	107.76
1	B	141	ASP	CA-C-N	6.68	124.47	119.66
1	B	141	ASP	C-N-CA	6.68	124.47	119.66
1	B	289	ILE	N-CA-C	-6.60	105.40	111.67
1	A	289	ILE	N-CA-C	-6.36	105.62	111.67
2	F	43	GLU	N-CA-C	-6.30	101.78	110.35
1	A	357	GLU	N-CA-C	6.28	121.73	111.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	43	GLU	N-CA-C	-6.15	101.82	110.50
1	B	100	GLU	N-CA-C	6.04	117.94	111.36
1	C	289	ILE	N-CA-C	-5.95	106.02	111.67
1	A	352	ASN	N-CA-C	5.89	120.54	113.23
1	A	168	LYS	N-CA-C	5.88	118.74	109.39
1	C	100	GLU	N-CA-C	5.86	117.66	111.28
1	A	199	ALA	N-CA-C	-5.72	99.97	109.46
1	B	357	GLU	N-CA-C	5.67	120.72	111.37
1	A	341	VAL	N-CA-C	5.66	116.31	110.36
1	C	94	ALA	N-CA-C	5.65	119.47	112.58
1	C	321	TRP	N-CA-C	5.63	118.14	111.33
1	C	177	ASN	N-CA-C	-5.58	104.83	111.69
1	A	94	ALA	N-CA-C	5.57	119.38	112.24
1	B	372	HIS	N-CA-C	5.55	120.04	112.88
1	A	372	HIS	N-CA-C	5.53	120.73	112.94
2	F	70	GLY	N-CA-C	-5.51	105.48	112.54
1	B	94	ALA	N-CA-C	5.48	119.25	112.24
1	B	2	ALA	N-CA-C	5.39	118.67	111.24
1	C	199	ALA	N-CA-C	-5.38	99.00	108.20
2	D	73	ASN	N-CA-C	-5.36	100.44	108.96
1	C	95	TRP	N-CA-C	-5.35	102.37	110.28
1	B	341	VAL	N-CA-C	5.20	115.82	110.36
1	B	168	LYS	N-CA-C	5.18	117.58	108.56
1	B	177	ASN	N-CA-C	-5.16	105.73	112.23
1	A	100	GLU	N-CA-C	5.16	116.90	111.28
1	C	372	HIS	N-CA-C	5.12	119.48	112.88

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	210	ARG	Sidechain

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	3136	0	3042	42	0
1	B	3136	0	3042	54	0
1	C	3136	0	3042	57	0
2	D	768	0	745	10	0
2	E	794	0	770	21	0
2	F	768	0	745	8	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	A	4	0	0	1	0
4	B	4	0	0	0	0
4	C	4	0	0	1	0
4	D	4	0	0	1	0
4	E	4	0	0	1	0
4	F	4	0	0	0	0
5	B	13	0	9	1	0
5	C	13	0	9	6	0
6	A	339	0	0	3	0
6	B	291	0	0	4	0
6	C	299	0	0	3	0
6	D	39	0	0	0	0
6	E	45	0	0	0	0
6	F	60	0	0	1	0
All	All	12864	0	11404	189	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 8.

All (189) 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:1:MET:H2	1:A:378:THR:HG22	1.34	0.90
2:E:92:GLU:HB2	2:E:107:LYS:HE3	1.55	0.87
1:A:289:ILE:HB	1:A:293:THR:HG23	1.57	0.86
1:C:262:ILE:HD11	5:C:2002:9CA:C7	2.12	0.79
1:C:262:ILE:HG23	1:C:270:LEU:HD11	1.67	0.77
1:C:262:ILE:CG2	1:C:270:LEU:HD11	2.19	0.72
1:A:290:ASP:OD1	1:A:293:THR:HG22	1.90	0.71
1:A:1:MET:H1	1:A:376:ILE:CG2	2.03	0.71
1:A:65:LEU:HD23	1:A:123:THR:HG22	1.73	0.71
2:E:108:LEU:O	2:E:109:GLU:HB3	1.92	0.70
1:A:1:MET:N	1:A:378:THR:HG22	2.04	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:184:ILE:HD11	1:A:200:LEU:CD1	2.20	0.70
1:B:65:LEU:HD23	1:B:123:THR:HG22	1.74	0.70
1:C:6:GLU:O	1:C:10:LYS:HG2	1.92	0.69
1:C:65:LEU:HD23	1:C:123:THR:HG22	1.73	0.69
2:D:7:LYS:HD2	2:D:99:GLU:OE2	1.92	0.69
1:B:1:MET:HG3	1:B:376:ILE:HB	1.75	0.68
1:A:194:LYS:HB2	6:A:547:HOH:O	1.95	0.66
1:C:210:ARG:HG2	6:C:932:HOH:O	1.96	0.66
1:B:209:ASP:OD2	1:B:212:GLN:HG3	1.97	0.65
1:C:94:ALA:HB1	1:C:108:LEU:HB2	1.79	0.65
2:D:90:VAL:HG23	2:D:107:LYS:HG2	1.79	0.65
2:D:19:ILE:HG21	2:D:54:SER:HA	1.79	0.65
1:C:262:ILE:CD1	5:C:2002:9CA:C7	2.75	0.64
1:B:311:LYS:O	1:B:315:GLN:HG3	1.98	0.63
2:E:14:MET:HE3	2:E:20:ARG:HB2	1.80	0.63
1:A:94:ALA:HB1	1:A:108:LEU:HB2	1.81	0.63
1:B:206:PRO:HG2	1:B:357:GLU:HB3	1.81	0.62
1:B:1:MET:H3	1:B:376:ILE:HG21	1.64	0.62
1:A:184:ILE:HD11	1:A:200:LEU:HD11	1.80	0.62
1:C:1:MET:HB2	1:C:378:THR:HG21	1.81	0.62
1:C:233:GLU:HG2	6:C:918:HOH:O	2.00	0.62
1:A:1:MET:H1	1:A:376:ILE:HG21	1.62	0.61
1:C:306:ASN:O	1:C:310:ARG:HG3	1.99	0.61
1:B:317:PHE:HA	1:B:321:TRP:HB2	1.83	0.61
2:E:4:ILE:HG23	2:E:5:TRP:N	2.15	0.60
1:B:361:ALA:HB3	6:B:1067:HOH:O	2.02	0.60
1:B:1:MET:HB2	1:B:378:THR:CG2	2.31	0.60
2:E:96:LYS:HB3	2:E:96:LYS:NZ	2.16	0.59
1:B:262:ILE:CG2	1:B:270:LEU:HD11	2.33	0.59
1:C:177:ASN:HD22	1:C:337:ARG:HE	1.51	0.58
1:A:118:ARG:HG3	6:F:903:HOH:O	2.03	0.58
2:D:7:LYS:HG2	2:D:101:TYR:CE2	2.39	0.57
1:C:262:ILE:CD1	5:C:2002:9CA:C8	2.83	0.57
1:C:48:LEU:HD23	1:C:137:LEU:HD23	1.84	0.57
1:C:220:ASP:OD1	1:C:222:VAL:HG22	2.05	0.57
1:A:287:VAL:HB	1:A:295:TYR:HB2	1.87	0.56
1:C:287:VAL:HB	1:C:295:TYR:HB2	1.88	0.56
1:C:177:ASN:ND2	1:C:337:ARG:HE	2.04	0.56
1:B:186:ILE:HG22	1:C:91:TRP:O	2.06	0.55
1:C:177:ASN:HD21	1:C:333:ASP:HB3	1.72	0.55
1:C:38:ILE:HA	1:C:44:LYS:HZ2	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:10:ALA:HB3	2:D:13:ASP:OD2	2.06	0.54
1:A:367:LYS:O	1:A:371:GLU:HG3	2.07	0.54
2:F:53:LEU:CD1	2:F:88:LEU:HD11	2.38	0.54
1:C:262:ILE:HD13	5:C:2002:9CA:C8	2.37	0.54
1:C:210:ARG:O	1:C:214:THR:HG22	2.08	0.53
1:C:311:LYS:O	1:C:315:GLN:HG3	2.09	0.53
2:E:3:GLN:HG3	2:E:4:ILE:N	2.23	0.53
1:A:91:TRP:O	1:C:186:ILE:HG22	2.09	0.53
1:B:48:LEU:HD23	1:B:137:LEU:HD23	1.89	0.53
1:B:196:ASN:HD22	1:B:198:LEU:HD21	1.74	0.53
1:B:171:TRP:CE2	1:B:172:ARG:HG3	2.44	0.52
1:B:287:VAL:HB	1:B:295:TYR:HB2	1.92	0.52
1:B:333:ASP:O	1:B:337:ARG:HG3	2.09	0.52
1:B:177:ASN:HD21	1:B:333:ASP:HB3	1.74	0.52
1:A:290:ASP:CG	1:A:293:THR:HG22	2.35	0.52
1:C:262:ILE:HD13	6:C:1073:HOH:O	2.09	0.52
2:F:7:LYS:HB2	2:F:7:LYS:NZ	2.25	0.51
1:C:38:ILE:HG23	1:C:57:ARG:HH21	1.74	0.51
2:D:14:MET:HE3	2:D:20:ARG:HB2	1.93	0.51
1:C:209:ASP:CG	1:C:212:GLN:HE21	2.19	0.51
1:B:191:ILE:O	1:B:195:ASP:HB2	2.11	0.50
1:C:231:ILE:HD13	1:C:262:ILE:HD12	1.92	0.50
1:B:312:LYS:O	1:B:312:LYS:HD3	2.11	0.50
2:E:4:ILE:CG2	2:E:5:TRP:N	2.75	0.50
1:B:118:ARG:NH2	6:B:931:HOH:O	2.45	0.50
1:C:210:ARG:HG3	1:C:211:LYS:N	2.27	0.50
1:C:171:TRP:CE2	1:C:172:ARG:HG3	2.47	0.49
1:B:210:ARG:O	1:B:214:THR:HG22	2.12	0.49
1:A:23:LEU:HD22	6:A:649:HOH:O	2.13	0.49
1:A:171:TRP:CE2	1:A:172:ARG:HG3	2.48	0.49
1:B:38:ILE:HG23	1:B:57:ARG:HH21	1.78	0.49
1:A:1:MET:N	1:A:378:THR:CG2	2.74	0.48
1:B:171:TRP:CG	1:B:288:PRO:HG3	2.48	0.48
1:C:167:ILE:HB	1:C:294:HIS:CE1	2.48	0.48
2:E:68:HIS:HB2	4:E:201:FES:S2	2.53	0.48
2:E:10:ALA:HB3	2:E:13:ASP:OD2	2.13	0.48
2:E:51:ALA:HB2	2:E:67:PHE:CG	2.48	0.48
1:A:94:ALA:CB	1:A:108:LEU:HB2	2.43	0.48
1:A:215:ARG:HB2	1:A:230:LEU:HD11	1.96	0.48
1:C:231:ILE:CD1	1:C:262:ILE:HD12	2.44	0.47
2:F:19:ILE:HG21	2:F:54:SER:HA	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:ARG:HG3	1:A:250:GLU:N	2.29	0.47
1:C:16:ALA:HB3	1:C:17:PRO:HD3	1.96	0.47
1:C:262:ILE:HD11	5:C:2002:9CA:C8	2.44	0.47
2:D:68:HIS:HB2	4:D:201:FES:S2	2.54	0.47
1:A:184:ILE:HD11	1:A:200:LEU:HD12	1.96	0.47
1:A:260:ASN:HD22	1:A:274:PRO:HD2	1.79	0.47
1:B:1:MET:N	1:B:376:ILE:HG21	2.30	0.47
1:B:90:CYS:O	1:B:94:ALA:HA	2.14	0.47
1:C:209:ASP:CG	1:C:212:GLN:HG3	2.40	0.47
2:E:96:LYS:HB3	2:E:96:LYS:HZ2	1.79	0.47
2:F:51:ALA:HB2	2:F:67:PHE:CG	2.49	0.47
1:C:93:HIS:HB2	4:C:401:FES:S1	2.55	0.47
2:E:108:LEU:O	2:E:109:GLU:CB	2.62	0.47
1:A:183:HIS:O	1:A:186:ILE:HG12	2.14	0.46
1:C:39:ASN:HB2	1:C:42:GLU:CD	2.40	0.46
1:C:233:GLU:HG3	1:C:234:HIS:CD2	2.50	0.46
2:E:47:THR:HG23	2:E:88:LEU:HD23	1.97	0.46
1:C:262:ILE:HD11	5:C:2002:9CA:C6	2.46	0.46
1:A:220:ASP:HB2	1:A:224:ARG:HB2	1.97	0.46
1:C:311:LYS:HE3	1:C:314:GLU:OE1	2.15	0.46
2:F:4:ILE:HG13	2:F:5:TRP:H	1.80	0.46
1:B:332:ASP:HA	1:B:335:TRP:NE1	2.30	0.45
1:C:285:TRP:HB2	1:C:297:PHE:HB3	1.99	0.45
2:D:51:ALA:HB2	2:D:67:PHE:CG	2.51	0.45
1:B:273:ASN:HA	1:B:274:PRO:HA	1.84	0.45
1:A:64:CYS:HB2	1:A:126:VAL:HG21	1.97	0.45
1:A:168:LYS:HE2	6:A:864:HOH:O	2.16	0.45
2:D:31:ALA:O	2:D:41:ALA:HA	2.17	0.45
1:B:1:MET:HB2	1:B:378:THR:HG21	1.97	0.45
2:E:4:ILE:HG23	2:E:5:TRP:CD1	2.52	0.45
1:B:220:ASP:OD1	1:B:222:VAL:N	2.48	0.45
1:B:167:ILE:HB	1:B:294:HIS:CE1	2.52	0.44
1:B:224:ARG:NH2	6:B:976:HOH:O	2.49	0.44
2:E:92:GLU:OE1	2:E:107:LYS:NZ	2.50	0.44
1:B:167:ILE:HD12	1:B:334:ILE:HG23	1.98	0.44
1:C:275:PHE:CG	1:C:276:PRO:HA	2.53	0.44
1:B:289:ILE:HB	1:B:293:THR:OG1	2.18	0.43
1:A:158:MET:HE1	1:A:281:MET:CE	2.49	0.43
1:A:175:VAL:HG11	1:A:365:TRP:CE2	2.53	0.43
1:C:118:ARG:HH11	1:C:118:ARG:HG3	1.83	0.43
1:A:93:HIS:HB2	4:A:401:FES:S1	2.58	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:321:TRP:HA	1:B:321:TRP:CE3	2.54	0.43
1:A:287:VAL:HA	1:A:288:PRO:HD3	1.88	0.43
1:A:143:PRO:HG3	1:A:147:ARG:CZ	2.49	0.43
1:B:94:ALA:CB	1:B:108:LEU:HB2	2.48	0.43
1:B:312:LYS:O	1:B:316:GLU:HG3	2.18	0.43
1:B:177:ASN:HD22	1:B:337:ARG:HE	1.65	0.43
1:B:283:PHE:HB2	1:B:299:THR:OG1	2.19	0.43
1:B:228:TYR:CD1	1:B:263:SER:HB3	2.54	0.43
1:A:136:TYR:CE1	1:A:143:PRO:HD2	2.54	0.43
1:C:333:ASP:O	1:C:337:ARG:HG3	2.19	0.43
1:B:94:ALA:HB1	1:B:108:LEU:HB2	2.01	0.42
1:B:207:GLY:HA3	1:B:234:HIS:O	2.19	0.42
1:B:312:LYS:HD2	1:B:316:GLU:OE2	2.18	0.42
1:C:160:ILE:HG23	1:C:299:THR:HB	2.01	0.42
1:A:285:TRP:HB2	1:A:297:PHE:HB3	2.01	0.42
1:A:308:GLU:H	1:A:308:GLU:CD	2.27	0.42
1:B:306:ASN:OD1	1:B:308:GLU:HB2	2.19	0.42
1:C:39:ASN:HB2	1:C:42:GLU:OE2	2.19	0.42
1:C:206:PRO:HG2	1:C:357:GLU:HB3	2.01	0.42
1:A:1:MET:HA	1:A:20:ASP:OD2	2.18	0.42
1:C:183:HIS:O	1:C:186:ILE:HG12	2.19	0.42
2:F:18:THR:C	2:F:19:ILE:HG13	2.44	0.42
1:B:1:MET:H3	1:B:376:ILE:CG2	2.30	0.42
1:C:228:TYR:CD1	1:C:263:SER:HB3	2.54	0.42
2:F:4:ILE:HG13	2:F:5:TRP:CD1	2.54	0.42
1:C:110:ASN:C	1:C:110:ASN:HD22	2.28	0.42
2:E:99:GLU:OE1	2:E:99:GLU:HA	2.18	0.42
1:B:168:LYS:HE2	6:B:906:HOH:O	2.18	0.42
2:E:3:GLN:HG3	2:E:4:ILE:H	1.83	0.42
2:E:19:ILE:HG21	2:E:54:SER:HA	2.02	0.42
1:A:64:CYS:HB2	1:A:126:VAL:CG2	2.50	0.42
1:B:207:GLY:HA2	1:B:236:VAL:HG23	2.01	0.41
1:C:346:ASP:O	1:C:347:ASP:HB2	2.20	0.41
1:A:260:ASN:HD22	1:A:274:PRO:CD	2.34	0.41
1:B:306:ASN:C	1:B:310:ARG:NH1	2.79	0.41
1:C:94:ALA:CB	1:C:108:LEU:HB2	2.47	0.41
1:B:196:ASN:O	1:B:197:ASP:C	2.64	0.41
2:E:25:VAL:O	2:E:25:VAL:HG13	2.19	0.41
2:F:5:TRP:CH2	2:F:92:GLU:HG3	2.56	0.41
1:B:346:ASP:O	1:B:347:ASP:HB2	2.21	0.41
1:C:215:ARG:HB3	1:C:228:TYR:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:ASP:OD1	1:A:223:GLY:HA2	2.20	0.41
1:B:39:ASN:HB2	1:B:42:GLU:OE1	2.20	0.41
1:C:1:MET:SD	1:C:1:MET:N	2.94	0.41
2:E:5:TRP:CZ3	2:E:94:GLU:HG2	2.57	0.40
1:A:38:ILE:CD1	1:A:55:VAL:HG12	2.52	0.40
1:A:73:GLY:HA3	1:C:343:PHE:CG	2.56	0.40
1:B:18:TYR:CE2	1:B:366:ARG:HG2	2.55	0.40
1:B:215:ARG:HB3	1:B:228:TYR:HB2	2.03	0.40
2:D:99:GLU:HB3	2:D:101:TYR:CE1	2.57	0.40
1:C:183:HIS:C	1:C:185:TYR:H	2.29	0.40
2:E:91:PHE:CD1	2:E:91:PHE:N	2.88	0.40
1:B:262:ILE:HD13	5:B:2001:9CA:C8	2.52	0.40
1:B:389:HIS:ND1	1:B:389:HIS:C	2.79	0.40
1:C:143:PRO:HG3	1:C:147:ARG:CZ	2.52	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	387/392 (99%)	372 (96%)	13 (3%)	2 (0%)	24	16
1	B	387/392 (99%)	366 (95%)	19 (5%)	2 (0%)	24	16
1	C	387/392 (99%)	366 (95%)	19 (5%)	2 (0%)	24	16
2	D	102/115 (89%)	98 (96%)	4 (4%)	0	100	100
2	E	105/115 (91%)	100 (95%)	5 (5%)	0	100	100
2	F	102/115 (89%)	101 (99%)	1 (1%)	0	100	100
All	All	1470/1521 (97%)	1403 (95%)	61 (4%)	6 (0%)	30	21

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	268	GLY
1	A	71	HIS
1	A	268	GLY
1	B	268	GLY
1	C	71	HIS
1	B	71	HIS

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	336/339 (99%)	331 (98%)	5 (2%)	57	54
1	B	336/339 (99%)	330 (98%)	6 (2%)	51	47
1	C	336/339 (99%)	330 (98%)	6 (2%)	51	47
2	D	82/93 (88%)	82 (100%)	0	100	100
2	E	85/93 (91%)	83 (98%)	2 (2%)	43	36
2	F	82/93 (88%)	80 (98%)	2 (2%)	43	36
All	All	1257/1296 (97%)	1236 (98%)	21 (2%)	53	49

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	LEU
1	A	137	LEU
1	A	269	VAL
1	A	356	PHE
1	A	368	LEU
1	B	46	LEU
1	B	137	LEU
1	B	210	ARG
1	B	312	LYS
1	B	356	PHE
1	B	368	LEU
1	C	46	LEU
1	C	137	LEU

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Mol	Chain	Res	Type
1	C	307	ASP
1	C	356	PHE
1	C	368	LEU
1	C	379	GLN
2	E	6	LEU
2	E	96	LYS
2	F	6	LEU
2	F	7	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	110	ASN
1	A	119	GLN
1	A	213	GLN
1	A	234	HIS
1	A	260	ASN
1	A	282	GLN
1	A	315	GLN
1	A	330	ASN
1	B	52	ASN
1	B	110	ASN
1	B	152	ASN
1	B	177	ASN
1	B	282	GLN
1	B	330	ASN
1	B	372	HIS
1	C	110	ASN
1	C	177	ASN
1	C	212	GLN
1	C	282	GLN
1	C	315	GLN
1	C	330	ASN
1	C	388	HIS
2	E	3	GLN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 3 are monoatomic - leaving 8 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	$\# Z > 2$
4	FES	D	201	2	0,4,4	-	-	-		
4	FES	A	401	1	0,4,4	-	-	-		
5	9CA	C	2002	-	15,15,15	1.37	3 (20%)	21,21,21	1.01	2 (9%)
4	FES	F	201	2	0,4,4	-	-	-		
4	FES	B	401	1	0,4,4	-	-	-		
4	FES	C	401	1	0,4,4	-	-	-		
5	9CA	B	2001	-	15,15,15	1.37	3 (20%)	21,21,21	0.99	0
4	FES	E	201	2	0,4,4	-	-	-		

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
4	FES	D	201	2	-	-	0/1/1/1
4	FES	A	401	1	-	-	0/1/1/1
5	9CA	C	2002	-	-	-	0/3/3/3
4	FES	F	201	2	-	-	0/1/1/1
4	FES	B	401	1	-	-	0/1/1/1
4	FES	C	401	1	-	-	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	9CA	B	2001	-	-	-	0/3/3/3
4	FES	E	201	2	-	-	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	2002	9CA	C9A-N9	-2.82	1.33	1.38
5	C	2002	9CA	C4B-C4A	-2.81	1.39	1.46
5	B	2001	9CA	C4B-C4A	-2.80	1.39	1.46
5	B	2001	9CA	C9A-N9	-2.79	1.33	1.38
5	B	2001	9CA	C8A-N9	-2.77	1.33	1.38
5	C	2002	9CA	C8A-N9	-2.75	1.33	1.38

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	2002	9CA	C1-C9A-N9	2.02	131.95	129.07
5	C	2002	9CA	C8-C8A-N9	2.01	131.94	129.07

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	201	FES	1	0
4	A	401	FES	1	0
5	C	2002	9CA	6	0
4	C	401	FES	1	0
5	B	2001	9CA	1	0
4	E	201	FES	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	389/392 (99%)	0.19	18 (4%) 37 43	15, 23, 39, 47	0
1	B	389/392 (99%)	0.47	29 (7%) 20 23	15, 27, 43, 49	0
1	C	389/392 (99%)	0.37	16 (4%) 41 48	18, 27, 40, 49	0
2	D	104/115 (90%)	1.08	17 (16%) 4 5	20, 37, 46, 48	0
2	E	107/115 (93%)	0.86	10 (9%) 14 16	22, 33, 42, 47	0
2	F	104/115 (90%)	0.94	9 (8%) 16 18	23, 34, 41, 45	0
All	All	1482/1521 (97%)	0.47	99 (6%) 24 27	15, 27, 43, 49	0

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	MET	9.1
1	C	1	MET	8.9
1	A	1	MET	8.0
1	A	221	VAL	5.6
1	B	389	HIS	5.2
1	A	389	HIS	4.7
1	A	220	ASP	4.7
2	F	4	ILE	4.3
2	D	104	GLY	4.3
1	C	262	ILE	4.1
1	B	2	ALA	4.0
1	A	2	ALA	4.0
2	D	4	ILE	3.9
1	B	222	VAL	3.9
1	B	221	VAL	3.8
1	C	209	ASP	3.8
1	B	211	LYS	3.7
1	C	307	ASP	3.7
1	B	209	ASP	3.6

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Mol	Chain	Res	Type	RSRZ
1	C	2	ALA	3.4
1	B	210	ARG	3.4
1	C	221	VAL	3.4
1	C	308	GLU	3.3
1	B	262	ILE	3.3
1	C	211	LYS	3.3
1	A	118	ARG	3.3
1	A	261	ASP	3.2
2	F	25	VAL	3.2
2	D	92	GLU	3.1
1	A	388	HIS	3.1
1	B	204	PHE	3.0
1	B	388	HIS	3.0
1	A	218	ASP	3.0
1	C	315	GLN	2.9
1	B	308	GLU	2.9
2	F	105	GLU	2.9
1	B	305	ALA	2.9
2	E	4	ILE	2.9
1	A	210	ARG	2.8
1	C	311	LYS	2.8
2	D	6	LEU	2.8
1	B	314	GLU	2.8
2	D	99	GLU	2.7
2	D	25	VAL	2.7
2	E	27	ALA	2.6
2	E	11	ALA	2.6
2	D	97	GLU	2.6
1	B	252	ALA	2.5
1	B	157	ASP	2.5
2	E	37	ASP	2.5
2	D	39	PHE	2.5
2	E	97	GLU	2.5
1	C	207	GLY	2.4
2	D	105	GLU	2.4
1	B	219	ASP	2.4
1	C	157	ASP	2.4
1	B	206	PRO	2.4
1	A	260	ASN	2.4
1	C	389	HIS	2.4
2	D	103	ALA	2.4
2	D	23	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	236	VAL	2.4
2	E	26	GLY	2.3
2	D	12	SER	2.3
2	E	13	ASP	2.3
1	A	311	LYS	2.3
1	B	44	LYS	2.2
1	A	314	GLU	2.2
2	D	98	GLY	2.2
1	C	210	ARG	2.2
1	B	230	LEU	2.2
2	E	3	GLN	2.2
1	B	218	ASP	2.2
2	D	89	GLY	2.2
1	B	212	GLN	2.2
1	B	304	CYS	2.2
2	F	92	GLU	2.1
1	A	222	VAL	2.1
1	C	236	VAL	2.1
2	D	5	TRP	2.1
2	F	100	VAL	2.1
2	E	96	LYS	2.1
2	F	106	LYS	2.1
2	D	101	TYR	2.1
1	B	312	LYS	2.1
1	C	388	HIS	2.1
2	F	6	LEU	2.1
1	A	308	GLU	2.0
2	D	7	LYS	2.0
1	B	313	TYR	2.0
1	B	306	ASN	2.0
1	A	305	ALA	2.0
2	F	103	ALA	2.0
1	A	209	ASP	2.0
1	B	307	ASP	2.0
1	B	255	GLU	2.0
1	A	310	ARG	2.0
2	E	35	VAL	2.0
2	F	90	VAL	2.0

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

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

6.3 Carbohydrates [i](#)

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, 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($\text{\AA}^2$)	Q<0.9
5	9CA	C	2002	13/13	0.90	0.10	27,28,28,28	0
5	9CA	B	2001	13/13	0.93	0.08	29,29,29,29	0
4	FES	C	401	4/4	0.99	0.04	21,22,24,24	0
4	FES	D	201	4/4	0.99	0.04	18,20,21,23	0
4	FES	E	201	4/4	0.99	0.04	22,22,23,24	0
4	FES	F	201	4/4	0.99	0.03	22,23,23,24	0
3	FE2	B	501	1/1	0.99	0.03	29,29,29,29	0
4	FES	A	401	4/4	0.99	0.04	20,20,20,21	0
3	FE2	C	501	1/1	1.00	0.03	25,25,25,25	0
3	FE2	A	501	1/1	1.00	0.03	23,23,23,23	0
4	FES	B	401	4/4	1.00	0.03	18,19,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.