



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 06:34 PM UTC

PDB ID : 4NBH / pdb_00004nbh
Title : Carbazole-bound Oxygenase with Gln282 replaced by Tyr and ferredoxin complex of carbazole 1,9a-dioxygenase
Authors : Ashikawa, Y.; Usami, Y.; Inoue, K.; Nojiri, H.
Deposited on : 2013-10-23
Resolution : 2.15 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

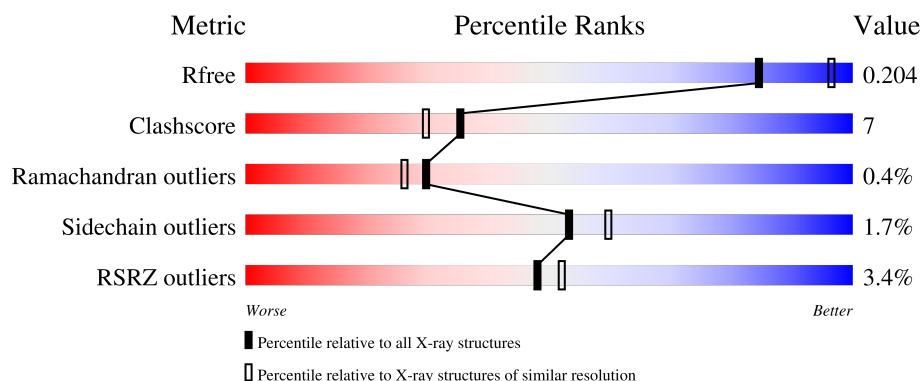
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	2% 80% 18% ..
1	B	392	4% 80% 18% ..
1	C	392	3% 77% 21% ..
2	D	115	3% 78% 12% 10%
2	E	115	7% 81% 16% ..

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Mol	Chain	Length	Quality of chain
2	F	115	3% 76% 14% • 10%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 12842 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	388	Total	C	N	O	S	0	0	0
			3131	2003	533	582	13			
1	B	390	Total	C	N	O	S	0	0	0
			3151	2015	539	584	13			
1	C	388	Total	C	N	O	S	0	0	0
			3131	2003	533	582	13			

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	282	TYR	GLN	engineered mutation	UNP Q84II6
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	282	TYR	GLN	engineered mutation	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	282	TYR	GLN	engineered mutation	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

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Chain	Residue	Modelled	Actual	Comment	Reference
C	389	HIS	-	expression tag	UNP Q84II6
C	390	HIS	-	expression tag	UNP Q84II6
C	391	HIS	-	expression tag	UNP Q84II6
C	392	HIS	-	expression tag	UNP Q84II6

- Molecule 2 is a protein called Ferredoxin CarAc.

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	112	Total	C	N	O	S	0	0	0
			844	529	148	160	7			
2	F	104	Total	C	N	O	S	0	0	0
			768	483	129	149	7			

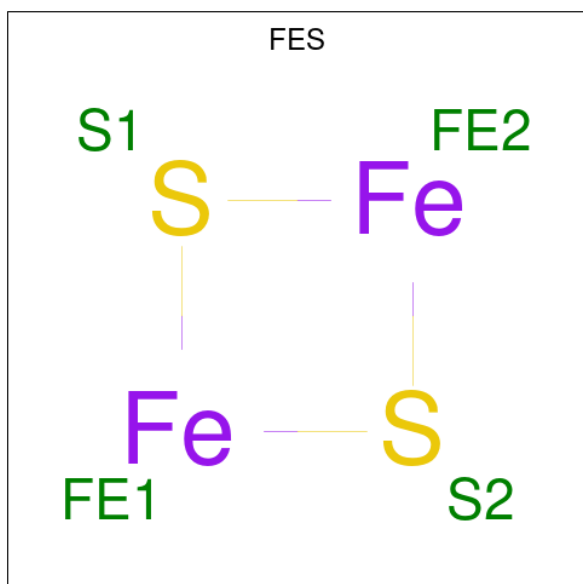
There are 24 discrepancies between the modelled and reference sequences:

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).

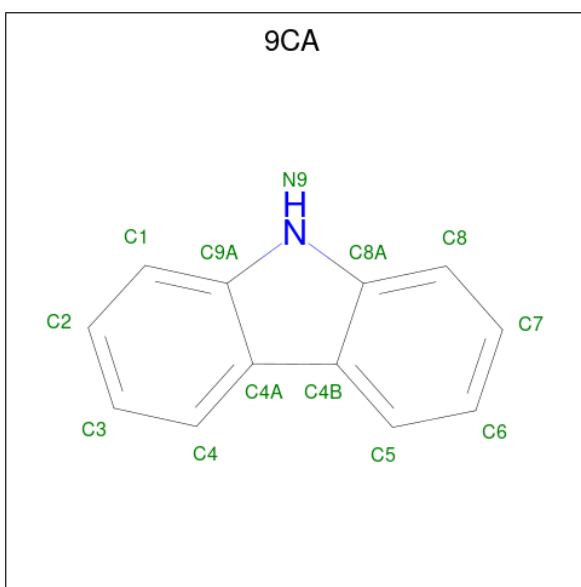
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Fe 1	0	0
3	B	1	Total 1	Fe 1	0	0
3	C	1	Total 1	Fe 1	0	0

- Molecule 4 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe₂S₂).



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

- Molecule 5 is 9H-CARBAZOLE (CCD ID: 9CA) (formula: C₁₂H₉N).



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		

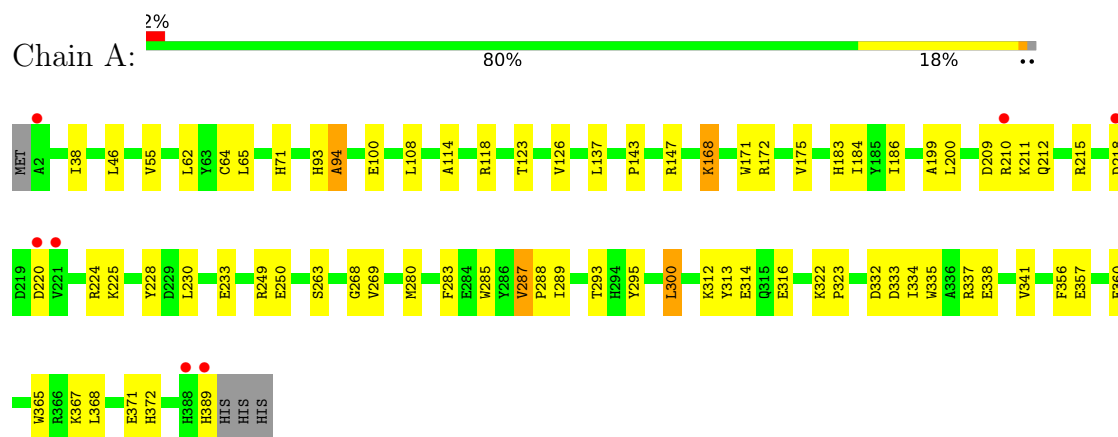
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	330	Total	O	0	0
			330	330		
6	B	263	Total	O	0	0
			263	263		
6	C	278	Total	O	0	0
			278	278		
6	D	30	Total	O	0	0
			30	30		
6	E	46	Total	O	0	0
			46	46		
6	F	49	Total	O	0	0
			49	49		

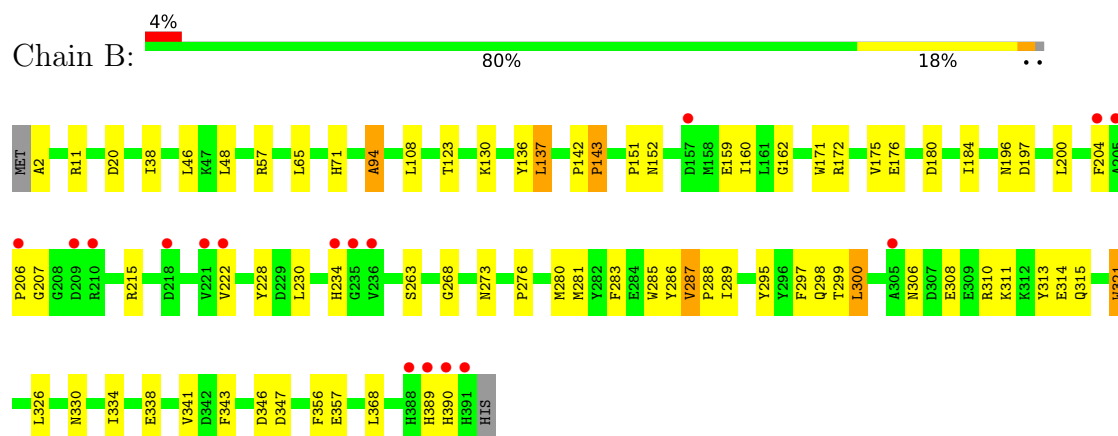
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

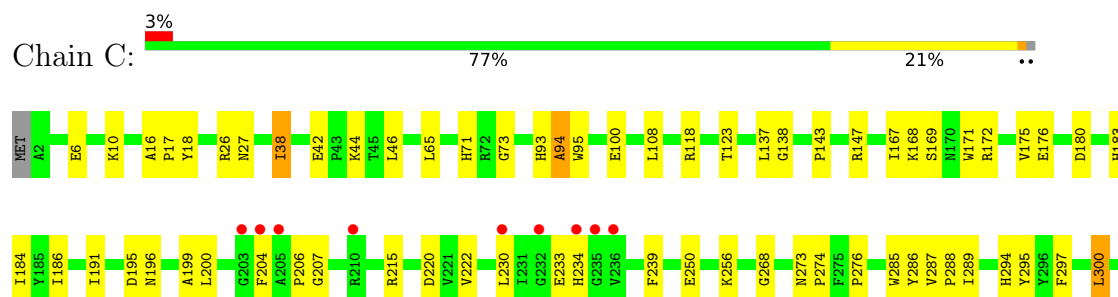
- Molecule 1: Terminal oxygenase component of carbazole



- Molecule 1: Terminal oxygenase component of carbazole

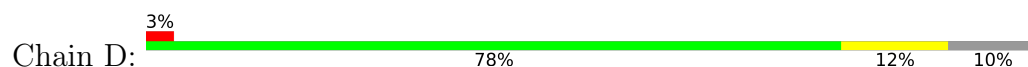


- Molecule 1: Terminal oxygenase component of carbazole

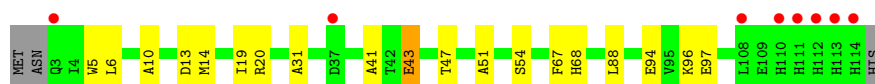
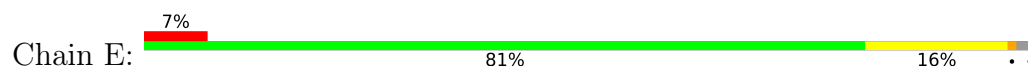




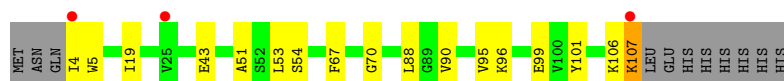
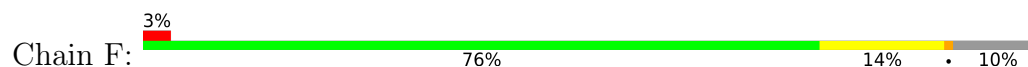
• Molecule 2: Ferredoxin CarAc



• Molecule 2: Ferredoxin CarAc



• Molecule 2: Ferredoxin CarAc



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	98.18Å 89.95Å 105.19Å 90.00° 104.29° 90.00°	Depositor
Resolution (Å)	40.92 – 2.15 40.92 – 2.15	Depositor EDS
% Data completeness (in resolution range)	99.6 (40.92-2.15) 99.7 (40.92-2.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.35 (at 2.16Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.183 , 0.211 0.178 , 0.204	Depositor DCC
R_{free} test set	4844 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	24.6	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 41.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12842	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% 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: 9CA, FES, FE2

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	0/3217	0.87	9/4368 (0.2%)
1	B	0.34	0/3239	0.86	7/4398 (0.2%)
1	C	0.36	0/3217	0.86	7/4368 (0.2%)
2	D	0.33	0/784	0.85	0/1066
2	E	0.33	0/865	0.84	1/1176 (0.1%)
2	F	0.34	0/784	0.87	2/1066 (0.2%)
All	All	0.36	0/12106	0.86	26/16442 (0.2%)

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	357	GLU	N-CA-C	8.74	121.60	111.11
1	C	289	ILE	N-CA-C	-7.52	105.34	111.81
1	A	287	VAL	N-CA-C	7.16	115.00	107.76
1	B	289	ILE	N-CA-C	-7.11	104.92	111.67
1	B	287	VAL	N-CA-C	7.04	114.87	107.76
1	A	357	GLU	N-CA-C	6.59	122.24	111.37
2	F	43	GLU	N-CA-C	-6.46	101.39	110.50
2	E	43	GLU	N-CA-C	-6.38	101.50	110.50
1	A	289	ILE	N-CA-C	-6.13	105.75	111.45
1	B	94	ALA	N-CA-C	6.12	120.08	112.24
1	A	341	VAL	N-CA-C	5.62	116.26	110.36
1	B	341	VAL	N-CA-C	5.58	116.22	110.36
1	C	199	ALA	N-CA-C	-5.48	99.80	108.73
1	A	199	ALA	N-CA-C	-5.47	100.38	109.46
1	C	95	TRP	N-CA-C	-5.42	101.82	109.96
1	A	100	GLU	N-CA-C	5.42	116.87	111.07
1	C	94	ALA	N-CA-C	5.37	119.11	112.24
1	A	94	ALA	N-CA-C	5.32	119.35	112.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	142	PRO	N-CA-C	5.24	116.19	110.58
1	A	372	HIS	N-CA-C	5.23	120.32	112.94
1	C	100	GLU	N-CA-C	5.21	116.96	111.28
1	C	352	ASN	N-CA-C	5.10	119.51	113.28
1	B	143	PRO	N-CA-C	-5.09	104.49	110.70
1	B	321	TRP	N-CA-C	5.08	117.22	111.02
2	F	70	GLY	N-CA-C	-5.03	106.10	112.54
1	A	168	LYS	N-CA-C	5.00	117.27	108.56

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3131	0	3031	46	0
1	B	3151	0	3045	50	0
1	C	3131	0	3031	50	0
2	D	768	0	745	7	0
2	E	844	0	805	10	0
2	F	768	0	745	12	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	0	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	0	0
6	A	330	0	0	3	0
6	B	263	0	0	4	0
6	C	278	0	0	3	0
6	D	30	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	E	46	0	0	0	0
6	F	49	0	0	3	0
All	All	12842	0	11420	173	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 (173) 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:B:2:ALA:HB1	6:B:779:HOH:O	1.44	1.13
1:A:280:MET:HE3	1:A:313:TYR:HE1	1.44	0.83
1:B:280:MET:HE2	1:B:300:LEU:HD22	1.61	0.82
1:C:6:GLU:O	1:C:10:LYS:HG2	1.88	0.74
1:C:65:LEU:HD23	1:C:123:THR:HG22	1.69	0.74
1:A:65:LEU:HD23	1:A:123:THR:HG22	1.72	0.72
1:C:220:ASP:OD1	1:C:222:VAL:HG22	1.92	0.68
1:B:230:LEU:HA	6:B:799:HOH:O	1.95	0.66
1:A:118:ARG:HD2	6:A:765:HOH:O	1.95	0.65
1:B:2:ALA:HB3	1:B:20:ASP:OD2	1.97	0.64
1:A:280:MET:HE3	1:A:313:TYR:CE1	2.30	0.64
2:F:106:LYS:NZ	6:F:349:HOH:O	2.20	0.64
1:B:38:ILE:HG23	1:B:57:ARG:HH21	1.63	0.63
1:B:65:LEU:HD23	1:B:123:THR:HG22	1.79	0.63
1:A:220:ASP:HB2	1:A:224:ARG:HB2	1.81	0.63
1:C:306:ASN:O	1:C:310:ARG:HG3	2.00	0.62
1:C:38:ILE:HD12	1:C:44:LYS:HD3	1.81	0.62
1:C:94:ALA:HB1	1:C:108:LEU:HB2	1.83	0.61
1:A:94:ALA:HB1	1:A:108:LEU:HB2	1.84	0.60
1:B:215:ARG:HB2	1:B:230:LEU:HD11	1.84	0.59
1:B:48:LEU:HD23	1:B:137:LEU:HD23	1.82	0.59
1:B:215:ARG:HB3	1:B:228:TYR:HB2	1.85	0.59
1:A:287:VAL:HB	1:A:295:TYR:HB2	1.84	0.59
1:A:38:ILE:HD13	1:A:62:LEU:HD22	1.83	0.59
1:C:333:ASP:O	1:C:337:ARG:HG3	2.03	0.58
2:F:96:LYS:HG3	6:F:321:HOH:O	2.04	0.58
1:A:280:MET:HE2	1:A:300:LEU:HD22	1.84	0.57
1:B:283:PHE:HB2	1:B:299:THR:OG1	2.05	0.57
2:D:51:ALA:HB2	2:D:67:PHE:CG	2.40	0.56
2:F:53:LEU:CD1	2:F:88:LEU:HD11	2.35	0.56
2:F:4:ILE:HG23	2:F:5:TRP:H	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:333:ASP:O	1:A:337:ARG:HG3	2.06	0.56
1:A:114:ALA:O	1:A:118:ARG:HD3	2.07	0.55
1:A:230:LEU:HB3	1:A:233:GLU:HG3	1.88	0.55
2:D:92:GLU:HG2	2:D:107:LYS:HE2	1.88	0.55
1:B:306:ASN:OD1	1:B:308:GLU:HB2	2.06	0.55
1:B:326:LEU:O	1:B:330:ASN:HB2	2.06	0.55
2:D:19:ILE:HG21	2:D:54:SER:HA	1.89	0.55
2:F:107:LYS:HB2	2:F:107:LYS:NZ	2.22	0.55
1:A:184:ILE:HD11	1:A:200:LEU:CD1	2.36	0.55
1:B:287:VAL:HB	1:B:295:TYR:HB2	1.90	0.54
1:B:334:ILE:O	1:B:338:GLU:HG3	2.08	0.54
1:A:209:ASP:OD1	1:A:212:GLN:HG3	2.08	0.54
1:B:276:PRO:HD2	1:B:300:LEU:HD21	1.90	0.53
1:B:11:ARG:HH21	1:C:388:HIS:CE1	2.26	0.53
2:E:47:THR:HG23	2:E:88:LEU:HD23	1.91	0.53
1:A:367:LYS:O	1:A:371:GLU:HG3	2.09	0.53
1:A:312:LYS:O	1:A:316:GLU:HG3	2.09	0.52
2:E:19:ILE:HG21	2:E:54:SER:HA	1.92	0.52
1:A:280:MET:CE	1:A:313:TYR:HE1	2.21	0.51
1:C:42:GLU:HG2	1:C:44:LYS:HZ2	1.75	0.51
1:A:215:ARG:HD2	6:A:687:HOH:O	2.09	0.51
1:C:94:ALA:CB	1:C:108:LEU:HB2	2.41	0.51
1:C:184:ILE:HD11	1:C:200:LEU:CD1	2.40	0.51
1:C:360:GLU:HG3	6:C:627:HOH:O	2.10	0.51
1:B:184:ILE:HD11	1:B:200:LEU:CD1	2.41	0.51
1:C:171:TRP:CE2	1:C:172:ARG:HG3	2.46	0.51
2:E:10:ALA:HB3	2:E:13:ASP:OD2	2.10	0.51
1:A:171:TRP:CE2	1:A:172:ARG:HG3	2.46	0.50
1:A:249:ARG:HG3	1:A:250:GLU:N	2.27	0.50
1:C:285:TRP:HB2	1:C:297:PHE:HB3	1.94	0.50
2:D:31:ALA:O	2:D:41:ALA:HA	2.12	0.50
1:C:176:GLU:O	1:C:180:ASP:HB2	2.11	0.49
5:B:503:9CA:H1	6:B:738:HOH:O	2.11	0.49
1:A:94:ALA:CB	1:A:108:LEU:HB2	2.42	0.49
1:C:143:PRO:HG3	1:C:147:ARG:CZ	2.42	0.49
1:C:276:PRO:CD	1:C:300:LEU:HD21	2.43	0.49
2:D:10:ALA:HB3	2:D:13:ASP:OD2	2.11	0.49
1:C:42:GLU:CG	1:C:44:LYS:HZ2	2.26	0.49
1:B:273:ASN:OD1	1:B:281:MET:HG3	2.13	0.49
2:E:51:ALA:HB2	2:E:67:PHE:CG	2.47	0.49
2:E:5:TRP:CZ3	2:E:94:GLU:HG2	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:222:VAL:O	1:B:222:VAL:HG12	2.12	0.48
1:B:171:TRP:CG	1:B:288:PRO:HG3	2.49	0.48
1:C:93:HIS:HB2	4:C:502:FES:S1	2.53	0.48
1:B:204:PHE:O	1:B:206:PRO:HD3	2.14	0.47
1:A:218:ASP:O	1:A:225:LYS:HD3	2.15	0.47
1:A:93:HIS:HB2	4:A:502:FES:S1	2.54	0.47
1:B:176:GLU:O	1:B:180:ASP:HB2	2.14	0.47
1:A:218:ASP:O	1:A:218:ASP:OD1	2.33	0.47
1:A:228:TYR:CD1	1:A:263:SER:HB3	2.49	0.47
2:F:51:ALA:HB2	2:F:67:PHE:CG	2.49	0.47
2:E:96:LYS:O	2:E:97:GLU:HB2	2.15	0.47
1:C:196:ASN:O	1:C:256:LYS:HD2	2.16	0.46
1:B:276:PRO:CD	1:B:300:LEU:HD21	2.44	0.46
1:B:311:LYS:O	1:B:315:GLN:HG3	2.15	0.46
1:B:2:ALA:HB3	1:B:20:ASP:CG	2.41	0.46
1:C:230:LEU:HB3	1:C:233:GLU:HG3	1.98	0.46
1:A:143:PRO:HG3	1:A:147:ARG:CZ	2.46	0.45
1:A:64:CYS:HB2	1:A:126:VAL:CG2	2.46	0.45
2:F:19:ILE:HG21	2:F:54:SER:HA	1.98	0.45
1:C:346:ASP:O	1:C:347:ASP:HB2	2.17	0.45
2:F:99:GLU:HB3	2:F:101:TYR:HE1	1.82	0.45
2:F:107:LYS:HB2	2:F:107:LYS:HZ2	1.82	0.45
1:A:210:ARG:HG3	1:A:360:GLU:CD	2.42	0.45
1:B:171:TRP:CE2	1:B:172:ARG:HG3	2.51	0.45
1:C:183:HIS:O	1:C:186:ILE:HG12	2.17	0.45
1:A:215:ARG:HB2	1:A:230:LEU:HD11	1.99	0.45
1:C:363:VAL:HG12	1:C:367:LYS:HE2	1.98	0.45
1:C:38:ILE:CD1	1:C:44:LYS:HD3	2.46	0.44
1:C:118:ARG:HH12	2:E:43:GLU:CD	2.26	0.44
1:C:287:VAL:HB	1:C:295:TYR:HB2	1.99	0.44
1:A:175:VAL:HG11	1:A:365:TRP:CE2	2.52	0.44
1:A:184:ILE:HD11	1:A:200:LEU:HD11	1.99	0.44
2:E:14:MET:HE3	2:E:20:ARG:HB2	1.99	0.44
1:C:27:ASN:HA	1:C:138:GLY:HA3	1.99	0.44
2:D:7:LYS:HG3	2:D:101:TYR:HE1	1.81	0.44
1:C:167:ILE:HB	1:C:294:HIS:CE1	2.53	0.44
1:C:239:PHE:O	1:C:250:GLU:HA	2.17	0.44
2:F:90:VAL:HG13	6:F:345:HOH:O	2.17	0.44
1:A:334:ILE:O	1:A:338:GLU:HG3	2.18	0.44
1:C:273:ASN:HA	1:C:274:PRO:HA	1.86	0.44
1:B:159:GLU:HG2	1:B:313:TYR:HD2	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:306:ASN:C	1:B:310:ARG:NH1	2.76	0.43
1:A:38:ILE:CD1	1:A:55:VAL:HG12	2.48	0.43
1:B:215:ARG:NH2	6:B:836:HOH:O	2.52	0.43
1:C:168:LYS:HE2	6:C:794:HOH:O	2.18	0.43
1:A:108:LEU:HD12	1:A:108:LEU:N	2.33	0.43
1:C:171:TRP:CG	1:C:288:PRO:HG3	2.53	0.43
1:A:332:ASP:HA	1:A:335:TRP:NE1	2.34	0.43
1:B:94:ALA:HB1	1:B:108:LEU:HB2	2.01	0.43
2:D:95:VAL:HA	2:D:99:GLU:O	2.19	0.43
1:A:287:VAL:HA	1:A:288:PRO:HD3	1.89	0.43
2:F:4:ILE:HG23	2:F:5:TRP:N	2.34	0.43
2:F:95:VAL:HA	2:F:99:GLU:O	2.19	0.43
1:B:321:TRP:HA	1:B:321:TRP:CE3	2.54	0.42
1:A:168:LYS:HA	1:A:293:THR:HG22	2.01	0.42
1:B:310:ARG:O	1:B:314:GLU:HG3	2.19	0.42
1:B:343:PHE:CG	1:C:73:GLY:HA3	2.55	0.42
1:A:322:LYS:HB3	1:A:323:PRO:HD3	2.02	0.42
1:A:183:HIS:O	1:A:186:ILE:HG12	2.19	0.42
1:B:38:ILE:HG23	1:B:57:ARG:NH2	2.31	0.42
1:A:64:CYS:HB2	1:A:126:VAL:HG21	2.01	0.42
1:B:151:PRO:O	1:B:152:ASN:HB2	2.19	0.42
1:C:204:PHE:O	1:C:206:PRO:HD3	2.19	0.42
1:C:215:ARG:HB2	1:C:230:LEU:HD11	2.02	0.42
1:C:276:PRO:HD3	1:C:300:LEU:HD21	2.02	0.42
1:C:317:PHE:HA	1:C:321:TRP:HB2	2.02	0.42
1:C:230:LEU:HA	6:C:873:HOH:O	2.20	0.42
1:B:207:GLY:HA3	1:B:234:HIS:O	2.19	0.41
1:B:389:HIS:HB3	1:B:390:HIS:H	1.59	0.41
1:C:108:LEU:N	1:C:108:LEU:HD12	2.34	0.41
1:C:322:LYS:HB3	1:C:323:PRO:HD3	2.02	0.41
1:A:218:ASP:OD1	1:A:218:ASP:C	2.63	0.41
1:A:312:LYS:HE2	1:A:316:GLU:OE2	2.20	0.41
1:B:285:TRP:HB2	1:B:297:PHE:HB3	2.02	0.41
1:B:390:HIS:ND1	1:B:390:HIS:C	2.79	0.41
1:C:276:PRO:HD2	1:C:300:LEU:HD21	2.03	0.41
1:B:162:GLY:HA3	1:B:298:GLN:O	2.20	0.41
1:C:184:ILE:HD11	1:C:200:LEU:HD11	2.01	0.41
2:E:68:HIS:HB2	4:E:201:FES:S2	2.60	0.41
1:A:322:LYS:HB3	1:A:323:PRO:CD	2.51	0.41
1:C:207:GLY:HA3	1:C:234:HIS:O	2.21	0.41
1:C:322:LYS:HB3	1:C:323:PRO:CD	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:357:GLU:H	1:B:357:GLU:CD	2.28	0.41
1:B:390:HIS:ND1	1:B:390:HIS:O	2.54	0.41
1:A:269:VAL:HB	1:A:285:TRP:CD1	2.55	0.41
1:B:65:LEU:CD2	1:B:123:THR:HG22	2.49	0.41
1:A:210:ARG:HG3	1:A:360:GLU:OE2	2.21	0.41
1:A:389:HIS:HB2	6:A:918:HOH:O	2.21	0.41
1:C:16:ALA:HB3	1:C:17:PRO:HD3	2.03	0.41
1:B:280:MET:HE3	1:B:313:TYR:CE1	2.56	0.40
1:B:175:VAL:HG22	1:B:286:TYR:CD2	2.56	0.40
1:B:228:TYR:CD1	1:B:263:SER:HB3	2.56	0.40
1:C:18:TYR:CE2	1:C:366:ARG:HG2	2.56	0.40
1:B:136:TYR:CE1	1:B:143:PRO:HD2	2.56	0.40
1:B:346:ASP:O	1:B:347:ASP:HB2	2.22	0.40
1:C:191:ILE:O	1:C:195:ASP:HB2	2.21	0.40
2:E:31:ALA:O	2:E:41:ALA:HA	2.22	0.40
1:B:130:LYS:HB3	1:B:160:ILE:HB	2.03	0.40
1:B:196:ASN:O	1:B:197:ASP:C	2.65	0.40
1:C:175:VAL:HG22	1:C:286:TYR:CD2	2.56	0.40
1:C:26:ARG:O	1:C:27:ASN:HB2	2.20	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	386/392 (98%)	369 (96%)	15 (4%)	2 (0%)	24	20
1	B	388/392 (99%)	361 (93%)	25 (6%)	2 (0%)	24	20
1	C	386/392 (98%)	367 (95%)	17 (4%)	2 (0%)	24	20
2	D	102/115 (89%)	96 (94%)	6 (6%)	0	100	100
2	E	110/115 (96%)	105 (96%)	5 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	F	102/115 (89%)	99 (97%)	3 (3%)	0	100	100
All	All	1474/1521 (97%)	1397 (95%)	71 (5%)	6 (0%)	30	26

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	268	GLY
1	B	268	GLY
1	C	268	GLY
1	C	71	HIS
1	A	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	335/339 (99%)	327 (98%)	8 (2%)	43	47
1	B	337/339 (99%)	332 (98%)	5 (2%)	57	64
1	C	335/339 (99%)	328 (98%)	7 (2%)	47	52
2	D	82/93 (88%)	82 (100%)	0	100	100
2	E	90/93 (97%)	89 (99%)	1 (1%)	65	72
2	F	82/93 (88%)	81 (99%)	1 (1%)	63	70
All	All	1261/1296 (97%)	1239 (98%)	22 (2%)	53	60

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	LEU
1	A	137	LEU
1	A	211	LYS
1	A	283	PHE
1	A	300	LEU

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Mol	Chain	Res	Type
1	A	314	GLU
1	A	356	PHE
1	A	368	LEU
1	B	46	LEU
1	B	137	LEU
1	B	300	LEU
1	B	356	PHE
1	B	368	LEU
1	C	38	ILE
1	C	46	LEU
1	C	137	LEU
1	C	169	SER
1	C	300	LEU
1	C	356	PHE
1	C	368	LEU
2	E	6	LEU
2	F	107	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	213	GLN
1	A	273	ASN
1	A	315	GLN
1	B	52	ASN
1	B	119	GLN
1	B	298	GLN
1	B	387	HIS
1	B	388	HIS
1	C	127	GLN
1	C	388	HIS
2	E	15	GLN
2	E	112	HIS

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	$\# Z > 2$
5	9CA	B	503	-	15,15,15	1.37	3 (20%)	21,21,21	1.00	0
4	FES	E	201	2	0,4,4	-	-	-		
4	FES	C	502	1	0,4,4	-	-	-		
5	9CA	C	503	-	15,15,15	1.38	3 (20%)	21,21,21	1.00	1 (4%)
4	FES	D	201	2	0,4,4	-	-	-		
4	FES	A	502	1	0,4,4	-	-	-		
4	FES	B	502	1	0,4,4	-	-	-		
4	FES	F	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
5	9CA	B	503	-	-	-	0/3/3/3
4	FES	E	201	2	-	-	0/1/1/1
4	FES	C	502	1	-	-	0/1/1/1
5	9CA	C	503	-	-	-	0/3/3/3
4	FES	D	201	2	-	-	0/1/1/1
4	FES	A	502	1	-	-	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	FES	B	502	1	-	-	0/1/1/1
4	FES	F	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	B	503	9CA	C4B-C4A	-2.84	1.39	1.46
5	C	503	9CA	C9A-N9	-2.83	1.33	1.38
5	C	503	9CA	C4B-C4A	-2.82	1.39	1.46
5	B	503	9CA	C9A-N9	-2.80	1.33	1.38
5	B	503	9CA	C8A-N9	-2.79	1.33	1.38
5	C	503	9CA	C8A-N9	-2.77	1.33	1.38

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	503	9CA	C8-C8A-N9	2.01	131.93	129.07

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	503	9CA	1	0
4	E	201	FES	1	0
4	C	502	FES	1	0
4	A	502	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	388/392 (98%)	-0.26	7 (1%) 67 71	11, 21, 38, 48	0
1	B	390/392 (99%)	0.14	17 (4%) 39 44	12, 26, 46, 53	0
1	C	388/392 (98%)	0.00	12 (3%) 51 55	14, 25, 41, 47	0
2	D	104/115 (90%)	0.63	4 (3%) 44 49	17, 37, 48, 54	0
2	E	112/115 (97%)	0.40	8 (7%) 22 25	20, 32, 44, 48	0
2	F	104/115 (90%)	0.55	3 (2%) 53 57	18, 32, 44, 51	0
All	All	1486/1521 (97%)	0.08	51 (3%) 48 52	11, 26, 45, 54	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	389	HIS	5.5
1	B	235	GLY	5.3
1	C	235	GLY	4.9
2	F	4	ILE	4.6
1	B	205	ALA	3.9
2	E	114	HIS	3.9
2	E	110	HIS	3.8
1	B	204	PHE	3.5
1	C	389	HIS	3.5
2	E	112	HIS	3.5
2	D	4	ILE	3.4
1	A	221	VAL	3.4
1	A	389	HIS	3.4
1	B	206	PRO	3.4
2	E	113	HIS	3.4
1	B	210	ARG	3.1
1	A	388	HIS	3.1
1	C	205	ALA	3.0
1	A	220	ASP	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	204	PHE	3.0
1	B	221	VAL	2.9
1	B	209	ASP	2.9
1	B	305	ALA	2.9
2	F	25	VAL	2.8
1	A	210	ARG	2.8
1	B	390	HIS	2.7
1	A	218	ASP	2.7
1	C	388	HIS	2.6
1	B	388	HIS	2.6
2	E	37	ASP	2.5
2	E	108	LEU	2.5
1	B	234	HIS	2.5
1	C	236	VAL	2.5
2	D	6	LEU	2.4
2	F	107	LYS	2.4
2	D	101	TYR	2.3
1	A	2	ALA	2.3
1	C	210	ARG	2.3
1	B	391	HIS	2.2
2	E	3	GLN	2.2
1	B	218	ASP	2.2
1	B	222	VAL	2.2
2	E	111	HIS	2.2
1	B	157	ASP	2.1
1	C	307	ASP	2.1
2	D	98	GLY	2.1
1	C	203	GLY	2.1
1	B	236	VAL	2.1
1	C	234	HIS	2.0
1	C	232	GLY	2.0
1	C	230	LEU	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 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(Å ²)	Q<0.9
5	9CA	B	503	13/13	0.79	0.14	39,40,42,42	0
5	9CA	C	503	13/13	0.82	0.14	36,39,41,41	0
4	FES	A	502	4/4	0.99	0.03	15,17,17,20	0
4	FES	B	502	4/4	0.99	0.03	11,12,13,19	0
4	FES	C	502	4/4	0.99	0.03	20,24,26,27	0
4	FES	D	201	4/4	0.99	0.03	16,19,19,20	0
4	FES	E	201	4/4	0.99	0.03	18,19,20,21	0
3	FE2	A	501	1/1	0.99	0.02	19,19,19,19	0
3	FE2	B	501	1/1	0.99	0.03	37,37,37,37	0
3	FE2	C	501	1/1	1.00	0.04	27,27,27,27	0
4	FES	F	201	4/4	1.00	0.03	18,18,18,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.