



# Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 12:36 pm BST

PDB ID : 3VMH  
Title : Oxygen-bound complex between oxygenase and ferredoxin in carbazole 1,9a-dioxygenase  
Authors : Ashikawa, Y.; Nojiri, H.  
Deposited on : 2011-12-12  
Resolution : 1.85 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

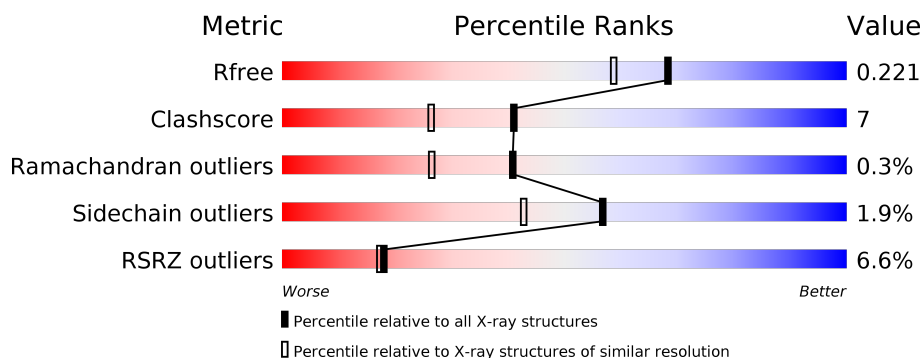
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.85 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))
$R_{free}$	130704	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

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  $\geq 3$ , 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	<div> <div>3%</div> <div>81%</div> <div>18%</div> <div>..</div> </div>
1	B	392	<div> <div>3%</div> <div>82%</div> <div>17%</div> <div>.</div> </div>
1	C	392	<div> <div>3%</div> <div>82%</div> <div>17%</div> <div>.</div> </div>
2	D	115	<div> <div>35%</div> <div>80%</div> <div>10%</div> <div>10%</div> </div>
2	E	115	<div> <div>5%</div> <div>83%</div> <div>10%</div> <div>7%</div> </div>
2	F	115	<div> <div>14%</div> <div>77%</div> <div>13%</div> <div>10%</div> </div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 13011 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	390	Total	C	N	O	S	0	0	0
			3146	2010	538	584	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			

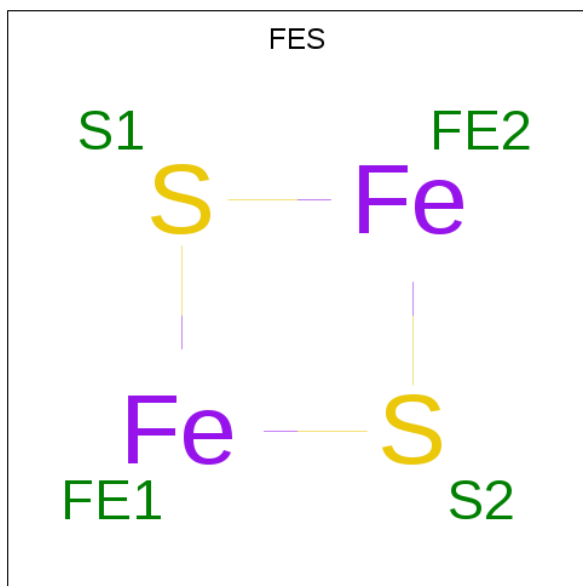
There are 24 discrepancies between the modelled and reference sequences:

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 (three-letter code: FE2) (formula: Fe).

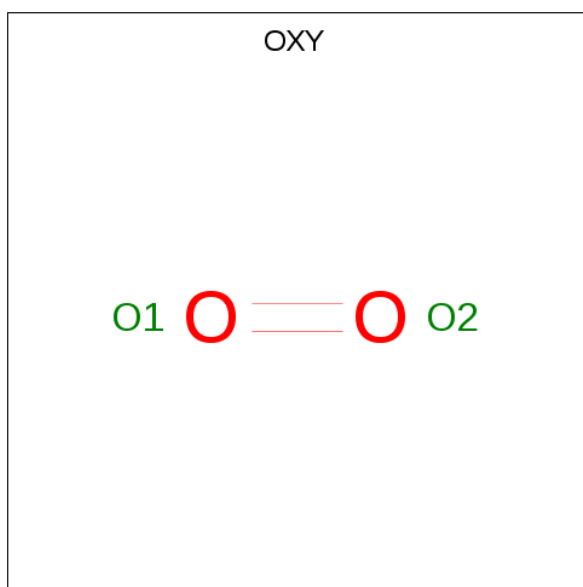
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Fe 1 1	0	0
3	A	1	Total Fe 1 1	0	0
3	C	1	Total Fe 1 1	0	0

- Molecule 4 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe<sub>2</sub>S<sub>2</sub>).



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 OXYGEN MOLECULE (three-letter code: OXY) (formula: O<sub>2</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total O 2 2	0	0
5	B	1	Total O 2 2	0	0
5	C	1	Total O 2 2	0	0

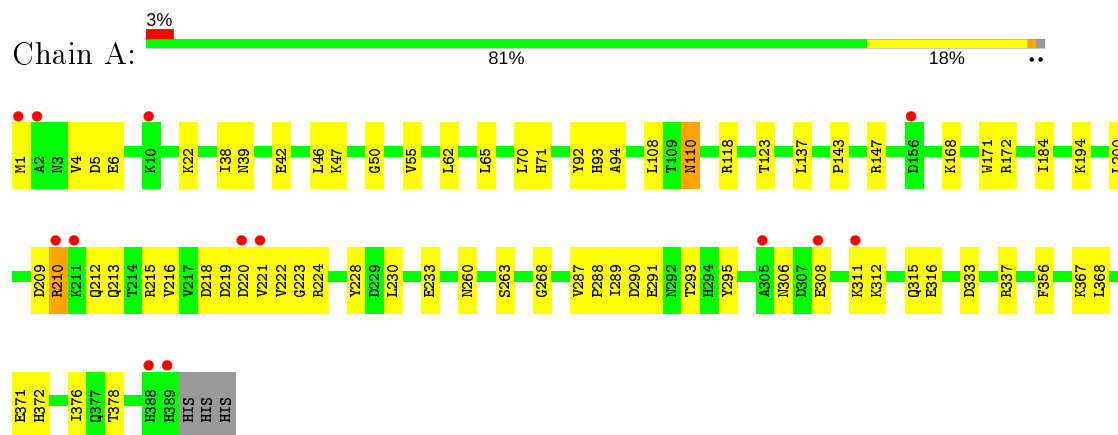
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	364	Total O 364 364	0	0
6	B	362	Total O 362 362	0	0
6	C	345	Total O 345 345	0	0
6	D	43	Total O 43 43	0	0
6	E	59	Total O 59 59	0	0
6	F	57	Total O 57 57	0	0

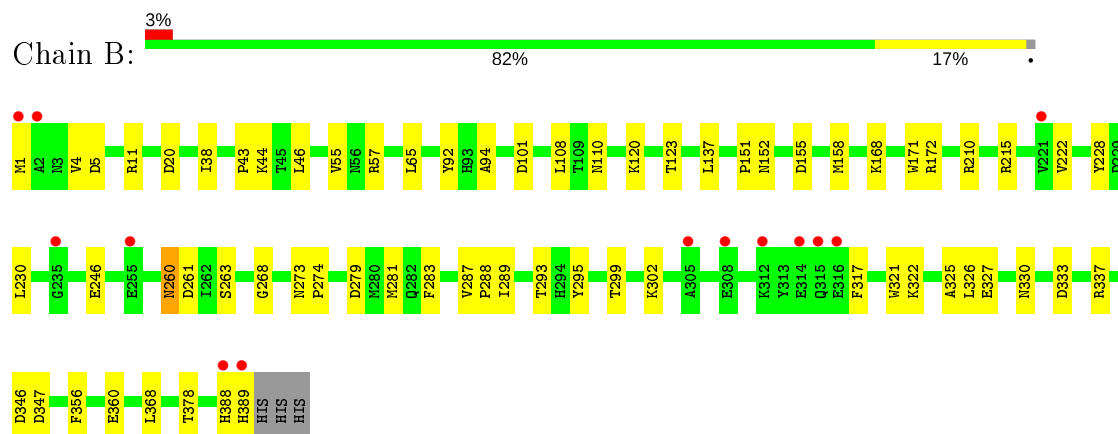
### 3 Residue-property plots [i](#)

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

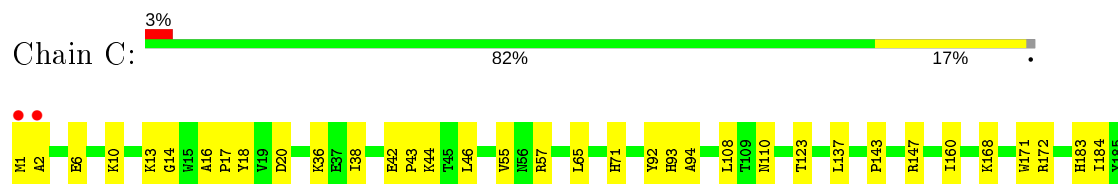
- Molecule 1: Terminal oxygenase component of carbazole

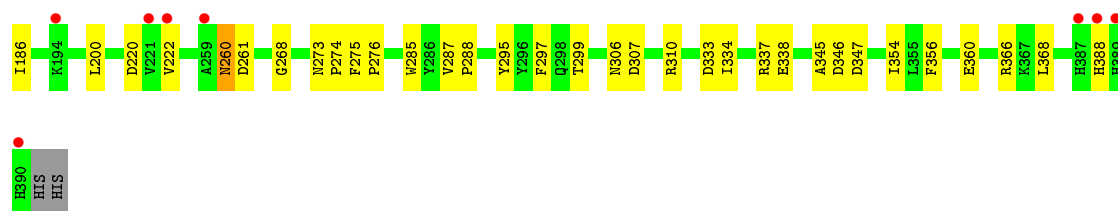


- Molecule 1: Terminal oxygenase component of carbazole

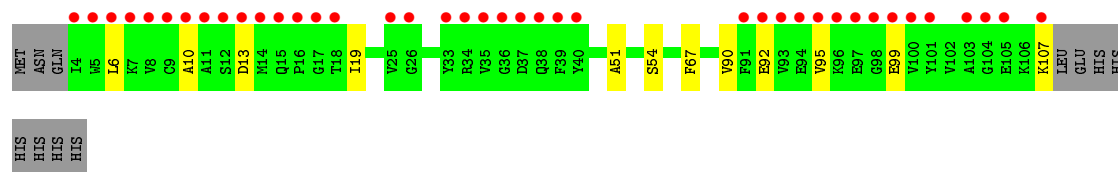
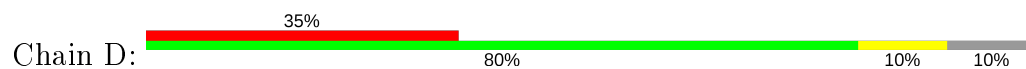


- Molecule 1: Terminal oxygenase component of carbazole

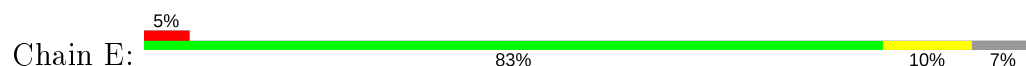




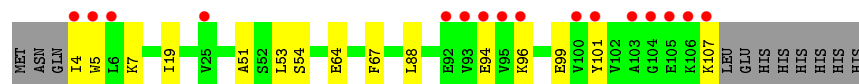
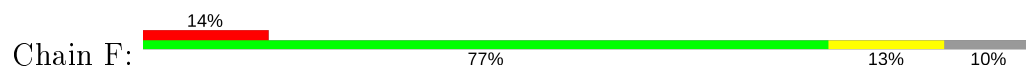
- Molecule 2: Ferredoxin component of carbazole



- 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, $\alpha$ , $\beta$ , $\gamma$	98.17Å 89.36Å 105.00Å 90.00° 104.10° 90.00°	Depositor
Resolution (Å)	47.61 – 1.85 47.61 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.9 (47.61-1.85) 99.8 (47.61-1.85)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.42 (at 1.84Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.196 , 0.225 0.192 , 0.221	Depositor DCC
$R_{free}$ test set	7432 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	28.2	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 46.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	13011	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.83% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.32	0/3221	0.60	0/4372
1	B	0.30	0/3221	0.60	0/4372
1	C	0.31	0/3232	0.60	0/4387
2	D	0.29	0/784	0.56	0/1066
2	E	0.30	0/810	0.58	0/1101
2	F	0.31	0/784	0.60	0/1066
All	All	0.31	0/12052	0.60	0/16364

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3136	0	3042	48	0
1	B	3136	0	3042	45	0
1	C	3146	0	3049	41	0
2	D	768	0	745	9	0
2	E	794	0	770	10	0
2	F	768	0	745	16	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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	0	0
4	F	4	0	0	0	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
5	C	2	0	0	0	0
6	A	364	0	0	8	0
6	B	362	0	0	5	0
6	C	345	0	0	5	0
6	D	43	0	0	0	0
6	E	59	0	0	0	0
6	F	57	0	0	1	0
All	All	13011	0	11393	165	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 (165) 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:230:LEU:HB3	1:A:233:GLU:HG3	1.58	0.86
1:C:6:GLU:O	1:C:10:LYS:HG2	1.76	0.85
2:F:94:GLU:HG2	2:F:96:LYS:HE2	1.60	0.84
1:A:289:ILE:HB	1:A:293:THR:HG23	1.63	0.81
1:C:1:MET:HA	1:C:20:ASP:OD2	1.82	0.79
1:C:65:LEU:HD23	1:C:123:THR:HG22	1.66	0.78
2:F:94:GLU:CG	2:F:96:LYS:HE2	2.14	0.77
1:A:118:ARG:HG3	6:A:1199:HOH:O	1.86	0.75
1:A:290:ASP:OD1	1:A:293:THR:HG22	1.88	0.71
1:C:220:ASP:OD1	1:C:222:VAL:HG22	1.90	0.71
1:B:120:LYS:HE2	6:B:437:HOH:O	1.91	0.69
2:F:4:ILE:HG12	2:F:5:TRP:N	2.07	0.67
1:A:65:LEU:HD23	1:A:123:THR:HG22	1.76	0.67
1:B:65:LEU:HD23	1:B:123:THR:HG22	1.78	0.66
1:B:168:LYS:HE2	6:B:1203:HOH:O	1.97	0.65
2:D:10:ALA:HB3	2:D:13:ASP:OD1	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:90:VAL:HG23	2:D:107:LYS:HG2	1.79	0.63
1:A:312:LYS:HE2	1:A:316:GLU:OE2	1.99	0.63
1:C:360:GLU:HG3	6:C:484:HOH:O	1.98	0.62
1:C:42:GLU:HG2	1:C:44:LYS:HE3	1.82	0.62
1:A:194:LYS:HB2	6:A:627:HOH:O	1.98	0.62
1:A:216:VAL:HG13	1:A:368:LEU:HD12	1.82	0.60
1:B:101:ASP:O	1:B:120:LYS:HE3	2.01	0.60
1:A:94:ALA:HB1	1:A:108:LEU:HB2	1.84	0.60
2:F:4:ILE:HG12	2:F:5:TRP:H	1.66	0.60
2:F:7:LYS:NZ	2:F:99:GLU:HG2	2.16	0.59
1:B:94:ALA:HB1	1:B:108:LEU:HB2	1.84	0.59
2:D:107:LYS:O	2:D:107:LYS:HG3	2.03	0.59
1:B:210:ARG:HH22	2:E:52:SER:HB2	1.68	0.59
1:B:360:GLU:HG3	6:B:491:HOH:O	2.02	0.58
1:B:260:ASN:HD22	1:B:261:ASP:N	2.01	0.58
2:F:4:ILE:HG23	2:F:5:TRP:H	1.70	0.57
1:B:38:ILE:HG23	1:B:57:ARG:HH21	1.70	0.56
1:C:36:LYS:HE3	6:C:1180:HOH:O	2.04	0.56
1:B:346:ASP:O	1:B:347:ASP:HB2	2.06	0.56
1:A:1:MET:N	1:A:378:THR:HG22	2.21	0.56
1:C:306:ASN:O	1:C:310:ARG:HG2	2.05	0.55
1:B:281:MET:HE3	1:B:283:PHE:CZ	2.41	0.55
2:D:51:ALA:HB2	2:D:67:PHE:CG	2.42	0.55
1:A:215:ARG:HB2	1:A:230:LEU:HD11	1.88	0.55
2:D:90:VAL:HG23	2:D:107:LYS:CG	2.37	0.54
1:A:220:ASP:HB2	1:A:224:ARG:HB2	1.90	0.54
2:E:4:ILE:HD12	2:E:4:ILE:N	2.22	0.54
1:A:287:VAL:HB	1:A:295:TYR:HB2	1.89	0.53
2:F:53:LEU:CD1	2:F:88:LEU:HD11	2.39	0.53
1:C:2:ALA:HA	6:C:1033:HOH:O	2.07	0.53
1:A:94:ALA:CB	1:A:108:LEU:HB2	2.39	0.52
1:A:1:MET:H2	1:A:378:THR:HG22	1.73	0.52
2:F:94:GLU:HG3	2:F:96:LYS:HE2	1.91	0.52
1:B:333:ASP:O	1:B:337:ARG:HG3	2.10	0.52
1:B:283:PHE:HB2	1:B:299:THR:OG1	2.10	0.51
1:C:260:ASN:C	1:C:260:ASN:HD22	2.12	0.51
1:B:260:ASN:C	1:B:260:ASN:HD22	2.14	0.51
1:A:1:MET:H1	1:A:376:ILE:HG21	1.76	0.51
1:B:287:VAL:HB	1:B:295:TYR:HB2	1.93	0.51
1:A:367:LYS:O	1:A:371:GLU:HG3	2.11	0.51
1:C:333:ASP:O	1:C:337:ARG:HG3	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:38:ILE:HG23	1:C:57:ARG:HH21	1.75	0.51
1:A:39:ASN:HB2	1:A:42:GLU:OE1	2.11	0.51
1:C:1:MET:SD	1:C:1:MET:N	2.83	0.51
2:E:4:ILE:H	2:E:4:ILE:CD1	2.24	0.51
1:A:168:LYS:HE2	6:A:1104:HOH:O	2.10	0.50
1:A:38:ILE:HD13	1:A:62:LEU:HD22	1.92	0.50
1:A:312:LYS:O	1:A:316:GLU:HG3	2.12	0.50
1:B:44:LYS:HE3	6:B:739:HOH:O	2.11	0.50
2:D:92:GLU:HG2	2:D:107:LYS:HE2	1.94	0.50
1:A:1:MET:H1	1:A:376:ILE:CG2	2.25	0.50
1:B:11:ARG:HH21	1:C:388:HIS:CE1	2.30	0.49
1:B:171:TRP:CE2	1:B:172:ARG:HG3	2.46	0.49
1:B:94:ALA:CB	1:B:108:LEU:HB2	2.42	0.49
1:B:171:TRP:CG	1:B:288:PRO:HG3	2.47	0.49
1:B:215:ARG:HB3	1:B:228:TYR:HB2	1.95	0.49
1:C:42:GLU:HG2	1:C:44:LYS:CE	2.43	0.48
2:F:4:ILE:HG23	2:F:5:TRP:N	2.28	0.48
1:A:171:TRP:CE2	1:A:172:ARG:HG3	2.49	0.48
2:E:4:ILE:HD12	2:E:4:ILE:H	1.78	0.48
2:F:107:LYS:HG2	2:F:107:LYS:O	2.13	0.48
1:A:118:ARG:HD2	6:A:1152:HOH:O	2.12	0.48
1:C:42:GLU:O	1:C:44:LYS:HE3	2.13	0.48
1:C:171:TRP:CE2	1:C:172:ARG:HG3	2.49	0.48
2:D:6:LEU:C	2:D:6:LEU:HD23	2.34	0.48
1:C:168:LYS:HE2	6:C:1202:HOH:O	2.14	0.47
1:C:18:TYR:CE2	1:C:366:ARG:HG2	2.50	0.47
1:C:261:ASP:HB3	1:C:273:ASN:O	2.13	0.47
1:A:184:ILE:HD11	1:A:200:LEU:CD1	2.45	0.47
1:B:222:VAL:HG12	1:B:222:VAL:O	2.13	0.47
1:C:287:VAL:HB	1:C:295:TYR:HB2	1.95	0.47
2:E:47:THR:HG23	2:E:88:LEU:HD23	1.97	0.47
2:E:7:LYS:HE3	2:E:99:GLU:OE2	2.15	0.47
1:A:47:LYS:HE3	1:A:50:GLY:HA2	1.97	0.47
1:A:118:ARG:HG3	1:A:118:ARG:O	2.15	0.47
1:B:215:ARG:HB2	1:B:230:LEU:HD11	1.97	0.46
1:A:1:MET:HG3	6:A:788:HOH:O	2.15	0.46
1:C:94:ALA:HB1	1:C:108:LEU:HB2	1.98	0.46
1:C:14:GLY:HA3	6:F:786:HOH:O	2.16	0.46
1:C:184:ILE:HD11	1:C:200:LEU:CD1	2.44	0.46
2:E:51:ALA:HB2	2:E:67:PHE:CG	2.50	0.46
2:F:19:ILE:HG21	2:F:54:SER:HA	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:ARG:HA	1:A:213:GLN:HG2	1.98	0.46
1:B:273:ASN:OD1	1:B:281:MET:HG2	2.15	0.46
1:B:279:ASP:OD2	1:B:302:LYS:NZ	2.49	0.46
1:C:108:LEU:N	1:C:108:LEU:HD12	2.31	0.46
1:C:13:LYS:NZ	2:F:64:GLU:OE2	2.45	0.46
1:C:43:PRO:HA	1:C:55:VAL:O	2.16	0.45
1:A:6:GLU:HG3	6:A:666:HOH:O	2.14	0.45
1:A:333:ASP:O	1:A:337:ARG:HG3	2.16	0.45
2:D:95:VAL:HA	2:D:99:GLU:O	2.16	0.45
1:C:171:TRP:CG	1:C:288:PRO:HG3	2.51	0.45
1:A:228:TYR:CD1	1:A:263:SER:HB3	2.51	0.45
1:B:322:LYS:O	1:B:327:GLU:HG3	2.17	0.45
2:F:99:GLU:HB3	2:F:101:TYR:HE1	1.81	0.45
1:A:218:ASP:OD1	1:A:372:HIS:CE1	2.69	0.45
1:A:290:ASP:CG	1:A:293:THR:HG22	2.37	0.45
2:F:51:ALA:HB2	2:F:67:PHE:CG	2.52	0.45
2:F:7:LYS:HZ2	2:F:99:GLU:HG2	1.80	0.44
1:B:326:LEU:O	1:B:330:ASN:HB2	2.18	0.44
1:C:307:ASP:HA	1:C:310:ARG:CG	2.47	0.44
1:B:321:TRP:O	1:B:325:ALA:HB3	2.18	0.44
1:A:222:VAL:HG23	1:A:224:ARG:HG3	1.98	0.44
1:B:302:LYS:HE2	6:B:840:HOH:O	2.17	0.44
1:C:346:ASP:O	1:C:347:ASP:HB2	2.18	0.44
1:A:143:PRO:HG3	1:A:147:ARG:CZ	2.48	0.44
1:A:311:LYS:O	1:A:315:GLN:HG3	2.17	0.44
1:A:38:ILE:CD1	1:A:55:VAL:HG12	2.48	0.44
1:B:289:ILE:HB	1:B:293:THR:OG1	2.18	0.44
1:A:219:ASP:OD1	1:A:223:GLY:HA2	2.18	0.44
1:C:143:PRO:HG3	1:C:147:ARG:CZ	2.48	0.43
2:E:99:GLU:OE1	2:E:99:GLU:HA	2.18	0.43
1:C:273:ASN:HA	1:C:274:PRO:HA	1.90	0.43
1:B:155:ASP:HB2	1:B:158:MET:HB2	2.01	0.43
1:C:183:HIS:O	1:C:186:ILE:HG12	2.19	0.43
1:B:287:VAL:HA	1:B:288:PRO:HD3	1.91	0.43
1:B:151:PRO:O	1:B:152:ASN:HB2	2.18	0.43
1:A:93:HIS:HB2	4:A:401:FES:S1	2.58	0.43
1:C:16:ALA:HB3	1:C:17:PRO:HD3	2.01	0.43
2:E:4:ILE:N	2:E:4:ILE:CD1	2.81	0.43
1:B:317:PHE:HA	1:B:321:TRP:HB2	2.00	0.42
1:C:160:ILE:HG23	1:C:299:THR:HB	2.01	0.42
1:B:38:ILE:HG23	1:B:38:ILE:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:306:ASN:OD1	1:A:308:GLU:HB2	2.19	0.42
1:B:273:ASN:HA	1:B:274:PRO:HA	1.88	0.42
1:B:261:ASP:HB3	1:B:273:ASN:O	2.20	0.42
1:B:1:MET:HB3	1:B:378:THR:HG21	2.01	0.42
1:A:220:ASP:OD1	1:A:221:VAL:N	2.48	0.42
2:F:99:GLU:HB3	2:F:101:TYR:CE1	2.55	0.42
1:B:43:PRO:HA	1:B:55:VAL:O	2.20	0.41
1:B:1:MET:N	1:B:20:ASP:OD2	2.34	0.41
1:C:285:TRP:HB2	1:C:297:PHE:HB3	2.02	0.41
1:A:287:VAL:HA	1:A:288:PRO:HD3	1.91	0.41
1:A:291:GLU:HG2	6:A:489:HOH:O	2.20	0.41
1:A:209:ASP:OD1	1:A:212:GLN:HG3	2.21	0.41
1:B:388:HIS:O	1:B:389:HIS:HB3	2.21	0.41
2:D:19:ILE:HG21	2:D:54:SER:HA	2.03	0.41
2:E:10:ALA:HB3	2:E:13:ASP:OD2	2.21	0.41
1:B:228:TYR:CD1	1:B:263:SER:HB3	2.56	0.40
1:A:4:VAL:HG12	1:A:5:ASP:N	2.35	0.40
6:A:520:HOH:O	1:B:388:HIS:HE1	2.04	0.40
1:C:345:ALA:HA	6:C:1178:HOH:O	2.20	0.40
1:A:70:LEU:HB3	1:C:354:ILE:HD12	2.02	0.40
1:A:110:ASN:C	1:A:110:ASN:HD22	2.25	0.40
1:B:4:VAL:HG12	1:B:5:ASP:N	2.35	0.40
1:C:275:PHE:CG	1:C:276:PRO:HA	2.56	0.40
1:C:334:ILE:O	1:C:338:GLU:HG3	2.22	0.40
1:C:93:HIS:HB2	4:C:401:FES:S1	2.61	0.40
1:B:260:ASN:ND2	1:B:260:ASN:H	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	387/392 (99%)	368 (95%)	17 (4%)	2 (0%)	29	15
1	B	387/392 (99%)	367 (95%)	19 (5%)	1 (0%)	41	26
1	C	388/392 (99%)	369 (95%)	17 (4%)	2 (0%)	29	15
2	D	102/115 (89%)	94 (92%)	8 (8%)	0	100	100
2	E	105/115 (91%)	100 (95%)	5 (5%)	0	100	100
2	F	102/115 (89%)	99 (97%)	3 (3%)	0	100	100
All	All	1471/1521 (97%)	1397 (95%)	69 (5%)	5 (0%)	41	26

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	268	GLY
1	C	268	GLY
1	A	268	GLY
1	C	71	HIS
1	A	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%)	328 (98%)	8 (2%)	49	33
1	B	336/339 (99%)	328 (98%)	8 (2%)	49	33
1	C	337/339 (99%)	330 (98%)	7 (2%)	53	38
2	D	82/93 (88%)	82 (100%)	0	100	100
2	E	85/93 (91%)	84 (99%)	1 (1%)	71	62
2	F	82/93 (88%)	82 (100%)	0	100	100
All	All	1258/1296 (97%)	1234 (98%)	24 (2%)	57	43

All (24) residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Type
1	A	22	LYS
1	A	46	LEU
1	A	92	TYR
1	A	110	ASN
1	A	137	LEU
1	A	210	ARG
1	A	260	ASN
1	A	356	PHE
1	B	46	LEU
1	B	92	TYR
1	B	110	ASN
1	B	137	LEU
1	B	246	GLU
1	B	260	ASN
1	B	356	PHE
1	B	368	LEU
1	C	46	LEU
1	C	92	TYR
1	C	110	ASN
1	C	137	LEU
1	C	260	ASN
1	C	356	PHE
1	C	368	LEU
2	E	6	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	110	ASN
1	A	165	GLN
1	A	260	ASN
1	B	110	ASN
1	B	165	GLN
1	B	260	ASN
1	C	110	ASN
1	C	165	GLN
1	C	260	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 ⓘ

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 12 ligands modelled in this entry, 3 are monoatomic - leaving 9 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	B	401	1	0,4,4	0.00	-	-		
5	OXY	B	602	3	1,1,1	1.18	0	-		
4	FES	C	401	1	0,4,4	0.00	-	-		
5	OXY	A	601	3	1,1,1	1.20	0	-		
5	OXY	C	603	3	1,1,1	1.17	0	-		
4	FES	A	401	1	0,4,4	0.00	-	-		
4	FES	D	201	2	0,4,4	0.00	-	-		
4	FES	F	201	2	0,4,4	0.00	-	-		
4	FES	E	201	2	0,4,4	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
4	FES	E	201	2	-	-	0/1/1/1

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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
4	FES	B	401	1	-	-	0/1/1/1
4	FES	F	201	2	-	-	0/1/1/1
4	FES	C	401	1	-	-	0/1/1/1

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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	401	FES	1	0
4	A	401	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	389/392 (99%)	0.02	13 (3%) 46 44	18, 27, 41, 48	0
1	B	389/392 (99%)	0.10	13 (3%) 46 44	19, 30, 43, 48	0
1	C	390/392 (99%)	-0.01	10 (2%) 56 54	21, 30, 40, 52	0
2	D	104/115 (90%)	1.68	40 (38%) 0 0	25, 40, 47, 50	0
2	E	107/115 (93%)	0.32	6 (5%) 24 23	25, 36, 43, 47	0
2	F	104/115 (90%)	0.68	16 (15%) 2 2	26, 37, 43, 48	0
All	All	1483/1521 (97%)	0.22	98 (6%) 18 17	18, 30, 44, 52	0

All (98) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	11.3
1	C	1	MET	11.1
2	D	4	ILE	7.1
1	B	1	MET	6.2
2	D	12	SER	6.1
1	A	221	VAL	6.1
2	D	101	TYR	6.0
2	D	95	VAL	5.9
2	D	5	TRP	5.8
1	C	390	HIS	5.8
2	D	93	VAL	5.8
2	D	98	GLY	5.7
2	D	104	GLY	5.6
2	F	6	LEU	5.6
2	D	97	GLU	5.2
2	D	105	GLU	4.8
2	D	25	VAL	4.8
2	D	37	ASP	4.7
2	D	99	GLU	4.7

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Mol	Chain	Res	Type	RSRZ
2	F	4	ILE	4.5
2	D	6	LEU	4.4
2	D	9	CYS	4.3
1	C	2	ALA	4.3
2	D	103	ALA	4.3
2	F	107	LYS	4.2
2	F	105	GLU	4.2
2	F	101	TYR	4.2
1	B	221	VAL	4.2
2	D	38	GLN	4.2
2	D	7	LYS	4.1
1	B	389	HIS	4.1
2	D	11	ALA	4.1
2	E	25	VAL	4.0
2	D	92	GLU	4.0
2	F	25	VAL	4.0
1	C	388	HIS	3.9
1	C	221	VAL	3.9
2	D	16	PRO	3.8
2	F	95	VAL	3.8
2	D	96	LYS	3.8
2	F	106	LYS	3.7
2	F	104	GLY	3.7
2	D	100	VAL	3.7
2	F	92	GLU	3.6
2	D	35	VAL	3.6
2	F	93	VAL	3.6
2	D	10	ALA	3.6
2	E	37	ASP	3.5
2	D	8	VAL	3.5
2	D	40	TYR	3.4
1	A	389	HIS	3.4
1	C	389	HIS	3.4
1	B	388	HIS	3.3
1	B	305	ALA	3.3
2	D	94	GLU	3.2
2	E	97	GLU	3.2
2	D	13	ASP	3.1
2	D	17	GLY	3.1
2	E	3	GLN	3.1
2	D	107	LYS	3.1
2	D	91	PHE	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	305	ALA	2.9
1	B	312	LYS	2.8
1	A	210	ARG	2.8
2	D	39	PHE	2.7
1	B	235	GLY	2.7
1	B	255	GLU	2.7
2	D	36	GLY	2.7
2	D	14	MET	2.6
1	B	308	GLU	2.6
2	D	26	GLY	2.5
1	A	2	ALA	2.5
2	F	94	GLU	2.5
2	F	5	TRP	2.5
1	A	388	HIS	2.4
2	D	15	GLN	2.4
2	F	100	VAL	2.4
1	A	220	ASP	2.4
1	B	2	ALA	2.4
2	F	103	ALA	2.4
2	D	34	ARG	2.3
1	A	311	LYS	2.3
1	A	156	ASP	2.3
1	B	314	GLU	2.3
2	E	98	GLY	2.3
1	A	308	GLU	2.3
1	B	315	GLN	2.3
2	E	12	SER	2.3
1	C	222	VAL	2.2
2	D	33	TYR	2.2
1	B	316	GLU	2.2
2	F	96	LYS	2.1
1	A	10	LYS	2.1
1	A	211	LYS	2.1
1	C	194	LYS	2.1
1	C	259	ALA	2.1
1	C	387	HIS	2.1
2	D	18	THR	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 carbohydrates in this entry.

### 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	OXY	B	602	2/2	0.91	0.08	40,40,40,41	0
5	OXY	A	601	2/2	0.93	0.13	31,31,31,34	0
5	OXY	C	603	2/2	0.97	0.09	34,34,34,35	0
4	FES	E	201	4/4	0.99	0.11	26,26,26,27	0
3	FE2	B	501	1/1	0.99	0.06	34,34,34,34	0
4	FES	A	401	4/4	0.99	0.12	21,22,22,24	0
4	FES	B	401	4/4	0.99	0.15	22,22,23,23	0
4	FES	F	201	4/4	0.99	0.08	23,25,26,26	0
4	FES	C	401	4/4	0.99	0.12	24,25,26,26	0
3	FE2	A	501	1/1	0.99	0.09	28,28,28,28	0
3	FE2	C	501	1/1	1.00	0.03	30,30,30,30	0
4	FES	D	201	4/4	1.00	0.09	25,26,26,27	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.