



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 10:02 PM GMT

PDB ID : 4ZGD  
Title : Mutant R157A of Fe-Type Nitrile Hydratase from Comamonas testosteroni Ni1  
Authors : Wu, R.; Martinez, S.; Holz, R.; Liu, D.  
Deposited on : 2015-04-22  
Resolution : 2.25 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

---

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

MolProbity : 4.02b-467  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026982  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20026982

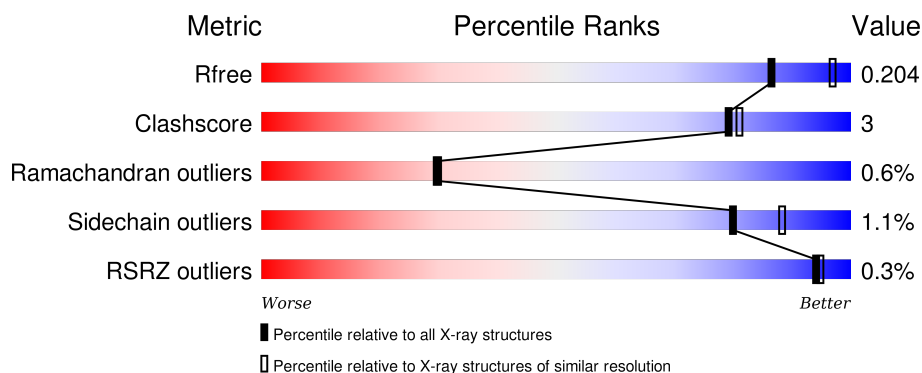
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.25 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	1640 (2.28-2.24)
Clashscore	102246	1095 (2.26-2.26)
Ramachandran outliers	100387	1063 (2.26-2.26)
Sidechain outliers	100360	1063 (2.26-2.26)
RSRZ outliers	91569	1647 (2.28-2.24)

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. 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	206	<div> <div>93%</div> <div>6%</div> </div>
1	C	206	<div> <div>93%</div> <div>6%</div> </div>
1	E	206	<div> <div>89%</div> <div>10%</div> </div>
1	G	206	<div> <div>%</div> <div>95%</div> <div>5%</div> </div>
1	I	206	<div> <div>92%</div> <div>8%</div> </div>

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Length	Quality of chain
1	K	206	 90% 8% .
1	M	206	 90% 10%
1	O	206	 90% 9% .
2	B	206	 96% .
2	D	206	 92% 7% .
2	F	206	 94% 6%
2	H	206	 95% .
2	J	206	 86% 13% .
2	L	206	 89% 10% .
2	N	206	 93% 7%
2	P	206	 92% 8%

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 28551 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 Nitrile hydratase alpha subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	206	Total	C	N	O	S	0	2	0
			1615	1031	266	309	9			
1	C	206	Total	C	N	O	S	0	0	0
			1603	1023	265	306	9			
1	E	206	Total	C	N	O	S	0	0	0
			1603	1023	265	306	9			
1	G	206	Total	C	N	O	S	0	0	0
			1603	1023	265	306	9			
1	I	206	Total	C	N	O	S	0	0	0
			1603	1023	265	306	9			
1	K	206	Total	C	N	O	S	0	0	0
			1603	1023	265	306	9			
1	M	206	Total	C	N	O	S	0	0	0
			1603	1023	265	306	9			
1	O	206	Total	C	N	O	S	0	0	0
			1603	1023	265	306	9			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	157	ALA	ARG	engineered mutation	UNP J9PBS0
C	157	ALA	ARG	engineered mutation	UNP J9PBS0
E	157	ALA	ARG	engineered mutation	UNP J9PBS0
G	157	ALA	ARG	engineered mutation	UNP J9PBS0
I	157	ALA	ARG	engineered mutation	UNP J9PBS0
K	157	ALA	ARG	engineered mutation	UNP J9PBS0
M	157	ALA	ARG	engineered mutation	UNP J9PBS0
O	157	ALA	ARG	engineered mutation	UNP J9PBS0

- Molecule 2 is a protein called Nitrile hydratase beta subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	206	Total 1589	C 1017	N 271	O 293	S 8	0	0	0
2	D	206	Total 1589	C 1017	N 271	O 293	S 8	0	0	0
2	F	206	Total 1589	C 1017	N 271	O 293	S 8	0	0	0
2	H	206	Total 1589	C 1017	N 271	O 293	S 8	0	0	0
2	J	206	Total 1589	C 1017	N 271	O 293	S 8	0	0	0
2	L	206	Total 1589	C 1017	N 271	O 293	S 8	0	0	0
2	N	206	Total 1589	C 1017	N 271	O 293	S 8	0	0	0
2	P	206	Total 1589	C 1017	N 271	O 293	S 8	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	112	PRO	ALA	conflict	UNP J9PBS1
D	112	PRO	ALA	conflict	UNP J9PBS1
F	112	PRO	ALA	conflict	UNP J9PBS1
H	112	PRO	ALA	conflict	UNP J9PBS1
J	112	PRO	ALA	conflict	UNP J9PBS1
L	112	PRO	ALA	conflict	UNP J9PBS1
N	112	PRO	ALA	conflict	UNP J9PBS1
P	112	PRO	ALA	conflict	UNP J9PBS1

- Molecule 3 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	1	Total 1	Fe 1	0	0
3	K	1	Total 1	Fe 1	0	0
3	E	1	Total 1	Fe 1	0	0
3	I	1	Total 1	Fe 1	0	0
3	C	1	Total 1	Fe 1	0	0
3	A	1	Total 1	Fe 1	0	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	O	1	Total 1	Fe 1	0	0
3	M	1	Total 1	Fe 1	0	0

- Molecule 4 is water.

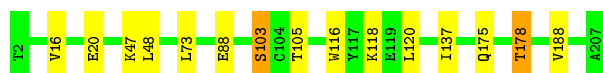
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	189	Total 189	O 189	0	0
4	B	176	Total 176	O 176	0	0
4	C	196	Total 196	O 196	0	0
4	D	191	Total 191	O 191	0	0
4	E	187	Total 187	O 187	0	0
4	F	208	Total 208	O 208	0	0
4	G	195	Total 195	O 195	0	0
4	H	182	Total 182	O 182	0	0
4	I	172	Total 172	O 172	0	0
4	J	172	Total 172	O 172	0	0
4	K	206	Total 206	O 206	0	0
4	L	181	Total 181	O 181	0	0
4	M	186	Total 186	O 186	0	0
4	N	194	Total 194	O 194	0	0
4	O	194	Total 194	O 194	0	0
4	P	166	Total 166	O 166	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 errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Nitrile hydratase alpha subunit

Chain A:  93% 6%




- Molecule 1: Nitrile hydratase alpha subunit

Chain C:  93% 6%



- Molecule 1: Nitrile hydratase alpha subunit

Chain E:  89% 10%



- Molecule 1: Nitrile hydratase alpha subunit

Chain G:  95% 5%




- Molecule 1: Nitrile hydratase alpha subunit

Chain I:  92% 8%

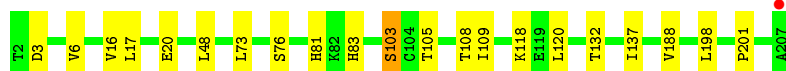


- Molecule 1: Nitrile hydratase alpha subunit

Chain K:  90% 8%



- Molecule 1: Nitrile hydratase alpha subunit



- Molecule 1: Nitrile hydratase alpha subunit



- Molecule 2: Nitrile hydratase beta subunit



- Molecule 2: Nitrile hydratase beta subunit



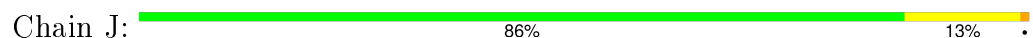
- Molecule 2: Nitrile hydratase beta subunit



- Molecule 2: Nitrile hydratase beta subunit



- Molecule 2: Nitrile hydratase beta subunit

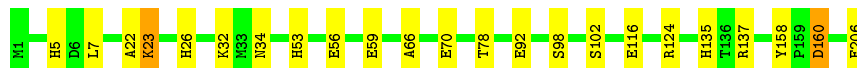






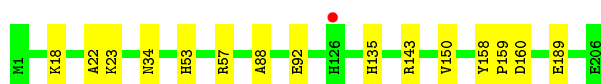
- Molecule 2: Nitrile hydratase beta subunit

Chain L: 89% 10%



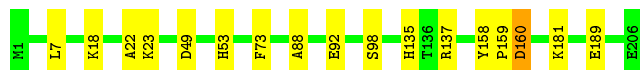
- Molecule 2: Nitrile hydratase beta subunit

Chain N: 93% 7%



- Molecule 2: Nitrile hydratase beta subunit

Chain P: 92% 8%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	111.59Å 111.59Å 474.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.49 – 2.25 50.56 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.9 (50.49-2.25) 95.3 (50.56-2.25)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.16	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.54 (at 2.25Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.172 , 0.203 0.173 , 0.204	Depositor DCC
$R_{free}$ test set	15125 reflections (5.06%)	DCC
Wilson B-factor (Å <sup>2</sup> )	22.7	Xtriage
Anisotropy	0.718	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 34.1	EDS
Estimated twinning fraction	0.467 for -h,-k,l 0.468 for h,-h-k,-l 0.469 for -k,-h,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 313637 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	28551	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.98% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CSD, FE

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.36	0/1642	0.52	0/2244
1	C	0.41	0/1624	0.54	0/2220
1	E	0.39	0/1624	0.54	0/2220
1	G	0.39	0/1624	0.54	0/2220
1	I	0.36	0/1624	0.48	0/2220
1	K	0.41	0/1624	0.57	0/2220
1	M	0.38	0/1624	0.54	0/2220
1	O	0.38	0/1624	0.51	0/2220
2	B	0.36	0/1636	0.50	0/2216
2	D	0.38	0/1636	0.50	0/2216
2	F	0.39	0/1636	0.51	0/2216
2	H	0.39	0/1636	0.52	0/2216
2	J	0.37	0/1636	0.52	0/2216
2	L	0.39	0/1636	0.52	0/2216
2	N	0.38	0/1636	0.50	0/2216
2	P	0.38	0/1636	0.50	0/2216
All	All	0.38	0/26098	0.52	0/35512

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

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	1615	0	1597	11	0
1	C	1603	0	1582	8	0
1	E	1603	0	1582	15	0
1	G	1603	0	1582	4	0
1	I	1603	0	1582	15	0
1	K	1603	0	1583	19	0
1	M	1603	0	1582	16	0
1	O	1603	0	1584	15	0
2	B	1589	0	1534	6	0
2	D	1589	0	1534	13	0
2	F	1589	0	1534	10	0
2	H	1589	0	1534	6	0
2	J	1589	0	1534	20	0
2	L	1589	0	1534	15	0
2	N	1589	0	1534	9	0
2	P	1589	0	1534	13	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
3	E	1	0	0	0	0
3	G	1	0	0	0	0
3	I	1	0	0	0	0
3	K	1	0	0	0	0
3	M	1	0	0	0	0
3	O	1	0	0	0	0
4	A	189	0	0	3	0
4	B	176	0	0	1	0
4	C	196	0	0	2	0
4	D	191	0	0	3	0
4	E	187	0	0	2	0
4	F	208	0	0	2	3
4	G	195	0	0	1	0
4	H	182	0	0	2	0
4	I	172	0	0	6	0
4	J	172	0	0	3	0
4	K	206	0	0	5	0
4	L	181	0	0	3	0
4	M	186	0	0	3	0
4	N	194	0	0	2	0
4	O	194	0	0	7	0
4	P	166	0	0	2	3
All	All	28551	0	24946	175	3

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 3.

All (175) 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:M:3:ASP:OD1	4:M:401:HOH:O	1.84	0.95
2:J:49:ASP:OD2	4:J:301:HOH:O	1.85	0.92
2:D:174:TYR:HE1	2:D:199:GLN:HG3	1.42	0.85
1:K:26:ASP:O	4:K:401:HOH:O	1.96	0.82
2:L:59:GLU:OE1	4:L:301:HOH:O	1.98	0.81
1:O:118:LYS:NZ	4:O:405:HOH:O	2.15	0.80
1:I:173:GLU:OE1	4:I:401:HOH:O	2.00	0.78
1:O:88:GLU:OE1	4:O:401:HOH:O	2.02	0.77
2:P:49:ASP:OD2	4:P:301:HOH:O	2.04	0.75
2:H:116:GLU:OE1	4:H:301:HOH:O	2.03	0.75
1:O:173:GLU:OE1	4:O:402:HOH:O	2.06	0.74
2:F:116:GLU:O	4:F:301:HOH:O	2.05	0.73
1:K:143:GLU:OE2	4:K:402:HOH:O	2.07	0.72
2:J:21:ASN:ND2	4:J:302:HOH:O	2.23	0.72
2:D:110:GLN:OE1	4:D:301:HOH:O	2.06	0.71
1:E:103:SER:HB2	1:E:118:LYS:HG2	1.74	0.69
1:O:26:ASP:OD1	4:O:403:HOH:O	2.09	0.69
2:D:174:TYR:CE1	2:D:199:GLN:HG3	2.26	0.68
2:L:116:GLU:O	4:L:302:HOH:O	2.12	0.68
2:H:116:GLU:O	4:H:302:HOH:O	2.13	0.67
1:K:103:SER:O	4:K:403:HOH:O	2.13	0.67
1:E:81:HIS:O	1:E:83:HIS:N	2.28	0.66
1:C:103:SER:HB2	1:C:118:LYS:HG2	1.79	0.65
1:E:76:SER:O	1:E:108:THR:OG1	2.11	0.64
2:J:92:GLU:HG2	2:J:98:SER:HA	1.79	0.64
1:C:2:THR:OG1	4:C:401:HOH:O	2.14	0.64
1:O:206:GLY:O	4:O:404:HOH:O	2.15	0.63
1:M:108:THR:HG23	1:M:109:ILE:HG12	1.81	0.62
2:N:23:LYS:NZ	4:N:302:HOH:O	2.33	0.61
1:E:3:ASP:HB2	1:E:6:VAL:HG23	1.82	0.61
2:P:92:GLU:HG2	2:P:98:SER:HA	1.83	0.61
2:D:186:ASP:OD1	4:D:302:HOH:O	2.15	0.60
1:O:76:SER:O	1:O:108:THR:HG21	2.02	0.60
1:I:119:GLU:HB3	4:I:461:HOH:O	2.03	0.59
2:B:92:GLU:HG2	2:B:98:SER:HA	1.85	0.59
1:E:51:LYS:HD3	1:E:58:PHE:CG	2.39	0.58
2:P:181:LYS:NZ	4:P:302:HOH:O	2.35	0.58
2:L:158:TYR:CE2	2:L:160:ASP:HB2	2.39	0.57
1:K:3:ASP:OD1	1:K:4:ASN:N	2.37	0.57
1:C:105:THR:HG22	4:C:539:HOH:O	2.03	0.57

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:198:LEU:HD12	1:I:198:LEU:HD12	1.86	0.57
1:A:103:SER:HB2	1:A:118:LYS:HG2	1.87	0.57
2:D:158:TYR:CE2	2:D:160:ASP:HB2	2.41	0.56
1:M:103:SER:HB2	1:M:118:LYS:HG2	1.87	0.55
1:E:108:THR:HG23	1:E:109:ILE:HG12	1.86	0.55
1:I:118:LYS:NZ	4:I:402:HOH:O	2.11	0.55
1:O:116:TRP:CD1	2:P:18:LYS:HG3	2.42	0.54
1:M:132:THR:OG1	2:N:189:GLU:OE1	2.23	0.54
1:C:108:THR:HG23	1:C:109:ILE:HG12	1.90	0.53
1:K:178:THR:HG21	1:M:201:PRO:CG	2.38	0.53
2:J:126:HIS:HD2	2:J:127:VAL:N	2.06	0.53
2:N:88:ALA:O	2:N:92:GLU:HG3	2.09	0.52
1:K:81:HIS:O	1:K:82:LYS:HG3	2.08	0.52
1:O:120:LEU:HD22	2:P:7:LEU:HD11	1.92	0.52
1:E:82:LYS:HE2	1:E:155:ASP:OD2	2.10	0.52
2:P:53:HIS:CG	2:P:135:HIS:HB2	2.45	0.52
2:D:206:GLU:N	4:D:306:HOH:O	2.38	0.51
2:L:92:GLU:HG2	2:L:98:SER:HA	1.93	0.51
1:K:108:THR:HG23	1:K:109:ILE:HG12	1.92	0.51
1:E:204:ALA:O	4:E:401:HOH:O	2.20	0.50
2:J:53:HIS:CG	2:J:135:HIS:HB2	2.47	0.49
1:A:88:GLU:OE1	4:A:401:HOH:O	2.20	0.49
1:G:82:LYS:HA	4:G:552:HOH:O	2.12	0.49
1:O:27:GLN:HG3	4:O:406:HOH:O	2.13	0.49
1:O:205:LEU:O	1:O:207:ALA:N	2.46	0.49
2:D:5:HIS:CG	2:D:56:GLU:HG2	2.48	0.48
1:M:3:ASP:HB3	1:M:6:VAL:HG23	1.95	0.48
2:B:53:HIS:CG	2:B:135:HIS:HB2	2.48	0.48
1:O:165:PRO:HB3	1:O:192:GLU:HA	1.96	0.48
1:I:120:LEU:HD22	2:J:7:LEU:HD11	1.95	0.48
1:A:47:LYS:NZ	4:A:408:HOH:O	2.47	0.48
2:J:126:HIS:CD2	2:J:127:VAL:N	2.81	0.48
2:F:10:LYS:HD2	4:F:348:HOH:O	2.14	0.47
1:C:206:GLY:HA2	1:C:207:ALA:HA	1.47	0.47
1:K:178:THR:HG21	1:M:201:PRO:HG3	1.95	0.47
2:D:22:ALA:O	2:D:23:LYS:HB2	2.14	0.47
1:M:76:SER:O	1:M:108:THR:OG1	2.16	0.47
2:J:179:LYS:HD3	2:J:181:LYS:HE2	1.97	0.47
2:D:1:MET:CE	2:D:56:GLU:HB3	2.45	0.47
2:J:158:TYR:CE2	2:J:160:ASP:HB2	2.50	0.46
1:M:16:VAL:O	1:M:20:GLU:HG3	2.15	0.46

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:81:HIS:O	1:M:83:HIS:N	2.48	0.46
1:E:174:ALA:HB1	1:I:203:VAL:HG11	1.96	0.46
2:J:61:ARG:O	2:J:65:THR:HB	2.15	0.46
1:K:198:LEU:HD12	1:M:198:LEU:HD12	1.95	0.46
1:C:116:TRP:CD1	2:D:18:LYS:HG3	2.51	0.46
1:A:48:LEU:HD23	1:A:73:LEU:HD12	1.98	0.46
1:K:81:HIS:O	1:K:82:LYS:O	2.34	0.46
1:E:116:TRP:CD1	2:F:18:LYS:HG3	2.50	0.46
2:L:124:ARG:HB2	2:L:206:GLU:HG3	1.98	0.46
1:M:137:ILE:HB	2:N:18:LYS:HD3	1.98	0.46
1:M:118:LYS:NZ	4:M:402:HOH:O	2.18	0.46
1:K:178:THR:HG21	1:M:201:PRO:HG2	1.98	0.46
1:O:132:THR:OG1	2:P:189:GLU:OE1	2.24	0.46
2:P:22:ALA:O	2:P:23:LYS:HB2	2.15	0.46
2:F:23:LYS:HE3	2:F:23:LYS:HB2	1.78	0.45
2:F:53:HIS:O	2:F:57:ARG:HG3	2.17	0.45
2:F:160:ASP:N	2:F:160:ASP:OD1	2.50	0.45
2:H:22:ALA:O	2:H:23:LYS:HB2	2.16	0.45
1:A:120:LEU:HD22	2:B:7:LEU:HD11	1.98	0.45
2:P:160:ASP:OD1	2:P:160:ASP:N	2.49	0.45
1:A:137:ILE:HD11	1:A:188:VAL:HG22	1.98	0.44
2:L:22:ALA:O	2:L:23:LYS:HB2	2.17	0.44
2:L:66:ALA:HB1	2:L:70:GLU:HB3	2.00	0.44
1:K:103:SER:OG	4:K:404:HOH:O	2.21	0.44
2:L:206:GLU:N	4:L:304:HOH:O	2.36	0.44
2:J:64:LEU:HD12	2:J:64:LEU:HA	1.87	0.44
2:J:22:ALA:O	2:J:23:LYS:HB2	2.18	0.44
1:O:119:GLU:HB3	4:O:440:HOH:O	2.17	0.44
1:A:116:TRP:CH2	1:A:188:VAL:HG21	2.53	0.43
2:D:78:THR:HG21	2:D:102:SER:HB2	2.00	0.43
1:E:78:PRO:HD2	1:E:81:HIS:HD2	1.81	0.43
1:I:116:TRP:CD1	2:J:18:LYS:HG3	2.53	0.43
1:G:116:TRP:CD1	2:H:18:LYS:HG3	2.52	0.43
2:J:88:ALA:O	2:J:92:GLU:HG3	2.19	0.43
1:A:73:LEU:HD23	1:A:73:LEU:HA	1.84	0.43
1:M:105:THR:HG22	4:M:517:HOH:O	2.18	0.43
2:L:23:LYS:HE3	2:L:26:HIS:CE1	2.54	0.43
1:E:164:ARG:HA	1:E:165:PRO:HD3	1.92	0.43
1:K:12:ASP:OD1	2:L:32:LYS:NZ	2.46	0.43
1:I:205:LEU:HD23	1:I:205:LEU:HA	1.86	0.43
1:I:48:LEU:HD23	1:I:73:LEU:HD12	2.01	0.42

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:28:GLU:HG3	4:B:385:HOH:O	2.18	0.42
2:J:78:THR:HG21	2:J:102:SER:HB2	2.00	0.42
1:I:10:ARG:NH1	4:I:410:HOH:O	2.39	0.42
1:K:120:LEU:HD22	2:L:7:LEU:HD11	2.01	0.42
2:F:22:ALA:O	2:F:23:LYS:HB2	2.19	0.42
1:I:108:THR:HG23	1:I:109:ILE:HG12	2.01	0.42
2:N:158:TYR:HA	2:N:159:PRO:HD3	1.88	0.42
1:K:164:ARG:HA	1:K:165:PRO:HD3	1.94	0.42
2:D:61:ARG:NH2	4:I:405:HOH:O	2.35	0.42
2:J:126:HIS:CD2	2:J:127:VAL:O	2.73	0.42
2:J:160:ASP:OD1	2:J:160:ASP:N	2.53	0.42
2:F:53:HIS:CG	2:F:135:HIS:HB2	2.55	0.41
1:K:17:LEU:HA	1:K:17:LEU:HD23	1.93	0.41
1:I:110:ILE:HG21	1:I:110:ILE:HD13	1.82	0.41
2:N:22:ALA:O	2:N:23:LYS:HB2	2.20	0.41
2:N:53:HIS:O	2:N:57:ARG:HG2	2.21	0.41
1:K:81:HIS:O	1:K:155:ASP:O	2.38	0.41
2:L:78:THR:HG21	2:L:102:SER:HB2	2.03	0.41
1:K:77:PHE:HA	1:K:108:THR:HG21	2.02	0.41
1:C:116:TRP:CH2	1:C:188:VAL:HG21	2.55	0.41
1:I:175:GLN:O	1:I:178:THR:HB	2.20	0.41
2:L:137:ARG:HD2	2:L:137:ARG:HA	1.86	0.41
2:J:66:ALA:HB1	2:J:70:GLU:HB3	2.03	0.41
1:E:81:HIS:O	1:E:155:ASP:O	2.38	0.41
1:E:51:LYS:NZ	4:E:404:HOH:O	2.38	0.41
2:P:137:ARG:HD2	2:P:137:ARG:HA	1.87	0.41
2:L:5:HIS:CG	2:L:56:GLU:HG2	2.56	0.41
2:D:158:TYR:HA	2:D:159:PRO:HD3	1.94	0.41
2:N:53:HIS:CG	2:N:135:HIS:HB2	2.55	0.41
1:G:78:PRO:HD2	1:G:81:HIS:HB2	2.02	0.41
1:G:108:THR:HG23	1:G:109:ILE:HG12	2.02	0.41
2:H:53:HIS:CG	2:H:135:HIS:HB2	2.56	0.41
2:J:197:VAL:HA	4:J:320:HOH:O	2.21	0.41
2:N:143:ARG:HD2	4:N:311:HOH:O	2.20	0.41
1:O:10:ARG:HB3	2:P:73:PHE:CZ	2.55	0.41
1:A:105:THR:HG22	4:A:526:HOH:O	2.21	0.40
2:H:160:ASP:N	2:H:160:ASP:OD1	2.54	0.40
2:P:88:ALA:O	2:P:92:GLU:HG3	2.21	0.40
1:A:175:GLN:O	1:A:178:THR:HG22	2.21	0.40
2:L:53:HIS:CG	2:L:135:HIS:HB2	2.56	0.40
1:I:73:LEU:HD23	1:I:73:LEU:HA	1.87	0.40

*Continued on next page...*



Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:137:ARG:HA	2:F:137:ARG:HD2	1.86	0.40
1:C:110:ILE:HD13	1:C:110:ILE:HG21	1.84	0.40
2:F:158:TYR:HA	2:F:159:PRO:HD3	1.94	0.40
2:B:158:TYR:CE2	2:B:160:ASP:HB2	2.56	0.40
2:B:160:ASP:OD1	2:B:160:ASP:N	2.54	0.40
1:K:203:VAL:HG13	4:K:454:HOH:O	2.21	0.40
1:M:48:LEU:HD23	1:M:73:LEU:HD12	2.03	0.40
1:O:76:SER:O	1:O:108:THR:CG2	2.69	0.40
1:I:119:GLU:HG3	4:I:405:HOH:O	2.21	0.40
2:P:158:TYR:HA	2:P:159:PRO:HD3	1.92	0.40
2:J:52:ARG:O	2:J:56:GLU:HG3	2.20	0.40
1:A:16:VAL:O	1:A:20:GLU:HG3	2.21	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:390:HOH:O	4:P:446:HOH:O[1_565]	1.74	0.46
4:F:432:HOH:O	4:P:401:HOH:O[1_565]	2.01	0.19
4:F:397:HOH:O	4:P:307:HOH:O[1_565]	2.11	0.09

## 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	204/206 (99%)	199 (98%)	4 (2%)	1 (0%)	34	34
1	C	202/206 (98%)	197 (98%)	3 (2%)	2 (1%)	19	16
1	E	202/206 (98%)	194 (96%)	5 (2%)	3 (2%)	13	8
1	G	202/206 (98%)	197 (98%)	3 (2%)	2 (1%)	19	16
1	I	202/206 (98%)	198 (98%)	3 (2%)	1 (0%)	34	34

Continued on next page...

*Continued from previous page...*

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	K	202/206 (98%)	194 (96%)	5 (2%)	3 (2%)	13	8
1	M	202/206 (98%)	196 (97%)	5 (2%)	1 (0%)	34	34
1	O	202/206 (98%)	197 (98%)	3 (2%)	2 (1%)	19	16
2	B	204/206 (99%)	201 (98%)	3 (2%)	0	100	100
2	D	204/206 (99%)	197 (97%)	6 (3%)	1 (0%)	34	34
2	F	204/206 (99%)	196 (96%)	8 (4%)	0	100	100
2	H	204/206 (99%)	193 (95%)	10 (5%)	1 (0%)	34	34
2	J	204/206 (99%)	199 (98%)	5 (2%)	0	100	100
2	L	204/206 (99%)	193 (95%)	10 (5%)	1 (0%)	34	34
2	N	204/206 (99%)	200 (98%)	4 (2%)	0	100	100
2	P	204/206 (99%)	199 (98%)	5 (2%)	0	100	100
All	All	3250/3296 (99%)	3150 (97%)	82 (2%)	18 (1%)	30	30

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	K	82	LYS
1	C	103	SER
1	E	205	LEU
1	G	103	SER
1	I	103	SER
1	K	103	SER
1	K	206	GLY
1	M	103	SER
1	A	103	SER
1	E	103	SER
1	G	83	HIS
1	O	103	SER
1	C	205	LEU
1	E	82	LYS
1	O	206	GLY
2	D	23	LYS
2	H	21	ASN
2	L	23	LYS

### 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	173/171 (101%)	172 (99%)	1 (1%)	90	94
1	C	171/171 (100%)	169 (99%)	2 (1%)	78	87
1	E	171/171 (100%)	170 (99%)	1 (1%)	90	94
1	G	171/171 (100%)	169 (99%)	2 (1%)	78	87
1	I	171/171 (100%)	170 (99%)	1 (1%)	90	94
1	K	171/171 (100%)	170 (99%)	1 (1%)	90	94
1	M	171/171 (100%)	168 (98%)	3 (2%)	66	77
1	O	171/171 (100%)	168 (98%)	3 (2%)	66	77
2	B	164/165 (99%)	163 (99%)	1 (1%)	90	94
2	D	164/165 (99%)	163 (99%)	1 (1%)	90	94
2	F	164/165 (99%)	162 (99%)	2 (1%)	78	87
2	H	164/165 (99%)	161 (98%)	3 (2%)	66	77
2	J	164/165 (99%)	161 (98%)	3 (2%)	66	77
2	L	164/165 (99%)	162 (99%)	2 (1%)	78	87
2	N	164/165 (99%)	161 (98%)	3 (2%)	66	77
2	P	164/165 (99%)	163 (99%)	1 (1%)	90	94
All	All	2682/2688 (100%)	2652 (99%)	30 (1%)	80	88

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	178	THR
2	B	160	ASP
1	C	178	THR
1	C	203	VAL
2	D	160	ASP
1	E	188	VAL
2	F	106	SER
2	F	160	ASP

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	G	178	THR
1	G	203	VAL
2	H	34	ASN
2	H	106	SER
2	H	160	ASP
1	I	178	THR
2	J	34	ASN
2	J	65	THR
2	J	160	ASP
1	K	178	THR
2	L	34	ASN
2	L	160	ASP
1	M	17	LEU
1	M	120	LEU
1	M	188	VAL
2	N	34	ASN
2	N	150	VAL
2	N	160	ASP
1	O	3	ASP
1	O	108	THR
1	O	178	THR
2	P	160	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	81	HIS
2	H	26	HIS
2	J	126	HIS
1	K	38	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

16 non-standard protein/DNA/RNA residues are modelled in this entry.

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	$\# Z  > 2$
1	CSD	A	102	1,3	2,7,8	0.56	0	2,8,10	4.10	2 (100%)
1	CSD	A	104	1,3	2,7,8	0.56	0	2,8,10	1.29	0
1	CSD	C	102	1,3	2,7,8	0.62	0	2,8,10	3.41	2 (100%)
1	CSD	C	104	1,3	2,7,8	0.52	0	2,8,10	1.59	1 (50%)
1	CSD	E	102	1,3	2,7,8	0.52	0	2,8,10	3.76	2 (100%)
1	CSD	E	104	1,3	2,7,8	0.55	0	2,8,10	1.31	0
1	CSD	G	102	1,3	2,7,8	0.55	0	2,8,10	4.16	2 (100%)
1	CSD	G	104	1,3	2,7,8	0.55	0	2,8,10	1.52	1 (50%)
1	CSD	I	102	1,3	2,7,8	0.57	0	2,8,10	4.15	2 (100%)
1	CSD	I	104	1,3	2,7,8	0.57	0	2,8,10	1.09	0
1	CSD	K	102	1,3	2,7,8	0.56	0	2,8,10	3.25	2 (100%)
1	CSD	K	104	1,3	2,7,8	0.64	0	2,8,10	2.31	1 (50%)
1	CSD	M	102	1,3	2,7,8	0.57	0	2,8,10	4.08	2 (100%)
1	CSD	M	104	1,3	2,7,8	0.56	0	2,8,10	1.35	0
1	CSD	O	102	1,3	2,7,8	0.53	0	2,8,10	3.82	2 (100%)
1	CSD	O	104	1,3	2,7,8	0.61	0	2,8,10	1.51	0

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
1	CSD	A	102	1,3	-	0/2/6/8	0/0/0/0
1	CSD	A	104	1,3	-	0/2/6/8	0/0/0/0
1	CSD	C	102	1,3	-	0/2/6/8	0/0/0/0
1	CSD	C	104	1,3	-	0/2/6/8	0/0/0/0
1	CSD	E	102	1,3	-	0/2/6/8	0/0/0/0
1	CSD	E	104	1,3	-	0/2/6/8	0/0/0/0
1	CSD	G	102	1,3	-	0/2/6/8	0/0/0/0
1	CSD	G	104	1,3	-	0/2/6/8	0/0/0/0
1	CSD	I	102	1,3	-	0/2/6/8	0/0/0/0
1	CSD	I	104	1,3	-	0/2/6/8	0/0/0/0
1	CSD	K	102	1,3	-	0/2/6/8	0/0/0/0

*Continued on next page...*

*Continued from previous page...*

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	CSD	K	104	1,3	-	0/2/6/8	0/0/0/0
1	CSD	M	102	1,3	-	0/2/6/8	0/0/0/0
1	CSD	M	104	1,3	-	0/2/6/8	0/0/0/0
1	CSD	O	102	1,3	-	0/2/6/8	0/0/0/0
1	CSD	O	104	1,3	-	0/2/6/8	0/0/0/0

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	102	CSD	O-C-CA	-3.16	117.24	125.72
1	A	102	CSD	O-C-CA	-3.12	117.36	125.72
1	G	102	CSD	O-C-CA	-2.99	117.70	125.72
1	O	102	CSD	O-C-CA	-2.88	118.00	125.72
1	M	102	CSD	O-C-CA	-2.81	118.18	125.72
1	E	102	CSD	O-C-CA	-2.72	118.42	125.72
1	I	102	CSD	O-C-CA	-2.68	118.53	125.72
1	C	102	CSD	O-C-CA	-2.66	118.59	125.72
1	C	104	CSD	O-C-CA	-2.13	120.01	125.72
1	G	104	CSD	O-C-CA	-2.05	120.22	125.72
1	K	104	CSD	OD1-SG-CB	2.78	110.00	105.43
1	K	102	CSD	OD1-SG-CB	3.33	110.91	105.43
1	C	102	CSD	OD1-SG-CB	4.02	112.05	105.43
1	O	102	CSD	OD1-SG-CB	4.57	112.96	105.43
1	E	102	CSD	OD1-SG-CB	4.57	112.96	105.43
1	A	102	CSD	OD1-SG-CB	4.88	113.46	105.43
1	M	102	CSD	OD1-SG-CB	5.04	113.72	105.43
1	G	102	CSD	OD1-SG-CB	5.06	113.75	105.43
1	I	102	CSD	OD1-SG-CB	5.22	114.02	105.43

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	204/206 (99%)	-0.48	0 100 100	18, 28, 43, 68	0
1	C	204/206 (99%)	-0.43	1 (0%) 91 92	18, 29, 47, 67	1 (0%)
1	E	204/206 (99%)	-0.43	1 (0%) 91 92	17, 29, 47, 62	1 (0%)
1	G	204/206 (99%)	-0.44	2 (0%) 84 85	17, 29, 46, 62	1 (0%)
1	I	204/206 (99%)	-0.47	0 100 100	19, 29, 43, 60	0
1	K	204/206 (99%)	-0.44	0 100 100	17, 29, 47, 73	1 (0%)
1	M	204/206 (99%)	-0.43	1 (0%) 91 92	18, 28, 44, 75	0
1	O	204/206 (99%)	-0.46	1 (0%) 91 92	19, 28, 44, 70	0
2	B	206/206 (100%)	-0.40	0 100 100	20, 29, 51, 75	0
2	D	206/206 (100%)	-0.40	1 (0%) 91 92	19, 29, 52, 67	0
2	F	206/206 (100%)	-0.40	0 100 100	18, 29, 53, 67	0
2	H	206/206 (100%)	-0.40	0 100 100	17, 28, 52, 69	0
2	J	206/206 (100%)	-0.43	1 (0%) 91 92	20, 29, 52, 69	0
2	L	206/206 (100%)	-0.40	0 100 100	18, 28, 52, 70	0
2	N	206/206 (100%)	-0.42	1 (0%) 91 92	20, 29, 51, 72	0
2	P	206/206 (100%)	-0.43	0 100 100	20, 29, 53, 74	0
All	All	3280/3296 (99%)	-0.43	9 (0%) 94 95	17, 29, 50, 75	4 (0%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	M	207	ALA	3.9
1	C	207	ALA	3.3
1	G	207	ALA	3.2
1	O	207	ALA	2.7
1	E	207	ALA	2.3

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
2	J	126	HIS	2.3
2	D	95	LEU	2.1
1	G	206	GLY	2.1
2	N	126	HIS	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
1	CSD	O	104	8/9	0.97	0.10	-	28,30,31,32	8
1	CSD	I	104	8/9	0.98	0.11	-	26,28,32,32	8
1	CSD	K	104	8/9	0.98	0.11	-	28,32,33,34	8
1	CSD	C	102	8/9	0.93	0.15	-	21,28,32,33	8
1	CSD	A	102	8/9	0.97	0.12	-	23,27,30,32	8
1	CSD	G	102	8/9	0.96	0.13	-	21,27,31,34	8
1	CSD	E	102	8/9	0.97	0.14	-	21,24,29,30	8
1	CSD	K	102	8/9	0.94	0.12	-	25,33,39,39	0
1	CSD	I	102	8/9	0.98	0.12	-	22,26,29,31	8
1	CSD	O	102	8/9	0.97	0.12	-	24,27,31,33	8
1	CSD	M	102	8/9	0.96	0.11	-	24,27,31,32	8
1	CSD	E	104	8/9	0.98	0.11	-	25,28,30,31	8
1	CSD	G	104	8/9	0.98	0.11	-	27,29,32,33	8
1	CSD	A	104	8/9	0.97	0.09	-	26,30,33,33	8
1	CSD	C	104	8/9	0.97	0.10	-	26,29,33,33	8
1	CSD	M	104	8/9	0.97	0.10	-	29,30,32,32	8

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains.

The B-factors column lists the minimum, median, 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	FE	O	301	1/1	0.98	0.11	-	35,35,35,35	1
3	FE	K	301	1/1	1.00	0.13	-	30,30,30,30	1
3	FE	A	301	1/1	0.99	0.13	-	29,29,29,29	1
3	FE	G	301	1/1	1.00	0.13	-	35,35,35,35	1
3	FE	M	301	1/1	0.99	0.14	-	33,33,33,33	1
3	FE	C	301	1/1	1.00	0.14	-	33,33,33,33	1
3	FE	E	301	1/1	0.99	0.14	-	31,31,31,31	1
3	FE	I	301	1/1	0.99	0.12	-	32,32,32,32	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.