



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2021 – 04:49 PM EDT

PDB ID : 3GPS  
Title : Crystal structure of the F87M/L110M mutant of human transthyretin at pH 5.5  
Authors : Palmieri, L.C.; Freire, J.B.B.; Foguel, D.; Lima, L.M.T.R.  
Deposited on : 2009-03-23  
Resolution : 1.78 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 : 2.23.2  
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) : 2.23.2

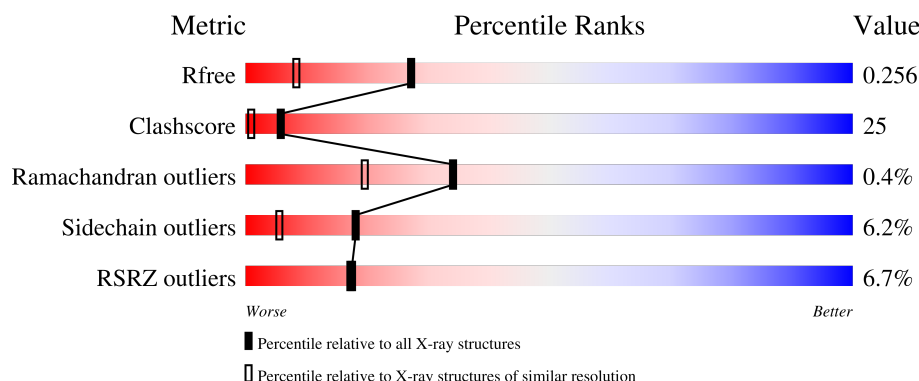
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


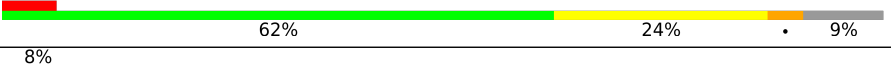
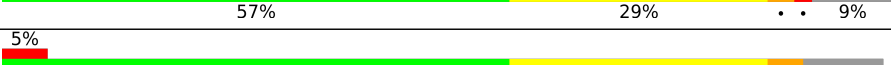

The reported resolution of this entry is 1.78 Å.

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	9185 (1.80-1.76)
Clashscore	141614	10184 (1.80-1.76)
Ramachandran outliers	138981	10051 (1.80-1.76)
Sidechain outliers	138945	10050 (1.80-1.76)
RSRZ outliers	127900	9032 (1.80-1.76)

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  $\geq 3$ , 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	127	
1	B	127	
1	C	127	
1	D	127	

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

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ACT	A	131[A]	-	-	X	-
3	ACT	A	131[B]	-	-	X	-
3	ACT	C	131	-	-	X	-
3	ACT	C	132	-	-	X	-
3	ACT	C	133	-	-	X	-
3	ACT	D	133[A]	-	-	X	-

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4167 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 Transthyretin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	115	Total	C	N	O	S	0	9	0
			969	614	161	190	4			
1	B	115	Total	C	N	O	S	0	6	0
			944	600	157	182	5			
1	C	116	Total	C	N	O	S	0	10	0
			983	624	161	194	4			
1	D	115	Total	C	N	O	S	0	9	0
			964	611	160	189	4			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	87	MET	PHE	engineered mutation	UNP P02766
A	110	MET	LEU	engineered mutation	UNP P02766
B	87	MET	PHE	engineered mutation	UNP P02766
B	110	MET	LEU	engineered mutation	UNP P02766
C	87	MET	PHE	engineered mutation	UNP P02766
C	110	MET	LEU	engineered mutation	UNP P02766
D	87	MET	PHE	engineered mutation	UNP P02766
D	110	MET	LEU	engineered mutation	UNP P02766

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Zn	0	0
			3	3		
2	B	3	Total	Zn	0	0
			3	3		
2	C	3	Total	Zn	0	0
			3	3		
2	D	3	Total	Zn	0	0
			3	3		

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula:  $C_2H_3O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	1
			8	4	4		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	1
			8	4	4		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		

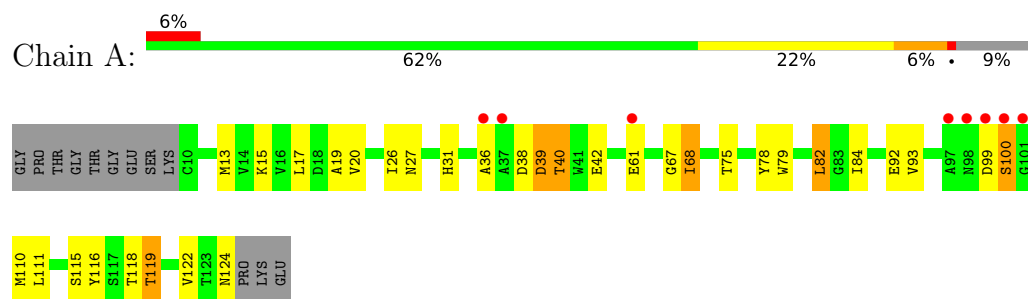
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	47	Total	O	0	0
			47	47		
5	B	51	Total	O	0	0
			51	51		
5	C	48	Total	O	0	0
			48	48		
5	D	59	Total	O	0	0
			59	59		

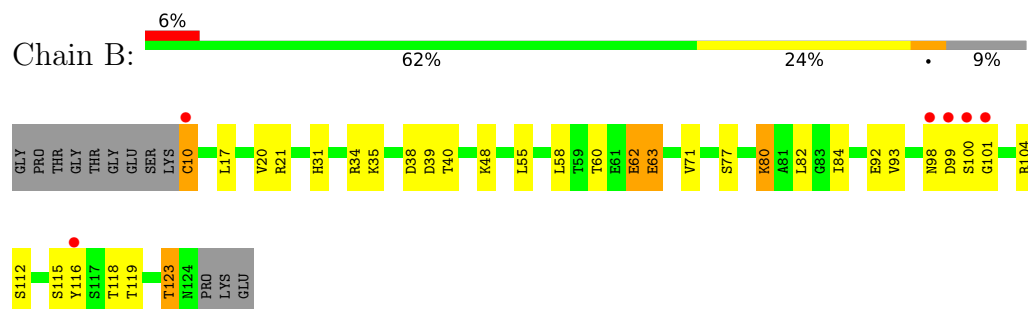
### 3 Residue-property plots [i](#)

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.

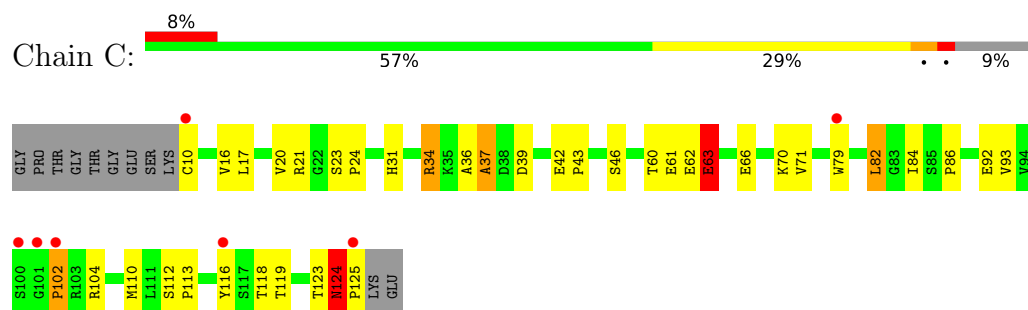
- Molecule 1: Transthyretin



- Molecule 1: Transthyretin

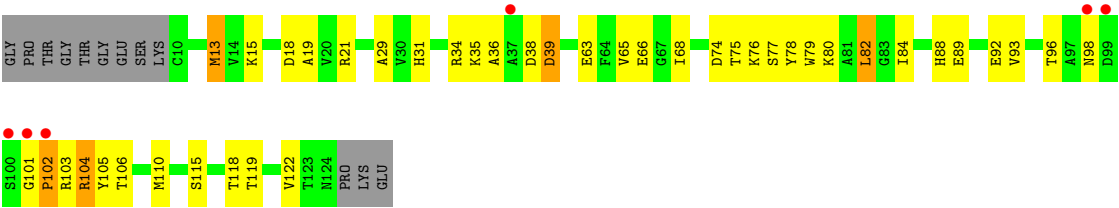


- Molecule 1: Transthyretin



- Molecule 1: Transthyretin







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	83.50Å 61.62Å 47.27Å 90.00° 90.01° 90.00°	Depositor
Resolution (Å)	27.90 – 1.78 27.90 – 1.78	Depositor EDS
% Data completeness (in resolution range)	99.2 (27.90-1.78) 99.2 (27.90-1.78)	Depositor EDS
$R_{merge}$	0.19	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.52 (at 1.78Å)	Xtriage
Refinement program	REFMAC 5.5.0059	Depositor
R, $R_{free}$	0.202 , 0.254 0.215 , 0.256	Depositor DCC
$R_{free}$ test set	2697 reflections (5.90%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	27.5	Xtriage
Anisotropy	0.176	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 43.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.480 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	4167	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.73% 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: ZN, ACT, GOL

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	1.05	2/994 (0.2%)	0.94	1/1355 (0.1%)
1	B	1.13	0/968	0.96	4/1317 (0.3%)
1	C	1.07	3/1008 (0.3%)	1.09	5/1374 (0.4%)
1	D	1.06	1/989 (0.1%)	0.97	2/1349 (0.1%)
All	All	1.08	6/3959 (0.2%)	0.99	12/5395 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	5
1	D	0	1
All	All	0	6

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	102	PRO	C-N	7.48	1.51	1.34
1	A	42	GLU	C-N	-6.60	1.21	1.34
1	A	78	TYR	CD1-CE1	-5.32	1.31	1.39
1	D	78	TYR	CD1-CE1	-5.21	1.31	1.39
1	C	63[A]	GLU	C-N	-5.07	1.22	1.34
1	C	63[B]	GLU	C-N	-5.07	1.22	1.34

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	124	ASN	C-N-CD	-11.31	95.72	120.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	102	PRO	O-C-N	-7.83	110.17	122.70
1	C	34	ARG	NE-CZ-NH2	7.24	123.92	120.30
1	A	82	LEU	CA-CB-CG	5.69	128.38	115.30
1	D	18	ASP	CB-CG-OD2	-5.41	113.43	118.30
1	B	21	ARG	NE-CZ-NH1	5.39	123.00	120.30
1	C	63[A]	GLU	C-N-CA	5.13	134.53	121.70
1	C	63[B]	GLU	C-N-CA	5.13	134.53	121.70
1	B	21	ARG	NE-CZ-NH2	-5.13	117.74	120.30
1	D	18	ASP	CB-CG-OD1	5.11	122.90	118.30
1	B	63	GLU	N-CA-CB	-5.05	101.52	110.60
1	B	34	ARG	NE-CZ-NH1	5.03	122.82	120.30

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	102	PRO	Mainchain
1	C	124	ASN	Mainchain
1	C	37	ALA	Peptide
1	C	61	GLU	Mainchain
1	D	101	GLY	Mainchain

## 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	969	0	922	47	0
1	B	944	0	912	48	0
1	C	983	0	939	86	0
1	D	964	0	918	43	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
2	C	3	0	0	0	0
2	D	3	0	0	0	0
3	A	8	0	6	5	0
3	B	12	0	9	2	0
3	C	12	0	9	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	16	0	12	4	0
4	A	6	0	8	0	0
4	B	18	0	24	0	0
4	C	6	0	8	2	0
4	D	12	0	16	0	0
5	A	47	0	0	7	0
5	B	51	0	0	2	0
5	C	48	0	0	2	0
5	D	59	0	0	9	0
All	All	4167	0	3783	197	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 25.

All (197) 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:119:THR:HG23	1:C:110:MET:SD	1.74	1.27
1:A:31[A]:HIS:HD2	5:A:245:HOH:O	1.13	1.25
1:A:119:THR:CG2	1:C:110:MET:SD	2.24	1.24
1:B:119:THR:HG22	1:D:110:MET:HE1	1.27	1.16
1:D:31[B]:HIS:CD2	5:D:231:HOH:O	2.02	1.12
1:C:124:ASN:HB3	1:C:125[B]:PRO:CD	1.80	1.11
1:C:124:ASN:HB3	1:C:125[B]:PRO:HD2	1.20	1.09
3:D:131:ACT:H1	5:D:243:HOH:O	1.51	1.08
1:C:124:ASN:N	1:C:125[A]:PRO:CD	2.18	1.07
1:C:124:ASN:N	1:C:125[A]:PRO:HD3	1.65	1.03
1:C:36:ALA:O	1:C:37:ALA:HB3	1.60	0.99
1:A:38:ASP:O	1:A:39:ASP:HB2	1.63	0.96
1:C:70:LYS:NZ	3:C:131:ACT:H3	1.81	0.95
1:C:36:ALA:HB2	1:C:42[B]:GLU:HG2	1.48	0.94
1:C:66:GLU:HG2	1:C:98:ASN:CB	1.98	0.93
1:D:93[A]:VAL:HG21	1:D:118[A]:THR:HG21	1.51	0.90
1:A:79[A]:TRP:HB2	5:A:227:HOH:O	1.70	0.90
1:B:119:THR:CG2	1:D:110:MET:HE1	2.00	0.90
1:D:31[A]:HIS:ND1	5:D:231:HOH:O	2.05	0.90
1:A:103[A]:ARG:HH11	1:A:103[A]:ARG:HG3	1.34	0.89
1:C:124:ASN:CB	1:C:125[B]:PRO:HD2	1.90	0.88
1:A:119:THR:HG21	1:C:110:MET:SD	2.11	0.86
1:A:103[A]:ARG:HH11	1:A:103[A]:ARG:CG	1.88	0.85
1:A:110:MET:SD	1:C:119[B]:THR:OG1	2.35	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:62[B]:GLU:HG3	1:C:62[B]:GLU:O	1.75	0.84
1:C:21[A]:ARG:HH21	1:C:82:LEU:CD2	1.90	0.84
1:A:15:LYS:HD2	5:A:259:HOH:O	1.77	0.84
1:A:31[B]:HIS:ND1	5:A:245:HOH:O	2.09	0.84
1:B:119:THR:HG22	1:D:110:MET:CE	2.06	0.84
1:C:66:GLU:HG2	1:C:98:ASN:HB3	1.57	0.84
1:C:36:ALA:HB2	1:C:42[B]:GLU:CG	2.08	0.83
1:B:119:THR:CG2	1:D:110:MET:CE	2.57	0.82
1:D:29:ALA:HB3	1:D:74[A]:ASP:OD2	1.80	0.81
1:C:70:LYS:HZ3	3:C:131:ACT:H3	1.46	0.80
1:C:66:GLU:HG2	1:C:98:ASN:O	1.81	0.79
1:C:36:ALA:O	1:C:37:ALA:CB	2.30	0.79
1:C:123:THR:C	1:C:125[A]:PRO:HD2	2.03	0.78
1:A:93:VAL:HG21	1:A:118[A]:THR:HG21	1.67	0.77
1:C:21[A]:ARG:NH2	1:C:82:LEU:CD2	2.47	0.77
1:B:82:LEU:HD11	1:C:82:LEU:HD12	1.65	0.76
1:C:31:HIS:NE2	4:C:134:GOL:H2	2.01	0.76
1:A:103[A]:ARG:HG3	1:A:103[A]:ARG:NH1	1.97	0.73
1:B:82:LEU:HD11	1:C:82:LEU:CD1	2.18	0.73
1:C:70:LYS:HZ2	3:C:131:ACT:H3	1.52	0.72
1:C:124:ASN:N	1:C:125[A]:PRO:HD2	2.03	0.72
1:C:123:THR:C	1:C:125[A]:PRO:CD	2.58	0.72
1:C:60:THR:OG1	1:C:63[B]:GLU:HG3	1.91	0.69
1:B:38:ASP:OD1	1:B:40:THR:HG22	1.93	0.68
1:C:66:GLU:HG2	1:C:98:ASN:HB2	1.77	0.66
1:C:116:TYR:HD1	1:C:118[B]:THR:HG23	1.60	0.66
1:D:15[A]:LYS:HE2	1:D:106:THR:HG21	1.78	0.66
1:C:21[A]:ARG:HH21	1:C:82:LEU:HD23	1.61	0.65
1:B:82:LEU:HG	1:C:82:LEU:HD11	1.78	0.65
1:A:36:ALA:HB3	1:A:40:THR:HG22	1.78	0.65
1:A:31[A]:HIS:CD2	5:A:245:HOH:O	2.03	0.65
1:A:102:PRO:O	1:A:102:PRO:HG2	1.97	0.64
1:C:66:GLU:HA	1:C:98:ASN:HB2	1.79	0.64
1:B:82:LEU:CD1	1:C:82:LEU:CD1	2.75	0.64
1:C:66:GLU:CG	1:C:98:ASN:O	2.45	0.64
1:A:116:TYR:HD1	1:A:118[B]:THR:HG23	1.63	0.63
1:D:35:LYS:HD2	1:D:39:ASP:O	1.97	0.63
1:B:108:ALA:HB3	1:B:119:THR:CG2	2.29	0.63
1:A:116:TYR:CD1	1:A:118[B]:THR:HG23	2.33	0.63
1:D:34:ARG:NH1	5:D:260:HOH:O	2.24	0.61
1:B:17:LEU:HD23	1:B:110[B]:MET:SD	2.41	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31[A]:HIS:HE1	5:A:244:HOH:O	1.84	0.60
1:B:35[B]:LYS:NZ	5:B:137:HOH:O	2.35	0.60
1:B:77:SER:HB3	1:B:80:LYS:HD3	1.83	0.60
1:B:10:CYS:O	1:B:104[A]:ARG:HG2	2.02	0.59
1:B:119:THR:HG21	1:D:110:MET:CE	2.32	0.59
1:D:115:SER:OG	3:D:133[B]:ACT:H1	2.02	0.59
1:C:124:ASN:HB3	1:C:125[B]:PRO:HD3	1.78	0.59
1:D:21:ARG:NH1	1:D:82:LEU:HD23	2.16	0.58
1:B:98:ASN:O	1:B:100:SER:N	2.36	0.58
1:C:66:GLU:CD	1:C:98:ASN:O	2.42	0.58
1:B:99:ASP:OD2	1:B:99:ASP:N	2.37	0.58
1:C:17[A]:LEU:N	1:C:17[A]:LEU:HD12	2.19	0.57
1:A:102:PRO:O	1:A:102:PRO:CG	2.53	0.57
1:D:102:PRO:O	1:D:102:PRO:HG2	2.04	0.57
1:C:116:TYR:CD1	1:C:118[B]:THR:HG23	2.40	0.57
1:A:75:THR:HG21	3:A:131[A]:ACT:CH3	2.34	0.56
1:B:119:THR:HG21	1:D:110:MET:HE2	1.87	0.56
1:D:68:ILE:HD12	1:D:96:THR:HG23	1.87	0.56
1:D:79[B]:TRP:HZ3	5:D:243:HOH:O	1.89	0.56
1:D:15[A]:LYS:HE2	1:D:106:THR:CG2	2.35	0.56
1:B:20:VAL:HG12	1:C:84:ILE:CD1	2.36	0.56
1:C:124:ASN:H	1:C:125[A]:PRO:HD3	1.63	0.55
1:D:31[B]:HIS:HD2	5:D:231:HOH:O	1.57	0.55
1:C:66:GLU:CG	1:C:98:ASN:HB3	2.32	0.55
1:B:104[B]:ARG:HB2	1:B:123:THR:HG22	1.87	0.55
1:B:31:HIS:HE1	5:B:232:HOH:O	1.89	0.55
1:C:93:VAL:HG22	1:C:118[B]:THR:HG21	1.89	0.55
1:C:10:CYS:N	1:C:104:ARG:NH1	2.55	0.54
1:B:84:ILE:HD12	1:C:20:VAL:HG12	1.90	0.54
1:B:84:ILE:CD1	1:C:20:VAL:HG12	2.38	0.54
1:A:115:SER:HB3	1:D:19:ALA:HB1	1.89	0.54
1:B:62:GLU:O	1:B:62:GLU:HG3	2.06	0.54
1:C:123:THR:C	1:C:125[A]:PRO:HD3	2.27	0.54
1:C:31:HIS:HE1	5:C:164:HOH:O	1.91	0.54
1:B:110[A]:MET:SD	1:D:119:THR:OG1	2.54	0.53
1:D:68:ILE:CD1	1:D:96:THR:HG23	2.38	0.53
1:A:38:ASP:OD1	1:A:40:THR:HB	2.08	0.53
1:C:21[A]:ARG:NH2	1:C:82:LEU:HD21	2.21	0.53
1:D:105:TYR:CD1	1:D:122:VAL:HG22	2.44	0.53
1:A:19:ALA:HB1	1:D:115:SER:HB3	1.91	0.53
1:B:93:VAL:HG11	1:B:118[A]:THR:HG21	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:124:ASN:CB	1:C:125[B]:PRO:CD	2.53	0.53
1:C:66:GLU:CG	1:C:98:ASN:CB	2.83	0.52
1:D:104:ARG:NH2	5:D:218:HOH:O	2.42	0.52
1:C:36:ALA:HB2	1:C:42[B]:GLU:HG3	1.91	0.52
1:B:82:LEU:CG	1:C:82:LEU:HD11	2.39	0.52
1:B:48:LYS:NZ	3:B:133:ACT:H1	2.25	0.51
1:C:36:ALA:CB	1:C:42[B]:GLU:HG2	2.32	0.51
1:D:77:SER:HA	1:D:80:LYS:HG2	1.93	0.51
1:A:75:THR:HG21	3:A:131[A]:ACT:H2	1.92	0.51
1:A:79[B]:TRP:CZ3	3:A:131[B]:ACT:OXT	2.63	0.51
1:C:37:ALA:C	1:C:39:ASP:H	2.14	0.51
1:B:115:SER:O	1:B:116:TYR:HB3	2.10	0.51
1:A:20:VAL:HG12	1:D:84:ILE:HD13	1.91	0.50
1:A:61[B]:GLU:OE2	1:A:61[B]:GLU:N	2.36	0.50
1:D:66:GLU:HG2	1:D:98:ASN:O	2.12	0.50
1:A:79[B]:TRP:HZ3	3:A:131[B]:ACT:OXT	1.94	0.50
1:D:115:SER:OG	3:D:133[A]:ACT:H3	2.12	0.49
1:C:46:SER:HB3	3:C:133:ACT:H2	1.94	0.49
1:C:60:THR:HG1	1:C:63[B]:GLU:HG3	1.76	0.49
1:C:70:LYS:HZ3	3:C:131:ACT:CH3	2.21	0.49
1:C:66:GLU:HG2	1:C:98:ASN:C	2.32	0.49
1:C:31:HIS:CE1	5:C:164:HOH:O	2.64	0.48
1:A:68:ILE:CD1	1:A:68:ILE:N	2.76	0.48
1:A:93:VAL:HG22	1:A:118[B]:THR:HG21	1.95	0.48
3:C:131:ACT:H1	5:D:148:HOH:O	2.12	0.48
1:B:93:VAL:CG2	1:B:118[B]:THR:HG21	2.43	0.48
1:B:116:TYR:HD1	1:B:118[B]:THR:HG23	1.79	0.48
1:D:21:ARG:NH1	1:D:82:LEU:CD2	2.76	0.48
1:D:36:ALA:O	1:D:38:ASP:O	2.32	0.48
1:B:93:VAL:HG22	1:B:118[B]:THR:HG21	1.96	0.47
1:B:104[A]:ARG:HB2	1:B:123:THR:HG22	1.96	0.47
1:C:112:SER:O	3:C:132:ACT:H2	2.15	0.47
1:B:17:LEU:HD23	1:B:110[B]:MET:CE	2.45	0.47
1:B:110[B]:MET:SD	1:B:110[B]:MET:N	2.89	0.46
1:C:113:PRO:HA	3:C:132:ACT:H2	1.96	0.46
1:C:123:THR:O	1:C:125[A]:PRO:HD2	2.14	0.46
1:C:124:ASN:OD1	1:C:124:ASN:O	2.33	0.46
1:A:17[B]:LEU:HD22	1:A:17[B]:LEU:N	2.30	0.46
1:C:16:VAL:C	1:C:17[A]:LEU:HD12	2.36	0.46
1:A:99:ASP:O	1:A:100:SER:OG	2.30	0.46
1:B:60:THR:OG1	1:B:63:GLU:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:115:SER:OG	3:D:133[A]:ACT:CH3	2.63	0.46
1:A:20:VAL:HG11	1:A:82:LEU:HD11	1.97	0.46
1:B:108:ALA:HB3	1:B:119:THR:HG22	1.99	0.45
1:C:124:ASN:O	1:C:125[A]:PRO:C	2.54	0.45
1:A:99:ASP:O	1:A:100:SER:CB	2.65	0.45
1:A:61[B]:GLU:H	1:A:61[B]:GLU:CD	2.15	0.45
1:B:118[A]:THR:HG22	1:B:119:THR:N	2.32	0.44
1:A:118[A]:THR:HG22	1:A:119:THR:N	2.33	0.44
1:B:119:THR:CG2	1:D:110:MET:HE2	2.42	0.44
1:A:105:TYR:CD1	1:A:122:VAL:HG22	2.53	0.44
1:A:19:ALA:HB1	1:D:115:SER:CB	2.48	0.43
1:C:96:THR:OG1	1:D:89[A]:GLU:OE2	2.27	0.43
1:C:79[A]:TRP:HE1	3:C:132:ACT:CH3	2.31	0.43
1:B:108:ALA:HB3	1:B:119:THR:HG23	2.00	0.43
1:C:79[A]:TRP:CD1	1:C:86:PRO:HB3	2.52	0.43
1:B:98:ASN:C	1:B:100:SER:N	2.72	0.43
1:B:55:LEU:HD21	1:B:58:LEU:HD21	2.00	0.43
1:C:93:VAL:CG2	1:C:118[B]:THR:HG21	2.48	0.43
1:A:68:ILE:N	1:A:68:ILE:HD12	2.34	0.42
1:C:37:ALA:N	1:C:39:ASP:H	2.17	0.42
1:B:98:ASN:O	1:B:101:GLY:N	2.51	0.42
1:C:31:HIS:NE2	4:C:134:GOL:C2	2.78	0.42
1:C:43:PRO:HB3	3:C:133:ACT:C	2.49	0.42
1:A:38:ASP:O	1:A:39:ASP:CB	2.42	0.42
1:B:112:SER:O	3:B:132:ACT:H3	2.20	0.42
1:C:34:ARG:O	1:C:42[B]:GLU:HG3	2.18	0.42
1:A:15:LYS:NZ	5:A:234:HOH:O	2.51	0.42
1:D:65:VAL:HG23	1:D:66:GLU:O	2.19	0.42
1:C:79[B]:TRP:HZ3	3:C:132:ACT:H3	1.85	0.42
1:D:13:MET:CE	1:D:15[A]:LYS:HD3	2.50	0.42
1:A:26:ILE:O	1:A:27:ASN:HB2	2.20	0.42
1:A:115:SER:CB	1:D:19:ALA:HB1	2.49	0.42
1:C:37:ALA:CA	1:C:39:ASP:H	2.33	0.42
1:A:75:THR:HG21	3:A:131[A]:ACT:H1	1.99	0.41
1:C:99:ASP:OD2	1:C:99:ASP:N	2.52	0.41
1:C:37:ALA:C	1:C:39:ASP:N	2.74	0.41
1:A:67:GLY:C	1:A:68:ILE:HD12	2.40	0.41
1:C:71:VAL:O	1:C:92[A]:GLU:HA	2.20	0.41
1:B:35[A]:LYS:C	1:B:35[A]:LYS:HD2	2.41	0.41
1:D:31[A]:HIS:CE1	5:D:231:HOH:O	2.64	0.41
1:D:75:THR:O	1:D:76:LYS:C	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:LEU:HD23	1:A:116:TYR:HB2	2.03	0.41
1:C:95:PHE:HA	1:D:88:HIS:HB2	2.03	0.41
1:B:20:VAL:HG12	1:C:84:ILE:HD13	2.03	0.41
1:C:23:SER:HB2	1:C:24:PRO:HD2	2.03	0.41
1:C:79[B]:TRP:CZ3	3:C:132:ACT:H3	2.57	0.41
1:C:17[B]:LEU:N	1:C:17[B]:LEU:HD22	2.36	0.40
1:B:71:VAL:O	1:B:92[A]:GLU:HA	2.22	0.40
1:B:82:LEU:HG	1:C:82:LEU:CD1	2.48	0.40
1:D:102:PRO:C	1:D:103:ARG:HG2	2.42	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	122/127 (96%)	114 (93%)	6 (5%)	2 (2%)	9	2
1	B	119/127 (94%)	115 (97%)	4 (3%)	0	100	100
1	C	123/127 (97%)	118 (96%)	5 (4%)	0	100	100
1	D	122/127 (96%)	117 (96%)	5 (4%)	0	100	100
All	All	486/508 (96%)	464 (96%)	20 (4%)	2 (0%)	34	19

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	100	SER
1	A	102	PRO

### 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	105/105 (100%)	94 (90%)	11 (10%)	7	1
1	B	102/105 (97%)	97 (95%)	5 (5%)	25	9
1	C	107/105 (102%)	104 (97%)	3 (3%)	43	27
1	D	105/105 (100%)	98 (93%)	7 (7%)	16	4
All	All	419/420 (100%)	393 (94%)	26 (6%)	18	5

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	MET
1	A	39	ASP
1	A	40	THR
1	A	68	ILE
1	A	84	ILE
1	A	92	GLU
1	A	102	PRO
1	A	103[A]	ARG
1	A	103[B]	ARG
1	A	119	THR
1	A	124	ASN
1	B	10	CYS
1	B	39	ASP
1	B	62	GLU
1	B	80	LYS
1	B	123	THR
1	C	63[A]	GLU
1	C	63[B]	GLU
1	C	82	LEU
1	D	13	MET
1	D	39	ASP
1	D	63	GLU
1	D	82	LEU
1	D	92	GLU

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Mol	Chain	Res	Type
1	D	102	PRO
1	D	104	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	B	98	ASN
1	C	124	ASN

### 5.3.3 RNA [i](#)

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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 31 ligands modelled in this entry, 12 are monoatomic - leaving 19 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	ACT	D	133[B]	-	1,3,3	1.21	0	0,3,3	-	-
3	ACT	D	131	-	1,3,3	1.39	0	0,3,3	-	-
4	GOL	D	134	-	5,5,5	0.34	0	5,5,5	0.41	0
4	GOL	C	134	-	5,5,5	0.42	0	5,5,5	0.45	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	ACT	B	133	-	1,3,3	2.16	1 (100%)	0,3,3	-	-
4	GOL	A	132	-	5,5,5	0.32	0	5,5,5	0.37	0
3	ACT	B	132	-	1,3,3	2.06	1 (100%)	0,3,3	-	-
4	GOL	D	135	-	5,5,5	0.40	0	5,5,5	0.21	0
4	GOL	B	135	-	5,5,5	0.38	0	5,5,5	0.35	0
3	ACT	B	131	-	1,3,3	1.32	0	0,3,3	-	-
4	GOL	B	134	-	5,5,5	0.41	0	5,5,5	0.53	0
3	ACT	D	132	-	1,3,3	0.95	0	0,3,3	-	-
3	ACT	C	132	-	1,3,3	1.32	0	0,3,3	-	-
3	ACT	C	133	-	1,3,3	4.18	1 (100%)	0,3,3	-	-
4	GOL	B	136	-	5,5,5	0.46	0	5,5,5	1.00	0
3	ACT	C	131	-	1,3,3	0.03	0	0,3,3	-	-
3	ACT	A	131[B]	-	1,3,3	4.26	1 (100%)	0,3,3	-	-
3	ACT	A	131[A]	-	1,3,3	1.54	0	0,3,3	-	-
3	ACT	D	133[A]	-	1,3,3	0.34	0	0,3,3	-	-

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	GOL	B	136	-	-	4/4/4/4	-
4	GOL	C	134	-	-	4/4/4/4	-
4	GOL	D	134	-	-	4/4/4/4	-
4	GOL	B	134	-	-	2/4/4/4	-
4	GOL	D	135	-	-	2/4/4/4	-
4	GOL	A	132	-	-	1/4/4/4	-
4	GOL	B	135	-	-	2/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	131[B]	ACT	CH3-C	4.26	1.54	1.48
3	C	133	ACT	CH3-C	4.18	1.54	1.48
3	B	133	ACT	CH3-C	2.16	1.51	1.48
3	B	132	ACT	CH3-C	2.06	1.51	1.48

There are no bond angle outliers.

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	134	GOL	O1-C1-C2-C3
4	B	135	GOL	C1-C2-C3-O3
4	B	136	GOL	O1-C1-C2-O2
4	B	136	GOL	O1-C1-C2-C3
4	B	136	GOL	C1-C2-C3-O3
4	C	134	GOL	O1-C1-C2-C3
4	C	134	GOL	C1-C2-C3-O3
4	C	134	GOL	O2-C2-C3-O3
4	D	134	GOL	O1-C1-C2-C3
4	B	135	GOL	O2-C2-C3-O3
4	D	134	GOL	C1-C2-C3-O3
4	D	135	GOL	O1-C1-C2-C3
4	D	134	GOL	O1-C1-C2-O2
4	D	134	GOL	O2-C2-C3-O3
4	D	135	GOL	O1-C1-C2-O2
4	B	136	GOL	O2-C2-C3-O3
4	C	134	GOL	O1-C1-C2-O2
4	A	132	GOL	O1-C1-C2-O2
4	B	134	GOL	O1-C1-C2-O2

There are no ring outliers.

11 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	133[B]	ACT	1	0
3	D	131	ACT	1	0
4	C	134	GOL	2	0
3	B	133	ACT	1	0
3	B	132	ACT	1	0
3	C	132	ACT	5	0
3	C	133	ACT	2	0
3	C	131	ACT	5	0
3	A	131[B]	ACT	2	0
3	A	131[A]	ACT	3	0
3	D	133[A]	ACT	2	0

## 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	115/127 (90%)	0.23	8 (6%) 16 16	21, 32, 55, 57	0
1	B	115/127 (90%)	0.27	7 (6%) 21 20	22, 35, 46, 53	0
1	C	116/127 (91%)	0.38	10 (8%) 10 10	22, 36, 48, 60	0
1	D	115/127 (90%)	0.22	6 (5%) 27 26	20, 31, 52, 54	0
All	All	461/508 (90%)	0.27	31 (6%) 17 17	20, 34, 52, 60	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	100	SER	12.1
1	B	100	SER	8.6
1	B	101	GLY	6.7
1	A	99	ASP	5.5
1	C	101	GLY	5.4
1	A	61[A]	GLU	4.9
1	A	98	ASN	4.7
1	D	98	ASN	4.4
1	D	37	ALA	3.7
1	D	100	SER	3.7
1	C	98	ASN	3.7
1	A	101	GLY	3.4
1	C	125[A]	PRO	3.0
1	A	37	ALA	3.0
1	C	97	ALA	2.9
1	D	102	PRO	2.9
1	A	100	SER	2.7
1	C	79[A]	TRP	2.7
1	A	97	ALA	2.5
1	B	10	CYS	2.5
1	D	101	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	10	CYS	2.4
1	D	99	ASP	2.4
1	B	99	ASP	2.3
1	C	99	ASP	2.3
1	B	116	TYR	2.3
1	B	98	ASN	2.2
1	A	36	ALA	2.2
1	C	102	PRO	2.1
1	C	116	TYR	2.1
1	B	110[A]	MET	2.1

## 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 monosaccharides 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(Å <sup>2</sup> )	Q<0.9
3	ACT	C	133	4/4	0.68	0.20	51,51,53,54	0
4	GOL	B	136	6/6	0.68	0.19	48,53,55,55	0
4	GOL	B	134	6/6	0.70	0.21	73,74,74,74	0
4	GOL	C	134	6/6	0.72	0.15	67,68,68,68	0
3	ACT	B	133	4/4	0.73	0.16	75,75,75,75	0
3	ACT	D	131	4/4	0.77	0.17	39,40,41,42	0
3	ACT	A	131[A]	4/4	0.77	0.31	30,32,32,33	4
3	ACT	C	132	4/4	0.77	0.20	51,51,52,52	0
3	ACT	A	131[B]	4/4	0.77	0.31	25,27,27,28	4
4	GOL	A	132	6/6	0.79	0.16	64,65,66,66	0
3	ACT	D	133[A]	4/4	0.82	0.29	31,34,34,34	4
3	ACT	D	133[B]	4/4	0.82	0.29	27,30,30,31	4
4	GOL	D	135	6/6	0.84	0.18	60,64,65,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	GOL	D	134	6/6	0.87	0.12	57,59,59,61	0
2	ZN	B	129	1/1	0.88	0.14	90,90,90,90	0
2	ZN	C	129	1/1	0.90	0.14	85,85,85,85	0
3	ACT	B	132	4/4	0.91	0.21	51,51,52,53	0
3	ACT	D	132	4/4	0.92	0.16	45,46,47,47	0
4	GOL	B	135	6/6	0.92	0.11	59,61,61,63	0
2	ZN	C	128	1/1	0.94	0.09	63,63,63,63	0
2	ZN	B	128	1/1	0.96	0.09	60,60,60,60	0
3	ACT	C	131	4/4	0.97	0.09	43,43,43,44	0
3	ACT	B	131	4/4	0.97	0.07	44,45,45,46	0
2	ZN	D	128	1/1	0.98	0.02	53,53,53,53	0
2	ZN	A	130	1/1	0.98	0.07	37,37,37,37	1
2	ZN	D	129	1/1	0.99	0.09	55,55,55,55	0
2	ZN	D	130	1/1	0.99	0.07	30,30,30,30	0
2	ZN	B	130	1/1	0.99	0.05	41,41,41,41	0
2	ZN	A	129	1/1	1.00	0.07	30,30,30,30	0
2	ZN	C	130	1/1	1.00	0.05	40,40,40,40	0
2	ZN	A	128	1/1	1.00	0.02	51,51,51,51	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.