



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:54 PM GMT

PDB ID : 1R6U
Title : Crystal structure of an active fragment of human tryptophanyl-tRNA synthetase with cytokine activity
Authors : Yang, X.-L.; Otero, F.J.; Skene, R.J.; McRee, D.E.; Ribas de Pouplana, L.; Schimmel, P.
Deposited on : 2003-10-16
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
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) : trunk26865

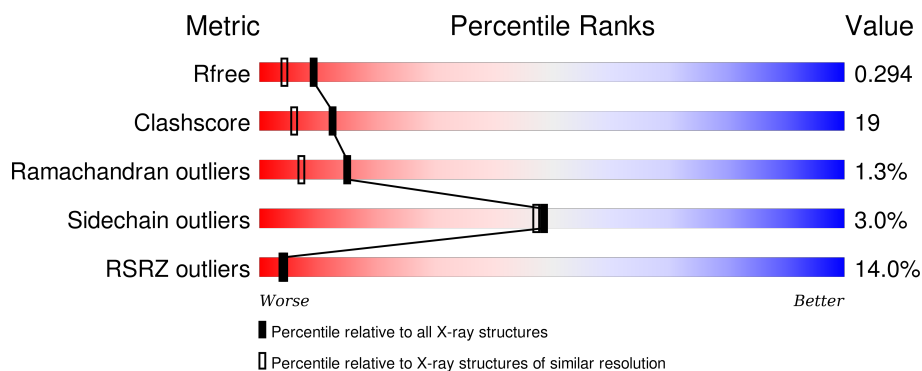
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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 ≥ 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	437	<div> <div>15%</div> <div>54%</div> <div>32%</div> <div>•</div> <div>11%</div> </div>
1	B	437	<div> <div>9%</div> <div>53%</div> <div>29%</div> <div>•</div> <div>17%</div> </div>

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	GOL	A	601	-	-	-	X
3	GOL	A	602	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6205 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 Tryptophanyl-tRNA synthetase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	388	Total	C	N	O	S	Se	0	0	0
			3129	2004	528	582	5	10			
1	B	362	Total	C	N	O	S	Se	0	0	0
			2927	1879	497	537	5	9			

There are 50 discrepancies between the modelled and reference sequences:

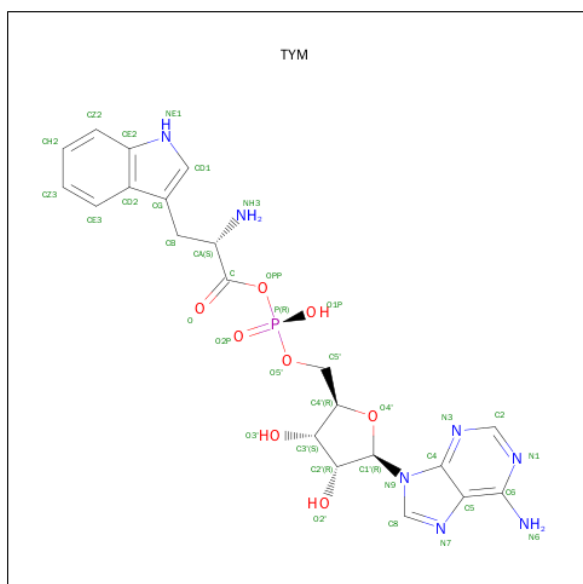
Chain	Residue	Modelled	Actual	Comment	Reference
A	143	MSE	MET	MODIFIED RESIDUE	UNP P23381
A	169	MSE	MET	MODIFIED RESIDUE	UNP P23381
A	195	MSE	MET	MODIFIED RESIDUE	UNP P23381
A	213	GLY	SER	ENGINEERED	UNP P23381
A	214	ASP	TYR	ENGINEERED	UNP P23381
A	241	MSE	MET	MODIFIED RESIDUE	UNP P23381
A	243	MSE	MET	MODIFIED RESIDUE	UNP P23381
A	319	MSE	MET	MODIFIED RESIDUE	UNP P23381
A	350	MSE	MET	MODIFIED RESIDUE	UNP P23381
A	401	MSE	MET	MODIFIED RESIDUE	UNP P23381
A	425	MSE	MET	MODIFIED RESIDUE	UNP P23381
A	461	MSE	MET	MODIFIED RESIDUE	UNP P23381
A	472	LYS	-	CLONING ARTIFACT	UNP P23381
A	473	LEU	-	CLONING ARTIFACT	UNP P23381
A	474	ALA	-	CLONING ARTIFACT	UNP P23381
A	475	ALA	-	CLONING ARTIFACT	UNP P23381
A	476	ALA	-	CLONING ARTIFACT	UNP P23381
A	477	LEU	-	CLONING ARTIFACT	UNP P23381
A	478	GLU	-	CLONING ARTIFACT	UNP P23381
A	479	HIS	-	EXPRESSION TAG	UNP P23381
A	480	HIS	-	EXPRESSION TAG	UNP P23381
A	481	HIS	-	EXPRESSION TAG	UNP P23381
A	482	HIS	-	EXPRESSION TAG	UNP P23381
A	483	HIS	-	EXPRESSION TAG	UNP P23381
A	484	HIS	-	EXPRESSION TAG	UNP P23381

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Chain	Residue	Modelled	Actual	Comment	Reference
B	143	MSE	MET	MODIFIED RESIDUE	UNP P23381
B	169	MSE	MET	MODIFIED RESIDUE	UNP P23381
B	195	MSE	MET	MODIFIED RESIDUE	UNP P23381
B	213	GLY	SER	ENGINEERED	UNP P23381
B	214	ASP	TYR	ENGINEERED	UNP P23381
B	241	MSE	MET	MODIFIED RESIDUE	UNP P23381
B	243	MSE	MET	MODIFIED RESIDUE	UNP P23381
B	319	MSE	MET	MODIFIED RESIDUE	UNP P23381
B	350	MSE	MET	MODIFIED RESIDUE	UNP P23381
B	401	MSE	MET	MODIFIED RESIDUE	UNP P23381
B	425	MSE	MET	MODIFIED RESIDUE	UNP P23381
B	461	MSE	MET	MODIFIED RESIDUE	UNP P23381
B	472	LYS	-	CLONING ARTIFACT	UNP P23381
B	473	LEU	-	CLONING ARTIFACT	UNP P23381
B	474	ALA	-	CLONING ARTIFACT	UNP P23381
B	475	ALA	-	CLONING ARTIFACT	UNP P23381
B	476	ALA	-	CLONING ARTIFACT	UNP P23381
B	477	LEU	-	CLONING ARTIFACT	UNP P23381
B	478	GLU	-	CLONING ARTIFACT	UNP P23381
B	479	HIS	-	EXPRESSION TAG	UNP P23381
B	480	HIS	-	EXPRESSION TAG	UNP P23381
B	481	HIS	-	EXPRESSION TAG	UNP P23381
B	482	HIS	-	EXPRESSION TAG	UNP P23381
B	483	HIS	-	EXPRESSION TAG	UNP P23381
B	484	HIS	-	EXPRESSION TAG	UNP P23381

- Molecule 2 is TRYPTOPHANYL-5'AMP (three-letter code: TYM) (formula: $C_{21}H_{24}N_7O_8P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			37	21	7	8	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

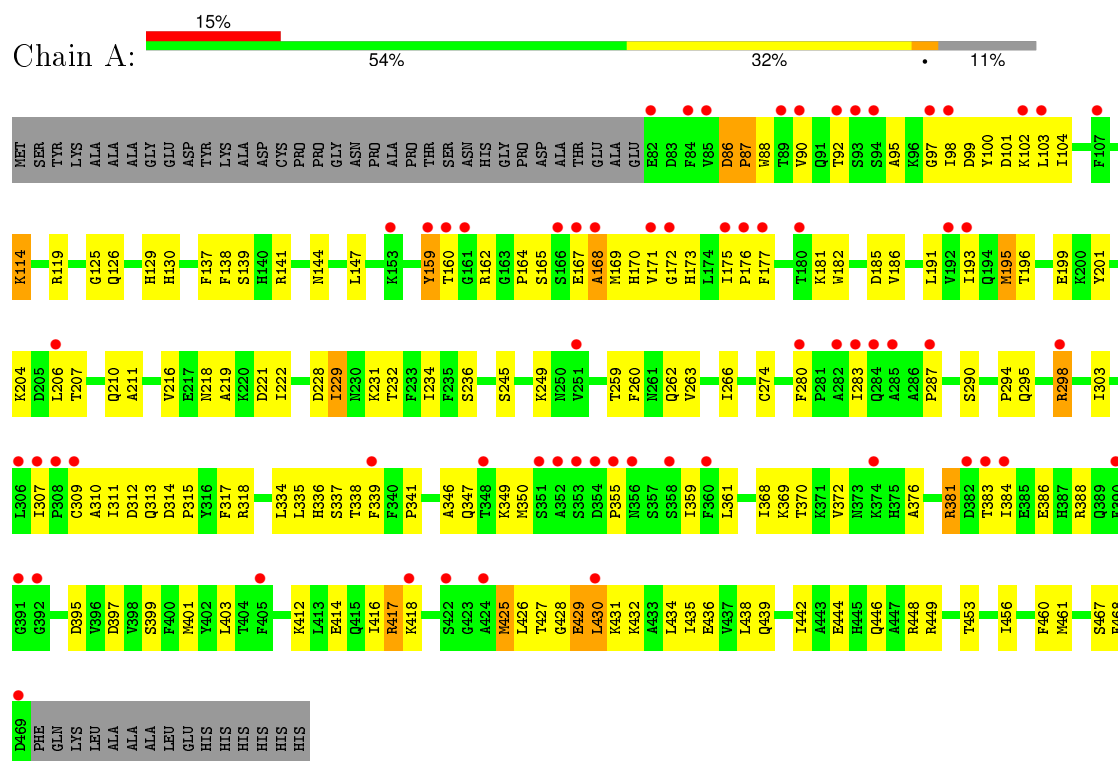
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	50	Total	O	0	0
			50	50		
4	B	50	Total	O	0	0
			50	50		

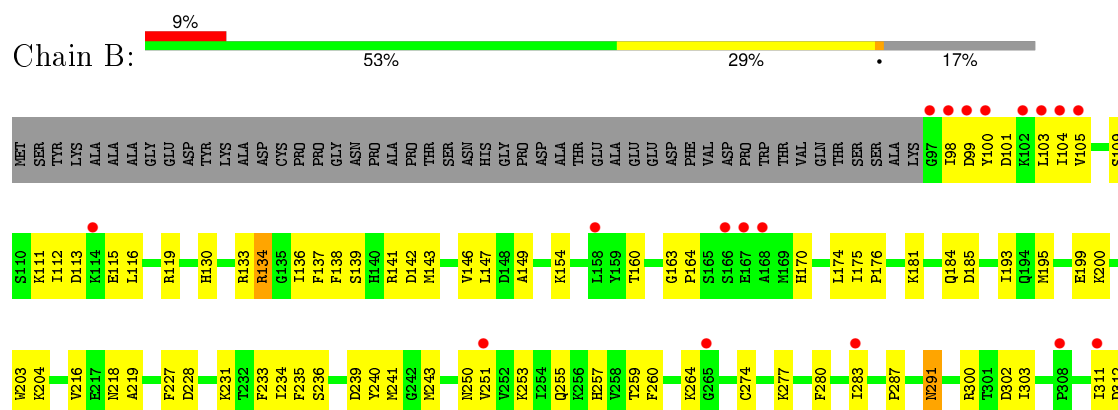
3 Residue-property plots [i](#)

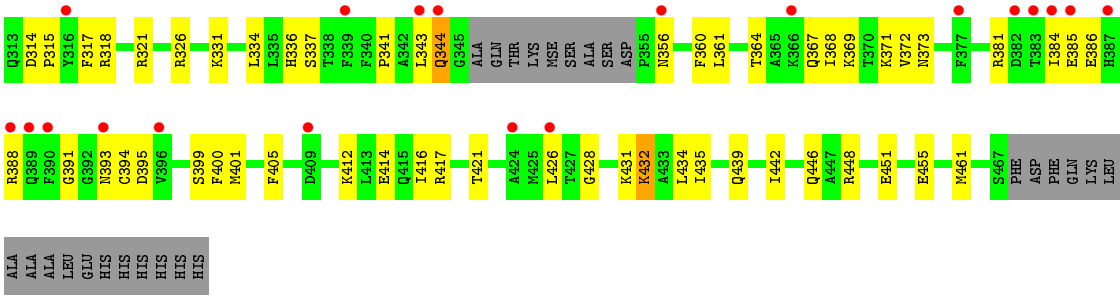
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Tryptophanyl-tRNA synthetase



• Molecule 1: Tryptophanyl-tRNA synthetase





ALA
ALA
ALA
LEU
GLU
HIS
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	134.66 Å 96.46 Å 97.13 Å 90.00° 129.90° 90.00°	Depositor
Resolution (Å)	30.00 – 2.00 29.88 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-2.00) 96.6 (29.88-2.00)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.60 (at 2.00 Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.254 , 0.297 0.252 , 0.294	Depositor DCC
R_{free} test set	3078 reflections (4.94%)	DCC
Wilson B-factor (Å ²)	37.7	Xtriage
Anisotropy	0.181	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 52.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 126480 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6205	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: GOL, TYM

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.38	0/3195	0.62	1/4297 (0.0%)
1	B	0.36	0/2988	0.58	0/4014
All	All	0.37	0/6183	0.60	1/8311 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	430	LEU	N-CA-C	-8.34	88.49	111.00

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	3129	0	3075	131	0
1	B	2927	0	2893	108	0
2	A	37	0	23	3	0
3	A	12	0	14	2	0
4	A	50	0	0	4	0
4	B	50	0	0	3	0
All	All	6205	0	6005	236	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 19.

All (236) 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:429:GLU:HB3	1:A:432:LYS:HB2	1.18	1.12
1:A:95:ALA:HA	1:A:347:GLN:NE2	1.88	0.89
1:A:416:ILE:HA	1:A:425:MSE:HE1	1.56	0.86
1:B:250:ASN:HD21	1:B:291:ASN:ND2	1.74	0.86
1:B:143:MSE:HE3	1:B:146:VAL:HB	1.57	0.84
1:A:245:SER:O	1:A:249:LYS:HE2	1.81	0.80
1:B:250:ASN:HD21	1:B:291:ASN:HD21	1.32	0.77
1:B:241:MSE:HE1	1:B:283:ILE:HD12	1.68	0.75
1:A:206:LEU:HB2	1:A:210:GLN:HE21	1.52	0.75
1:A:432:LYS:O	1:A:436:GLU:HG3	1.86	0.75
1:A:312:ASP:HB3	1:A:339:PHE:CZ	2.23	0.74
1:B:174:LEU:HD21	1:B:361:LEU:HD21	1.69	0.73
1:B:253:LYS:O	1:B:257:HIS:HD2	1.71	0.73
1:B:141:ARG:HD3	1:B:321:ARG:NH1	2.03	0.72
1:A:95:ALA:HA	1:A:347:GLN:HE22	1.54	0.71
1:A:312:ASP:HB3	1:A:339:PHE:HZ	1.54	0.71
1:B:428:GLY:O	1:B:432:LYS:HG2	1.91	0.71
1:B:141:ARG:HD2	1:B:334:LEU:HD12	1.73	0.70
1:B:373:ASN:HA	1:B:431:LYS:HD3	1.71	0.70
1:B:104:ILE:HG21	4:B:1080:HOH:O	1.92	0.69
1:A:137:PHE:CZ	1:A:337:SER:HB3	2.27	0.69
1:B:175:ILE:HB	1:B:176:PRO:HD3	1.74	0.69
1:B:448:ARG:O	1:B:451:GLU:HG2	1.93	0.69
1:A:169:MSE:HE3	1:A:221:ASP:HB2	1.74	0.68
1:A:141:ARG:HG3	1:A:141:ARG:HH11	1.59	0.67
1:A:309:CYS:SG	4:A:1087:HOH:O	2.52	0.67
1:B:251:VAL:O	1:B:255:GLN:HG3	1.95	0.66
1:A:90:VAL:HG23	1:A:349:LYS:HE3	1.78	0.65
1:A:181:LYS:HE2	1:A:185:ASP:OD2	1.96	0.64
1:A:229:ILE:HG12	1:A:460:PHE:CE1	2.32	0.64
1:B:112:ILE:HG23	1:B:116:LEU:HD23	1.80	0.64
1:A:310:ALA:HB3	1:A:312:ASP:OD1	1.99	0.63
1:B:300:ARG:HG2	1:B:300:ARG:HH11	1.62	0.63
1:B:104:ILE:HD11	1:B:111:LYS:N	2.14	0.62
1:A:429:GLU:HB3	1:A:432:LYS:CB	2.11	0.62
1:B:104:ILE:HD13	1:B:139:SER:HB3	1.82	0.62
1:B:321:ARG:O	1:B:331:LYS:HE3	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:336:HIS:CD2	4:A:1087:HOH:O	2.53	0.61
1:B:104:ILE:HD12	1:B:109:SER:OG	2.00	0.61
1:B:356:ASN:HA	1:B:371:LYS:HE2	1.82	0.61
1:B:369:LYS:HA	1:B:435:ILE:HD13	1.83	0.61
1:A:429:GLU:C	1:A:432:LYS:H	2.04	0.61
1:A:418:LYS:HB2	1:A:418:LYS:NZ	2.15	0.61
1:A:144:ASN:HB2	4:A:1034:HOH:O	2.01	0.61
1:B:112:ILE:HB	1:B:138:PHE:O	2.02	0.60
1:A:438:LEU:O	1:A:442:ILE:HG12	2.01	0.60
1:A:444:GLU:HB3	1:A:448:ARG:HH12	1.65	0.60
1:A:442:ILE:O	1:A:446:GLN:HG3	2.01	0.60
1:B:384:ILE:HG23	1:B:385:GLU:OE2	2.02	0.60
1:B:113:ASP:OD1	1:B:115:GLU:HB3	2.02	0.59
1:B:234:ILE:O	1:B:461:MSE:HA	2.02	0.59
1:A:201:TYR:HA	1:A:206:LEU:HD11	1.84	0.59
1:B:181:LYS:HD3	1:B:185:ASP:OD2	2.02	0.59
1:A:171:VAL:HG23	1:A:359:ILE:O	2.02	0.59
1:B:119:ARG:HD3	1:B:147:LEU:HD13	1.86	0.58
1:B:101:ASP:O	1:B:104:ILE:HG22	2.03	0.58
1:A:137:PHE:CE1	1:A:337:SER:HB3	2.38	0.58
1:A:431:LYS:O	1:A:435:ILE:HG13	2.04	0.57
1:A:206:LEU:CB	1:A:210:GLN:HE21	2.18	0.56
1:A:100:TYR:O	1:A:101:ASP:HB3	2.06	0.56
1:A:368:ILE:O	1:A:372:VAL:HG23	2.05	0.56
1:A:229:ILE:O	1:A:229:ILE:HD13	2.05	0.56
1:A:199:GLU:HB2	1:A:280:PHE:CZ	2.40	0.56
1:A:416:ILE:HG23	1:A:425:MSE:SE	2.55	0.55
1:A:427:THR:O	1:A:428:GLY:C	2.44	0.55
1:A:274:CYS:HA	1:B:259:THR:HA	1.87	0.55
1:A:169:MSE:HA	1:A:173:HIS:ND1	2.21	0.55
1:A:141:ARG:NH1	1:A:141:ARG:HG3	2.22	0.55
1:B:137:PHE:CE1	1:B:337:SER:HB3	2.42	0.55
1:B:442:ILE:O	1:B:446:GLN:HG3	2.07	0.55
1:A:234:ILE:O	1:A:461:MSE:HA	2.06	0.55
1:B:139:SER:OG	1:B:311:ILE:HD13	2.07	0.54
1:B:104:ILE:HG12	1:B:111:LYS:HB2	1.89	0.54
1:B:203:TRP:CD1	1:B:277:LYS:HE2	2.43	0.54
1:A:100:TYR:C	1:A:102:LYS:H	2.11	0.53
1:A:453:THR:O	1:A:456:ILE:HB	2.09	0.53
1:B:136:ILE:HD12	1:B:405:PHE:CE1	2.44	0.53
1:B:98:ILE:HG21	1:B:100:TYR:HE2	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:164:PRO:HG3	1:A:195:MSE:HG3	1.89	0.53
1:A:307:ILE:HB	1:A:334:LEU:HD23	1.90	0.53
1:B:318:ARG:HH11	1:B:318:ARG:HG2	1.73	0.53
1:B:199:GLU:HB2	1:B:280:PHE:CZ	2.44	0.53
1:B:98:ILE:HG21	1:B:100:TYR:CE2	2.43	0.53
1:B:130:HIS:O	1:B:134:ARG:HB2	2.08	0.53
1:A:429:GLU:C	1:A:431:LYS:N	2.53	0.53
1:A:260:PHE:O	1:A:263:VAL:HG12	2.09	0.52
1:A:294:PRO:HA	1:A:298:ARG:O	2.08	0.52
1:B:317:PHE:O	1:B:321:ARG:HG3	2.10	0.52
1:B:141:ARG:HD3	1:B:321:ARG:HH11	1.71	0.52
1:B:364:THR:O	1:B:368:ILE:HG13	2.09	0.52
1:A:162:ARG:HH11	1:A:162:ARG:HG3	1.74	0.52
1:B:432:LYS:HA	1:B:432:LYS:HE3	1.90	0.52
1:B:241:MSE:CE	1:B:283:ILE:HD12	2.38	0.52
1:A:98:ILE:N	1:A:98:ILE:HD12	2.25	0.52
1:A:99:ASP:OD1	1:A:102:LYS:HD2	2.10	0.52
1:A:453:THR:OG1	1:A:456:ILE:HG12	2.10	0.52
1:B:204:LYS:HE3	4:B:1078:HOH:O	2.08	0.51
1:A:159:TYR:CZ	1:A:287:PRO:HB2	2.45	0.51
1:B:341:PRO:HG3	1:B:401:MSE:HE2	1.92	0.51
1:B:368:ILE:CD1	1:B:442:ILE:HD12	2.40	0.51
1:B:112:ILE:HG12	1:B:116:LEU:HD23	1.91	0.51
1:B:341:PRO:HG2	1:B:401:MSE:HB3	1.91	0.51
1:B:233:PHE:CZ	1:B:235:PHE:HB3	2.45	0.51
1:A:129:HIS:CD2	3:A:602:GOL:H31	2.45	0.50
1:A:173:HIS:O	1:A:176:PRO:HD2	2.12	0.50
1:A:141:ARG:HG3	1:A:334:LEU:HD12	1.93	0.50
1:B:326:ARG:HH11	1:B:326:ARG:HG3	1.76	0.50
1:A:412:LYS:O	1:A:416:ILE:HG13	2.11	0.50
1:A:429:GLU:OE1	1:A:432:LYS:HG3	2.12	0.50
1:A:444:GLU:HB3	1:A:448:ARG:NH1	2.26	0.50
1:A:372:VAL:HG12	1:A:431:LYS:HB3	1.94	0.49
1:A:177:PHE:CE1	1:A:222:ILE:HD12	2.47	0.49
1:A:172:GLY:O	1:A:175:ILE:HG12	2.12	0.49
1:A:206:LEU:CD1	1:A:211:ALA:HB2	2.42	0.49
1:A:283:ILE:N	1:A:283:ILE:HD12	2.27	0.49
1:A:103:LEU:HD23	1:A:311:ILE:HD13	1.93	0.49
1:A:383:THR:HG22	1:A:386:GLU:OE1	2.12	0.49
1:B:143:MSE:HE3	1:B:146:VAL:CB	2.36	0.49
1:B:228:ASP:CB	1:B:231:LYS:HE2	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:231:LYS:HG2	1:A:468:PHE:CE1	2.47	0.49
1:B:139:SER:OG	1:B:336:HIS:HB2	2.13	0.49
1:B:133:ARG:O	1:B:134:ARG:HG3	2.13	0.49
1:A:86:ASP:OD2	1:A:88:TRP:HE3	1.96	0.49
1:A:182:TRP:O	1:A:186:VAL:HG22	2.13	0.49
1:A:449:ARG:HG2	1:A:449:ARG:HH11	1.78	0.49
1:A:262:GLN:O	1:A:266:ILE:HG12	2.13	0.48
1:B:321:ARG:HB3	1:B:331:LYS:HE2	1.95	0.48
1:B:112:ILE:CB	1:B:138:PHE:O	2.61	0.48
1:A:87:PRO:HD3	1:A:315:PRO:HG2	1.94	0.48
1:A:369:LYS:HG3	1:A:370:THR:N	2.28	0.48
1:B:149:ALA:HA	1:B:154:LYS:HE3	1.96	0.48
1:A:199:GLU:HB2	1:A:280:PHE:CE1	2.48	0.48
1:A:395:ASP:O	1:A:401:MSE:HE3	2.13	0.48
1:B:228:ASP:HB3	1:B:231:LYS:HE2	1.96	0.48
1:A:381:ARG:HG2	1:A:381:ARG:HH11	1.78	0.48
1:B:143:MSE:CE	1:B:146:VAL:HB	2.38	0.48
1:B:98:ILE:HG23	1:B:99:ASP:N	2.29	0.48
1:A:280:PHE:HA	1:A:283:ILE:HD13	1.95	0.48
1:A:97:GLY:HA2	1:A:346:ALA:O	2.14	0.48
1:A:95:ALA:HA	1:A:347:GLN:HE21	1.75	0.47
1:A:170:HIS:ND1	1:A:350:MSE:SE	2.97	0.47
1:A:384:ILE:O	1:A:388:ARG:HG2	2.14	0.47
1:B:388:ARG:HH21	1:B:426:LEU:HD11	1.79	0.47
1:A:318:ARG:HB3	4:A:1041:HOH:O	2.13	0.47
1:A:119:ARG:HG2	1:A:147:LEU:HD13	1.95	0.47
1:A:426:LEU:C	1:A:428:GLY:N	2.64	0.47
1:B:149:ALA:O	1:B:154:LYS:HB2	2.15	0.47
1:B:104:ILE:CG2	4:B:1080:HOH:O	2.57	0.47
1:A:196:THR:HB	1:A:199:GLU:HB3	1.96	0.47
1:A:314:ASP:N	1:A:315:PRO:CD	2.78	0.47
1:A:138:PHE:HD2	1:A:138:PHE:H	1.61	0.47
1:B:435:ILE:O	1:B:439:GLN:HG3	2.14	0.47
1:B:104:ILE:HG23	1:B:105:VAL:N	2.29	0.46
1:B:394:CYS:HB3	1:B:400:PHE:CD2	2.51	0.46
1:B:393:ASN:OD1	1:B:395:ASP:HB2	2.15	0.46
1:A:206:LEU:HD12	1:A:206:LEU:C	2.36	0.46
1:A:259:THR:OG1	1:A:262:GLN:HG3	2.15	0.46
1:B:184:GLN:HE22	1:B:227:PHE:HD1	1.64	0.46
1:A:169:MSE:HG3	1:A:361:LEU:HD22	1.97	0.46
1:B:283:ILE:O	1:B:287:PRO:HD3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:TYR:CD1	1:B:103:LEU:HD12	2.50	0.46
1:A:418:LYS:HB2	1:A:418:LYS:HZ2	1.78	0.46
1:B:381:ARG:HG3	1:B:386:GLU:O	2.16	0.45
1:A:100:TYR:CZ	1:A:338:THR:HB	2.51	0.45
1:A:207:THR:H	1:A:210:GLN:NE2	2.15	0.45
1:B:104:ILE:HG23	1:B:105:VAL:HG13	1.99	0.45
1:B:311:ILE:HG23	1:B:312:ASP:N	2.31	0.45
1:B:300:ARG:HG2	1:B:300:ARG:NH1	2.27	0.45
1:A:206:LEU:HD11	1:A:211:ALA:HB2	1.99	0.45
1:A:201:TYR:HE1	3:A:601:GOL:HO2	1.62	0.45
1:A:259:THR:HA	1:B:274:CYS:HA	1.99	0.45
1:A:372:VAL:O	1:A:376:ALA:HB3	2.16	0.45
1:B:184:GLN:NE2	1:B:227:PHE:HD1	2.15	0.45
1:B:163:GLY:HA3	1:B:200:LYS:HE2	1.99	0.45
1:A:435:ILE:O	1:A:439:GLN:HG3	2.17	0.45
1:B:98:ILE:HG23	1:B:99:ASP:H	1.82	0.45
1:B:414:GLU:OE2	1:B:417:ARG:NH2	2.51	0.44
1:A:139:SER:HB3	1:A:336:HIS:HB2	1.99	0.44
1:A:216:VAL:O	1:A:219:ALA:HB3	2.18	0.44
1:A:100:TYR:O	1:A:101:ASP:CB	2.66	0.44
1:A:114:LYS:HE3	1:A:114:LYS:HA	2.00	0.44
1:A:310:ALA:HB3	2:A:501:TYM:O2'	2.18	0.44
1:B:136:ILE:O	1:B:337:SER:HA	2.17	0.44
1:A:307:ILE:HD13	1:A:317:PHE:CE1	2.52	0.44
1:A:218:ASN:N	1:A:218:ASN:HD22	2.16	0.44
1:A:349:LYS:HB2	2:A:501:TYM:N6	2.32	0.44
1:B:104:ILE:CD1	1:B:139:SER:HB3	2.47	0.43
1:A:274:CYS:HB2	1:B:255:GLN:O	2.18	0.43
1:A:339:PHE:O	1:A:341:PRO:HD3	2.18	0.43
1:A:207:THR:OG1	1:A:210:GLN:HG3	2.19	0.43
1:A:403:LEU:HD21	1:A:434:LEU:HA	2.00	0.43
1:A:397:ASP:OD1	1:A:399:SER:HB2	2.18	0.43
1:A:130:HIS:H	1:A:130:HIS:CD2	2.37	0.43
1:B:260:PHE:CE2	1:B:264:LYS:HD2	2.54	0.43
1:B:142:ASP:OD1	1:B:321:ARG:NH2	2.52	0.43
1:A:228:ASP:OD1	1:A:231:LYS:HE3	2.19	0.43
1:B:314:ASP:N	1:B:315:PRO:CD	2.82	0.43
1:B:373:ASN:HD21	1:B:432:LYS:HE2	1.84	0.43
1:B:388:ARG:NH2	1:B:426:LEU:HD11	2.34	0.43
1:A:165:SER:CB	1:A:168:ALA:HB3	2.49	0.43
1:A:428:GLY:O	1:A:429:GLU:C	2.57	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:384:ILE:HG23	1:B:385:GLU:CD	2.38	0.43
1:A:103:LEU:CD2	1:A:311:ILE:HD13	2.49	0.43
1:A:165:SER:HB3	1:A:168:ALA:HB3	2.01	0.43
1:B:412:LYS:O	1:B:416:ILE:HG13	2.18	0.43
1:A:191:LEU:HD23	1:A:232:THR:HG23	2.00	0.42
1:A:162:ARG:HG3	2:A:501:TYM:O1P	2.18	0.42
1:B:326:ARG:HG3	1:B:326:ARG:NH1	2.34	0.42
1:B:104:ILE:CG2	1:B:105:VAL:N	2.82	0.42
1:A:101:ASP:HA	1:A:104:ILE:HD12	2.01	0.42
1:A:167:GLU:O	1:A:168:ALA:O	2.37	0.42
1:B:160:THR:O	1:B:193:ILE:HA	2.19	0.42
1:B:250:ASN:ND2	1:B:291:ASN:ND2	2.55	0.42
1:B:216:VAL:O	1:B:219:ALA:HB3	2.19	0.42
1:A:160:THR:O	1:A:193:ILE:HA	2.19	0.42
1:B:98:ILE:O	1:B:99:ASP:HB3	2.20	0.42
1:A:162:ARG:NH1	1:A:162:ARG:HG3	2.35	0.42
1:A:92:THR:HG23	1:A:347:GLN:HA	2.02	0.42
1:B:300:ARG:NH2	1:B:302:ASP:OD2	2.53	0.41
1:A:138:PHE:N	1:A:138:PHE:CD2	2.86	0.41
1:B:391:GLY:HA3	1:B:421:THR:O	2.20	0.41
1:B:164:PRO:HG3	1:B:195:MSE:HE3	2.01	0.41
1:A:162:ARG:NE	1:A:218:ASN:OD1	2.45	0.41
1:A:414:GLU:OE2	1:A:417:ARG:NH1	2.52	0.41
1:A:218:ASN:O	1:A:222:ILE:HG12	2.21	0.41
1:B:399:SER:HA	1:B:434:LEU:HD22	2.03	0.41
1:B:344:GLN:HB2	1:B:344:GLN:HE21	1.67	0.41
1:B:170:HIS:HA	1:B:360:PHE:HA	2.02	0.41
1:A:426:LEU:O	1:A:427:THR:C	2.59	0.41
1:B:240:TYR:HA	1:B:243:MSE:HG2	2.03	0.41
1:A:313:GLN:O	1:A:317:PHE:CD2	2.74	0.41
1:B:368:ILE:HD13	1:B:442:ILE:HD12	2.02	0.41
1:A:290:SER:HB3	1:A:303:ILE:HB	2.02	0.40
1:B:367:GLN:O	1:B:371:LYS:HG2	2.21	0.40
1:A:125:GLY:O	1:A:126:GLN:HG2	2.21	0.40
1:B:115:GLU:OE1	1:B:115:GLU:O	2.39	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	386/437 (88%)	360 (93%)	20 (5%)	6 (2%)	12	5
1	B	358/437 (82%)	336 (94%)	18 (5%)	4 (1%)	17	9
All	All	744/874 (85%)	696 (94%)	38 (5%)	10 (1%)	15	7

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	168	ALA
1	A	381	ARG
1	A	429	GLU
1	A	355	PRO
1	B	134	ARG
1	B	343	LEU
1	A	467	SER
1	B	303	ILE
1	B	372	VAL
1	A	87	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	343/369 (93%)	330 (96%)	13 (4%)	40	36
1	B	320/369 (87%)	313 (98%)	7 (2%)	60	62
All	All	663/738 (90%)	643 (97%)	20 (3%)	48	47

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	86	ASP
1	A	114	LYS
1	A	159	TYR
1	A	195	MSE
1	A	204	LYS
1	A	229	ILE
1	A	236	SER
1	A	295	GLN
1	A	298	ARG
1	A	335	LEU
1	A	417	ARG
1	A	425	MSE
1	A	430	LEU
1	B	218	ASN
1	B	236	SER
1	B	239	ASP
1	B	291	ASN
1	B	344	GLN
1	B	432	LYS
1	B	455	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	GLN
1	A	130	HIS
1	A	152	ASN
1	A	210	GLN
1	A	255	GLN
1	A	373	ASN
1	A	389	GLN
1	A	415	GLN
1	A	446	GLN
1	B	118	ASN
1	B	126	GLN
1	B	184	GLN
1	B	257	HIS
1	B	261	ASN
1	B	291	ASN
1	B	344	GLN
1	B	356	ASN

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 ⓘ

3 ligands 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$
2	TYM	A	501	-	33,41,41	0.96	3 (9%)	31,61,61	0.81	0
3	GOL	A	601	-	5,5,5	0.34	0	5,5,5	1.90	3 (60%)
3	GOL	A	602	-	5,5,5	0.28	0	5,5,5	1.88	3 (60%)

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
2	TYM	A	501	-	-	0/16/39/39	0/5/5/5
3	GOL	A	601	-	-	0/4/4/4	0/0/0/0
3	GOL	A	602	-	-	0/4/4/4	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	TYM	C2-N3	2.01	1.35	1.32
2	A	501	TYM	CH2-CZ3	2.11	1.43	1.38
2	A	501	TYM	CZ3-CE3	2.38	1.42	1.36

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	GOL	O2-C2-C3	2.19	118.69	108.65
3	A	602	GOL	O2-C2-C3	2.31	119.26	108.65
3	A	601	GOL	O2-C2-C1	2.33	119.32	108.65
3	A	602	GOL	O2-C2-C1	2.39	119.61	108.65
3	A	602	GOL	C3-C2-C1	2.55	121.13	111.12
3	A	601	GOL	C3-C2-C1	2.77	121.99	111.12

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	TYM	3	0
3	A	601	GOL	1	0
3	A	602	GOL	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	378/437 (86%)	0.90	64 (16%) 2 3	22, 49, 76, 84	11 (2%)
1	B	353/437 (80%)	0.63	38 (10%) 8 8	20, 51, 77, 88	6 (1%)
All	All	731/874 (83%)	0.77	102 (13%) 4 4	20, 50, 76, 88	17 (2%)

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	353	SER	5.6
1	A	390	PHE	5.5
1	B	390	PHE	5.5
1	A	167	GLU	5.5
1	B	384	ILE	5.2
1	A	355	PRO	5.2
1	B	98	ILE	5.0
1	A	352	ALA	4.9
1	B	100	TYR	4.6
1	A	90	VAL	4.5
1	A	85	VAL	4.4
1	A	168	ALA	4.4
1	B	166	SER	4.3
1	B	389	GLN	4.2
1	B	102	LYS	4.2
1	A	98	ILE	4.0
1	A	283	ILE	4.0
1	A	161	GLY	3.9
1	A	160	THR	3.9
1	B	97	GLY	3.9
1	B	103	LEU	3.8
1	A	391	GLY	3.7
1	A	469	ASP	3.7
1	A	374	LYS	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	104	ILE	3.5
1	B	168	ALA	3.5
1	B	385	GLU	3.5
1	A	159	TYR	3.5
1	A	360	PHE	3.4
1	B	426	LEU	3.4
1	A	93	SER	3.4
1	A	166	SER	3.4
1	B	311	ILE	3.4
1	B	167	GLU	3.4
1	A	175	ILE	3.4
1	A	172	GLY	3.3
1	A	94	SER	3.3
1	A	282	ALA	3.2
1	A	177	PHE	3.2
1	A	176	PRO	3.2
1	B	383	THR	3.2
1	A	192	VAL	3.2
1	A	107	PHE	3.1
1	B	344	GLN	3.1
1	A	84	PHE	3.0
1	A	251	VAL	3.0
1	A	354	ASP	3.0
1	A	307	ILE	3.0
1	A	392	GLY	2.9
1	A	308	PRO	2.9
1	A	351	SER	2.9
1	A	298	ARG	2.8
1	B	382	ASP	2.8
1	A	382	ASP	2.8
1	A	92	THR	2.8
1	A	103	LEU	2.8
1	A	82	GLU	2.8
1	A	384	ILE	2.8
1	A	97	GLY	2.7
1	B	283	ILE	2.7
1	B	396	VAL	2.7
1	B	343	LEU	2.7
1	B	251	VAL	2.7
1	A	383	THR	2.7
1	B	366	LYS	2.7
1	B	393	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	339	PHE	2.6
1	A	309	CYS	2.6
1	A	89	THR	2.5
1	B	356	ASN	2.5
1	B	409	ASP	2.5
1	A	285	ALA	2.5
1	B	377	PHE	2.5
1	A	430	LEU	2.5
1	B	388	ARG	2.5
1	A	171	VAL	2.5
1	B	114	LYS	2.5
1	A	422	SER	2.4
1	B	99	ASP	2.4
1	A	287	PRO	2.4
1	A	348	THR	2.4
1	A	193	ILE	2.4
1	B	387	HIS	2.4
1	B	316	TYR	2.4
1	B	265	GLY	2.3
1	A	180	THR	2.3
1	A	418	LYS	2.3
1	A	356	ASN	2.3
1	A	206	LEU	2.3
1	A	358	SER	2.2
1	A	102	LYS	2.2
1	A	153	LYS	2.2
1	B	424	ALA	2.2
1	B	308	PRO	2.2
1	B	105	VAL	2.2
1	A	405	PHE	2.2
1	B	158	LEU	2.1
1	A	306	LEU	2.1
1	A	339	PHE	2.1
1	A	280	PHE	2.1
1	A	284	GLN	2.0
1	A	424	ALA	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. 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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	GOL	A	602	6/6	0.62	0.33	4.33	55,57,57,57	0
3	GOL	A	601	6/6	0.80	0.29	3.53	52,54,54,55	0
2	TYM	A	501	37/37	0.87	0.24	0.32	45,53,60,60	8

6.5 Other polymers [i](#)

There are no such residues in this entry.