



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:58 AM UTC

PDB ID : 6FQB / pdb_00006fqb
Title : MurT/GatD peptidoglycan amidotransferase complex from *Streptococcus pneumoniae* R6
Authors : Morlot, C.; Contreras-Martel, C.; Leisico, F.; Straume, D.; Peters, K.; Hegnar, O.A.; Simon, N.; Villard, A.M.; Breukink, E.; Gravier-Pelletier, C.; Le Corre, L.; Vollmer, W.; Pietrancosta, N.; Havarstein, L.S.; Zapun, A.
Deposited on : 2018-02-13
Resolution : 3.00 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

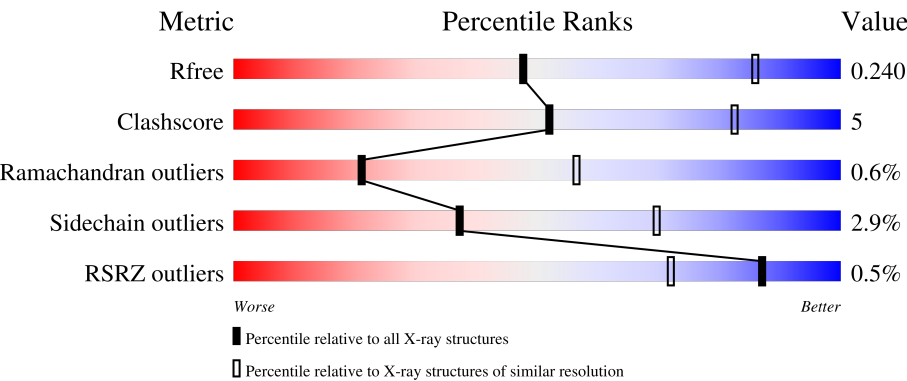
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	465	% 70% 10% • 18%
1	B	465	71% 9% • 19%
1	C	465	71% 9% • 19%
1	D	465	71% 9% • 18%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	E	260	85% 10%5%
2	F	260	84% 10%5%
2	G	260	86% 9%5%
2	H	260	83% 11%5%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 19727 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 Mur ligase family protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	383	Total	C	N	O	S	0	0	0
			2989	1906	497	573	13			
1	B	378	Total	C	N	O	S	0	0	0
			2952	1886	489	564	13			
1	C	375	Total	C	N	O	S	0	0	0
			2937	1876	485	563	13			
1	D	380	Total	C	N	O	S	0	0	0
			2977	1900	494	570	13			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-17	MET	-	initiating methionine	UNP A0A0B7LND9
A	-16	HIS	-	expression tag	UNP A0A0B7LND9
A	-15	HIS	-	expression tag	UNP A0A0B7LND9
A	-14	HIS	-	expression tag	UNP A0A0B7LND9
A	-13	HIS	-	expression tag	UNP A0A0B7LND9
A	-12	HIS	-	expression tag	UNP A0A0B7LND9
A	-11	HIS	-	expression tag	UNP A0A0B7LND9
A	-10	HIS	-	expression tag	UNP A0A0B7LND9
A	-9	HIS	-	expression tag	UNP A0A0B7LND9
A	-8	GLU	-	expression tag	UNP A0A0B7LND9
A	-7	ASN	-	expression tag	UNP A0A0B7LND9
A	-6	LEU	-	expression tag	UNP A0A0B7LND9
A	-5	TYR	-	expression tag	UNP A0A0B7LND9
A	-4	PHE	-	expression tag	UNP A0A0B7LND9
A	-3	GLN	-	expression tag	UNP A0A0B7LND9
A	-2	GLY	-	expression tag	UNP A0A0B7LND9
A	-1	SER	-	expression tag	UNP A0A0B7LND9
A	0	HIS	-	expression tag	UNP A0A0B7LND9
B	-17	MET	-	initiating methionine	UNP A0A0B7LND9
B	-16	HIS	-	expression tag	UNP A0A0B7LND9
B	-15	HIS	-	expression tag	UNP A0A0B7LND9

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP A0A0B7LND9
B	-13	HIS	-	expression tag	UNP A0A0B7LND9
B	-12	HIS	-	expression tag	UNP A0A0B7LND9
B	-11	HIS	-	expression tag	UNP A0A0B7LND9
B	-10	HIS	-	expression tag	UNP A0A0B7LND9
B	-9	HIS	-	expression tag	UNP A0A0B7LND9
B	-8	GLU	-	expression tag	UNP A0A0B7LND9
B	-7	ASN	-	expression tag	UNP A0A0B7LND9
B	-6	LEU	-	expression tag	UNP A0A0B7LND9
B	-5	TYR	-	expression tag	UNP A0A0B7LND9
B	-4	PHE	-	expression tag	UNP A0A0B7LND9
B	-3	GLN	-	expression tag	UNP A0A0B7LND9
B	-2	GLY	-	expression tag	UNP A0A0B7LND9
B	-1	SER	-	expression tag	UNP A0A0B7LND9
B	0	HIS	-	expression tag	UNP A0A0B7LND9
C	-17	MET	-	initiating methionine	UNP A0A0B7LND9
C	-16	HIS	-	expression tag	UNP A0A0B7LND9
C	-15	HIS	-	expression tag	UNP A0A0B7LND9
C	-14	HIS	-	expression tag	UNP A0A0B7LND9
C	-13	HIS	-	expression tag	UNP A0A0B7LND9
C	-12	HIS	-	expression tag	UNP A0A0B7LND9
C	-11	HIS	-	expression tag	UNP A0A0B7LND9
C	-10	HIS	-	expression tag	UNP A0A0B7LND9
C	-9	HIS	-	expression tag	UNP A0A0B7LND9
C	-8	GLU	-	expression tag	UNP A0A0B7LND9
C	-7	ASN	-	expression tag	UNP A0A0B7LND9
C	-6	LEU	-	expression tag	UNP A0A0B7LND9
C	-5	TYR	-	expression tag	UNP A0A0B7LND9
C	-4	PHE	-	expression tag	UNP A0A0B7LND9
C	-3	GLN	-	expression tag	UNP A0A0B7LND9
C	-2	GLY	-	expression tag	UNP A0A0B7LND9
C	-1	SER	-	expression tag	UNP A0A0B7LND9
C	0	HIS	-	expression tag	UNP A0A0B7LND9
D	-17	MET	-	initiating methionine	UNP A0A0B7LND9
D	-16	HIS	-	expression tag	UNP A0A0B7LND9
D	-15	HIS	-	expression tag	UNP A0A0B7LND9
D	-14	HIS	-	expression tag	UNP A0A0B7LND9
D	-13	HIS	-	expression tag	UNP A0A0B7LND9
D	-12	HIS	-	expression tag	UNP A0A0B7LND9
D	-11	HIS	-	expression tag	UNP A0A0B7LND9
D	-10	HIS	-	expression tag	UNP A0A0B7LND9
D	-9	HIS	-	expression tag	UNP A0A0B7LND9

Continued on next page...

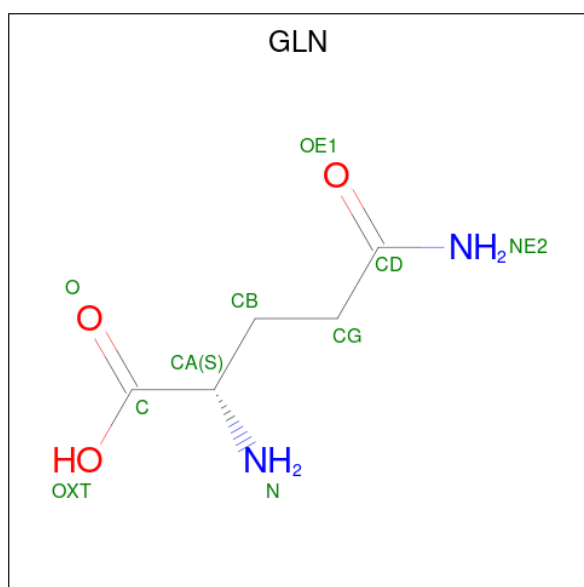
Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	-8	GLU	-	expression tag	UNP A0A0B7LND9
D	-7	ASN	-	expression tag	UNP A0A0B7LND9
D	-6	LEU	-	expression tag	UNP A0A0B7LND9
D	-5	TYR	-	expression tag	UNP A0A0B7LND9
D	-4	PHE	-	expression tag	UNP A0A0B7LND9
D	-3	GLN	-	expression tag	UNP A0A0B7LND9
D	-2	GLY	-	expression tag	UNP A0A0B7LND9
D	-1	SER	-	expression tag	UNP A0A0B7LND9
D	0	HIS	-	expression tag	UNP A0A0B7LND9

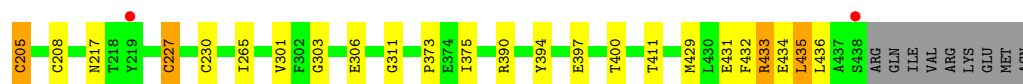
- Molecule 2 is a protein called Cobyric acid synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	247	Total	C	N	O	S	0	0	0
			1958	1242	326	385	5			
2	F	247	Total	C	N	O	S	0	0	0
			1958	1242	326	385	5			
2	G	247	Total	C	N	O	S	0	0	0
			1958	1242	326	385	5			
2	H	247	Total	C	N	O	S	0	0	0
			1958	1242	326	385	5			

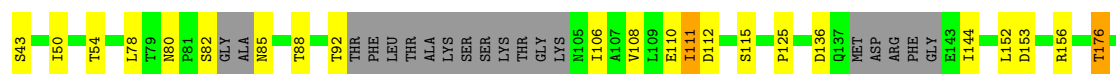
- Molecule 3 is GLUTAMINE (CCD ID: GLN) (formula: C₅H₁₀N₂O₃).



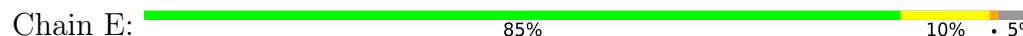
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	E	1	Total	C	N	O	0	0
			10	5	2	3		
3	F	1	Total	C	N	O	0	0
			10	5	2	3		
3	G	1	Total	C	N	O	0	0
			10	5	2	3		
3	H	1	Total	C	N	O	0	0
			10	5	2	3		



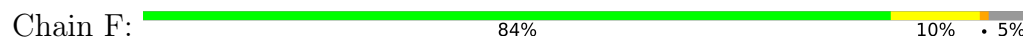
- Molecule 1: Mur ligase family protein



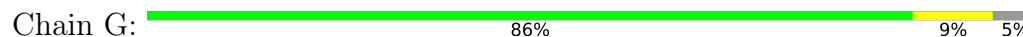
- Molecule 2: Cobyric acid synthase



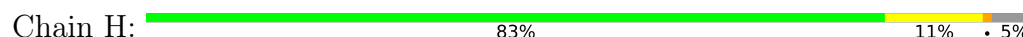
- Molecule 2: Cobyric acid synthase



- Molecule 2: Cobyric acid synthase



- Molecule 2: Cobyric acid synthase



ASP
VAL
LYS
SER
LYS
LYS
ALA
ASP
PHE
SER

4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	288.43Å 288.43Å 115.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 3.00 50.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	96.2 (50.00-3.00) 96.2 (50.00-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.16 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.189 , 0.228 0.205 , 0.240	Depositor DCC
R_{free} test set	6854 reflections (9.59%)	wwPDB-VP
Wilson B-factor (Å ²)	88.4	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 49.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.011 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	19727	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.35% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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	0.86	5/3043 (0.2%)	0.99	9/4133 (0.2%)
1	B	0.78	1/3005 (0.0%)	0.98	3/4080 (0.1%)
1	C	0.78	0/2990	0.98	3/4061 (0.1%)
1	D	0.84	5/3030 (0.2%)	0.96	1/4115 (0.0%)
2	E	0.83	1/2001 (0.0%)	0.96	2/2707 (0.1%)
2	F	0.76	1/2001 (0.0%)	0.96	2/2707 (0.1%)
2	G	0.82	1/2001 (0.0%)	0.99	3/2707 (0.1%)
2	H	0.81	1/2001 (0.0%)	1.01	6/2707 (0.2%)
All	All	0.81	15/20072 (0.1%)	0.98	29/27217 (0.1%)

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	136	ASP	CA-C	9.33	1.58	1.52
1	A	434	GLU	C-N	8.91	1.45	1.33
1	A	84	ALA	CA-C	6.67	1.60	1.52
1	A	85	ASN	CA-C	5.72	1.60	1.52
1	A	84	ALA	N-CA	5.36	1.52	1.46
1	A	407	GLN	CD-OE1	5.31	1.33	1.23
1	D	437	ALA	CA-C	5.25	1.59	1.52
2	G	58	HIS	CD2-NE2	-5.21	1.32	1.37
2	F	58	HIS	CD2-NE2	-5.17	1.32	1.37
1	D	436	LEU	CA-C	5.16	1.59	1.52
2	E	198	LYS	CA-C	-5.15	1.48	1.53
1	B	217	ASN	CG-OD1	5.15	1.33	1.23
2	H	196	HIS	CD2-NE2	-5.13	1.32	1.37
1	D	407	GLN	CD-OE1	5.08	1.33	1.23
1	D	217	ASN	CG-OD1	5.08	1.33	1.23

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	436	LEU	N-CA-C	-10.49	99.53	110.97

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	56	SER	N-CA-C	10.30	117.23	108.78
1	C	435	LEU	N-CA-C	-7.97	99.60	110.68
1	B	49	GLU	N-CA-C	-7.60	98.56	109.96
2	F	58	HIS	CB-CG-CD2	-6.68	122.52	131.20
2	G	58	HIS	CB-CG-CD2	-6.67	122.53	131.20
1	C	217	ASN	N-CA-C	6.59	119.64	108.90
2	H	196	HIS	CB-CG-CD2	-6.56	122.67	131.20
1	A	434	GLU	CA-C-N	6.42	130.98	122.06
1	A	434	GLU	C-N-CA	6.42	130.98	122.06
1	A	82	SER	CB-CA-C	-5.91	109.75	116.54
1	B	436	LEU	N-CA-C	-5.89	104.71	112.72
2	H	56	SER	CA-C-O	5.80	121.30	117.94
1	A	435	LEU	N-CA-C	-5.74	101.62	109.71
2	G	58	HIS	CB-CG-ND1	5.74	131.30	122.70
2	H	196	HIS	CB-CG-ND1	5.67	131.21	122.70
1	A	295	PHE	N-CA-C	-5.67	105.00	111.07
2	F	58	HIS	CB-CG-ND1	5.65	131.18	122.70
1	D	241	LEU	N-CA-C	-5.65	98.54	108.20
2	E	39	LEU	N-CA-C	-5.44	105.26	111.14
2	E	198	LYS	N-CA-C	-5.42	98.00	107.20
2	G	55	VAL	CB-CA-C	-5.40	104.92	111.88
1	A	80	ASN	CA-C-N	5.31	125.21	119.85
1	A	80	ASN	C-N-CA	5.31	125.21	119.85
1	A	105	ASN	N-CA-C	5.20	117.76	109.81
1	A	119	ILE	CB-CA-C	-5.18	105.25	111.88
2	H	55	VAL	CB-CA-C	-5.16	106.80	112.68
2	H	56	SER	CB-CA-C	-5.11	109.72	117.07
1	B	119	ILE	CB-CG1-CD1	5.01	124.33	113.80

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	2989	0	2986	48	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2952	0	2945	28	0
1	C	2937	0	2929	24	0
1	D	2977	0	2973	18	0
2	E	1958	0	1857	24	0
2	F	1958	0	1857	17	1
2	G	1958	0	1857	18	0
2	H	1958	0	1857	27	1
3	E	10	0	7	0	0
3	F	10	0	7	1	0
3	G	10	0	7	0	0
3	H	10	0	7	3	0
All	All	19727	0	19289	198	1

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 5.

All (198) 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:86:MET:HE1	1:A:122:TYR:CD2	1.94	1.02
1:A:116:LEU:HD12	1:A:116:LEU:O	1.70	0.91
2:H:219:ARG:HG3	2:H:219:ARG:HH11	1.34	0.91
1:B:115:SER:OG	1:B:119:ILE:CD1	2.22	0.88
1:B:81:PRO:O	1:B:83:GLY:N	2.07	0.88
2:F:25:ASN:C	2:F:26:LEU:HD23	1.98	0.87
1:C:431:GLU:O	1:C:434:GLU:HG2	1.75	0.86
1:A:87:ILE:HG22	1:A:87:ILE:O	1.76	0.83
1:A:86:MET:CE	1:A:122:TYR:CE2	2.65	0.80
1:A:86:MET:HE1	1:A:122:TYR:CE2	2.22	0.75
1:A:205:CYS:SG	1:A:208:CYS:N	2.60	0.74
1:C:306:GLU:OE2	1:C:433:ARG:NH1	2.21	0.74
1:D:205:CYS:SG	1:D:208:CYS:N	2.61	0.74
1:B:205:CYS:SG	1:B:208:CYS:N	2.60	0.73
1:A:86:MET:HE1	1:A:122:TYR:HD2	1.48	0.72
1:A:46:LYS:O	1:A:46:LYS:HG2	1.88	0.72
2:H:219:ARG:HG3	2:H:219:ARG:NH1	2.02	0.72
1:A:44:LEU:C	1:A:44:LEU:HD12	2.16	0.71
1:B:115:SER:OG	1:B:119:ILE:HD11	1.88	0.71
1:A:116:LEU:HD12	1:A:116:LEU:C	2.13	0.71
1:B:88:THR:O	1:B:89:GLY:C	2.34	0.70
1:A:130:ILE:HD11	1:A:151:ILE:HD13	1.72	0.70

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:23:TYR:HA	2:G:56:SER:HB2	1.73	0.70
1:C:205:CYS:SG	1:C:208:CYS:N	2.62	0.69
2:F:130:VAL:HG23	2:F:131:MET:HG3	1.74	0.69
1:A:296:ASP:HA	1:A:299:ARG:HD3	1.74	0.69
2:E:198:LYS:O	2:E:198:LYS:HG3	1.93	0.69
2:E:116:TYR:OH	2:G:116:TYR:OH	1.92	0.69
2:G:23:TYR:HA	2:G:56:SER:CB	2.23	0.68
2:E:23:TYR:CD1	2:E:82:ILE:HD13	2.28	0.67
2:H:5:SER:HB3	2:H:54:ILE:HG13	1.75	0.67
1:B:115:SER:OG	1:B:119:ILE:HD12	1.93	0.65
1:D:296:ASP:HA	1:D:299:ARG:HD3	1.77	0.65
2:H:170:LEU:HD23	2:H:195:VAL:HG23	1.78	0.65
2:H:198:LYS:O	2:H:198:LYS:HG3	1.96	0.65
2:F:26:LEU:HD23	2:F:26:LEU:N	2.06	0.65
1:A:44:LEU:HD12	1:A:44:LEU:O	1.97	0.65
1:A:115:SER:O	1:A:119:ILE:HD12	1.97	0.64
1:A:211:ILE:HD13	2:E:25:ASN:HD22	1.63	0.63
2:G:128:LEU:HB3	2:G:130:VAL:HG13	1.80	0.63
2:H:131:MET:HE2	2:H:174:GLN:NE2	2.14	0.62
2:E:167:ARG:HD2	2:E:188:GLU:OE1	1.99	0.62
1:B:87:ILE:HG22	1:B:87:ILE:O	2.00	0.61
1:A:116:LEU:HD11	1:A:120:CYS:SG	2.39	0.61
1:C:135:ARG:NH1	1:C:136:ASP:O	2.34	0.60
1:B:88:THR:O	1:B:90:ILE:N	2.34	0.60
1:B:147:THR:O	1:B:148:TYR:C	2.43	0.60
2:H:107:CYS:SG	3:H:301:GLN:OE1	2.61	0.59
2:E:167:ARG:HH11	2:E:188:GLU:CD	2.10	0.59
2:G:165:GLN:HG3	2:G:166:GLY:N	2.18	0.59
2:F:27:MET:HE1	2:F:79:GLN:HG3	1.84	0.58
2:F:38:MET:HE2	2:F:210:LEU:O	2.03	0.58
1:A:156:ARG:NH2	1:A:176:THR:O	2.37	0.58
2:H:107:CYS:HG	3:H:301:GLN:CD	2.11	0.58
2:G:54:ILE:HG22	2:G:54:ILE:O	2.01	0.58
1:D:50:ILE:HG23	1:D:125:PRO:HA	1.86	0.57
1:D:156:ARG:NH2	1:D:176:THR:O	2.37	0.57
1:B:156:ARG:NH2	1:B:176:THR:O	2.37	0.57
1:B:50:ILE:HG23	1:B:125:PRO:HA	1.85	0.57
1:B:115:SER:O	1:B:119:ILE:HD12	2.04	0.57
1:C:156:ARG:NH2	1:C:176:THR:O	2.37	0.57
2:H:55:VAL:O	2:H:55:VAL:HG12	2.04	0.57
2:H:170:LEU:HD11	2:H:186:ASN:ND2	2.20	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:ILE:O	1:A:122:TYR:N	2.37	0.57
1:A:50:ILE:HG23	1:A:125:PRO:HA	1.86	0.56
2:F:165:GLN:HG3	2:F:166:GLY:N	2.20	0.56
2:G:42:VAL:HG12	2:G:221:VAL:HG11	1.86	0.56
1:D:80:ASN:O	1:D:82:SER:N	2.38	0.56
2:H:107:CYS:SG	3:H:301:GLN:CD	2.88	0.56
1:C:50:ILE:HG23	1:C:125:PRO:HA	1.86	0.56
1:D:227:CYS:SG	1:D:230:CYS:N	2.79	0.56
2:E:42:VAL:HG12	2:E:221:VAL:HG11	1.88	0.56
1:A:227:CYS:SG	1:A:230:CYS:N	2.78	0.56
2:F:42:VAL:HG12	2:F:221:VAL:HG11	1.88	0.56
2:G:5:SER:HB3	2:G:54:ILE:HG13	1.88	0.55
1:C:429:MET:O	1:C:433:ARG:HG3	2.06	0.55
2:G:130:VAL:HG23	2:G:131:MET:HG3	1.89	0.55
1:D:54:THR:HG22	1:D:111:ILE:O	2.07	0.55
2:H:38:MET:HE2	2:H:210:LEU:O	2.06	0.55
1:A:86:MET:HE2	1:A:122:TYR:HE2	1.72	0.55
2:F:27:MET:HE2	2:F:82:ILE:HG13	1.89	0.55
1:C:227:CYS:SG	1:C:230:CYS:N	2.79	0.54
2:E:27:MET:HE1	2:E:79:GLN:HG3	1.88	0.54
1:B:227:CYS:SG	1:B:230:CYS:N	2.78	0.54
2:E:167:ARG:NH1	2:E:188:GLU:OE1	2.37	0.54
2:E:40:LYS:HG3	2:E:50:VAL:HG21	1.90	0.54
2:F:25:ASN:O	2:F:26:LEU:HD23	2.07	0.54
1:C:434:GLU:HA	1:C:434:GLU:OE1	2.07	0.54
2:F:128:LEU:HB3	2:F:130:VAL:HG13	1.89	0.53
1:A:435:LEU:C	1:A:437:ALA:H	2.16	0.53
2:H:40:LYS:HG3	2:H:50:VAL:HG21	1.88	0.53
2:G:197:TYR:CD2	2:G:198:LYS:HG2	2.44	0.53
2:H:8:SER:OG	2:H:50:VAL:HG22	2.09	0.53
2:E:165:GLN:HG3	2:E:166:GLY:H	1.74	0.53
2:F:167:ARG:HD2	2:F:188:GLU:OE1	2.09	0.52
1:A:130:ILE:CD1	1:A:151:ILE:HD13	2.39	0.52
1:C:431:GLU:O	1:C:434:GLU:CG	2.52	0.52
1:A:119:ILE:HD12	1:A:119:ILE:H	1.75	0.52
1:A:119:ILE:O	1:A:121:ASP:N	2.42	0.52
2:E:165:GLN:HG3	2:E:166:GLY:N	2.25	0.52
1:A:211:ILE:HD13	2:E:25:ASN:ND2	2.24	0.52
2:E:8:SER:OG	2:E:50:VAL:HG22	2.09	0.52
1:A:148:TYR:O	1:A:151:ILE:HG13	2.10	0.51
1:A:46:LYS:O	1:A:46:LYS:CG	2.58	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:74:TYR:CD1	1:B:106:ILE:HG21	2.45	0.51
2:E:167:ARG:NH1	2:E:188:GLU:CD	2.68	0.51
2:H:219:ARG:NH1	2:H:219:ARG:CG	2.73	0.51
2:E:26:LEU:HB2	2:E:82:ILE:HD11	1.92	0.51
1:B:80:ASN:ND2	1:B:84:ALA:HB3	2.26	0.51
2:E:198:LYS:O	2:E:198:LYS:CG	2.56	0.50
1:C:44:LEU:HD13	1:C:46:LYS:HG3	1.92	0.50
2:H:106:ILE:O	2:H:107:CYS:C	2.53	0.50
1:B:81:PRO:C	1:B:83:GLY:N	2.69	0.50
2:G:196:HIS:C	2:G:196:HIS:HD1	2.20	0.50
2:F:55:VAL:HG12	2:F:55:VAL:O	2.10	0.49
1:B:80:ASN:HD22	1:B:84:ALA:HB3	1.76	0.49
1:B:88:THR:C	1:B:90:ILE:N	2.71	0.49
1:C:85:ASN:O	1:C:88:THR:HG23	2.12	0.49
2:G:23:TYR:HA	2:G:56:SER:HB3	1.95	0.49
1:A:87:ILE:O	1:A:87:ILE:CG2	2.49	0.49
1:A:115:SER:O	1:A:119:ILE:CD1	2.60	0.48
2:H:57:LEU:C	2:H:59:ASP:H	2.21	0.48
1:A:82:SER:OG	1:A:83:GLY:N	2.41	0.48
1:D:85:ASN:O	1:D:88:THR:HG23	2.13	0.48
1:A:80:ASN:OD1	1:A:80:ASN:N	2.40	0.47
1:A:208:CYS:HB3	1:A:230:CYS:HB3	1.71	0.47
2:E:23:TYR:CD1	2:E:82:ILE:CD1	2.97	0.47
2:H:57:LEU:O	2:H:59:ASP:N	2.48	0.47
2:H:198:LYS:O	2:H:198:LYS:CG	2.63	0.47
2:E:128:LEU:HB3	2:E:130:VAL:HG13	1.96	0.47
1:A:119:ILE:HG22	1:A:123:ILE:HG12	1.97	0.47
2:H:201:PHE:CD1	2:H:220:LEU:HD21	2.50	0.47
1:A:116:LEU:CD1	1:A:120:CYS:SG	3.02	0.46
2:G:23:TYR:CD1	2:G:82:ILE:HD13	2.50	0.46
2:E:143:PHE:HD2	2:E:164:HIS:CE1	2.32	0.46
1:C:301:VAL:O	1:C:301:VAL:HG12	2.15	0.46
1:D:188:LEU:N	1:D:188:LEU:HD12	2.31	0.46
1:A:265:ILE:HD13	1:A:303:GLY:O	2.15	0.46
1:D:111:ILE:CG2	1:D:112:ASP:N	2.77	0.46
1:A:390:ARG:NH1	2:E:245:GLU:OE1	2.49	0.45
2:E:1:MET:N	2:E:59:ASP:OD1	2.39	0.45
1:B:147:THR:O	1:B:150:MET:N	2.50	0.45
1:A:86:MET:CE	1:A:122:TYR:CD2	2.77	0.45
2:F:23:TYR:CD1	2:F:82:ILE:HD13	2.52	0.45
1:D:188:LEU:N	1:D:188:LEU:CD1	2.80	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:265:ILE:HD13	1:D:303:GLY:O	2.17	0.45
2:H:23:TYR:CD1	2:H:82:ILE:HD13	2.52	0.45
1:B:265:ILE:HD13	1:B:303:GLY:O	2.16	0.44
2:G:56:SER:OG	2:G:57:LEU:N	2.50	0.44
2:H:5:SER:HB3	2:H:54:ILE:CG1	2.44	0.44
1:A:67:VAL:HG22	1:A:108:VAL:CG2	2.47	0.44
1:A:116:LEU:HG	1:A:154:ALA:CB	2.47	0.44
1:C:265:ILE:HD13	1:C:303:GLY:O	2.17	0.44
1:A:86:MET:CE	1:A:122:TYR:HE2	2.18	0.44
2:H:130:VAL:HG23	2:H:131:MET:HG3	1.99	0.44
2:H:131:MET:HE2	2:H:174:GLN:HE21	1.82	0.44
2:E:57:LEU:HD23	2:E:57:LEU:HA	1.73	0.44
1:C:432:PHE:CD2	1:C:432:PHE:O	2.70	0.44
1:C:81:PRO:O	1:C:82:SER:HB2	2.18	0.44
1:D:186:PHE:CD2	1:D:218:THR:CG2	3.01	0.43
1:B:208:CYS:HB3	1:B:230:CYS:HB3	1.72	0.43
1:A:151:ILE:HD12	1:A:152:LEU:N	2.33	0.43
1:D:78:LEU:HD21	1:D:92:THR:HG22	2.01	0.43
1:A:108:VAL:O	1:A:108:VAL:HG23	2.19	0.42
1:C:208:CYS:HB3	1:C:230:CYS:HB3	1.70	0.42
1:B:390:ARG:NH1	2:F:245:GLU:OE1	2.52	0.42
2:F:41:TYR:CZ	2:F:45:LYS:HE2	2.54	0.42
1:C:153:ASP:OD1	1:C:156:ARG:NH1	2.53	0.42
1:A:119:ILE:C	1:A:121:ASP:N	2.77	0.42
2:F:57:LEU:C	2:F:59:ASP:H	2.26	0.42
2:E:40:LYS:CG	2:E:50:VAL:HG21	2.50	0.42
1:C:81:PRO:O	1:C:82:SER:CB	2.68	0.42
2:G:57:LEU:HD12	2:G:57:LEU:HA	1.92	0.42
2:H:43:ALA:HB3	2:H:50:VAL:HG11	2.02	0.42
1:A:409:LEU:O	1:A:409:LEU:HD23	2.20	0.41
1:C:375:ILE:HD12	1:C:394:TYR:CE2	2.55	0.41
1:B:54:THR:HG22	1:B:111:ILE:O	2.20	0.41
1:C:433:ARG:C	1:C:435:LEU:H	2.29	0.41
1:D:111:ILE:HG22	1:D:112:ASP:N	2.32	0.41
1:D:375:ILE:HD12	1:D:394:TYR:CE2	2.55	0.41
1:B:153:ASP:OD1	1:B:156:ARG:NH1	2.54	0.41
1:D:153:ASP:OD1	1:D:156:ARG:NH1	2.53	0.41
2:H:40:LYS:CG	2:H:50:VAL:HG21	2.50	0.41
1:A:153:ASP:OD1	1:A:156:ARG:NH1	2.54	0.41
1:C:54:THR:HG22	1:C:111:ILE:O	2.21	0.41
1:C:390:ARG:NH1	2:G:245:GLU:OE1	2.54	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:319:LEU:CA	1:D:429:MET:HE1	2.51	0.41
1:A:54:THR:HG22	1:A:111:ILE:O	2.20	0.41
1:A:165:LEU:HD13	1:A:172:PHE:HB3	2.02	0.41
1:B:165:LEU:HD13	1:B:172:PHE:HB3	2.02	0.41
1:B:375:ILE:HD12	1:B:394:TYR:CE2	2.56	0.41
2:F:165:GLN:HG3	3:F:301:GLN:OXT	2.21	0.41
1:B:319:LEU:CA	1:B:429:MET:HE1	2.51	0.41
2:G:46:LEU:HD11	2:G:241:ILE:CD1	2.51	0.41
1:B:409:LEU:HD23	1:B:409:LEU:O	2.20	0.40
2:G:198:LYS:HG3	2:G:198:LYS:O	2.21	0.40
1:C:434:GLU:OE1	1:C:434:GLU:CA	2.69	0.40
2:H:57:LEU:C	2:H:59:ASP:N	2.78	0.40

All (1) 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)
2:F:116:TYR:OH	2:H:116:TYR:OH[2_555]	2.05	0.15

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	377/465 (81%)	342 (91%)	32 (8%)	3 (1%)	16	50
1	B	370/465 (80%)	334 (90%)	33 (9%)	3 (1%)	16	50
1	C	367/465 (79%)	332 (90%)	32 (9%)	3 (1%)	16	50
1	D	372/465 (80%)	338 (91%)	32 (9%)	2 (0%)	24	60
2	E	245/260 (94%)	230 (94%)	15 (6%)	0	100	100
2	F	245/260 (94%)	228 (93%)	16 (6%)	1 (0%)	30	65
2	G	245/260 (94%)	229 (94%)	16 (6%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	245/260 (94%)	230 (94%)	13 (5%)	2 (1%)	16	50
All	All	2466/2900 (85%)	2263 (92%)	189 (8%)	14 (1%)	21	56

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	58	HIS
1	B	89	GLY
2	F	58	HIS
1	A	120	CYS
1	D	373	PRO
2	H	197	TYR
1	A	373	PRO
1	B	373	PRO
1	C	86	MET
1	C	311	GLY
1	C	373	PRO
1	D	311	GLY
1	A	311	GLY
1	B	311	GLY

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	323/395 (82%)	310 (96%)	13 (4%)	28	62
1	B	318/395 (80%)	310 (98%)	8 (2%)	42	72
1	C	319/395 (81%)	307 (96%)	12 (4%)	29	63
1	D	323/395 (82%)	308 (95%)	15 (5%)	24	58
2	E	207/219 (94%)	205 (99%)	2 (1%)	68	84
2	F	207/219 (94%)	202 (98%)	5 (2%)	43	73
2	G	207/219 (94%)	204 (99%)	3 (1%)	59	80
2	H	207/219 (94%)	203 (98%)	4 (2%)	50	76

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	2111/2456 (86%)	2049 (97%)	62 (3%)	37 70

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	88	THR
1	A	106	ILE
1	A	115	SER
1	A	116	LEU
1	A	144	ILE
1	A	152	LEU
1	A	176	THR
1	A	205	CYS
1	A	227	CYS
1	A	354	ILE
1	A	397	GLU
1	A	400	THR
1	A	411	THR
2	E	82	ILE
2	E	97	ILE
1	B	147	THR
1	B	152	LEU
1	B	188	LEU
1	B	205	CYS
1	B	227	CYS
1	B	397	GLU
1	B	400	THR
1	B	411	THR
2	F	26	LEU
2	F	82	ILE
2	F	97	ILE
2	F	139	THR
2	F	227	LYS
1	C	43	SER
1	C	106	ILE
1	C	115	SER
1	C	144	ILE
1	C	152	LEU
1	C	176	THR
1	C	205	CYS
1	C	227	CYS
1	C	397	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	400	THR
1	C	411	THR
1	C	433	ARG
2	G	82	ILE
2	G	97	ILE
2	G	139	THR
1	D	43	SER
1	D	106	ILE
1	D	108	VAL
1	D	110	GLU
1	D	111	ILE
1	D	115	SER
1	D	144	ILE
1	D	152	LEU
1	D	176	THR
1	D	204	LEU
1	D	205	CYS
1	D	227	CYS
1	D	397	GLU
1	D	400	THR
1	D	411	THR
2	H	54	ILE
2	H	82	ILE
2	H	97	ILE
2	H	139	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	GLN
1	A	85	ASN
1	A	132	ASN
1	A	305	GLN
1	A	404	ASN
2	E	25	ASN
2	E	141	ASN
1	B	132	ASN
1	B	149	ASN
1	B	259	GLN
1	B	305	GLN
1	B	328	GLN
1	B	414	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	F	75	GLN
1	C	132	ASN
1	C	137	GLN
1	C	250	ASN
1	C	259	GLN
1	C	264	GLN
1	C	404	ASN
2	G	16	GLN
2	G	49	HIS
2	G	98	GLN
2	G	111	GLN
2	G	141	ASN
1	D	132	ASN
1	D	137	GLN
1	D	199	ASN
1	D	259	GLN
1	D	305	GLN
2	H	16	GLN
2	H	49	HIS
2	H	111	GLN
2	H	141	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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
3	GLN	G	301	-	8,9,9	0.88	1 (12%)	8,11,11	1.10	1 (12%)
3	GLN	E	301	-	8,9,9	0.90	1 (12%)	8,11,11	1.51	1 (12%)
3	GLN	F	301	-	8,9,9	0.90	1 (12%)	8,11,11	1.47	1 (12%)
3	GLN	H	301	-	8,9,9	0.83	1 (12%)	8,11,11	1.09	1 (12%)

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
3	GLN	G	301	-	-	6/9/9/9	-
3	GLN	E	301	-	-	6/9/9/9	-
3	GLN	F	301	-	-	6/9/9/9	-
3	GLN	H	301	-	-	4/9/9/9	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	301	GLN	OXT-C	-2.29	1.23	1.30
3	E	301	GLN	OXT-C	-2.19	1.23	1.30
3	H	301	GLN	OXT-C	-2.18	1.23	1.30
3	G	301	GLN	OXT-C	-2.00	1.24	1.30

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	301	GLN	OXT-C-O	-4.13	114.71	124.08
3	F	301	GLN	OXT-C-O	-4.12	114.72	124.08
3	G	301	GLN	OXT-C-O	-3.05	117.16	124.08
3	H	301	GLN	OXT-C-O	-3.04	117.18	124.08

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	E	301	GLN	O-C-CA-N
3	F	301	GLN	O-C-CA-N
3	G	301	GLN	O-C-CA-N
3	H	301	GLN	O-C-CA-N
3	E	301	GLN	OXT-C-CA-N
3	F	301	GLN	OXT-C-CA-N
3	H	301	GLN	OXT-C-CA-N
3	G	301	GLN	OXT-C-CA-N
3	F	301	GLN	OE1-CD-CG-CB
3	E	301	GLN	NE2-CD-CG-CB
3	G	301	GLN	OXT-C-CA-CB
3	F	301	GLN	NE2-CD-CG-CB
3	E	301	GLN	O-C-CA-CB
3	E	301	GLN	OXT-C-CA-CB
3	E	301	GLN	OE1-CD-CG-CB
3	G	301	GLN	O-C-CA-CB
3	G	301	GLN	OE1-CD-CG-CB
3	G	301	GLN	NE2-CD-CG-CB
3	F	301	GLN	O-C-CA-CB
3	F	301	GLN	OXT-C-CA-CB
3	H	301	GLN	OXT-C-CA-CB
3	H	301	GLN	NE2-CD-CG-CB

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	301	GLN	1	0
3	H	301	GLN	3	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	383/465 (82%)	-0.34	4 (1%) 79 59	30, 83, 130, 162	0
1	B	378/465 (81%)	-0.28	1 (0%) 90 79	66, 94, 132, 161	0
1	C	375/465 (80%)	-0.29	2 (0%) 87 72	30, 90, 134, 165	0
1	D	380/465 (81%)	-0.33	2 (0%) 87 72	58, 88, 131, 162	0
2	E	247/260 (95%)	-0.53	1 (0%) 88 76	46, 67, 95, 143	0
2	F	247/260 (95%)	-0.36	1 (0%) 88 76	56, 85, 124, 143	0
2	G	247/260 (95%)	-0.44	1 (0%) 88 76	51, 78, 122, 146	0
2	H	247/260 (95%)	-0.44	1 (0%) 88 76	51, 82, 117, 143	0
All	All	2504/2900 (86%)	-0.36	13 (0%) 87 72	30, 85, 127, 165	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	82	SER	5.9
1	D	441	ILE	3.4
2	F	1	MET	2.7
2	G	1	MET	2.7
1	C	438	SER	2.7
2	H	1	MET	2.6
1	A	81	PRO	2.6
1	A	436	LEU	2.3
2	E	25	ASN	2.3
1	C	219	TYR	2.2
1	D	439	ARG	2.2
1	A	142	GLY	2.1
1	B	90	ILE	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GLN	F	301	10/10	0.79	0.20	111,140,151,153	0
3	GLN	H	301	10/10	0.83	0.16	97,120,132,139	0
3	GLN	G	301	10/10	0.84	0.16	102,124,137,140	0
3	GLN	E	301	10/10	0.90	0.13	88,105,125,125	0

6.5 Other polymers [i](#)

There are no such residues in this entry.