



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 23, 2021 – 04:10 PM EST

PDB ID : 1YBD
Title : Crystal structure analysis of uridylylate kinase from *Neisseria meningitidis*
Authors : Kumaran, D.; Swaminathan, S.; Burley, S.K.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2004-12-20
Resolution : 2.60 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 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 : 2.16
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) : 2.16

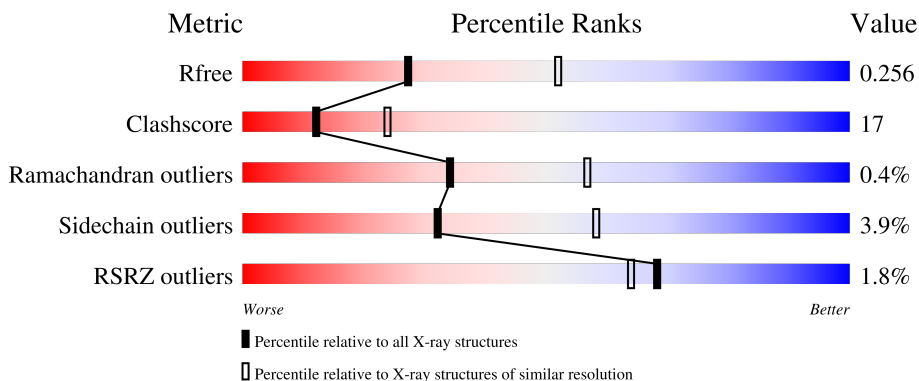
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	239	
1	B	239	
1	C	239	

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	250	-	X	-	-
3	GOL	B	251	-	X	-	-
3	GOL	C	252	-	X	-	-

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 5446 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 Uridylate kinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	236	1771	1115	302	340	3	11	0	0	0
1	B	236	1771	1115	302	340	3	11	0	0	0
1	C	236	1771	1115	302	340	3	11	0	0	0

There are 36 discrepancies between the modelled and reference sequences:

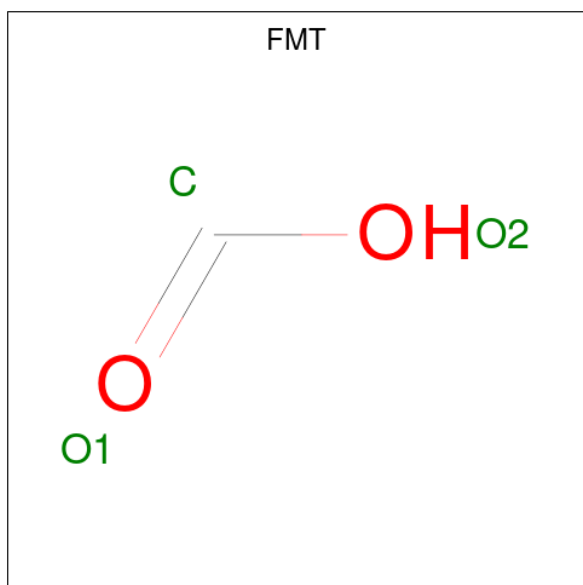
Chain	Residue	Modelled	Actual	Comment	Reference
A	3	MSE	MET	modified residue	UNP P65932
A	22	MSE	MET	modified residue	UNP P65932
A	47	MSE	MET	modified residue	UNP P65932
A	71	MSE	MET	modified residue	UNP P65932
A	79	MSE	MET	modified residue	UNP P65932
A	81	MSE	MET	modified residue	UNP P65932
A	82	MSE	MET	modified residue	UNP P65932
A	86	MSE	MET	modified residue	UNP P65932
A	110	MSE	MET	modified residue	UNP P65932
A	156	MSE	MET	modified residue	UNP P65932
A	161	MSE	MET	modified residue	UNP P65932
A	200	MSE	MET	modified residue	UNP P65932
B	3	MSE	MET	modified residue	UNP P65932
B	22	MSE	MET	modified residue	UNP P65932
B	47	MSE	MET	modified residue	UNP P65932
B	71	MSE	MET	modified residue	UNP P65932
B	79	MSE	MET	modified residue	UNP P65932
B	81	MSE	MET	modified residue	UNP P65932
B	82	MSE	MET	modified residue	UNP P65932
B	86	MSE	MET	modified residue	UNP P65932
B	110	MSE	MET	modified residue	UNP P65932
B	156	MSE	MET	modified residue	UNP P65932
B	161	MSE	MET	modified residue	UNP P65932

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Chain	Residue	Modelled	Actual	Comment	Reference
B	200	MSE	MET	modified residue	UNP P65932
C	3	MSE	MET	modified residue	UNP P65932
C	22	MSE	MET	modified residue	UNP P65932
C	47	MSE	MET	modified residue	UNP P65932
C	71	MSE	MET	modified residue	UNP P65932
C	79	MSE	MET	modified residue	UNP P65932
C	81	MSE	MET	modified residue	UNP P65932
C	82	MSE	MET	modified residue	UNP P65932
C	86	MSE	MET	modified residue	UNP P65932
C	110	MSE	MET	modified residue	UNP P65932
C	156	MSE	MET	modified residue	UNP P65932
C	161	MSE	MET	modified residue	UNP P65932
C	200	MSE	MET	modified residue	UNP P65932

- Molecule 2 is FORMIC ACID (three-letter code: FMT) (formula: CH_2O_2).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	3	1	2	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $\text{C}_3\text{H}_8\text{O}_3$).



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

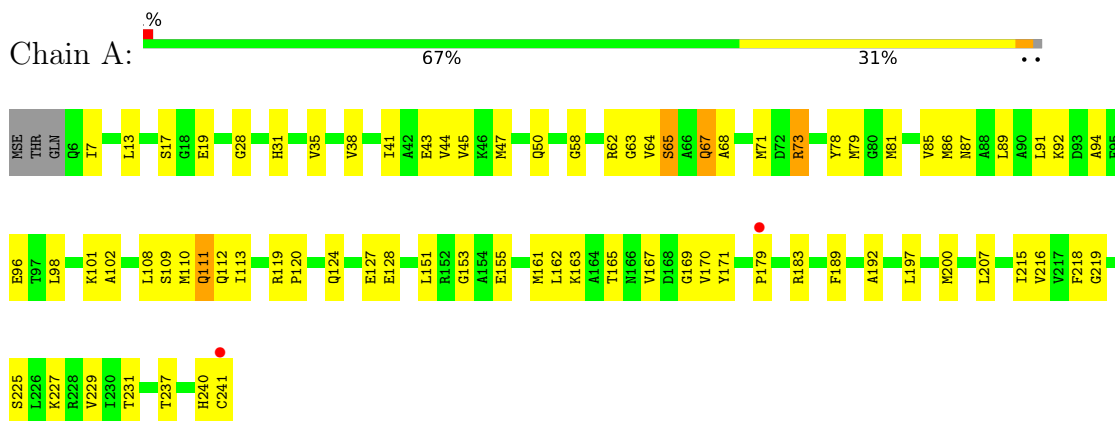
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	40	Total O 40 40	0	0
4	B	38	Total O 38 38	0	0
4	C	34	Total O 34 34	0	0

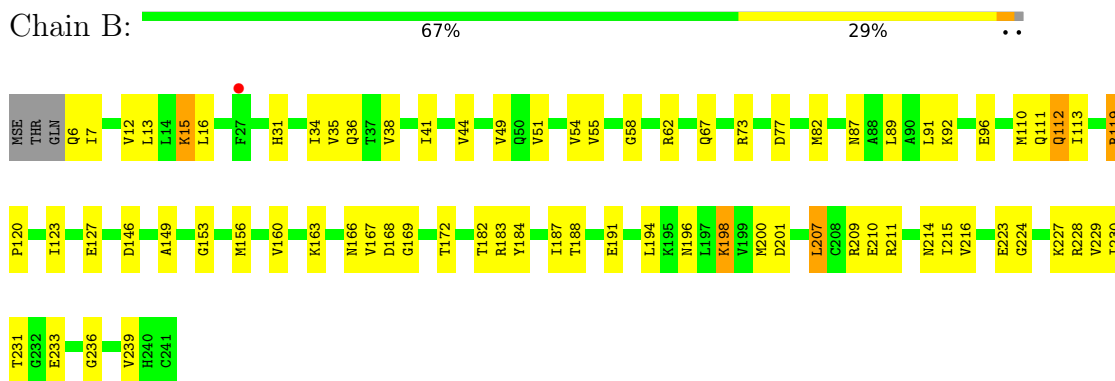
3 Residue-property plots

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.

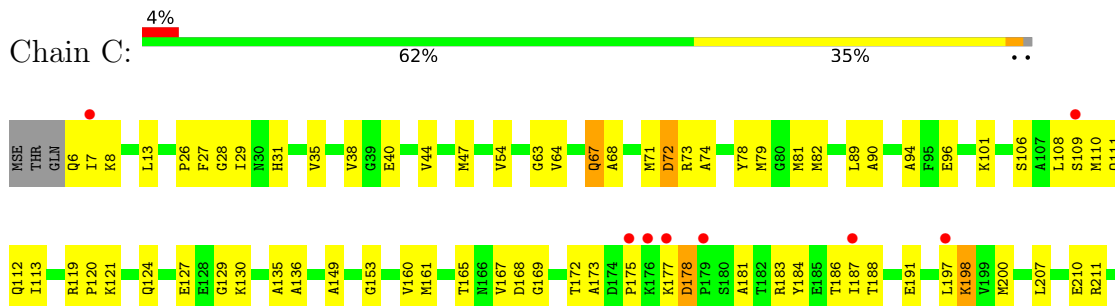
- Molecule 1: Uridylate kinase

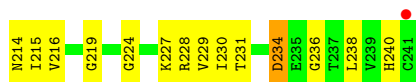


- Molecule 1: Uridylate kinase



- Molecule 1: Uridylate kinase





4 Data and refinement statistics i

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	158.29Å 158.29Å 58.74Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.17 – 2.60 47.17 – 2.55	Depositor EDS
% Data completeness (in resolution range)	96.9 (47.17-2.60) 95.3 (47.17-2.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.61 (at 2.54Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.206 , 0.259 0.205 , 0.256	Depositor DCC
R_{free} test set	1247 reflections (4.56%)	wwPDB-VP
Wilson B-factor (Å ²)	38.7	Xtrriage
Anisotropy	0.256	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 42.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.034 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5446	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.71% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.37	0/1781	0.61	0/2381
1	B	0.35	0/1781	0.62	0/2381
1	C	0.36	0/1781	0.60	0/2381
All	All	0.36	0/5343	0.61	0/7143

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1771	0	1810	71	0
1	B	1771	0	1810	53	0
1	C	1771	0	1810	83	0
2	A	3	0	1	0	0
3	A	6	0	4	1	0
3	B	6	0	4	0	0
3	C	6	0	4	0	0
4	A	40	0	0	0	0
4	B	38	0	0	0	0
4	C	34	0	0	2	0
All	All	5446	0	5443	184	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 17.

All (184) 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:109:SER:H	1:C:111:GLN:HE21	1.03	1.00
1:C:111:GLN:HG2	1:C:112:GLN:HG3	1.50	0.92
1:A:7:ILE:HD11	1:A:47:MSE:HE2	1.54	0.90
1:C:68:ALA:HA	1:C:71:MSE:HE2	1.53	0.88
1:A:109:SER:H	1:C:111:GLN:NE2	1.73	0.86
1:C:200:MSE:HE3	1:C:215:ILE:HD13	1.67	0.77
1:A:109:SER:N	1:C:111:GLN:HE21	1.81	0.75
1:C:119:ARG:HB3	1:C:120:PRO:HD3	1.71	0.72
1:C:165:THR:HG23	1:C:167:VAL:H	1.54	0.71
1:B:15:LYS:C	1:B:15:LYS:HD3	2.11	0.71
1:C:168:ASP:HA	1:C:219:GLY:H	1.56	0.70
1:A:108:LEU:HD13	1:C:113:ILE:HD11	1.73	0.69
1:A:67:GLN:NE2	1:C:31:HIS:ND1	2.42	0.67
1:A:179:PRO:HG2	1:B:187:ILE:HG22	1.76	0.67
1:C:89:LEU:HD11	1:C:113:ILE:HG23	1.75	0.67
1:A:31:HIS:ND1	1:C:67:GLN:NE2	2.42	0.67
1:B:196:ASN:HD21	1:B:198:LYS:HE3	1.59	0.67
1:A:108:LEU:HD13	1:C:113:ILE:CD1	2.26	0.66
1:B:111:GLN:HG2	1:B:112:GLN:HG2	1.78	0.66
1:A:124:GLN:O	1:A:128:GLU:HG3	1.96	0.65
1:C:13:LEU:HD13	1:C:153:GLY:N	2.12	0.65
1:C:173:ALA:HB3	1:C:181:ALA:HB2	1.79	0.65
1:C:26:PRO:HG2	1:C:27:PHE:CD1	2.32	0.64
1:A:113:ILE:HG12	1:C:82:MSE:HE3	1.80	0.64
1:A:108:LEU:HA	1:C:111:GLN:NE2	2.13	0.64
1:C:198:LYS:HD2	1:C:198:LYS:H	1.64	0.63
1:A:43:GLU:CD	1:A:227:LYS:HD2	2.19	0.63
1:C:121:LYS:O	1:C:124:GLN:HB3	1.98	0.63
1:B:188:THR:OG1	1:B:191:GLU:HG3	1.99	0.63
1:C:68:ALA:HA	1:C:71:MSE:CE	2.28	0.63
1:C:96:GLU:HG3	4:C:280:HOH:O	1.98	0.63
1:B:196:ASN:ND2	1:B:198:LYS:HE3	2.13	0.63
1:A:161:MSE:HE2	1:A:163:LYS:HB2	1.81	0.62
1:C:113:ILE:N	1:C:113:ILE:HD12	2.15	0.61
1:B:123:ILE:O	1:B:127:GLU:HG3	1.99	0.61
1:A:111:GLN:HE21	1:C:109:SER:H	1.49	0.61
1:B:119:ARG:HB3	1:B:120:PRO:HD3	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:68:ALA:CA	1:C:71:MSE:HE2	2.29	0.61
1:B:13:LEU:HD13	1:B:153:GLY:N	2.17	0.60
1:A:108:LEU:CD1	1:C:113:ILE:HD11	2.32	0.59
1:A:189:PHE:CD2	1:A:241:CYS:HB3	2.37	0.59
1:C:227:LYS:O	1:C:231:THR:HG23	2.02	0.59
1:A:86:MSE:HB3	1:C:79:MSE:HE3	1.83	0.59
1:A:87:ASN:O	1:A:91:LEU:HB2	2.03	0.58
1:C:101:LYS:HD3	1:C:130:LYS:NZ	2.19	0.58
1:C:82:MSE:HE1	1:C:110:MSE:HE1	1.86	0.58
1:C:187:ILE:HG22	1:C:191:GLU:HB2	1.86	0.57
1:A:200:MSE:HE3	1:A:215:ILE:HD13	1.86	0.57
1:C:230:ILE:C	1:C:230:ILE:HD12	2.24	0.57
1:C:64:VAL:HG21	1:C:79:MSE:HE2	1.86	0.56
1:C:210:GLU:OE2	1:C:211:ARG:NH1	2.37	0.56
1:A:7:ILE:CD1	1:A:47:MSE:HE2	2.33	0.56
1:A:179:PRO:CG	1:B:187:ILE:HG22	2.36	0.56
1:A:64:VAL:HG13	1:C:29:ILE:HD12	1.87	0.55
1:A:200:MSE:CE	1:A:215:ILE:HD13	2.36	0.55
1:C:54:VAL:HG21	1:C:149:ALA:HA	1.89	0.55
1:A:63:GLY:O	1:C:28:GLY:HA3	2.07	0.54
1:B:200:MSE:CE	1:B:215:ILE:HD13	2.37	0.54
1:A:13:LEU:HD13	1:A:153:GLY:N	2.22	0.54
1:B:7:ILE:HD12	1:B:7:ILE:N	2.24	0.53
1:B:167:VAL:HG12	1:B:169:GLY:H	1.73	0.53
1:B:215:ILE:HB	1:B:239:VAL:HB	1.89	0.53
1:C:187:ILE:HG22	1:C:188:THR:H	1.73	0.53
1:A:92:LYS:HB3	1:A:92:LYS:NZ	2.23	0.53
1:B:200:MSE:HE3	1:B:215:ILE:HD13	1.91	0.53
1:C:167:VAL:HG12	1:C:169:GLY:H	1.73	0.53
1:C:13:LEU:HD13	1:C:153:GLY:CA	2.39	0.52
1:C:64:VAL:HG11	1:C:79:MSE:HE1	1.90	0.52
1:A:151:LEU:O	1:A:155:GLU:HG3	2.10	0.52
1:B:209:ARG:HH11	1:B:209:ARG:HG2	1.74	0.52
1:C:188:THR:OG1	1:C:191:GLU:HG3	2.09	0.52
1:A:225:SER:O	1:A:229:VAL:HG23	2.09	0.52
1:B:82:MSE:HE2	1:B:82:MSE:HA	1.92	0.51
1:A:73:ARG:HD3	1:A:73:ARG:O	2.10	0.51
1:A:113:ILE:H	1:A:113:ILE:HD12	1.75	0.51
1:B:31:HIS:O	1:B:35:VAL:HG23	2.11	0.51
1:B:87:ASN:O	1:B:91:LEU:HB2	2.11	0.50
1:C:38:VAL:HG21	1:C:94:ALA:CB	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:183:ARG:NH2	1:C:234:ASP:OD1	2.44	0.50
1:A:58:GLY:O	1:A:62:ARG:HG3	2.11	0.50
1:A:170:VAL:HG23	1:A:237:THR:HB	1.92	0.50
1:A:108:LEU:HB3	1:C:111:GLN:HB3	1.94	0.50
1:A:89:LEU:HD11	1:A:113:ILE:CG2	2.42	0.50
1:B:227:LYS:O	1:B:231:THR:HG23	2.11	0.50
1:B:228:ARG:HB3	1:B:233:GLU:HB2	1.92	0.50
1:C:7:ILE:CD1	1:C:47:MSE:HE2	2.41	0.50
1:A:171:TYR:CZ	1:A:183:ARG:HB2	2.47	0.49
1:A:179:PRO:HB2	1:B:191:GLU:HB3	1.93	0.49
1:B:6:GLN:C	1:B:7:ILE:HD12	2.32	0.49
1:A:38:VAL:HG12	1:A:98:LEU:HD12	1.94	0.49
1:A:31:HIS:O	1:A:35:VAL:HG23	2.13	0.49
1:B:54:VAL:HG21	1:B:149:ALA:HA	1.94	0.49
1:B:15:LYS:O	1:B:15:LYS:HD3	2.12	0.49
1:B:207:LEU:HD23	1:B:211:ARG:HG2	1.94	0.49
1:A:227:LYS:O	1:A:231:THR:HG23	2.13	0.49
1:B:224:GLY:O	1:B:228:ARG:HG3	2.13	0.49
1:B:41:ILE:O	1:B:44:VAL:HG22	2.13	0.49
1:C:172:THR:HG22	1:C:184:TYR:CZ	2.47	0.48
1:B:168:ASP:O	1:B:236:GLY:HA2	2.14	0.48
1:C:200:MSE:CE	1:C:215:ILE:HD13	2.39	0.48
1:C:7:ILE:HD11	1:C:47:MSE:HE2	1.95	0.48
1:A:85:VAL:CG2	1:A:110:MSE:HE1	2.44	0.47
1:B:216:VAL:HG11	1:B:229:VAL:HG11	1.97	0.47
1:C:44:VAL:CG1	1:C:230:ILE:HG12	2.44	0.47
1:A:64:VAL:HG12	1:A:79:MSE:HE1	1.97	0.47
1:B:13:LEU:HD13	1:B:153:GLY:CA	2.45	0.47
1:C:165:THR:O	1:C:219:GLY:HA2	2.15	0.47
1:C:113:ILE:H	1:C:113:ILE:HD12	1.79	0.47
1:B:207:LEU:CD2	1:B:211:ARG:HG2	2.44	0.47
1:B:34:ILE:O	1:B:38:VAL:HG23	2.15	0.47
1:C:68:ALA:HA	1:C:71:MSE:HG3	1.98	0.46
1:C:135:ALA:O	1:C:136:ALA:HB3	2.15	0.46
1:B:210:GLU:OE1	1:B:211:ARG:HD2	2.16	0.46
1:C:101:LYS:HD3	1:C:130:LYS:HZ2	1.81	0.46
1:A:192:ALA:HA	1:A:197:LEU:HD12	1.97	0.46
1:C:177:LYS:O	1:C:178:ASP:HB2	2.15	0.46
1:A:119:ARG:HB3	1:A:120:PRO:HD3	1.96	0.46
1:A:165:THR:HG23	1:A:167:VAL:H	1.81	0.46
1:B:216:VAL:HG11	1:B:229:VAL:CG1	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:VAL:HG12	1:A:64:VAL:O	2.15	0.46
1:C:127:GLU:C	1:C:129:GLY:H	2.20	0.45
1:C:78:TYR:HA	1:C:81:MSE:HE3	1.98	0.45
1:A:111:GLN:NE2	1:C:108:LEU:HA	2.31	0.45
1:B:167:VAL:HG12	1:B:168:ASP:N	2.30	0.45
1:B:172:THR:HA	1:B:184:TYR:CE2	2.51	0.45
1:C:224:GLY:O	1:C:228:ARG:HG3	2.16	0.45
1:A:85:VAL:HG21	1:A:110:MSE:HE1	1.99	0.45
1:B:13:LEU:HD21	1:B:54:VAL:HG23	1.97	0.45
1:A:38:VAL:HG21	1:A:94:ALA:HB3	1.99	0.45
1:C:216:VAL:HG11	1:C:229:VAL:CG1	2.47	0.45
1:A:112:GLN:HG3	1:C:108:LEU:HD22	1.98	0.45
1:A:19:GLU:OE1	1:A:19:GLU:N	2.44	0.45
1:B:36:GLN:NE2	1:B:223:GLU:HG3	2.32	0.45
1:A:38:VAL:HG21	1:A:94:ALA:CB	2.47	0.45
1:A:96:GLU:OE1	1:A:101:LYS:HA	2.17	0.44
1:A:165:THR:O	1:A:219:GLY:HA2	2.16	0.44
1:A:17:SER:HB2	3:A:250:GOL:H12	1.98	0.44
1:B:182:THR:CG2	1:B:183:ARG:N	2.80	0.44
1:B:73:ARG:NH1	1:B:77:ASP:OD2	2.51	0.44
1:B:194:LEU:C	1:B:196:ASN:H	2.21	0.44
1:C:31:HIS:O	1:C:35:VAL:HG23	2.18	0.44
1:B:92:LYS:O	1:B:96:GLU:HG2	2.18	0.44
1:B:228:ARG:HH21	1:B:233:GLU:HB3	1.83	0.43
1:A:78:TYR:HA	1:A:81:MSE:HE3	2.00	0.43
1:A:50:GLN:NE2	1:A:127:GLU:HA	2.33	0.43
1:B:15:LYS:NZ	1:B:146:ASP:OD1	2.41	0.43
1:C:8:LYS:HB3	1:C:230:ILE:O	2.18	0.43
1:A:162:LEU:HD23	1:A:216:VAL:HB	2.00	0.43
1:A:28:GLY:HA3	1:C:63:GLY:O	2.19	0.43
1:A:113:ILE:HD11	1:C:108:LEU:CD1	2.48	0.43
1:A:41:ILE:O	1:A:44:VAL:HG22	2.19	0.43
1:B:49:VAL:HG21	1:B:230:ILE:HG21	1.99	0.43
1:B:89:LEU:HD11	1:B:113:ILE:HG23	2.00	0.43
1:C:6:GLN:NE2	4:C:272:HOH:O	2.52	0.43
1:B:119:ARG:HA	1:B:156:MSE:HE1	2.01	0.43
1:A:92:LYS:O	1:A:96:GLU:HG2	2.19	0.43
1:A:162:LEU:HB3	1:A:218:PHE:HE2	1.85	0.42
1:C:106:SER:OG	1:C:110:MSE:HE3	2.20	0.42
1:C:72:ASP:OD1	1:C:74:ALA:HB3	2.18	0.42
1:A:43:GLU:OE1	1:A:227:LYS:HD2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:MSE:HE3	1:C:90:ALA:HA	2.01	0.42
1:C:175:PRO:HA	1:C:178:ASP:O	2.20	0.42
1:C:47:MSE:SE	1:C:227:LYS:HE3	2.70	0.42
1:C:160:VAL:HG22	1:C:161:MSE:N	2.34	0.42
1:A:92:LYS:HG3	1:A:102:ALA:HB3	2.01	0.42
1:B:12:VAL:CG2	1:B:51:VAL:HG22	2.49	0.42
1:B:160:VAL:HG23	1:B:214:ASN:O	2.20	0.41
1:C:44:VAL:HG12	1:C:230:ILE:HG12	2.01	0.41
1:C:188:THR:HA	1:C:240:HIS:O	2.21	0.41
1:C:67:GLN:O	1:C:71:MSE:HG3	2.20	0.41
1:A:189:PHE:CE2	1:A:241:CYS:HB3	2.56	0.41
1:A:65:SER:O	1:A:68:ALA:HB3	2.19	0.41
1:B:16:LEU:O	1:B:55:VAL:HA	2.19	0.41
1:C:187:ILE:HD13	1:C:197:LEU:CD1	2.50	0.41
1:C:215:ILE:O	1:C:238:LEU:HD12	2.21	0.41
1:A:167:VAL:HG12	1:A:169:GLY:H	1.86	0.41
1:B:58:GLY:O	1:B:62:ARG:HG3	2.20	0.41
1:C:216:VAL:HG11	1:C:229:VAL:HG11	2.03	0.41
1:B:163:LYS:HD2	1:B:200:MSE:HE2	2.01	0.41
1:A:41:ILE:O	1:A:45:VAL:HG23	2.21	0.41
1:C:160:VAL:HG23	1:C:214:ASN:O	2.21	0.41
1:A:165:THR:HG23	1:A:167:VAL:O	2.21	0.40
1:C:215:ILE:HG22	1:C:216:VAL:N	2.35	0.40
1:C:64:VAL:HG11	1:C:79:MSE:CE	2.51	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	234/239 (98%)	225 (96%)	8 (3%)	1 (0%)	34 57

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	234/239 (98%)	220 (94%)	14 (6%)	0	100	100
1	C	234/239 (98%)	224 (96%)	8 (3%)	2 (1%)	17	35
All	All	702/717 (98%)	669 (95%)	30 (4%)	3 (0%)	34	57

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	178	ASP
1	C	236	GLY
1	A	111	GLN

5.3.2 Protein sidechains [i](#)

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	186/177 (105%)	181 (97%)	5 (3%)	44	71
1	B	186/177 (105%)	177 (95%)	9 (5%)	25	49
1	C	186/177 (105%)	178 (96%)	8 (4%)	29	54
All	All	558/531 (105%)	536 (96%)	22 (4%)	32	58

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	SER
1	A	67	GLN
1	A	73	ARG
1	A	207	LEU
1	A	240	HIS
1	B	15	LYS
1	B	67	GLN
1	B	110	MSE
1	B	112	GLN
1	B	119	ARG
1	B	166	ASN

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Mol	Chain	Res	Type
1	B	198	LYS
1	B	201	ASP
1	B	207	LEU
1	C	40	GLU
1	C	67	GLN
1	C	72	ASP
1	C	73	ARG
1	C	186	THR
1	C	198	LYS
1	C	207	LEU
1	C	234	ASP

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

Mol	Chain	Res	Type
1	A	67	GLN
1	A	111	GLN
1	A	240	HIS
1	B	6	GLN
1	B	111	GLN
1	B	166	ASN
1	B	196	ASN
1	C	36	GLN
1	C	67	GLN
1	C	111	GLN
1	C	157	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

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	GOL	B	251	-	5,5,5	4.46	5 (100%)	5,5,5	5.77	3 (60%)
3	GOL	C	252	-	5,5,5	4.41	5 (100%)	5,5,5	5.80	3 (60%)
2	FMT	A	253	-	0,2,2	0.00	-	0,1,1	0.00	-
3	GOL	A	250	-	5,5,5	4.42	5 (100%)	5,5,5	5.79	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
3	GOL	C	252	-	-	3/4/4/4	-
3	GOL	B	251	-	-	2/4/4/4	-
3	GOL	A	250	-	-	2/4/4/4	-

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	251	GOL	C3-C2	-7.29	1.21	1.51
3	A	250	GOL	C3-C2	-7.19	1.22	1.51
3	C	252	GOL	C3-C2	-6.92	1.23	1.51
3	C	252	GOL	O1-C1	4.70	1.62	1.42
3	B	251	GOL	O1-C1	4.60	1.61	1.42
3	A	250	GOL	O1-C1	4.57	1.61	1.42
3	C	252	GOL	O3-C3	3.74	1.58	1.42
3	A	250	GOL	O3-C3	3.49	1.57	1.42
3	B	251	GOL	O3-C3	3.48	1.57	1.42
3	A	250	GOL	C1-C2	-2.70	1.40	1.51
3	C	252	GOL	O2-C2	-2.61	1.35	1.43
3	B	251	GOL	C1-C2	-2.55	1.41	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	251	GOL	O2-C2	-2.55	1.35	1.43
3	C	252	GOL	C1-C2	-2.50	1.41	1.51
3	A	250	GOL	O2-C2	-2.36	1.36	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	252	GOL	O3-C3-C2	10.55	160.78	110.20
3	A	250	GOL	O3-C3-C2	10.50	160.56	110.20
3	B	251	GOL	O3-C3-C2	10.50	160.55	110.20
3	A	250	GOL	O2-C2-C3	6.75	138.85	109.12
3	C	252	GOL	O2-C2-C3	6.72	138.73	109.12
3	B	251	GOL	O2-C2-C3	6.68	138.53	109.12
3	C	252	GOL	O1-C1-C2	3.39	126.45	110.20
3	A	250	GOL	O1-C1-C2	3.31	126.08	110.20
3	B	251	GOL	O1-C1-C2	3.29	125.96	110.20

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	251	GOL	O1-C1-C2-C3
3	B	251	GOL	C1-C2-C3-O3
3	C	252	GOL	O1-C1-C2-C3
3	C	252	GOL	C1-C2-C3-O3
3	A	250	GOL	C1-C2-C3-O3
3	A	250	GOL	O1-C1-C2-C3
3	C	252	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	250	GOL	1	0

5.7 Other polymers [i](#)

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	225/239 (94%)	-0.23	2 (0%) 84 82	12, 37, 58, 68	0
1	B	225/239 (94%)	-0.25	1 (0%) 92 91	14, 38, 57, 64	0
1	C	225/239 (94%)	-0.08	9 (4%) 38 31	18, 44, 66, 79	0
All	All	675/717 (94%)	-0.19	12 (1%) 68 64	12, 39, 60, 79	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	175	PRO	5.1
1	C	176	LYS	4.5
1	C	179	PRO	3.7
1	C	177	LYS	3.2
1	A	241	CYS	2.6
1	C	197	LEU	2.4
1	A	179	PRO	2.3
1	C	7	ILE	2.2
1	C	187	ILE	2.2
1	C	109	SER	2.1
1	C	241	CYS	2.0
1	B	27	PHE	2.0

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, 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(Å ²)	Q<0.9
3	GOL	C	252	6/6	0.56	0.27	52,54,55,56	0
3	GOL	A	250	6/6	0.77	0.29	42,47,48,50	0
3	GOL	B	251	6/6	0.82	0.17	41,42,44,45	0
2	FMT	A	253	3/3	0.95	0.39	31,31,33,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.