



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 12:43 am BST

PDB ID : 1B7B
Title : Carbamate kinase from *Enterococcus faecalis*
Authors : Marina, A.; Alzari, P.M.; Bravo, J.; Uriarte, M.; Barcelona, B.; Fita, I.; Rubio, V.
Deposited on : 1999-01-20
Resolution : 2.80 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 : 2.11
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) : 2.11

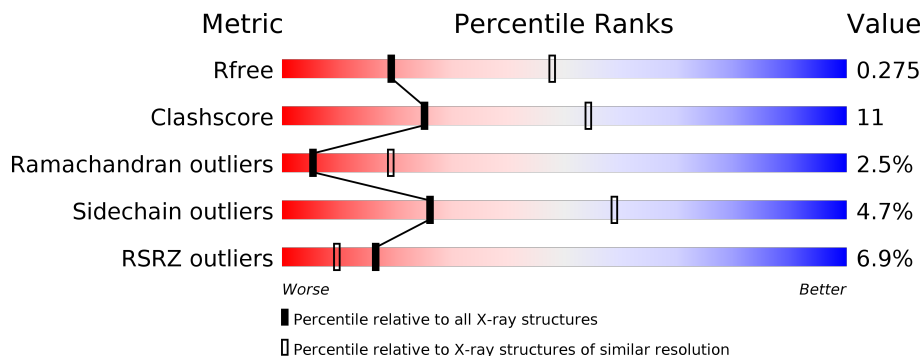
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 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	310	 4% 75% 21% ..
1	B	310	 11% 74% 22% ..
1	C	310	 3% 76% 19% ..
1	D	310	 10% 73% 23% ..

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 9512 atoms, of which 280 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 CARBAMATE KINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
1	A	307	Total 2356	C 1441	H 70	N 387	O 449	S 9	0	0	0
1	B	307	Total 2356	C 1441	H 70	N 387	O 449	S 9	0	0	0
1	C	307	Total 2356	C 1441	H 70	N 387	O 449	S 9	0	0	0
1	D	307	Total 2356	C 1441	H 70	N 387	O 449	S 9	0	0	0

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	Total 5	O 4	S 1	0	0
2	B	1	Total 5	O 4	S 1	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

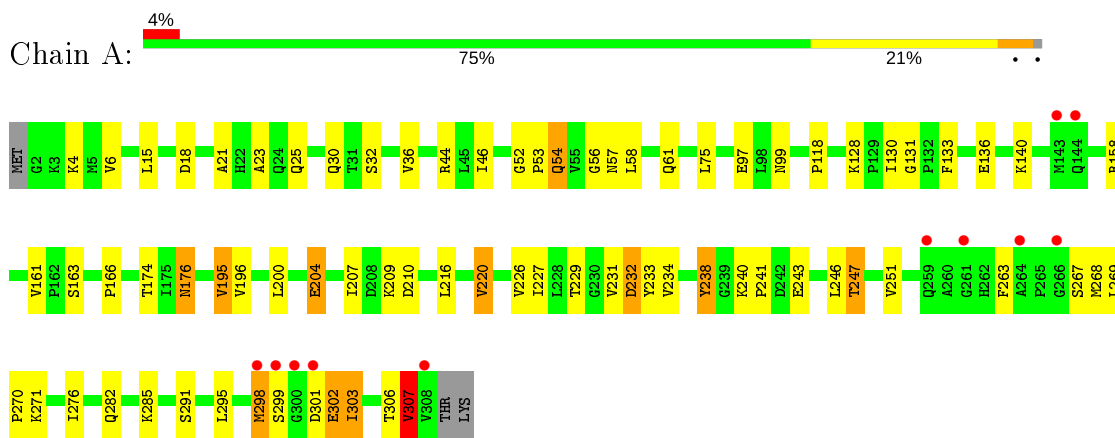
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	16	Total	O	0	0
			16	16		
3	B	17	Total	O	0	0
			17	17		
3	C	18	Total	O	0	0
			18	18		
3	D	17	Total	O	0	0
			17	17		

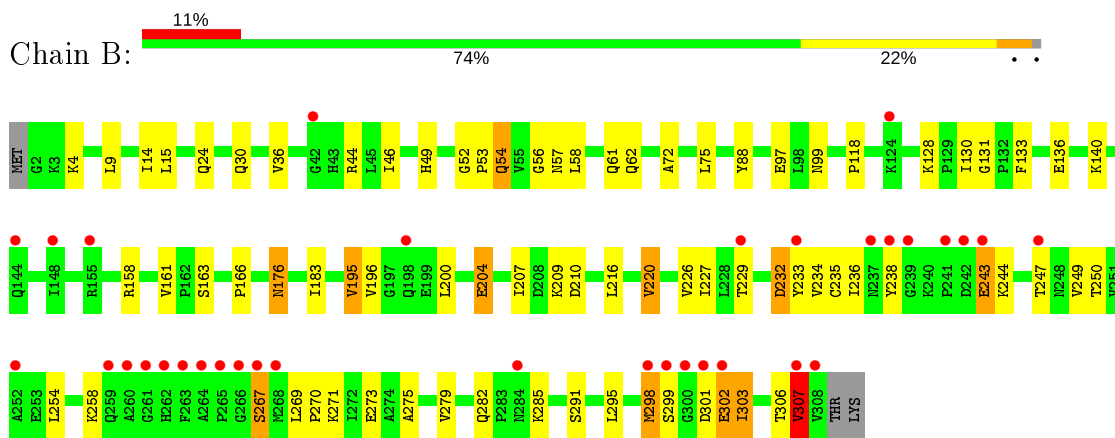
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

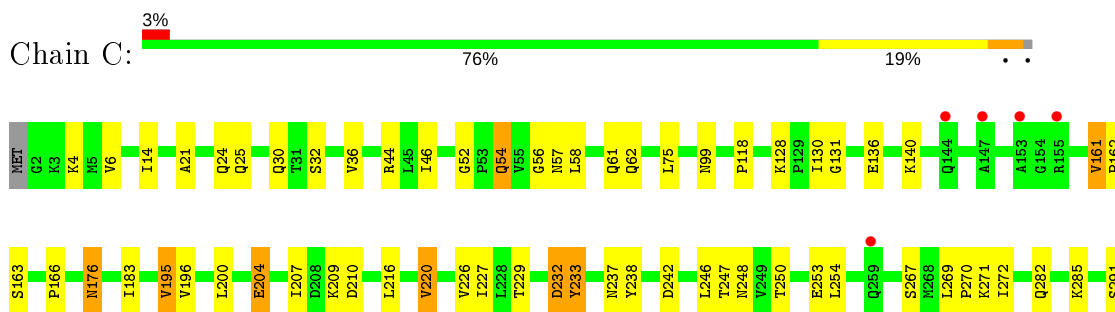
- Molecule 1: CARBAMATE KINASE

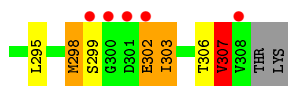


- Molecule 1: CARBAMATE KINASE

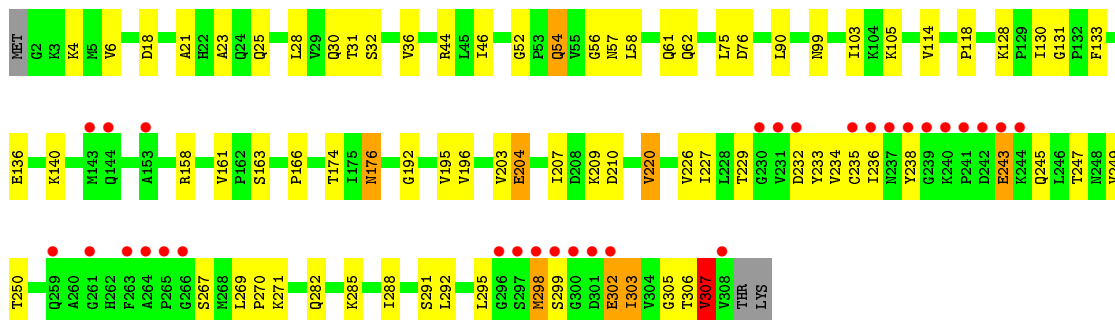
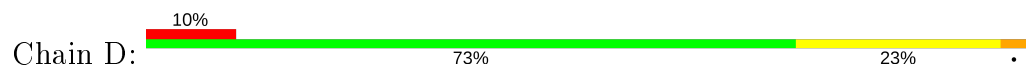


- Molecule 1: CARBAMATE KINASE





• Molecule 1: CARBAMATE KINASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.86Å 172.92Å 98.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.80 14.91 – 2.80	Depositor EDS
% Data completeness (in resolution range)	89.5 (15.00-2.80) 82.0 (14.91-2.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtrriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.224 , 0.283 0.222 , 0.275	Depositor DCC
R_{free} test set	2893 reflections (10.20%)	wwPDB-VP
Wilson B-factor (Å ²)	37.6	Xtrriage
Anisotropy	0.132	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 50.5	EDS
L-test for twinning ¹	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	9512	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 28.17 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.9298e-03.*

¹Theoretical 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: SO4

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	0.30	0/2321	0.57	0/3151
1	B	0.30	0/2321	0.57	0/3151
1	C	0.30	0/2321	0.58	0/3151
1	D	0.30	0/2321	0.57	0/3151
All	All	0.30	0/9284	0.57	0/12604

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	2286	70	2317	51	0
1	B	2286	70	2317	52	0
1	C	2286	70	2317	43	0
1	D	2286	70	2317	54	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
3	A	16	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	17	0	0	0	0
3	C	18	0	0	0	0
3	D	17	0	0	0	0
All	All	9232	280	9268	198	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 11.

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:B:232:ASP:HB2	1:B:302:GLU:HG3	1.50	0.91
1:C:232:ASP:HB2	1:C:302:GLU:HG3	1.69	0.74
1:C:75:LEU:HB3	1:C:204:GLU:HG3	1.70	0.74
1:A:75:LEU:HB3	1:A:204:GLU:HG3	1.70	0.73
1:B:235:CYS:SG	1:B:243:GLU:HG3	2.29	0.73
1:D:75:LEU:HB3	1:D:204:GLU:HG3	1.73	0.70
1:A:234:VAL:HG23	1:A:303:ILE:HG13	1.74	0.70
1:B:234:VAL:HG23	1:B:303:ILE:HG13	1.73	0.70
1:A:246:LEU:HB2	1:A:303:ILE:HD12	1.74	0.69
1:B:75:LEU:HB3	1:B:204:GLU:HG3	1.75	0.68
1:D:269:LEU:HB3	1:D:270:PRO:HD3	1.77	0.67
1:B:234:VAL:H	1:B:303:ILE:HD11	1.61	0.64
1:A:128:LYS:HB2	1:A:163:SER:OG	1.97	0.64
1:C:269:LEU:HB3	1:C:270:PRO:HD3	1.79	0.64
1:B:269:LEU:HB3	1:B:270:PRO:HD3	1.80	0.64
1:D:128:LYS:HB2	1:D:163:SER:OG	1.98	0.64
1:A:75:LEU:CB	1:A:204:GLU:HG3	2.28	0.63
1:C:128:LYS:HB2	1:C:163:SER:OG	1.99	0.63
1:A:233:TYR:CD1	1:A:302:GLU:HB3	2.34	0.62
1:A:269:LEU:HB3	1:A:270:PRO:HD3	1.79	0.62
1:A:267:SER:O	1:A:270:PRO:HD2	1.99	0.61
1:B:249:VAL:HG21	1:B:254:LEU:HD21	1.82	0.61
1:B:128:LYS:HB2	1:B:163:SER:OG	2.00	0.61
1:B:4:LYS:HA	1:B:44:ARG:HB2	1.83	0.60
1:A:232:ASP:HB2	1:A:302:GLU:HG3	1.83	0.60
1:A:233:TYR:HA	1:A:303:ILE:HG12	1.83	0.59
1:B:136:GLU:HG2	1:B:140:LYS:HE3	1.83	0.59
1:B:46:ILE:HD13	1:B:220:VAL:HG11	1.84	0.59
1:B:58:LEU:HA	1:B:61:GLN:HE21	1.66	0.59
1:D:267:SER:O	1:D:271:LYS:HG2	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:227:ILE:CG2	1:C:271:LYS:HD2	2.32	0.59
1:C:250:THR:HB	1:C:253:GLU:HG3	1.84	0.59
1:D:176:ASN:HD21	1:D:220:VAL:HA	1.65	0.59
1:D:54:GLN:H	1:D:54:GLN:NE2	2.01	0.59
1:C:246:LEU:HB2	1:C:303:ILE:HD12	1.85	0.58
1:A:136:GLU:HG2	1:A:140:LYS:HE3	1.85	0.58
1:C:233:TYR:CD1	1:C:302:GLU:HB3	2.39	0.58
1:A:226:VAL:HG11	1:A:295:LEU:HD11	1.85	0.58
1:C:247:THR:O	1:C:303:ILE:HA	2.03	0.57
1:D:4:LYS:HA	1:D:44:ARG:HB2	1.85	0.57
1:D:136:GLU:HG2	1:D:140:LYS:HE3	1.88	0.56
1:D:250:THR:HA	1:D:306:THR:O	2.05	0.56
1:C:226:VAL:HG11	1:C:295:LEU:HD11	1.85	0.56
1:A:130:ILE:HD12	1:A:161:VAL:HB	1.87	0.56
1:D:176:ASN:ND2	1:D:220:VAL:HA	2.21	0.56
1:D:227:ILE:CG2	1:D:271:LYS:HD2	2.36	0.55
1:B:233:TYR:CD1	1:B:302:GLU:HB3	2.41	0.55
1:B:226:VAL:HG11	1:B:295:LEU:HD11	1.88	0.55
1:C:267:SER:O	1:C:270:PRO:HD2	2.06	0.55
1:C:118:PRO:HD3	1:C:196:VAL:HG13	1.89	0.54
1:C:306:THR:O	1:C:307:VAL:HG13	2.07	0.54
1:A:18:ASP:HB3	1:A:23:ALA:CB	2.38	0.54
1:C:46:ILE:HD13	1:C:220:VAL:HG11	1.89	0.54
1:A:227:ILE:CG2	1:A:271:LYS:HD2	2.38	0.53
1:D:46:ILE:HD13	1:D:220:VAL:HG11	1.90	0.53
1:C:58:LEU:HA	1:C:61:GLN:HE21	1.74	0.53
1:D:235:CYS:SG	1:D:245:GLN:HG2	2.48	0.53
1:D:232:ASP:O	1:D:233:TYR:HB2	2.08	0.53
1:B:209:LYS:HG3	1:B:210:ASP:N	2.24	0.53
1:B:227:ILE:CG2	1:B:271:LYS:HD2	2.40	0.52
1:D:118:PRO:HD3	1:D:196:VAL:HG13	1.92	0.52
1:D:58:LEU:HA	1:D:61:GLN:HE21	1.74	0.52
1:B:54:GLN:H	1:B:54:GLN:NE2	2.07	0.52
1:C:75:LEU:CB	1:C:204:GLU:HG3	2.37	0.52
1:A:58:LEU:HA	1:A:61:GLN:HE21	1.75	0.52
1:A:54:GLN:H	1:A:54:GLN:NE2	2.07	0.52
1:D:226:VAL:HG11	1:D:295:LEU:HD11	1.91	0.52
1:C:232:ASP:O	1:C:233:TYR:HB2	2.10	0.52
1:A:166:PRO:HG3	1:A:207:ILE:HG21	1.91	0.52
1:B:232:ASP:O	1:B:233:TYR:HB2	2.09	0.51
1:B:166:PRO:HG3	1:B:207:ILE:HG21	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:ASN:HD21	1:A:220:VAL:HA	1.75	0.51
1:A:4:LYS:HA	1:A:44:ARG:HB2	1.91	0.51
1:C:229:THR:O	1:C:291:SER:HA	2.10	0.51
1:A:32:SER:O	1:A:36:VAL:HG23	2.11	0.51
1:C:237:ASN:CB	1:C:242:ASP:HB2	2.41	0.51
1:D:76:ASP:HB3	1:D:203:VAL:O	2.11	0.51
1:D:75:LEU:CB	1:D:204:GLU:HG3	2.38	0.50
1:A:306:THR:O	1:A:307:VAL:HG13	2.12	0.50
1:D:267:SER:O	1:D:270:PRO:HD2	2.11	0.50
1:D:234:VAL:HG23	1:D:303:ILE:HG13	1.92	0.50
1:B:56:GLY:HA3	1:B:131:GLY:CA	2.42	0.50
1:A:46:ILE:HD13	1:A:220:VAL:HG11	1.94	0.50
1:B:75:LEU:CB	1:B:204:GLU:HG3	2.41	0.50
1:D:166:PRO:HG3	1:D:207:ILE:HG21	1.94	0.50
1:A:216:LEU:O	1:A:220:VAL:HB	2.12	0.49
1:B:195:VAL:HG11	1:B:200:LEU:HD22	1.94	0.49
1:B:229:THR:O	1:B:291:SER:HA	2.12	0.49
1:B:130:ILE:HD12	1:B:161:VAL:HB	1.94	0.49
1:A:267:SER:O	1:A:271:LYS:HG2	2.12	0.49
1:B:306:THR:O	1:B:307:VAL:HG13	2.12	0.49
1:D:114:VAL:HG23	1:D:192:GLY:HA3	1.95	0.49
1:A:251:VAL:HG13	1:A:276:ILE:HG23	1.94	0.49
1:C:166:PRO:HG3	1:C:207:ILE:HG21	1.94	0.49
1:B:118:PRO:HD3	1:B:196:VAL:HG13	1.95	0.49
1:C:136:GLU:HG2	1:C:140:LYS:HE3	1.94	0.49
1:D:247:THR:O	1:D:303:ILE:HA	2.13	0.49
1:B:258:LYS:NZ	1:B:273:GLU:HG2	2.27	0.49
1:C:14:ILE:HG22	1:C:24:GLN:HE21	1.77	0.49
1:B:250:THR:HA	1:B:306:THR:O	2.12	0.49
1:D:18:ASP:HB3	1:D:23:ALA:CB	2.43	0.49
1:B:233:TYR:HA	1:B:303:ILE:HG12	1.94	0.48
1:A:234:VAL:H	1:A:303:ILE:HD11	1.78	0.48
1:A:246:LEU:HB2	1:A:303:ILE:CD1	2.44	0.48
1:A:118:PRO:HD3	1:A:196:VAL:HG13	1.96	0.48
1:A:133:PHE:CE2	1:A:158:ARG:HD2	2.49	0.48
1:A:232:ASP:O	1:A:233:TYR:HB2	2.14	0.48
1:C:130:ILE:HD12	1:C:161:VAL:HB	1.96	0.47
1:D:130:ILE:HD12	1:D:161:VAL:HB	1.95	0.47
1:C:282:GLN:HB2	1:C:285:LYS:HG3	1.95	0.47
1:C:54:GLN:H	1:C:54:GLN:NE2	2.13	0.47
1:B:176:ASN:HD21	1:B:220:VAL:HA	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:62:GLN:HB3	1:B:72:ALA:HB1	1.96	0.47
1:D:233:TYR:CD1	1:D:302:GLU:HB3	2.50	0.47
1:D:236:ILE:O	1:D:243:GLU:HA	2.13	0.47
1:D:57:ASN:O	1:D:61:GLN:HG3	2.14	0.47
1:A:57:ASN:O	1:A:61:GLN:HG3	2.14	0.47
1:D:58:LEU:O	1:D:62:GLN:HG2	2.15	0.47
1:A:176:ASN:ND2	1:A:220:VAL:HA	2.30	0.46
1:A:15:LEU:HD13	1:A:53:PRO:HB2	1.98	0.46
1:C:21:ALA:O	1:C:25:GLN:HG3	2.15	0.46
1:D:232:ASP:HB2	1:D:302:GLU:HG3	1.97	0.46
1:D:229:THR:O	1:D:291:SER:HA	2.15	0.46
1:B:58:LEU:O	1:B:62:GLN:HG2	2.15	0.46
1:A:195:VAL:HG11	1:A:200:LEU:HD22	1.98	0.46
1:C:195:VAL:HG11	1:C:200:LEU:HD22	1.98	0.46
1:B:57:ASN:O	1:B:61:GLN:HG3	2.15	0.46
1:D:56:GLY:HA3	1:D:131:GLY:CA	2.46	0.45
1:B:195:VAL:HG22	1:B:200:LEU:O	2.16	0.45
1:D:118:PRO:HB3	1:D:196:VAL:HG11	1.98	0.45
1:D:282:GLN:HB2	1:D:285:LYS:HG3	1.97	0.45
1:B:216:LEU:O	1:B:220:VAL:HB	2.16	0.45
1:B:282:GLN:HB2	1:B:285:LYS:HG3	1.99	0.45
1:B:236:ILE:HD11	1:B:244:LYS:HD3	1.98	0.45
1:C:209:LYS:HG3	1:C:210:ASP:N	2.30	0.45
1:B:14:ILE:HG22	1:B:24:GLN:HE21	1.82	0.45
1:C:216:LEU:O	1:C:220:VAL:HB	2.16	0.45
1:A:229:THR:O	1:A:291:SER:HA	2.17	0.45
1:A:282:GLN:HB2	1:A:285:LYS:CG	2.47	0.45
1:C:233:TYR:HA	1:C:303:ILE:HG12	1.98	0.45
1:C:57:ASN:O	1:C:61:GLN:HG3	2.17	0.45
1:D:28:LEU:HD22	1:D:90:LEU:HD13	1.98	0.45
1:B:267:SER:O	1:B:271:LYS:HG2	2.16	0.45
1:D:306:THR:O	1:D:307:VAL:HG13	2.17	0.45
1:B:133:PHE:CE2	1:B:158:ARG:HD2	2.52	0.44
1:B:267:SER:O	1:B:270:PRO:HD2	2.17	0.44
1:B:176:ASN:ND2	1:B:220:VAL:HA	2.31	0.44
1:B:295:LEU:HD23	1:B:301:ASP:OD2	2.18	0.44
1:C:176:ASN:HD21	1:C:220:VAL:HA	1.82	0.44
1:A:282:GLN:HB2	1:A:285:LYS:HG3	1.99	0.44
1:C:4:LYS:HA	1:C:44:ARG:HB2	2.00	0.44
1:C:254:LEU:HD13	1:C:272:ILE:HG23	1.99	0.44
1:C:58:LEU:O	1:C:62:GLN:HG2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:56:GLY:HA3	1:A:131:GLY:CA	2.47	0.44
1:C:118:PRO:HB3	1:C:196:VAL:HG11	2.00	0.44
1:A:295:LEU:HD23	1:A:301:ASP:OD2	2.18	0.43
1:D:31:THR:HG22	1:D:292:LEU:HD23	2.00	0.43
1:A:247:THR:O	1:A:303:ILE:HA	2.18	0.43
1:C:56:GLY:HA3	1:C:131:GLY:CA	2.48	0.43
1:A:238:TYR:O	1:A:243:GLU:HB2	2.19	0.43
1:C:6:VAL:HA	1:C:46:ILE:O	2.17	0.43
1:D:235:CYS:HB3	1:D:243:GLU:HG3	2.01	0.43
1:D:32:SER:O	1:D:36:VAL:HG23	2.18	0.43
1:A:240:LYS:HB3	1:A:241:PRO:HD2	1.99	0.43
1:D:209:LYS:HG3	1:D:210:ASP:N	2.34	0.43
1:B:118:PRO:HB3	1:B:196:VAL:HG11	2.00	0.43
1:B:279:VAL:HG11	1:B:306:THR:HA	2.01	0.43
1:D:249:VAL:HG23	1:D:305:GLY:HA2	2.01	0.43
1:A:209:LYS:HG3	1:A:210:ASP:N	2.33	0.42
1:A:263:PHE:HB3	1:A:268:MET:HB3	2.01	0.42
1:B:282:GLN:HB2	1:B:285:LYS:CG	2.49	0.42
1:B:15:LEU:HD13	1:B:53:PRO:HB2	2.01	0.42
1:A:174:THR:HG23	1:C:183:ILE:HD13	2.00	0.42
1:D:133:PHE:CE2	1:D:158:ARG:HD2	2.55	0.42
1:C:282:GLN:HB2	1:C:285:LYS:CG	2.49	0.42
1:D:103:ILE:HG22	1:D:105:LYS:HG2	2.01	0.42
1:B:136:GLU:O	1:B:140:LYS:HG3	2.20	0.42
1:B:183:ILE:HD13	1:D:174:THR:HG23	2.02	0.42
1:B:36:VAL:HG21	1:B:97:GLU:HB3	2.02	0.41
1:A:36:VAL:HG21	1:A:97:GLU:HB3	2.01	0.41
1:B:275:ALA:O	1:B:279:VAL:HG23	2.21	0.41
1:D:136:GLU:O	1:D:140:LYS:HG3	2.21	0.41
1:C:161:VAL:HG13	1:C:162:PRO:HD2	2.02	0.41
1:D:233:TYR:HA	1:D:303:ILE:HG12	2.02	0.41
1:C:267:SER:O	1:C:271:LYS:HG2	2.21	0.41
1:A:229:THR:HG22	1:A:231:VAL:H	1.86	0.41
1:A:232:ASP:CB	1:A:302:GLU:HG3	2.51	0.41
1:D:6:VAL:HA	1:D:46:ILE:O	2.20	0.41
1:D:114:VAL:CG2	1:D:192:GLY:HA3	2.50	0.40
1:D:233:TYR:CE1	1:D:302:GLU:HB3	2.55	0.40
1:A:6:VAL:HA	1:A:46:ILE:O	2.21	0.40
1:A:21:ALA:O	1:A:25:GLN:HG3	2.22	0.40
1:B:9:LEU:O	1:B:49:HIS:HA	2.21	0.40
1:D:234:VAL:H	1:D:303:ILE:HD11	1.87	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:288:ILE:HA	1:D:303:ILE:O	2.21	0.40
1:D:28:LEU:HA	1:D:31:THR:OG1	2.22	0.40
1:C:32:SER:O	1:C:36:VAL:HG23	2.21	0.40
1:D:21:ALA:O	1:D:25:GLN:HG3	2.21	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	305/310 (98%)	282 (92%)	16 (5%)	7 (2%)	6	21
1	B	305/310 (98%)	282 (92%)	14 (5%)	9 (3%)	4	15
1	C	305/310 (98%)	284 (93%)	14 (5%)	7 (2%)	6	21
1	D	305/310 (98%)	278 (91%)	20 (7%)	7 (2%)	6	21
All	All	1220/1240 (98%)	1126 (92%)	64 (5%)	30 (2%)	5	19

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	238	TYR
1	A	299	SER
1	A	303	ILE
1	B	247	THR
1	B	299	SER
1	B	303	ILE
1	C	299	SER
1	C	303	ILE
1	D	238	TYR
1	D	299	SER
1	D	303	ILE

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Mol	Chain	Res	Type
1	A	247	THR
1	A	307	VAL
1	B	267	SER
1	B	307	VAL
1	C	238	TYR
1	D	243	GLU
1	D	307	VAL
1	B	298	MET
1	C	307	VAL
1	A	298	MET
1	B	243	GLU
1	C	298	MET
1	D	52	GLY
1	B	52	GLY
1	C	233	TYR
1	D	298	MET
1	B	238	TYR
1	C	52	GLY
1	A	52	GLY

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	246/249 (99%)	235 (96%)	11 (4%)	27	60
1	B	246/249 (99%)	234 (95%)	12 (5%)	25	57
1	C	246/249 (99%)	233 (95%)	13 (5%)	22	54
1	D	246/249 (99%)	236 (96%)	10 (4%)	30	64
All	All	984/996 (99%)	938 (95%)	46 (5%)	26	59

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	GLN
1	A	54	GLN

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Mol	Chain	Res	Type
1	A	99	ASN
1	A	176	ASN
1	A	195	VAL
1	A	204	GLU
1	A	220	VAL
1	A	232	ASP
1	A	298	MET
1	A	302	GLU
1	A	307	VAL
1	B	30	GLN
1	B	54	GLN
1	B	88	TYR
1	B	99	ASN
1	B	176	ASN
1	B	195	VAL
1	B	204	GLU
1	B	220	VAL
1	B	232	ASP
1	B	298	MET
1	B	302	GLU
1	B	307	VAL
1	C	30	GLN
1	C	54	GLN
1	C	99	ASN
1	C	161	VAL
1	C	176	ASN
1	C	195	VAL
1	C	204	GLU
1	C	220	VAL
1	C	232	ASP
1	C	248	ASN
1	C	298	MET
1	C	302	GLU
1	C	307	VAL
1	D	30	GLN
1	D	54	GLN
1	D	99	ASN
1	D	176	ASN
1	D	195	VAL
1	D	204	GLU
1	D	220	VAL
1	D	298	MET

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Mol	Chain	Res	Type
1	D	302	GLU
1	D	307	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	54	GLN
1	A	61	GLN
1	A	63	GLN
1	A	96	GLN
1	A	176	ASN
1	B	24	GLN
1	B	54	GLN
1	B	61	GLN
1	B	63	GLN
1	B	170	HIS
1	B	176	ASN
1	B	248	ASN
1	C	24	GLN
1	C	54	GLN
1	C	61	GLN
1	C	63	GLN
1	C	176	ASN
1	D	24	GLN
1	D	54	GLN
1	D	61	GLN
1	D	63	GLN
1	D	176	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 carbohydrates 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
2	SO4	B	2375	-	4,4,4	0.75	0	6,6,6	0.54	0
2	SO4	A	1375	-	4,4,4	0.70	0	6,6,6	0.63	0
2	SO4	D	4375	-	4,4,4	0.72	0	6,6,6	0.56	0
2	SO4	C	3375	-	4,4,4	0.71	0	6,6,6	0.67	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [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	307/310 (99%)	-0.18	11 (3%) 42 32	5, 29, 68, 94	0
1	B	307/310 (99%)	0.28	34 (11%) 5 3	11, 37, 86, 96	0
1	C	307/310 (99%)	-0.33	10 (3%) 46 36	4, 26, 63, 88	0
1	D	307/310 (99%)	0.24	30 (9%) 7 4	7, 31, 93, 100	0
All	All	1228/1240 (99%)	0.01	85 (6%) 16 10	4, 31, 82, 100	0

All (85) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	261	GLY	10.7
1	A	308	VAL	9.1
1	A	300	GLY	9.0
1	D	230	GLY	7.1
1	A	299	SER	7.1
1	D	300	GLY	7.0
1	B	299	SER	6.7
1	D	241	PRO	6.6
1	D	266	GLY	6.6
1	C	308	VAL	6.5
1	B	308	VAL	6.0
1	D	299	SER	6.0
1	D	308	VAL	5.9
1	D	238	TYR	5.5
1	B	265	PRO	5.2
1	D	236	ILE	5.0
1	D	239	GLY	5.0
1	D	244	LYS	5.0
1	B	298	MET	4.9
1	D	263	PHE	4.8
1	B	301	ASP	4.8

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Mol	Chain	Res	Type	RSRZ
1	C	144	GLN	4.6
1	D	298	MET	4.5
1	D	231	VAL	4.4
1	D	144	GLN	4.4
1	D	232	ASP	4.4
1	B	266	GLY	4.4
1	D	237	ASN	4.3
1	D	264	ALA	4.2
1	B	259	GLN	4.1
1	B	300	GLY	4.1
1	B	238	TYR	4.0
1	D	297	SER	4.0
1	B	261	GLY	4.0
1	D	259	GLN	3.9
1	D	243	GLU	3.9
1	D	143	MET	3.7
1	B	144	GLN	3.7
1	A	298	MET	3.7
1	D	301	ASP	3.7
1	C	153	ALA	3.7
1	B	42	GLY	3.6
1	B	239	GLY	3.6
1	D	240	LYS	3.4
1	B	237	ASN	3.4
1	D	265	PRO	3.4
1	C	299	SER	3.4
1	A	261	GLY	3.3
1	B	242	ASP	3.2
1	A	144	GLN	3.2
1	C	302	GLU	3.1
1	B	262	HIS	3.1
1	B	155	ARG	3.0
1	A	266	GLY	2.9
1	D	153	ALA	2.9
1	B	264	ALA	2.9
1	B	263	PHE	2.9
1	B	267	SER	2.8
1	B	307	VAL	2.8
1	B	241	PRO	2.7
1	B	260	ALA	2.7
1	B	252	ALA	2.7
1	B	284	ASN	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	302	GLU	2.7
1	A	143	MET	2.6
1	C	301	ASP	2.6
1	D	242	ASP	2.6
1	B	233	TYR	2.5
1	B	268	MET	2.4
1	A	301	ASP	2.4
1	A	259	GLN	2.4
1	B	198	GLN	2.3
1	A	264	ALA	2.3
1	D	235	CYS	2.3
1	B	229	THR	2.2
1	B	243	GLU	2.2
1	D	302	GLU	2.2
1	B	247	THR	2.2
1	D	296	GLY	2.1
1	B	124	LYS	2.1
1	C	155	ARG	2.1
1	C	259	GLN	2.1
1	C	147	ALA	2.1
1	C	300	GLY	2.1
1	B	148	ILE	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 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. 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(\AA^2)	Q<0.9
2	SO4	C	3375	5/5	0.97	0.13	19,23,27,31	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	SO4	B	2375	5/5	0.98	0.10	33,37,40,42	0
2	SO4	D	4375	5/5	0.98	0.12	43,44,47,47	0
2	SO4	A	1375	5/5	0.99	0.08	13,14,20,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.