



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:50 AM GMT

PDB ID : 3G7K
Title : Crystal Structure of Methylitaconate-delta-isomerase
Authors : Messerschmidt, A.; Macieira, S.; Velarde, M.
Deposited on : 2009-02-10
Resolution : 2.70 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

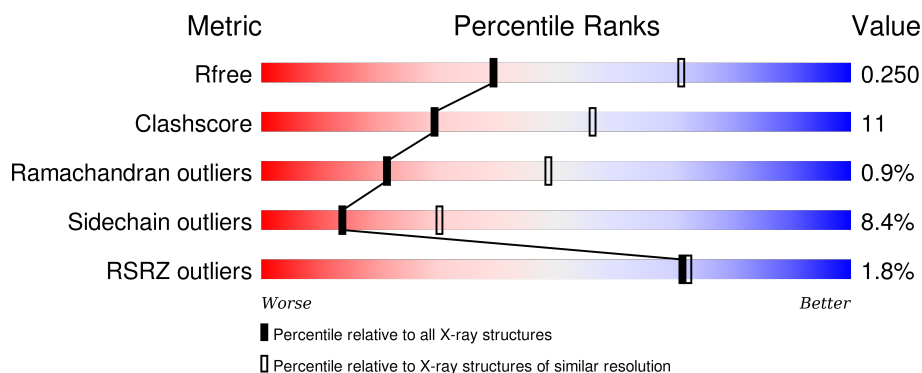
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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	391	<div> <div></div> <div>80% 13% . .</div> </div>
1	B	391	<div> <div>2%</div> <div>76% 20% . .</div> </div>
1	C	391	<div> <div>4%</div> <div>41% 13% . 43%</div> </div>
1	D	391	<div> <div></div> <div>75% 19% . .</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10592 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 3-methylitaconate isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	379	Total	C	N	O	S	0	0	0
			2807	1757	474	558	18			
1	B	387	Total	C	N	O	S	0	0	0
			2868	1795	486	569	18			
1	C	224	Total	C	N	O	S	0	0	0
			1645	1033	284	319	9			
1	D	379	Total	C	N	O	S	0	0	0
			2807	1757	474	558	18			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	381	GLY	-	EXPRESSION TAG	UNP Q0QLE6
A	382	SER	-	EXPRESSION TAG	UNP Q0QLE6
A	383	ALA	-	EXPRESSION TAG	UNP Q0QLE6
A	384	TRP	-	EXPRESSION TAG	UNP Q0QLE6
A	385	SER	-	EXPRESSION TAG	UNP Q0QLE6
A	386	HIS	-	EXPRESSION TAG	UNP Q0QLE6
A	387	PRO	-	EXPRESSION TAG	UNP Q0QLE6
A	388	GLN	-	EXPRESSION TAG	UNP Q0QLE6
A	389	PHE	-	EXPRESSION TAG	UNP Q0QLE6
A	390	GLU	-	EXPRESSION TAG	UNP Q0QLE6
A	391	LYS	-	EXPRESSION TAG	UNP Q0QLE6
B	381	GLY	-	EXPRESSION TAG	UNP Q0QLE6
B	382	SER	-	EXPRESSION TAG	UNP Q0QLE6
B	383	ALA	-	EXPRESSION TAG	UNP Q0QLE6
B	384	TRP	-	EXPRESSION TAG	UNP Q0QLE6
B	385	SER	-	EXPRESSION TAG	UNP Q0QLE6
B	386	HIS	-	EXPRESSION TAG	UNP Q0QLE6
B	387	PRO	-	EXPRESSION TAG	UNP Q0QLE6
B	388	GLN	-	EXPRESSION TAG	UNP Q0QLE6
B	389	PHE	-	EXPRESSION TAG	UNP Q0QLE6
B	390	GLU	-	EXPRESSION TAG	UNP Q0QLE6

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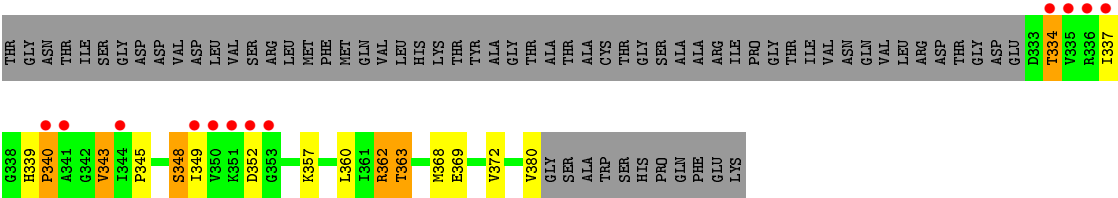
Chain	Residue	Modelled	Actual	Comment	Reference
B	391	LYS	-	EXPRESSION TAG	UNP Q0QLE6
C	381	GLY	-	EXPRESSION TAG	UNP Q0QLE6
C	382	SER	-	EXPRESSION TAG	UNP Q0QLE6
C	383	ALA	-	EXPRESSION TAG	UNP Q0QLE6
C	384	TRP	-	EXPRESSION TAG	UNP Q0QLE6
C	385	SER	-	EXPRESSION TAG	UNP Q0QLE6
C	386	HIS	-	EXPRESSION TAG	UNP Q0QLE6
C	387	PRO	-	EXPRESSION TAG	UNP Q0QLE6
C	388	GLN	-	EXPRESSION TAG	UNP Q0QLE6
C	389	PHE	-	EXPRESSION TAG	UNP Q0QLE6
C	390	GLU	-	EXPRESSION TAG	UNP Q0QLE6
C	391	LYS	-	EXPRESSION TAG	UNP Q0QLE6
D	381	GLY	-	EXPRESSION TAG	UNP Q0QLE6
D	382	SER	-	EXPRESSION TAG	UNP Q0QLE6
D	383	ALA	-	EXPRESSION TAG	UNP Q0QLE6
D	384	TRP	-	EXPRESSION TAG	UNP Q0QLE6
D	385	SER	-	EXPRESSION TAG	UNP Q0QLE6
D	386	HIS	-	EXPRESSION TAG	UNP Q0QLE6
D	387	PRO	-	EXPRESSION TAG	UNP Q0QLE6
D	388	GLN	-	EXPRESSION TAG	UNP Q0QLE6
D	389	PHE	-	EXPRESSION TAG	UNP Q0QLE6
D	390	GLU	-	EXPRESSION TAG	UNP Q0QLE6
D	391	LYS	-	EXPRESSION TAG	UNP Q0QLE6

- Molecule 2 is water.

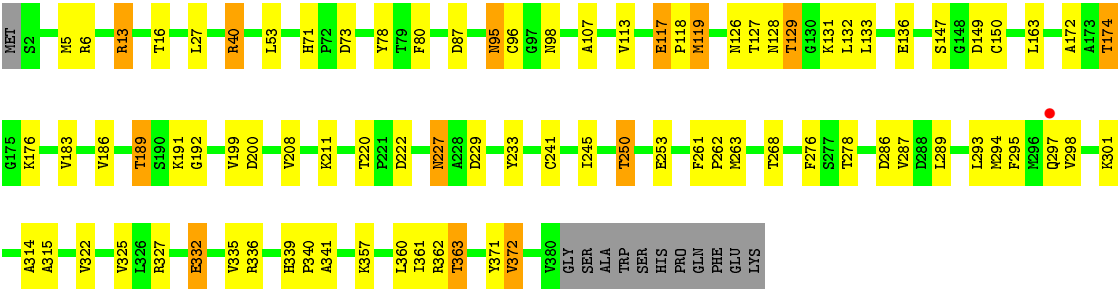
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	129	Total O 129 129	0	0
2	B	107	Total O 107 107	0	0
2	C	77	Total O 77 77	0	0
2	D	152	Total O 152 152	0	0

- Molecule 1: 3-methylitaconate isomerase





● Molecule 1: 3-methylitaconate isomerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	53.08Å 142.08Å 228.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.70 49.72 – 2.70	Depositor EDS
% Data completeness (in resolution range)	97.8 (50.00-2.70) 97.7 (49.72-2.70)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.65 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.200 , 0.256 0.199 , 0.250	Depositor DCC
R_{free} test set	2402 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	50.3	Xtriage
Anisotropy	0.354	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 53.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 47452 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10592	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.98% 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

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	0.63	1/2849 (0.0%)	0.74	3/3871 (0.1%)
1	B	0.58	0/2914	0.70	1/3961 (0.0%)
1	C	0.60	0/1669	0.70	0/2265
1	D	0.65	1/2849 (0.0%)	0.73	2/3871 (0.1%)
All	All	0.62	2/10281 (0.0%)	0.72	6/13968 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	C	0	1
1	D	0	1
All	All	0	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	332	GLU	CG-CD	6.73	1.62	1.51
1	A	9	CYS	CB-SG	-6.25	1.71	1.82

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	13	ARG	NE-CZ-NH2	-6.76	116.92	120.30
1	B	365	ARG	NE-CZ-NH2	-5.82	117.39	120.30
1	D	13	ARG	NE-CZ-NH2	-5.59	117.50	120.30
1	A	372	VAL	CB-CA-C	-5.33	101.28	111.40
1	A	117	GLU	C-N-CD	-5.27	109.00	120.60
1	D	117	GLU	C-N-CD	-5.11	109.37	120.60

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	117	GLU	Peptide
1	C	117	GLU	Peptide
1	D	117	GLU	Peptide

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	2807	0	2819	53	0
1	B	2868	0	2869	77	0
1	C	1645	0	1652	44	0
1	D	2807	0	2819	66	0
2	A	129	0	0	5	0
2	B	107	0	0	5	0
2	C	77	0	0	5	0
2	D	152	0	0	8	0
All	All	10592	0	10159	229	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 (229) 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:D:13:ARG:NH2	1:D:363:THR:HG22	1.45	1.31
1:D:297:GLN:HG3	2:D:434:HOH:O	1.34	1.28
1:A:13:ARG:NH2	1:A:363:THR:HG22	1.51	1.25
1:B:13:ARG:NH2	1:B:363:THR:HG22	1.53	1.21
1:D:13:ARG:HH22	1:D:363:THR:CG2	1.56	1.18
1:A:13:ARG:HH22	1:A:363:THR:CG2	1.60	1.12
1:C:13:ARG:NH2	1:C:363:THR:HG23	1.66	1.10
1:C:13:ARG:HH21	1:C:363:THR:HG23	1.09	1.08
1:B:13:ARG:HH22	1:B:363:THR:CG2	1.68	1.06
1:D:13:ARG:HH22	1:D:363:THR:HG22	0.92	1.04

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:96:CYS:SG	2:D:448:HOH:O	2.14	1.04
1:B:13:ARG:HH22	1:B:363:THR:HG22	1.03	1.03
1:B:287:VAL:O	1:B:327:ARG:NH1	1.96	0.97
1:A:189:THR:HG22	1:A:191:LYS:H	1.28	0.97
1:D:95:ASN:ND2	1:D:96:CYS:H	1.62	0.96
1:B:189:THR:HG22	1:B:191:LYS:H	1.30	0.95
1:D:95:ASN:HD22	1:D:96:CYS:H	1.12	0.94
1:B:126:ASN:ND2	1:B:129:THR:HB	1.86	0.91
1:B:126:ASN:HD22	1:B:129:THR:HB	1.37	0.88
1:D:126:ASN:HD22	1:D:129:THR:HB	1.39	0.88
1:B:80:PHE:H	1:B:128:ASN:HD21	1.14	0.87
1:C:71:HIS:HD2	1:C:73:ASP:H	1.20	0.87
1:C:148:GLY:O	1:C:158:THR:HG23	1.72	0.86
1:D:71:HIS:HD2	1:D:73:ASP:H	1.22	0.86
1:A:148:GLY:O	1:A:158:THR:HG23	1.76	0.85
1:A:80:PHE:H	1:A:128:ASN:HD21	1.19	0.85
1:B:189:THR:HB	1:B:192:GLY:O	1.77	0.84
1:D:80:PHE:H	1:D:128:ASN:HD21	1.26	0.83
1:C:126:ASN:HD22	1:C:129:THR:HB	1.43	0.83
1:D:250:THR:HG22	1:D:253:GLU:H	1.46	0.81
1:D:129:THR:HG22	1:D:131:LYS:H	1.47	0.80
1:B:250:THR:HG22	1:B:253:GLU:H	1.48	0.79
1:D:363:THR:HG21	2:D:408:HOH:O	1.83	0.79
1:B:129:THR:HG21	2:B:477:HOH:O	1.85	0.77
1:D:126:ASN:ND2	1:D:129:THR:HB	2.00	0.77
1:D:189:THR:CG2	1:D:191:LYS:H	1.97	0.77
1:D:95:ASN:HD22	1:D:96:CYS:N	1.83	0.76
1:C:129:THR:HG22	1:C:131:LYS:H	1.48	0.76
1:C:334:THR:HG22	1:C:345:PRO:HB2	1.70	0.74
1:B:371:TYR:OH	1:C:369:GLU:OE2	2.06	0.73
1:D:174:THR:HG22	1:D:176:LYS:H	1.53	0.73
1:B:278:THR:O	1:B:280:ASN:N	2.18	0.72
1:D:189:THR:HG22	1:D:191:LYS:H	1.55	0.71
1:B:189:THR:HG22	1:B:191:LYS:N	2.06	0.70
1:B:40:ARG:NH2	1:B:378:LYS:HD2	2.07	0.69
1:A:275:ASP:HB2	1:A:282:ILE:CD1	2.24	0.67
1:A:363:THR:HG21	2:A:451:HOH:O	1.93	0.67
1:A:189:THR:HB	1:A:192:GLY:O	1.94	0.67
1:B:189:THR:HG21	1:B:233:TYR:CE1	2.30	0.66
1:B:189:THR:CG2	1:B:233:TYR:HE1	2.09	0.66
1:B:129:THR:HG22	1:B:131:LYS:H	1.61	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:95:ASN:ND2	1:D:96:CYS:N	2.40	0.65
1:A:360:LEU:HD12	1:A:362:ARG:HD3	1.77	0.65
1:A:129:THR:HG23	1:A:131:LYS:HG2	1.80	0.64
1:C:134:TYR:HB2	1:C:165:ASP:HB3	1.79	0.63
1:B:339:HIS:HB2	1:B:340:PRO:CD	2.28	0.63
1:B:357:LYS:HE2	2:B:462:HOH:O	1.98	0.63
1:B:331:ASP:HA	2:B:441:HOH:O	1.98	0.63
1:A:352:ASP:HA	2:A:514:HOH:O	1.97	0.62
1:D:13:ARG:NH2	1:D:363:THR:CG2	2.30	0.62
1:B:189:THR:HG21	1:B:233:TYR:OH	1.99	0.62
1:B:329:THR:HG22	1:B:332:GLU:HB3	1.81	0.62
1:D:119:MET:HE1	1:D:136:GLU:HA	1.81	0.62
1:C:80:PHE:H	1:C:128:ASN:HD21	1.46	0.62
1:B:53:LEU:HD11	1:B:372:VAL:HG13	1.82	0.62
1:B:220:THR:HG22	1:B:222:ASP:H	1.63	0.62
1:B:303:TYR:HD2	1:B:308:THR:HG1	1.46	0.61
1:A:13:ARG:HH22	1:A:363:THR:HG22	0.66	0.61
1:A:339:HIS:HB2	1:A:340:PRO:CD	2.30	0.61
1:C:50:ILE:O	1:D:13:ARG:HG3	2.01	0.61
1:B:278:THR:C	1:B:280:ASN:H	2.05	0.60
1:C:168:ASP:HB3	2:C:467:HOH:O	2.01	0.60
1:C:169:THR:HG22	2:C:465:HOH:O	2.02	0.60
1:D:53:LEU:HD11	1:D:372:VAL:HG13	1.83	0.59
1:A:189:THR:CG2	1:A:191:LYS:H	2.11	0.59
1:D:189:THR:HG22	1:D:191:LYS:N	2.17	0.59
1:A:168:ASP:O	1:A:168:ASP:OD1	2.19	0.59
1:D:136:GLU:HB2	1:D:163:LEU:HB3	1.83	0.59
1:A:361:ILE:O	1:A:362:ARG:HD2	2.03	0.59
1:C:115:ALA:HB1	1:C:139:VAL:HG11	1.84	0.59
1:D:189:THR:HB	1:D:192:GLY:O	2.03	0.58
1:A:275:ASP:HB2	1:A:282:ILE:HD13	1.85	0.58
1:B:339:HIS:HB2	1:B:340:PRO:HD3	1.84	0.58
1:A:189:THR:HG22	1:A:191:LYS:N	2.10	0.58
1:D:16:THR:OG1	1:D:98:ASN:ND2	2.37	0.58
1:C:363:THR:HG21	2:C:407:HOH:O	2.03	0.57
1:B:363:THR:HG21	2:B:396:HOH:O	2.04	0.57
1:B:13:ARG:NH2	1:B:363:THR:CG2	2.40	0.57
1:A:50:ILE:O	1:B:13:ARG:HG3	2.04	0.57
1:B:119:MET:HE1	1:B:136:GLU:HG2	1.86	0.56
1:D:189:THR:HG23	1:D:191:LYS:H	1.71	0.56
1:D:189:THR:HG21	1:D:233:TYR:OH	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:119:MET:HE1	1:D:136:GLU:CA	2.36	0.56
1:D:339:HIS:HB2	1:D:340:PRO:CD	2.36	0.56
1:B:148:GLY:O	1:B:158:THR:HG23	2.06	0.56
1:C:339:HIS:HB2	1:C:340:PRO:HD2	1.87	0.55
1:C:126:ASN:ND2	1:C:129:THR:HB	2.18	0.55
1:C:128:ASN:ND2	1:C:128:ASN:H	2.05	0.55
1:B:189:THR:CG2	1:B:233:TYR:CE1	2.88	0.54
1:D:332:GLU:HG2	2:D:470:HOH:O	2.08	0.54
1:B:129:THR:HG23	1:B:131:LYS:HG2	1.89	0.53
1:A:189:THR:HG21	1:A:233:TYR:OH	2.07	0.53
1:A:126:ASN:ND2	1:A:129:THR:HB	2.24	0.53
1:A:189:THR:CG2	1:A:233:TYR:HE1	2.21	0.53
1:C:32:GLU:HG2	2:C:464:HOH:O	2.09	0.53
1:B:269:LYS:N	1:B:325:VAL:HG22	2.24	0.53
1:A:129:THR:HG22	1:A:131:LYS:H	1.73	0.53
1:B:199:VAL:HG21	1:B:314:ALA:HB2	1.91	0.53
1:B:172:ALA:HB3	2:B:455:HOH:O	2.08	0.53
1:A:250:THR:HG22	1:A:252:LYS:H	1.72	0.52
1:D:339:HIS:HB2	1:D:340:PRO:HD2	1.90	0.52
1:D:287:VAL:O	1:D:327:ARG:NH1	2.42	0.52
1:B:327:ARG:HG2	1:B:328:ASP:O	2.10	0.52
1:D:129:THR:CG2	1:D:131:LYS:HG2	2.39	0.52
1:C:129:THR:HG22	1:C:131:LYS:N	2.22	0.52
1:B:70:THR:HB	1:B:77:ASP:OD1	2.09	0.52
1:D:276:PHE:HE1	1:D:341:ALA:HB2	1.74	0.52
1:A:183:VAL:HG23	1:A:184:VAL:HG23	1.92	0.52
1:B:204:PRO:HB2	1:B:262:PRO:HB3	1.91	0.52
1:C:369:GLU:OE2	1:D:6:ARG:NH1	2.44	0.51
1:B:189:THR:HG23	1:B:233:TYR:HE1	1.74	0.51
1:B:80:PHE:H	1:B:128:ASN:ND2	1.95	0.51
1:A:361:ILE:O	1:A:362:ARG:CD	2.59	0.51
1:B:115:ALA:HB1	1:B:139:VAL:HG11	1.92	0.51
1:C:13:ARG:NH2	1:C:363:THR:CG2	2.57	0.51
1:D:128:ASN:H	1:D:128:ASN:HD22	1.58	0.51
1:A:12:MET:HG2	1:A:366:ARG:HA	1.93	0.51
1:A:115:ALA:HB1	1:A:139:VAL:HG11	1.93	0.51
1:D:263:MET:HE3	1:D:295:PHE:CD1	2.46	0.51
1:B:189:THR:HG21	1:B:233:TYR:HE1	1.73	0.51
1:C:334:THR:CG2	1:C:345:PRO:HB2	2.38	0.51
1:B:9:CYS:SG	1:B:372:VAL:HG22	2.51	0.51
1:C:77:ASP:OD2	1:C:125:HIS:HD2	1.94	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:189:THR:HG21	1:B:233:TYR:CZ	2.46	0.50
1:C:96:CYS:HB3	1:C:99:ILE:HB	1.94	0.50
1:A:31:GLN:HE21	1:A:31:GLN:HA	1.77	0.50
1:D:5:MET:HE1	2:D:421:HOH:O	2.12	0.49
1:C:348:SER:O	1:C:349:ILE:HG13	2.12	0.49
1:D:119:MET:HE1	1:D:136:GLU:CB	2.42	0.49
1:D:276:PHE:CE1	1:D:341:ALA:HB2	2.48	0.49
1:A:31:GLN:NE2	1:A:34:ARG:HD2	2.27	0.49
1:D:71:HIS:CD2	1:D:73:ASP:H	2.14	0.49
1:D:53:LEU:HD11	1:D:372:VAL:CG1	2.43	0.49
1:A:77:ASP:OD2	1:A:125:HIS:HD2	1.95	0.49
1:D:129:THR:HG22	1:D:131:LYS:N	2.22	0.49
1:C:337:ILE:O	1:C:343:VAL:HA	2.13	0.48
1:C:71:HIS:CD2	1:C:73:ASP:H	2.12	0.48
1:A:80:PHE:H	1:A:128:ASN:ND2	2.00	0.48
1:B:236:GLU:O	1:B:240:LYS:HG3	2.14	0.48
1:B:9:CYS:CB	1:B:372:VAL:HG22	2.44	0.48
1:C:95:ASN:HD21	1:C:133:LEU:HD11	1.78	0.48
1:B:70:THR:O	1:B:70:THR:CG2	2.61	0.48
1:B:303:TYR:HD2	1:B:308:THR:OG1	1.97	0.48
1:B:235:GLU:HG2	1:B:297:GLN:HG3	1.96	0.47
1:B:276:PHE:HE1	1:B:341:ALA:HB2	1.80	0.47
1:D:13:ARG:HH22	1:D:363:THR:HG23	1.64	0.47
1:A:189:THR:HG21	1:A:233:TYR:CE1	2.50	0.47
1:C:368:MET:HG3	1:D:371:TYR:O	2.15	0.47
1:D:199:VAL:HG21	1:D:314:ALA:HB2	1.97	0.47
1:D:107:ALA:HB1	1:D:113:VAL:HG23	1.97	0.46
1:A:245:ILE:HG23	1:A:245:ILE:O	2.16	0.46
1:C:171:GLY:O	1:C:172:ALA:C	2.53	0.46
1:A:131:LYS:NZ	1:A:170:ALA:O	2.48	0.46
1:C:98:ASN:O	1:C:101:ALA:HB3	2.16	0.46
1:C:362:ARG:HD2	1:C:362:ARG:HA	1.42	0.45
1:B:80:PHE:N	1:B:128:ASN:HD21	1.97	0.45
1:D:40:ARG:CZ	2:D:519:HOH:O	2.64	0.45
1:B:276:PHE:CE1	1:B:341:ALA:HB2	2.52	0.45
1:B:114:LYS:N	1:C:68:PRO:HG2	2.31	0.45
1:C:80:PHE:CE1	1:C:96:CYS:HB2	2.52	0.45
1:D:263:MET:HE1	1:D:293:LEU:HD21	1.98	0.45
1:B:69:SER:OG	1:B:74:ALA:HB3	2.17	0.44
1:D:361:ILE:O	1:D:362:ARG:HD2	2.17	0.44
1:B:361:ILE:O	1:B:362:ARG:HD2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:70:THR:HG22	1:C:125:HIS:CD2	2.52	0.44
1:A:362:ARG:HA	1:A:362:ARG:HD2	1.60	0.44
1:B:235:GLU:HA	1:B:235:GLU:OE1	2.17	0.44
1:D:211:LYS:HD3	2:D:513:HOH:O	2.18	0.44
1:C:131:LYS:HD3	1:C:131:LYS:HA	1.73	0.44
1:C:51:ASP:OD1	1:D:13:ARG:HD3	2.18	0.44
1:A:166:PHE:CE1	1:A:360:LEU:HD21	2.53	0.44
1:C:14:ALA:HB3	1:C:99:ILE:HD13	2.00	0.44
1:A:371:TYR:O	1:B:368:MET:HG3	2.18	0.44
1:B:302:THR:O	1:B:339:HIS:HB2	2.18	0.44
1:C:20:ILE:HD12	1:C:42:PHE:CZ	2.52	0.44
1:B:349:ILE:HD12	1:B:357:LYS:HD2	2.01	0.43
1:C:150:CYS:HB3	1:C:159:ASN:O	2.19	0.43
1:A:129:THR:CG2	1:A:131:LYS:H	2.31	0.43
1:A:126:ASN:HD22	1:A:129:THR:HB	1.83	0.43
1:A:220:THR:HG22	1:A:223:VAL:H	1.83	0.43
1:B:129:THR:CG2	1:B:131:LYS:H	2.26	0.43
1:D:227:ASN:C	1:D:227:ASN:HD22	2.21	0.43
1:A:240:LYS:HE3	2:A:461:HOH:O	2.19	0.43
1:A:198:ILE:HG21	1:A:245:ILE:HD13	2.00	0.43
1:B:293:LEU:O	1:B:299:LEU:HD12	2.19	0.43
1:D:132:LEU:O	1:D:133:LEU:HD23	2.19	0.43
1:B:220:THR:HB	1:B:223:VAL:HG23	2.01	0.42
1:B:326:LEU:HG	1:B:327:ARG:O	2.18	0.42
1:B:263:MET:HE3	1:B:295:PHE:HD1	1.84	0.42
1:A:80:PHE:N	1:A:128:ASN:HD21	2.01	0.42
1:C:167:SER:HB3	1:C:357:LYS:HA	2.01	0.42
1:A:9:CYS:SG	1:A:372:VAL:HG22	2.59	0.42
1:A:119:MET:HA	1:A:137:VAL:O	2.20	0.42
1:D:301:LYS:HB2	1:D:301:LYS:HE2	1.82	0.42
1:B:322:VAL:O	1:B:325:VAL:HG12	2.20	0.42
1:C:3:ASP:OD2	1:D:147:SER:N	2.48	0.42
1:D:286:ASP:O	1:D:336:ARG:HD2	2.20	0.41
1:A:339:HIS:HB2	1:A:340:PRO:HD3	2.02	0.41
1:A:189:THR:HG23	1:A:233:TYR:HE1	1.84	0.41
1:B:126:ASN:HD22	1:B:129:THR:CB	2.20	0.41
1:A:365:ARG:NH2	1:B:373:TYR:O	2.53	0.41
1:D:294:MET:HB3	1:D:294:MET:HE2	1.83	0.41
1:D:241:CYS:O	1:D:245:ILE:HG12	2.19	0.41
1:B:9:CYS:HB2	1:B:372:VAL:HG22	2.01	0.41
1:D:129:THR:HG21	2:D:540:HOH:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:128:ASN:ND2	1:B:128:ASN:H	2.18	0.41
1:A:259:PRO:HB2	1:A:297:GLN:HE22	1.86	0.41
1:A:47:VAL:HG22	2:A:504:HOH:O	2.21	0.41
1:B:16:THR:OG1	1:B:98:ASN:ND2	2.52	0.41
1:C:172:ALA:HB3	2:C:458:HOH:O	2.21	0.41
1:D:278:THR:O	1:D:278:THR:CG2	2.69	0.41
1:B:286:ASP:O	1:B:336:ARG:HD2	2.21	0.41
1:B:58:PRO:O	1:B:62:LYS:HE2	2.21	0.41
1:D:174:THR:HB	1:D:200:ASP:OD2	2.22	0.40
1:D:315:ALA:HA	1:D:322:VAL:HG12	2.02	0.40
1:A:51:ASP:OD1	1:B:13:ARG:HD3	2.21	0.40
1:A:59:LEU:HD13	2:A:429:HOH:O	2.21	0.40
1:B:348:SER:HA	1:B:357:LYS:O	2.21	0.40
1:D:261:PHE:HA	1:D:262:PRO:C	2.41	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	377/391 (96%)	361 (96%)	14 (4%)	2 (0%)	34 63
1	B	385/391 (98%)	360 (94%)	22 (6%)	3 (1%)	24 51
1	C	220/391 (56%)	200 (91%)	15 (7%)	5 (2%)	8 20
1	D	377/391 (96%)	356 (94%)	19 (5%)	2 (0%)	34 63
All	All	1359/1564 (87%)	1277 (94%)	70 (5%)	12 (1%)	21 49

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	118	PRO

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Mol	Chain	Res	Type
1	B	93	ASN
1	B	279	GLY
1	C	118	PRO
1	C	172	ALA
1	D	118	PRO
1	A	46	ASP
1	D	172	ALA
1	C	352	ASP
1	B	46	ASP
1	C	340	PRO
1	C	88	ALA

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	308/319 (97%)	285 (92%)	23 (8%)	17	38
1	B	314/319 (98%)	291 (93%)	23 (7%)	17	39
1	C	175/319 (55%)	157 (90%)	18 (10%)	9	20
1	D	308/319 (97%)	279 (91%)	29 (9%)	11	25
All	All	1105/1276 (87%)	1012 (92%)	93 (8%)	14	30

All (93) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	LEU
1	A	31	GLN
1	A	47	VAL
1	A	48	ARG
1	A	70	THR
1	A	76	VAL
1	A	119	MET
1	A	127	THR
1	A	129	THR
1	A	133	LEU

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Mol	Chain	Res	Type
1	A	158	THR
1	A	220	THR
1	A	245	ILE
1	A	255	SER
1	A	289	LEU
1	A	325	VAL
1	A	331	ASP
1	A	335	VAL
1	A	360	LEU
1	A	362	ARG
1	A	363	THR
1	A	365	ARG
1	A	372	VAL
1	B	5	MET
1	B	27	LEU
1	B	70	THR
1	B	86	THR
1	B	119	MET
1	B	128	ASN
1	B	129	THR
1	B	136	GLU
1	B	161	PRO
1	B	208	VAL
1	B	222	ASP
1	B	229	ASP
1	B	235	GLU
1	B	250	THR
1	B	285	ASP
1	B	289	LEU
1	B	303	TYR
1	B	325	VAL
1	B	329	THR
1	B	331	ASP
1	B	335	VAL
1	B	363	THR
1	B	372	VAL
1	C	27	LEU
1	C	31	GLN
1	C	32	GLU
1	C	40	ARG
1	C	50	ILE
1	C	62	LYS

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Mol	Chain	Res	Type
1	C	70	THR
1	C	128	ASN
1	C	158	THR
1	C	176	LYS
1	C	334	THR
1	C	343	VAL
1	C	348	SER
1	C	360	LEU
1	C	362	ARG
1	C	363	THR
1	C	372	VAL
1	C	380	VAL
1	D	27	LEU
1	D	40	ARG
1	D	78	TYR
1	D	87	ASP
1	D	95	ASN
1	D	119	MET
1	D	127	THR
1	D	129	THR
1	D	149	ASP
1	D	150	CYS
1	D	174	THR
1	D	183	VAL
1	D	186	VAL
1	D	189	THR
1	D	208	VAL
1	D	220	THR
1	D	222	ASP
1	D	227	ASN
1	D	229	ASP
1	D	250	THR
1	D	268	THR
1	D	289	LEU
1	D	298	VAL
1	D	325	VAL
1	D	335	VAL
1	D	357	LYS
1	D	360	LEU
1	D	363	THR
1	D	372	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	125	HIS
1	A	126	ASN
1	A	128	ASN
1	A	297	GLN
1	B	125	HIS
1	B	126	ASN
1	B	128	ASN
1	C	31	GLN
1	C	71	HIS
1	C	125	HIS
1	C	126	ASN
1	C	128	ASN
1	D	71	HIS
1	D	95	ASN
1	D	125	HIS
1	D	126	ASN
1	D	128	ASN
1	D	227	ASN
1	D	280	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 ⓘ

There are no ligands in this entry.

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	379/391 (96%)	-0.08	4 (1%) 82 83	26, 43, 59, 68	0
1	B	387/391 (98%)	0.10	6 (1%) 74 75	28, 48, 71, 84	0
1	C	224/391 (57%)	0.20	14 (6%) 23 22	32, 51, 100, 108	0
1	D	379/391 (96%)	-0.17	1 (0%) 94 95	24, 39, 55, 61	0
All	All	1369/1564 (87%)	-0.01	25 (1%) 71 72	24, 45, 69, 108	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	341	ALA	5.4
1	C	337	ILE	5.2
1	C	344	ILE	5.1
1	C	340	PRO	5.0
1	C	352	ASP	4.6
1	C	335	VAL	4.4
1	C	353	GLY	4.4
1	C	350	VAL	4.2
1	C	334	THR	3.8
1	B	223	VAL	3.6
1	C	336	ARG	3.6
1	C	349	ILE	3.4
1	C	176	LYS	3.0
1	C	177	VAL	2.9
1	A	352	ASP	2.9
1	A	351	LYS	2.9
1	A	185	ASP	2.8
1	A	215	MET	2.7
1	B	386	HIS	2.6
1	D	297	GLN	2.5
1	B	226	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	351	LYS	2.4
1	B	351	LYS	2.3
1	B	383	ALA	2.2
1	B	218	THR	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 carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.