



# Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 03:31 pm BST

PDB ID : 3EOM  
Title : 2.4 Å crystal structure of native glutaryl-coa dehydrogenase from Burkholderia pseudomallei  
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)  
Deposited on : 2008-09-28  
Resolution : 2.40 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](mailto: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.

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

MolProbity	:	4.02b-467
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

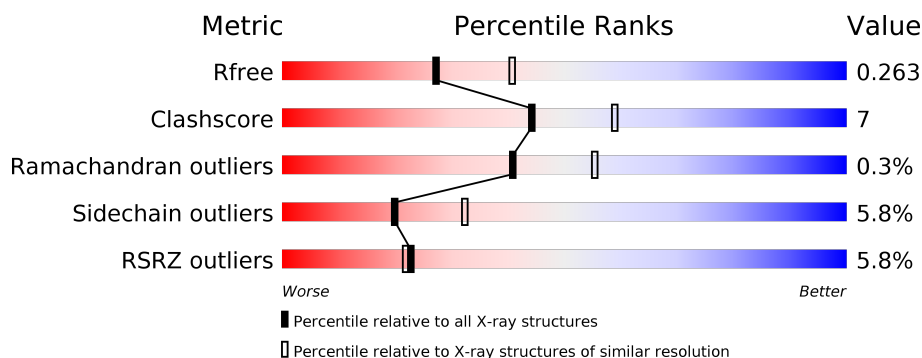
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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  $\geq 3$ , 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	396	<div> <div>4%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div>• •</div> </div> </div>
1	B	396	<div> <div>8%</div> <div> <div></div> <div>79%</div> <div>15%</div> <div>• 5%</div> </div> </div>
1	C	396	<div> <div>6%</div> <div> <div></div> <div>80%</div> <div>15%</div> <div>• •</div> </div> </div>
1	D	396	<div> <div>4%</div> <div> <div></div> <div>79%</div> <div>16%</div> <div>• •</div> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11908 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 Glutaryl-CoA dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	387	Total	C	N	O	S	0	0	0
			2976	1881	525	554	16			
1	B	378	Total	C	N	O	S	0	0	0
			2909	1839	515	539	16			
1	C	381	Total	C	N	O	S	0	0	0
			2924	1846	517	545	16			
1	D	381	Total	C	N	O	S	0	0	0
			2932	1854	519	543	16			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	EXPRESSION TAG	UNP Q3JP94
B	0	SER	-	EXPRESSION TAG	UNP Q3JP94
C	0	SER	-	EXPRESSION TAG	UNP Q3JP94
D	0	SER	-	EXPRESSION TAG	UNP Q3JP94

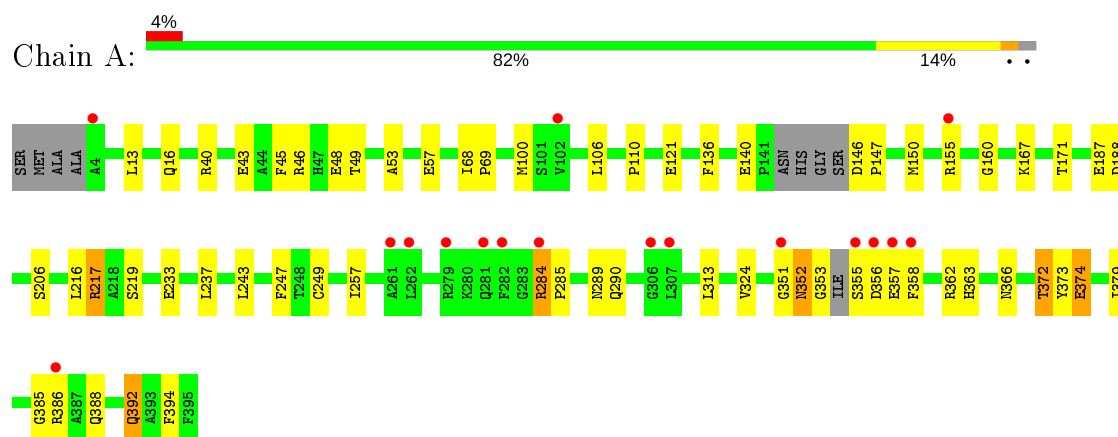
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	50	Total	O	0	0
			50	50		
2	B	34	Total	O	0	0
			34	34		
2	C	36	Total	O	0	0
			36	36		
2	D	47	Total	O	0	0
			47	47		

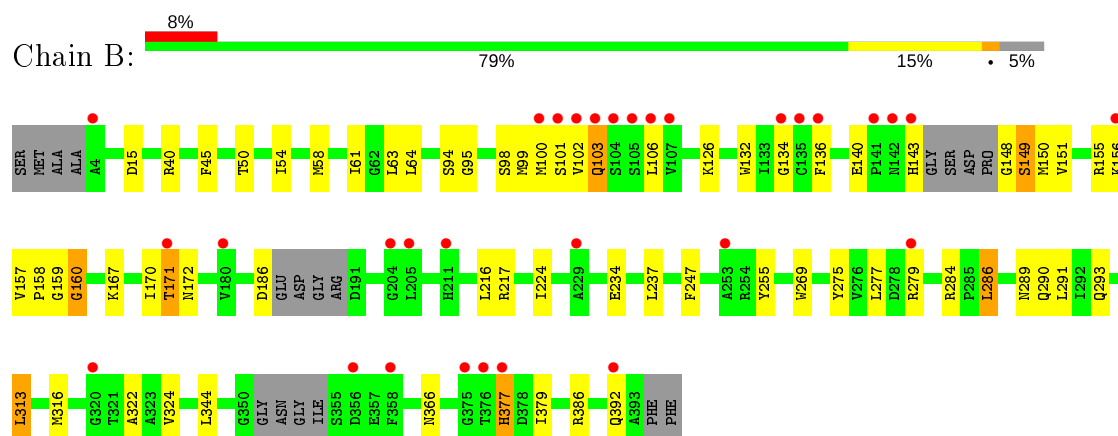
### 3 Residue-property plots [i](#)

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.

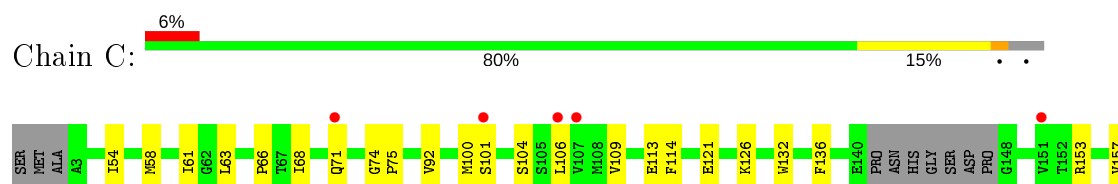
- Molecule 1: Glutaryl-CoA dehydrogenase

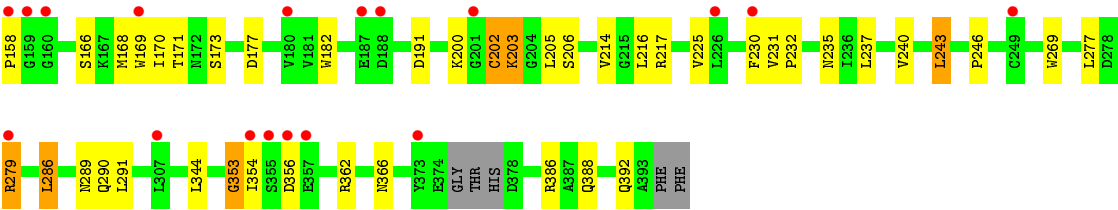


- Molecule 1: Glutaryl-CoA dehydrogenase

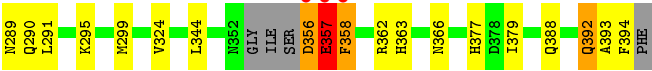
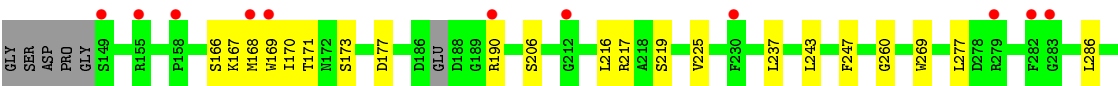
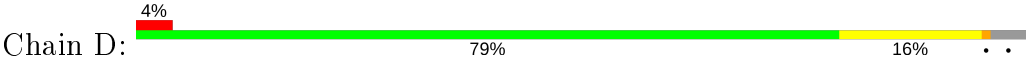


- Molecule 1: Glutaryl-CoA dehydrogenase





● Molecule 1: Glutaryl-CoA dehydrogenase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	98.11Å 106.77Å 145.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.40 29.21 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.3 (30.00-2.40) 99.3 (29.21-2.40)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.59 (at 2.39Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.211 , 0.264 0.208 , 0.263	Depositor DCC
$R_{free}$ test set	3035 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	53.2	Xtriage
Anisotropy	0.023	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 35.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	11908	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.09% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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](#)

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	0.54	0/3034	0.63	0/4098
1	B	0.55	0/2964	0.62	0/4003
1	C	0.51	0/2977	0.66	1/4019 (0.0%)
1	D	0.54	0/2987	0.62	0/4033
All	All	0.53	0/11962	0.63	1/16153 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	353	GLY	N-CA-C	12.67	144.77	113.10

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	2976	0	2961	47	0
1	B	2909	0	2906	40	0
1	C	2924	0	2924	42	0
1	D	2932	0	2928	40	0
2	A	50	0	0	2	0
2	B	34	0	0	0	0
2	C	36	0	0	0	0
2	D	47	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	11908	0	11719	155	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 7.

All (155) 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:355:SER:N	1:A:358:PHE:HD2	1.63	0.97
1:A:355:SER:N	1:A:358:PHE:CD2	2.38	0.92
1:A:146:ASP:HB3	1:A:147:PRO:HD3	1.55	0.86
1:A:290:GLN:HE22	1:D:289:ASN:HD21	1.26	0.83
1:B:377:HIS:CE1	1:B:379:ILE:HG22	2.15	0.82
1:A:355:SER:O	1:A:358:PHE:N	2.16	0.79
1:B:275:TYR:O	1:B:279:ARG:HG2	1.85	0.76
1:D:356:ASP:C	1:D:358:PHE:H	1.90	0.75
1:A:247:PHE:HZ	1:A:324:VAL:HG11	1.53	0.74
1:A:290:GLN:NE2	1:D:289:ASN:HD21	1.84	0.74
1:B:377:HIS:HE1	1:B:379:ILE:HG22	1.52	0.73
1:B:159:GLY:N	1:B:160:GLY:HA2	2.03	0.71
1:B:157:VAL:HB	1:B:158:PRO:HD2	1.70	0.71
1:B:140:GLU:OE2	1:B:150:MET:HB2	1.91	0.70
1:B:103:GLN:HE22	1:B:134:GLY:H	1.39	0.70
1:B:50:THR:HG21	1:B:99:MET:CE	2.22	0.70
1:D:43:GLU:OE2	1:D:46:ARG:NH1	2.25	0.69
1:D:216:LEU:H	1:D:366:ASN:HD22	1.38	0.68
1:B:279:ARG:HB2	1:B:286:LEU:HD22	1.75	0.67
1:A:290:GLN:HE22	1:D:289:ASN:ND2	1.92	0.67
1:A:353:GLY:C	1:A:355:SER:N	2.49	0.66
1:C:279:ARG:HB3	1:C:286:LEU:HD22	1.77	0.66
1:D:21:ASP:O	1:D:25:VAL:HG23	1.96	0.66
1:C:243:LEU:O	1:C:246:PRO:HD2	1.96	0.65
1:D:356:ASP:OD2	1:D:357:GLU:N	2.30	0.65
1:A:247:PHE:HZ	1:A:324:VAL:CG1	2.09	0.64
1:A:355:SER:O	1:A:357:GLU:N	2.30	0.64
1:A:355:SER:O	1:A:356:ASP:C	2.35	0.64
1:B:50:THR:HG21	1:B:99:MET:HE3	1.81	0.62
1:D:92:VAL:O	1:D:362:ARG:NH1	2.33	0.62
1:C:203:LYS:HE2	1:C:230:PHE:HB3	1.82	0.61
1:C:54:ILE:HG22	1:C:58:MET:CE	2.31	0.61
1:D:377:HIS:HD2	2:D:400:HOH:O	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:388:GLN:HE22	1:B:269:TRP:HE1	1.48	0.60
1:A:388:GLN:NE2	1:B:269:TRP:HE1	2.01	0.59
1:B:247:PHE:HZ	1:B:324:VAL:HG11	1.67	0.59
1:A:43:GLU:OE2	1:A:46:ARG:NH1	2.35	0.58
1:C:206:SER:HB2	1:C:225:VAL:HB	1.85	0.58
1:C:54:ILE:CG2	1:C:58:MET:HE2	2.33	0.58
1:D:216:LEU:H	1:D:366:ASN:ND2	2.02	0.58
1:A:13:LEU:HB3	1:A:16:GLN:HG3	1.86	0.58
1:D:99:MET:HE2	1:D:219:SER:HA	1.86	0.58
1:B:159:GLY:H	1:B:160:GLY:HA2	1.69	0.57
1:C:269:TRP:HE1	1:D:388:GLN:HE22	1.49	0.57
1:D:356:ASP:C	1:D:358:PHE:N	2.56	0.57
1:B:98:SER:O	1:B:102:VAL:HG23	2.05	0.56
1:D:7:HIS:CD2	1:D:13:LEU:HD21	2.41	0.56
1:A:362:ARG:NH1	2:A:420:HOH:O	2.34	0.56
1:D:393:ALA:O	1:D:394:PHE:HB2	2.06	0.56
1:C:353:GLY:C	1:C:354:ILE:HD12	2.27	0.55
1:C:54:ILE:HG22	1:C:58:MET:HE2	1.89	0.55
1:A:216:LEU:H	1:A:366:ASN:HD22	1.54	0.54
1:D:168:MET:HG3	1:D:169:TRP:CD1	2.43	0.54
1:A:216:LEU:H	1:A:366:ASN:ND2	2.05	0.54
1:B:40:ARG:HH21	1:B:54:ILE:HG12	1.72	0.54
1:C:388:GLN:HE22	1:D:269:TRP:HE1	1.55	0.54
1:A:187:GLU:O	1:A:188:ASP:C	2.45	0.54
1:A:257:ILE:CD1	1:A:374:GLU:HG3	2.38	0.53
1:B:45:PHE:O	1:B:217:ARG:NH2	2.41	0.53
1:C:354:ILE:O	1:C:354:ILE:CG2	2.57	0.53
1:D:206:SER:HB3	1:D:225:VAL:HB	1.90	0.53
1:A:363:HIS:HD2	2:A:399:HOH:O	1.92	0.52
1:D:81:SER:O	1:D:85:ILE:HG13	2.09	0.52
1:A:247:PHE:CZ	1:A:324:VAL:HG11	2.37	0.52
1:A:45:PHE:O	1:A:217:ARG:NH2	2.42	0.52
1:A:385:GLY:HA3	1:B:293:GLN:HB3	1.91	0.52
1:B:247:PHE:CZ	1:B:324:VAL:HG11	2.45	0.52
1:D:45:PHE:O	1:D:217:ARG:NH2	2.43	0.51
1:B:316:MET:HB3	1:B:322:ALA:HB2	1.92	0.51
1:A:140:GLU:OE1	1:A:150:MET:HG3	2.10	0.51
1:C:386:ARG:HH11	1:C:392:GLN:HE21	1.59	0.51
1:A:40:ARG:NH2	1:A:57:GLU:OE2	2.44	0.51
1:A:386:ARG:HG3	1:A:392:GLN:HA	1.92	0.51
1:B:216:LEU:H	1:B:366:ASN:HD22	1.59	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:103:GLN:OE1	1:D:134:GLY:N	2.29	0.51
1:D:140:GLU:HG2	1:D:166:SER:O	2.12	0.50
1:C:170:ILE:HG22	1:C:173:SER:HB3	1.93	0.50
1:B:149:SER:O	1:B:149:SER:OG	2.30	0.50
1:D:106:LEU:O	1:D:136:PHE:HD2	1.95	0.50
1:D:356:ASP:O	1:D:358:PHE:N	2.45	0.49
1:B:95:GLY:O	1:B:99:MET:HG3	2.12	0.49
1:C:354:ILE:HD12	1:C:354:ILE:N	2.28	0.49
1:D:392:GLN:OE1	1:D:394:PHE:HA	2.13	0.48
1:A:353:GLY:HA2	1:C:214:VAL:HG23	1.95	0.48
1:B:106:LEU:HB3	1:B:136:PHE:CG	2.49	0.48
1:C:114:PHE:CD1	1:C:240:VAL:HB	2.48	0.48
1:B:50:THR:HG21	1:B:99:MET:HE1	1.95	0.48
1:C:61:ILE:HG13	1:C:63:LEU:HG	1.94	0.48
1:A:140:GLU:HG2	1:A:167:LYS:HD3	1.96	0.48
1:B:64:LEU:HD22	1:B:103:GLN:HG2	1.96	0.48
1:C:168:MET:HG3	1:C:169:TRP:CD1	2.49	0.48
1:C:66:PRO:HA	1:C:75:PRO:HG2	1.95	0.48
1:C:177:ASP:HA	1:C:200:LYS:HB2	1.96	0.47
1:A:372:THR:HG22	1:A:373:TYR:CD1	2.50	0.47
1:B:126:LYS:HD2	1:B:132:TRP:CE2	2.50	0.47
1:A:160:GLY:HA2	1:A:233:GLU:HG2	1.97	0.47
1:D:356:ASP:C	1:D:356:ASP:OD2	2.53	0.47
1:C:153:ARG:NH1	1:C:191:ASP:OD1	2.47	0.47
1:C:216:LEU:H	1:C:366:ASN:HD22	1.61	0.47
1:C:104:SER:O	1:C:109:VAL:HG23	2.14	0.46
1:B:61:ILE:HG13	1:B:63:LEU:HG	1.96	0.46
1:D:295:LYS:O	1:D:299:MET:HG3	2.16	0.46
1:C:54:ILE:HG22	1:C:58:MET:HE3	1.96	0.46
1:B:216:LEU:H	1:B:366:ASN:ND2	2.14	0.46
1:A:146:ASP:HB3	1:A:147:PRO:CD	2.38	0.45
1:B:157:VAL:HB	1:B:158:PRO:CD	2.44	0.45
1:B:171:THR:O	1:B:172:ASN:HB2	2.17	0.45
1:C:170:ILE:CG2	1:C:173:SER:HB3	2.46	0.45
1:C:136:PHE:CE2	1:C:182:TRP:NE1	2.81	0.45
1:A:106:LEU:HB3	1:A:136:PHE:CD2	2.52	0.45
1:A:284:ARG:HG3	1:A:285:PRO:HD2	1.99	0.45
1:A:355:SER:C	1:A:357:GLU:N	2.69	0.45
1:C:354:ILE:O	1:C:354:ILE:HG22	2.17	0.44
1:B:54:ILE:O	1:B:58:MET:HG3	2.16	0.44
1:D:363:HIS:HD2	2:D:397:HOH:O	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:386:ARG:NH1	1:C:392:GLN:HE21	2.14	0.44
1:C:232:PRO:HB2	1:C:235:ASN:HD22	1.83	0.44
1:D:70:GLU:HA	1:D:74:GLY:O	2.18	0.44
1:D:13:LEU:HB3	1:D:16:GLN:HG3	2.00	0.43
1:A:351:GLY:O	1:A:352:ASN:HB2	2.18	0.43
1:D:392:GLN:HB3	1:D:392:GLN:HE21	1.67	0.43
1:A:392:GLN:HG2	1:A:394:PHE:HD2	1.84	0.43
1:C:92:VAL:O	1:C:362:ARG:NH1	2.50	0.43
1:C:68:ILE:O	1:C:74:GLY:HA3	2.18	0.43
1:C:106:LEU:HB3	1:C:136:PHE:CG	2.53	0.42
1:C:203:LYS:CE	1:C:230:PHE:HB3	2.48	0.42
1:A:110:PRO:HB3	1:A:249:CYS:SG	2.59	0.42
1:B:148:GLY:O	1:B:150:MET:N	2.53	0.42
1:A:386:ARG:NE	1:A:392:GLN:HB2	2.34	0.42
1:A:289:ASN:HD21	1:D:290:GLN:HE22	1.68	0.42
1:B:143:HIS:CE1	1:B:151:VAL:HG13	2.55	0.42
1:C:388:GLN:NE2	1:D:269:TRP:HE1	2.16	0.42
1:D:56:ARG:O	1:D:60:GLU:HG2	2.20	0.42
1:B:170:ILE:HG21	1:B:224:ILE:HD11	2.01	0.42
1:C:202:CYS:O	1:C:205:LEU:HB3	2.20	0.42
1:A:48:GLU:HB2	1:A:217:ARG:NH2	2.35	0.41
1:B:290:GLN:HE22	1:C:289:ASN:HB3	1.85	0.41
1:D:167:LYS:HB3	1:D:170:ILE:HD11	2.02	0.41
1:A:68:ILE:HA	1:A:69:PRO:HD3	1.94	0.41
1:A:106:LEU:HB3	1:A:136:PHE:CG	2.55	0.41
1:B:140:GLU:HG2	1:B:167:LYS:HD3	2.02	0.41
1:C:216:LEU:H	1:C:366:ASN:ND2	2.19	0.41
1:C:231:VAL:HG12	1:C:235:ASN:HB2	2.03	0.41
1:C:157:VAL:HB	1:C:158:PRO:HD2	2.03	0.41
1:B:289:ASN:HB3	1:C:290:GLN:HE22	1.84	0.41
1:D:132:TRP:HA	1:D:177:ASP:OD1	2.21	0.41
1:B:255:TYR:HD2	1:B:313:LEU:HD13	1.85	0.41
1:B:94:SER:HB2	1:B:366:ASN:HB3	2.03	0.41
1:C:126:LYS:HD2	1:C:132:TRP:CE2	2.55	0.41
1:A:352:ASN:HB3	1:C:169:TRP:CH2	2.57	0.40
1:A:372:THR:HG22	1:A:373:TYR:HD1	1.85	0.40
1:D:247:PHE:CZ	1:D:324:VAL:HG11	2.56	0.40
1:A:53:ALA:O	1:A:57:GLU:HG3	2.21	0.40
1:D:90:GLU:OE1	1:D:260:GLY:HA3	2.21	0.40
1:D:68:ILE:HA	1:D:69:PRO:HD3	1.95	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	381/396 (96%)	363 (95%)	17 (4%)	1 (0%)	41	55
1	B	370/396 (93%)	357 (96%)	11 (3%)	2 (0%)	29	41
1	C	375/396 (95%)	362 (96%)	13 (4%)	0	100	100
1	D	373/396 (94%)	357 (96%)	15 (4%)	1 (0%)	41	55
All	All	1499/1584 (95%)	1439 (96%)	56 (4%)	4 (0%)	41	55

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	149	SER
1	A	352	ASN
1	D	357	GLU
1	B	160	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	307/313 (98%)	291 (95%)	16 (5%)	23	38
1	B	301/313 (96%)	282 (94%)	19 (6%)	18	28
1	C	301/313 (96%)	283 (94%)	18 (6%)	19	31
1	D	302/313 (96%)	285 (94%)	17 (6%)	21	34
All	All	1211/1252 (97%)	1141 (94%)	70 (6%)	20	32

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	THR
1	A	100	MET
1	A	121	GLU
1	A	155	ARG
1	A	171	THR
1	A	206	SER
1	A	217	ARG
1	A	219	SER
1	A	237	LEU
1	A	243	LEU
1	A	284	ARG
1	A	313	LEU
1	A	372	THR
1	A	374	GLU
1	A	379	ILE
1	A	392	GLN
1	B	15	ASP
1	B	100	MET
1	B	101	SER
1	B	103	GLN
1	B	155	ARG
1	B	156	LYS
1	B	171	THR
1	B	186	ASP
1	B	234	GLU
1	B	237	LEU
1	B	277	LEU
1	B	284	ARG
1	B	286	LEU
1	B	291	LEU
1	B	313	LEU
1	B	344	LEU
1	B	377	HIS
1	B	386	ARG
1	B	392	GLN
1	C	71	GLN
1	C	100	MET
1	C	101	SER
1	C	113	GLU
1	C	121	GLU
1	C	166	SER
1	C	171	THR

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Mol	Chain	Res	Type
1	C	202	CYS
1	C	203	LYS
1	C	217	ARG
1	C	237	LEU
1	C	243	LEU
1	C	277	LEU
1	C	279	ARG
1	C	286	LEU
1	C	291	LEU
1	C	344	LEU
1	C	356	ASP
1	D	98	SER
1	D	100	MET
1	D	117	ASP
1	D	171	THR
1	D	173	SER
1	D	190	ARG
1	D	237	LEU
1	D	243	LEU
1	D	277	LEU
1	D	286	LEU
1	D	291	LEU
1	D	344	LEU
1	D	356	ASP
1	D	357	GLU
1	D	358	PHE
1	D	379	ILE
1	D	392	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	281	GLN
1	A	290	GLN
1	A	293	GLN
1	A	366	ASN
1	A	388	GLN
1	B	16	GLN
1	B	47	HIS
1	B	103	GLN
1	B	143	HIS
1	B	290	GLN

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Mol	Chain	Res	Type
1	B	293	GLN
1	B	366	ASN
1	B	377	HIS
1	B	388	GLN
1	C	235	ASN
1	C	290	GLN
1	C	366	ASN
1	C	388	GLN
1	C	392	GLN
1	D	7	HIS
1	D	281	GLN
1	D	293	GLN
1	D	366	ASN
1	D	388	GLN

### 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](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	387/396 (97%)	0.18	17 (4%) 34 33	41, 56, 78, 90	0
1	B	378/396 (95%)	0.29	31 (8%) 11 10	39, 61, 77, 90	0
1	C	381/396 (96%)	0.31	23 (6%) 21 20	40, 66, 87, 92	0
1	D	381/396 (96%)	0.23	17 (4%) 33 31	38, 60, 79, 90	0
All	All	1527/1584 (96%)	0.25	88 (5%) 23 22	38, 60, 81, 92	0

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	3	ALA	5.2
1	B	102	VAL	4.4
1	C	356	ASP	4.2
1	B	205	LEU	4.2
1	B	101	SER	3.8
1	B	377	HIS	3.7
1	A	357	GLU	3.6
1	A	358	PHE	3.6
1	B	143	HIS	3.6
1	D	282	PHE	3.5
1	A	282	PHE	3.5
1	C	230	PHE	3.5
1	B	180	VAL	3.4
1	C	169	TRP	3.4
1	C	151	VAL	3.3
1	C	158	PRO	3.3
1	B	376	THR	3.2
1	B	358	PHE	3.2
1	B	106	LEU	3.2
1	C	373	TYR	3.1
1	B	4	ALA	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	149	SER	3.1
1	B	253	ALA	3.1
1	D	158	PRO	3.1
1	B	107	VAL	3.0
1	D	279	ARG	3.0
1	D	141	PRO	2.9
1	B	229	ALA	2.9
1	C	355	SER	2.8
1	A	102	VAL	2.8
1	B	356	ASP	2.8
1	B	392	GLN	2.8
1	C	159	GLY	2.8
1	B	375	GLY	2.8
1	B	156	LYS	2.7
1	C	201	GLY	2.6
1	C	71	GLN	2.6
1	B	204	GLY	2.6
1	C	106	LEU	2.6
1	B	135	CYS	2.5
1	A	351	GLY	2.5
1	C	101	SER	2.5
1	A	281	GLN	2.5
1	A	356	ASP	2.5
1	B	211	HIS	2.5
1	A	306	GLY	2.5
1	C	279	ARG	2.5
1	D	358	PHE	2.5
1	D	5	THR	2.5
1	C	107	VAL	2.4
1	D	190	ARG	2.4
1	C	187	GLU	2.4
1	B	103	GLN	2.4
1	C	226	LEU	2.4
1	C	180	VAL	2.4
1	C	354	ILE	2.4
1	D	356	ASP	2.4
1	D	169	TRP	2.3
1	B	100	MET	2.3
1	A	307	LEU	2.2
1	D	283	GLY	2.2
1	D	230	PHE	2.2
1	B	141	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	4	ALA	2.2
1	D	212	GLY	2.2
1	B	279	ARG	2.2
1	C	307	LEU	2.2
1	C	357	GLU	2.2
1	D	357	GLU	2.2
1	A	261	ALA	2.2
1	A	262	LEU	2.2
1	A	386	ARG	2.2
1	D	168	MET	2.2
1	B	142	ASN	2.2
1	D	155	ARG	2.2
1	C	160	GLY	2.2
1	A	279	ARG	2.1
1	B	136	PHE	2.1
1	B	105	SER	2.1
1	B	171	THR	2.1
1	C	188	ASP	2.1
1	A	155	ARG	2.1
1	B	134	GLY	2.1
1	B	320	GLY	2.1
1	B	104	SER	2.1
1	C	249	CYS	2.0
1	A	355	SER	2.0
1	A	284	ARG	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.