



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:58 PM UTC

PDB ID : 3E49 / pdb_00003e49
Title : Crystal structure of a prokaryotic domain of unknown function (duf849) with a tim barrel fold (bx_e_c0966) from burkholderia xenovorans lb400 at 1.75 Å resolution
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2008-08-11
Resolution : 1.75 Å (reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

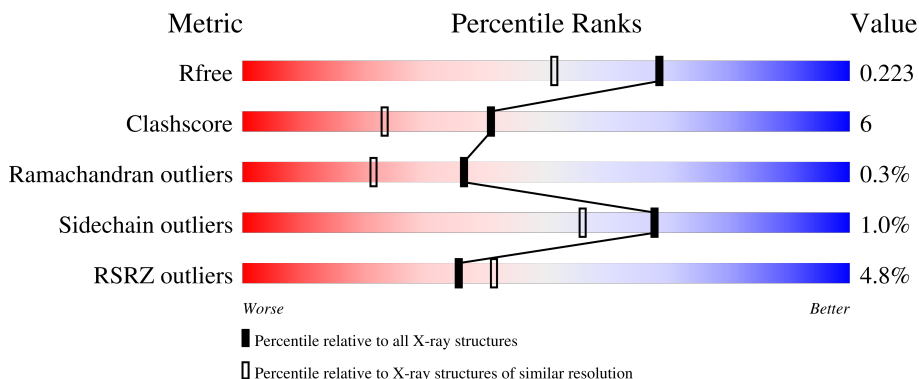
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

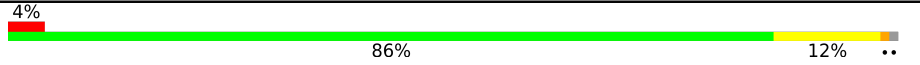



The reported resolution of this entry is 1.75 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3183 (1.76-1.76)
Clashscore	190562	3299 (1.76-1.76)
Ramachandran outliers	187476	3274 (1.76-1.76)
Sidechain outliers	187428	3274 (1.76-1.76)
RSRZ outliers	180081	3183 (1.76-1.76)

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	311	 4% 86% 12% ..
1	B	311	 8% 83% 15% ..
1	C	311	 3% 85% 12% ..
1	D	311	 4% 84% 13% ..

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
5	IMD	D	501	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 10552 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 uncharacterized protein DUF849 with a TIM barrel fold.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	307	2376	1510	423	431	3	9	0	6	0
1	B	307	2362	1503	416	431	3	9	0	6	0
1	C	307	2372	1503	423	434	3	9	0	7	0
1	D	307	2408	1534	425	437	3	9	0	10	0

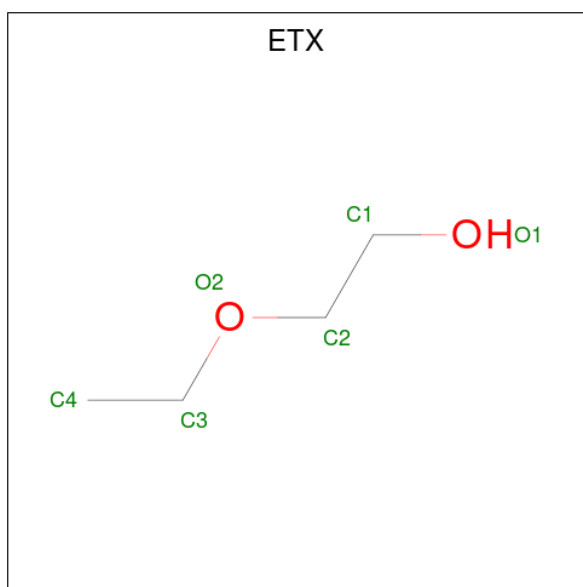
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP Q13GE9
B	0	GLY	-	expression tag	UNP Q13GE9
C	0	GLY	-	expression tag	UNP Q13GE9
D	0	GLY	-	expression tag	UNP Q13GE9

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Zn 1	0	0
2	B	1	Total 1	Zn 1	0	0
2	C	1	Total 1	Zn 1	0	0
2	D	1	Total 1	Zn 1	0	0

- Molecule 3 is 2-ETHOXYETHANOL (CCD ID: ETX) (formula: C₄H₁₀O₂).



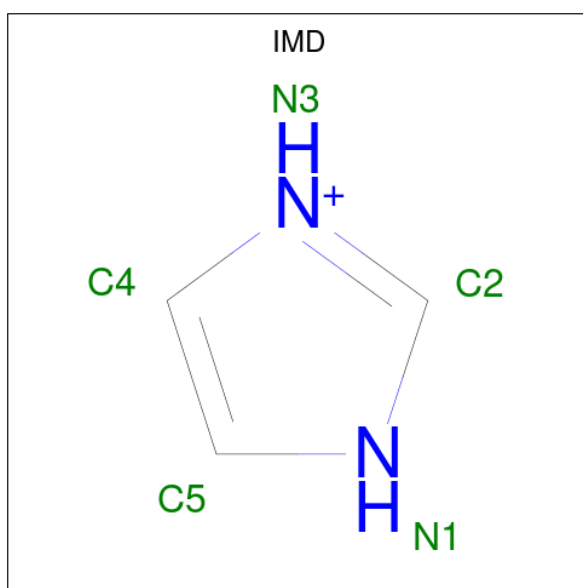
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 4 2	0	0
3	B	1	Total C O 6 4 2	0	0
3	B	1	Total C O 6 4 2	0	0
3	B	1	Total C O 6 4 2	0	0
3	C	1	Total C O 6 4 2	0	0
3	C	1	Total C O 6 4 2	0	0
3	C	1	Total C O 6 4 2	0	0

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is IMIDAZOLE (CCD ID: IMD) (formula: C₃H₅N₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	N	0	0
			5	3	2		
5	C	1	Total	C	N	0	0
			5	3	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	D	1	Total	C	N	0	0
			5	3	2		
5	D	1	Total	C	N	0	0
			5	3	2		
5	D	1	Total	C	N	0	0
			5	3	2		

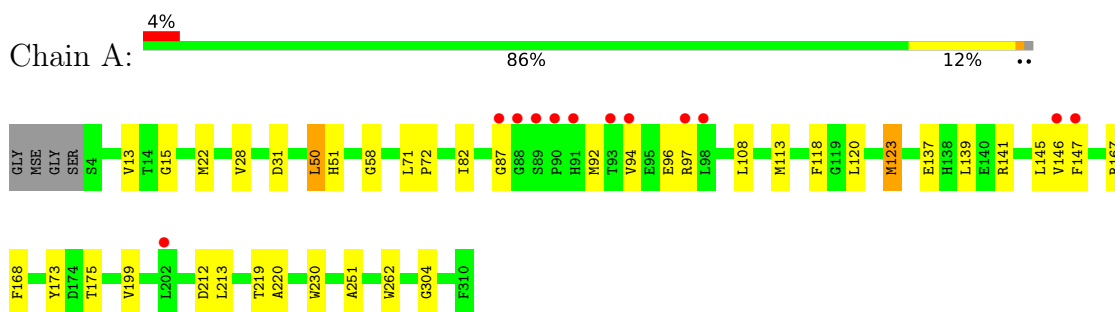
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	225	Total	O	0	0
			225	225		
6	B	195	Total	O	0	4
			199	199		
6	C	254	Total	O	0	3
			257	257		
6	D	272	Total	O	0	2
			274	274		

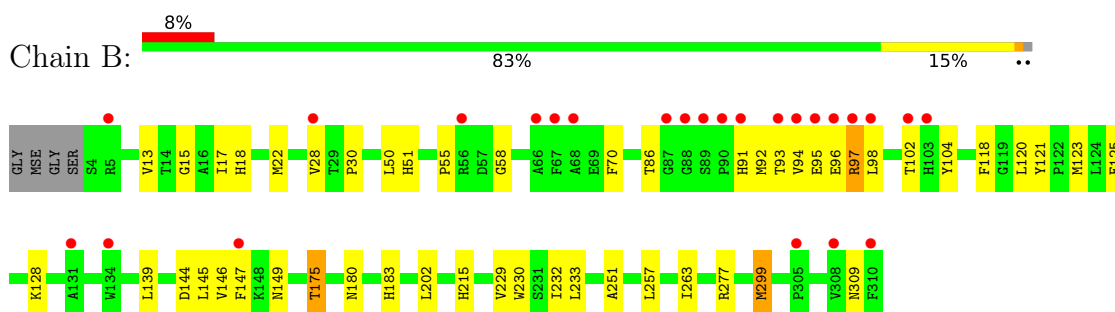
3 Residue-property plots [i](#)

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.

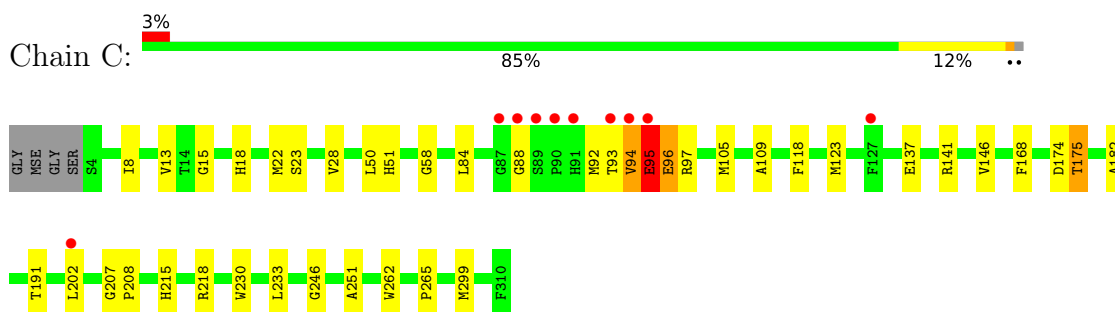
- Molecule 1: uncharacterized protein DUF849 with a TIM barrel fold



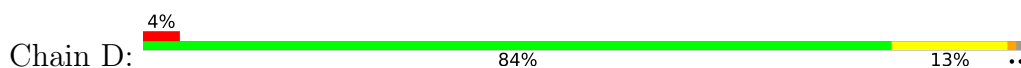
- Molecule 1: uncharacterized protein DUF849 with a TIM barrel fold

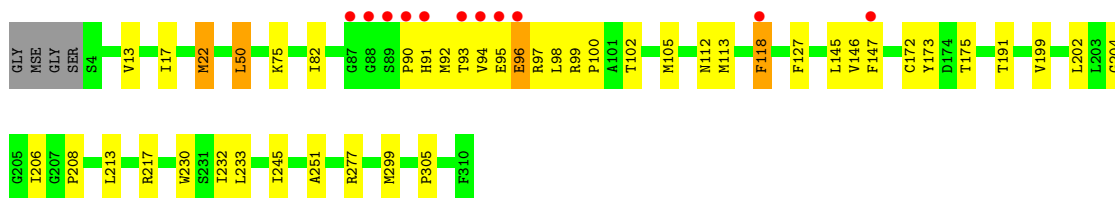


- Molecule 1: uncharacterized protein DUF849 with a TIM barrel fold



- Molecule 1: uncharacterized protein DUF849 with a TIM barrel fold





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	59.49Å 68.28Å 83.95Å 69.99° 88.07° 72.84°	Depositor
Resolution (Å)	29.66 – 1.75 29.66 – 1.75	Depositor EDS
% Data completeness (in resolution range)	97.2 (29.66-1.75) 97.2 (29.66-1.75)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.08 (at 1.75Å)	Xtrriage
Refinement program	REFMAC 5.2.0019, PHENIX	Depositor
R, R_{free}	0.177 , 0.218 0.184 , 0.223	Depositor DCC
R_{free} test set	5804 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	20.7	Xtrriage
Anisotropy	0.176	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 54.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10552	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.49% 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: ACT, ZN, ETX, IMD

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	0.96	5/2445 (0.2%)	1.02	10/3307 (0.3%)
1	B	0.97	2/2431 (0.1%)	1.04	8/3292 (0.2%)
1	C	1.09	7/2443 (0.3%)	1.07	8/3307 (0.2%)
1	D	1.06	6/2489 (0.2%)	1.07	5/3366 (0.1%)
All	All	1.02	20/9808 (0.2%)	1.05	31/13272 (0.2%)

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	299	MSE	SE-CE	-7.91	1.71	1.95
1	C	299	MSE	SE-CE	-7.61	1.72	1.95
1	C	22	MSE	SE-CE	-7.55	1.72	1.95
1	D	113	MSE	SE-CE	-6.84	1.75	1.95
1	A	22	MSE	SE-CE	-6.43	1.76	1.95
1	B	22	MSE	SE-CE	-6.31	1.76	1.95
1	C	94	VAL	N-CA	6.22	1.54	1.46
1	D	105	MSE	SE-CE	-6.05	1.77	1.95
1	A	113	MSE	SE-CE	-5.96	1.77	1.95
1	C	105	MSE	C-O	5.82	1.26	1.23
1	B	299	MSE	SE-CE	5.79	2.12	1.95
1	C	95	GLU	N-CA	5.72	1.53	1.46
1	C	174	ASP	N-CA	-5.72	1.39	1.46
1	D	206	ILE	CA-CB	-5.70	1.48	1.54
1	D	172	CYS	CA-CB	-5.64	1.45	1.53
1	A	219	THR	CA-CB	-5.34	1.44	1.53
1	A	123	MSE	SE-CE	-5.28	1.79	1.95
1	C	182	ALA	N-CA	5.21	1.52	1.46
1	D	22	MSE	SE-CE	-5.18	1.79	1.95
1	A	220	ALA	N-CA	5.04	1.52	1.46

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	94	VAL	N-CA-C	9.57	121.44	111.00
1	C	13	VAL	N-CA-C	8.53	121.75	111.09
1	C	88	GLY	N-CA-C	-7.38	100.25	111.24
1	B	70	PHE	N-CA-C	7.08	120.03	111.82
1	C	50	LEU	N-CA-C	7.07	120.42	108.90
1	D	13	VAL	N-CA-C	6.98	120.79	111.44
1	D	50	LEU	N-CA-C	6.78	120.44	109.40
1	D	96	GLU	N-CA-C	-6.46	103.52	111.40
1	A	113	MSE	N-CA-C	6.44	120.06	112.72
1	B	50	LEU	N-CA-C	6.28	119.70	109.59
1	A	13	VAL	N-CA-C	6.14	119.67	111.44
1	B	86	THR	N-CA-C	-6.13	105.09	113.30
1	B	96	GLU	N-CA-C	-6.10	104.26	111.03
1	B	13	VAL	N-CA-C	5.86	118.42	111.09
1	A	50	LEU	N-CA-C	5.68	118.66	109.40
1	C	23	SER	CA-C-N	-5.60	113.90	120.12
1	C	23	SER	C-N-CA	-5.60	113.90	120.12
1	B	121	TYR	CA-C-N	-5.55	113.90	119.56
1	B	121	TYR	C-N-CA	-5.55	113.90	119.56
1	D	245	ILE	N-CA-C	-5.54	104.97	110.62
1	A	219	THR	O-C-N	5.54	128.07	122.09
1	A	304	GLY	CA-C-N	5.43	125.10	119.56
1	A	304	GLY	C-N-CA	5.43	125.10	119.56
1	A	212	ASP	N-CA-C	-5.38	105.49	111.36
1	A	168	PHE	N-CA-C	5.33	118.26	109.72
1	C	8	ILE	CB-CA-C	-5.24	104.95	111.08
1	C	96	GLU	N-CA-C	-5.10	105.81	111.36
1	B	263	ILE	N-CA-C	-5.08	108.56	113.53
1	D	204	GLY	N-CA-C	-5.05	108.03	114.85
1	A	262	TRP	N-CA-C	5.04	117.96	109.95
1	C	168	PHE	N-CA-C	5.01	117.65	109.59

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2376	0	2337	19	0
1	B	2362	0	2311	39	0
1	C	2372	0	2320	26	0
1	D	2408	0	2386	36	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	6	0	10	0	0
3	B	18	0	30	3	0
3	C	18	0	30	2	0
4	B	4	0	3	1	0
4	C	4	0	3	1	0
5	B	5	0	5	0	0
5	C	5	0	5	1	0
5	D	15	0	15	5	0
6	A	225	0	0	2	0
6	B	199	0	0	5	0
6	C	257	0	0	0	0
6	D	274	0	0	6	0
All	All	10552	0	9455	123	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 6.

All (123) 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:92[B]:MSE:SE	1:A:96:GLU:CB	2.50	1.09
1:C:94:VAL:HG12	1:C:95:GLU:H	1.20	1.02
1:D:90:PRO:HB2	1:D:145[A]:LEU:HD21	1.49	0.92
1:B:123:MSE:HE3	1:B:139:LEU:HD22	1.54	0.89
1:C:94:VAL:HG12	1:C:95:GLU:N	1.89	0.86
1:C:95:GLU:CB	4:C:504:ACT:H3	2.06	0.84
1:D:92[B]:MSE:SE	1:D:96:GLU:CB	2.80	0.79
1:C:94:VAL:CG1	1:C:95:GLU:H	1.96	0.78
1:A:123:MSE:HE3	1:A:139:LEU:HD22	1.67	0.77
1:C:94:VAL:HG22	1:C:97:ARG:NH1	2.00	0.76
1:B:299:MSE:O	3:B:502:ETX:H42	1.91	0.71
1:B:98:LEU:O	1:B:102:THR:HG23	1.91	0.70
1:D:94:VAL:HG12	1:D:95:GLU:HG3	1.74	0.70
1:B:92[B]:MSE:HE2	1:B:93:THR:H	1.56	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:94:VAL:O	1:C:96:GLU:N	2.22	0.67
1:D:94:VAL:HG22	1:D:97[A]:ARG:NH1	2.09	0.67
1:A:31:ASP:CB	6:A:720:HOH:O	2.42	0.66
1:D:91:HIS:ND1	1:D:145[B]:LEU:HD12	2.11	0.66
1:B:149[A]:ASN:OD1	6:B:661:HOH:O	2.15	0.65
1:C:93:THR:O	1:C:97:ARG:HG3	1.96	0.65
5:D:501:IMD:C2	6:D:759:HOH:O	2.44	0.64
1:B:94:VAL:HG22	1:B:97:ARG:NH1	2.13	0.64
1:A:123:MSE:HE3	1:A:139:LEU:CD2	2.27	0.63
1:C:94:VAL:HG22	1:C:97:ARG:HH11	1.63	0.63
1:D:118:PHE:CE1	1:D:147:PHE:CD2	2.87	0.63
1:B:94:VAL:HG12	1:B:95:GLU:N	2.14	0.62
1:C:265:PRO:HA	3:C:503:ETX:H22	1.82	0.62
1:D:277[A]:ARG:HG2	6:D:616:HOH:O	1.97	0.62
1:B:120:LEU:HB3	1:B:123:MSE:CE	2.30	0.62
1:D:112:ASN:ND2	1:D:147:PHE:CE2	2.69	0.61
1:B:230:TRP:CE3	1:B:251:ALA:HB2	2.36	0.61
1:A:120:LEU:HB3	1:A:123:MSE:CE	2.32	0.59
1:D:118:PHE:CZ	1:D:202[B]:LEU:HD21	2.37	0.59
1:D:92[B]:MSE:HE2	1:D:93:THR:H	1.66	0.59
1:B:118:PHE:CE2	1:B:202[B]:LEU:HD21	2.38	0.59
1:D:91:HIS:ND1	1:D:145[B]:LEU:CD1	2.68	0.57
1:A:145:LEU:O	6:A:670:HOH:O	2.17	0.56
1:C:215[B]:HIS:HD2	1:C:218:ARG:HH11	1.52	0.56
1:D:118:PHE:CE1	1:D:147:PHE:HD2	2.22	0.56
1:A:137:GLU:CB	1:A:141:ARG:NH2	2.69	0.55
1:D:118:PHE:CE2	1:D:202[B]:LEU:HD21	2.42	0.55
1:C:118:PHE:CE2	1:C:202:LEU:HD21	2.42	0.55
1:B:28:VAL:HG13	1:B:58:GLY:HA2	1.89	0.54
1:A:87:GLY:O	1:A:97:ARG:HG2	2.08	0.54
1:A:146:VAL:HG11	1:B:146:VAL:HG11	1.90	0.54
1:B:118:PHE:CE1	1:B:147:PHE:HD2	2.27	0.53
1:B:118:PHE:CD2	1:B:202[B]:LEU:HD21	2.44	0.52
1:D:91:HIS:CE1	1:D:145[B]:LEU:HD12	2.45	0.52
1:D:94:VAL:HG22	1:D:97[A]:ARG:HH11	1.73	0.52
5:D:501:IMD:H2	6:D:759:HOH:O	2.06	0.51
1:B:232:ILE:CD1	1:B:251:ALA:HB1	2.41	0.51
1:D:94:VAL:HG12	1:D:95:GLU:N	2.26	0.51
1:D:94:VAL:C	1:D:96:GLU:H	2.19	0.51
1:A:50:LEU:HD12	1:A:82:ILE:HG23	1.93	0.50
5:D:502:IMD:H4	6:D:507:HOH:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:217:ARG:CZ	5:D:501:IMD:H5	2.42	0.50
1:D:230:TRP:CE2	1:D:232[A]:ILE:HD11	2.47	0.50
1:C:262:TRP:CE2	3:C:503:ETX:H42	2.47	0.50
1:C:146:VAL:HG11	1:D:146:VAL:HG11	1.92	0.49
1:B:118:PHE:CE1	1:B:147:PHE:CD2	3.00	0.49
1:D:232[A]:ILE:CD1	1:D:251:ALA:HB1	2.41	0.49
1:D:22:MSE:HG2	1:D:127:PHE:CG	2.48	0.49
1:C:230:TRP:CE3	1:C:251:ALA:HB2	2.48	0.49
1:B:202[B]:LEU:HD12	6:B:653:HOH:O	2.13	0.48
1:D:230:TRP:CZ2	1:D:232[B]:ILE:HD11	2.47	0.48
5:D:501:IMD:N3	6:D:759:HOH:O	2.35	0.48
1:D:99:ARG:HG3	6:D:756:HOH:O	2.12	0.48
1:B:94:VAL:CG1	1:B:95:GLU:N	2.76	0.48
1:A:118:PHE:CE1	1:A:147:PHE:CD2	3.02	0.47
1:C:175:THR:OG1	1:C:215[A]:HIS:HE1	1.97	0.47
1:B:232:ILE:CD1	1:B:251:ALA:CB	2.93	0.47
1:B:94:VAL:HG12	1:B:95:GLU:H	1.78	0.47
4:B:504:ACT:H1	6:B:609:HOH:O	2.14	0.47
1:B:95:GLU:HA	1:B:98:LEU:HD12	1.97	0.47
1:B:229:VAL:CG2	3:B:502:ETX:H43	2.45	0.47
1:A:230:TRP:CE3	1:A:251:ALA:HB2	2.50	0.46
1:C:18:HIS:HD2	1:C:123:MSE:HE1	1.80	0.46
1:A:213:LEU:C	1:A:213:LEU:HD23	2.41	0.46
1:C:94:VAL:HA	1:C:97:ARG:HD3	1.96	0.46
1:D:99:ARG:HB2	1:D:100:PRO:HD3	1.98	0.46
1:C:92[B]:MSE:HG3	1:C:93:THR:N	2.31	0.46
1:A:15:GLY:HA2	1:A:51:HIS:CD2	2.51	0.46
1:C:93:THR:O	1:C:97:ARG:CG	2.64	0.46
1:D:213:LEU:C	1:D:213:LEU:HD23	2.41	0.46
1:B:15:GLY:HA2	1:B:51:HIS:CD2	2.51	0.46
1:D:17:ILE:HD11	1:D:118:PHE:HE2	1.81	0.46
1:B:202[A]:LEU:HD11	6:B:657:HOH:O	2.16	0.45
1:D:98:LEU:O	1:D:102:THR:HG23	2.16	0.45
1:B:229:VAL:HG22	3:B:502:ETX:H11	1.97	0.45
1:D:95:GLU:HA	1:D:98:LEU:HB2	1.97	0.45
1:B:91:HIS:ND1	1:B:145:LEU:HD12	2.31	0.45
1:B:277:ARG:HG2	6:B:686:HOH:O	2.17	0.45
1:A:28:VAL:HG13	1:A:58:GLY:HA2	1.98	0.45
1:B:94:VAL:CG1	1:B:95:GLU:H	2.30	0.45
1:D:75:LYS:HE3	1:D:305:PRO:O	2.17	0.45
1:D:90:PRO:O	1:D:145[A]:LEU:HD11	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:91:HIS:ND1	1:B:145:LEU:CD1	2.80	0.44
1:C:207:GLY:HA2	5:C:505:IMD:H2	1.99	0.44
1:B:94:VAL:HG22	1:B:97:ARG:HH11	1.79	0.44
1:B:104:TYR:HA	1:B:309:ASN:HB3	1.98	0.44
1:B:125:GLU:O	1:B:128:LYS:NZ	2.49	0.44
1:D:50:LEU:HD12	1:D:82:ILE:HG23	1.99	0.44
1:B:175:THR:CG2	1:B:215[B]:HIS:NE2	2.81	0.44
1:A:173:TYR:CE2	1:A:199:VAL:HG11	2.53	0.43
1:B:17:ILE:HG23	1:B:120:LEU:HD22	2.00	0.43
1:B:18:HIS:ND1	1:B:257:LEU:O	2.51	0.43
1:A:71:LEU:HB2	1:A:72:PRO:HD3	2.00	0.43
1:C:246:GLY:C	1:C:251:ALA:HB3	2.43	0.43
1:D:230:TRP:CZ2	1:D:232[A]:ILE:HD11	2.54	0.43
1:A:108:LEU:HA	1:A:167:ARG:O	2.20	0.42
1:C:137:GLU:HG3	1:C:141:ARG:HG2	2.01	0.42
1:D:232[A]:ILE:CD1	1:D:251:ALA:CB	2.98	0.42
1:C:15:GLY:HA2	1:C:51:HIS:CD2	2.55	0.42
1:C:28:VAL:HG13	1:C:58:GLY:HA2	2.02	0.42
1:C:215[B]:HIS:CD2	1:C:218:ARG:HH11	2.35	0.41
1:C:84:LEU:O	1:C:109:ALA:HA	2.21	0.41
1:B:30:PRO:HD3	1:B:55:PRO:HG3	2.03	0.41
1:B:180:ASN:O	1:B:183:HIS:HB3	2.20	0.41
1:D:173:TYR:CE2	1:D:199:VAL:HG11	2.56	0.41
1:A:118:PHE:HE1	1:A:147:PHE:HD2	1.69	0.41
1:B:123:MSE:HE3	1:B:139:LEU:CD2	2.37	0.40
1:D:22:MSE:HG2	1:D:127:PHE:CD1	2.56	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	311/311 (100%)	304 (98%)	7 (2%)	0	100	100
1	B	311/311 (100%)	299 (96%)	11 (4%)	1 (0%)	36	21
1	C	312/311 (100%)	300 (96%)	10 (3%)	2 (1%)	21	8
1	D	315/311 (101%)	305 (97%)	9 (3%)	1 (0%)	36	21
All	All	1249/1244 (100%)	1208 (97%)	37 (3%)	4 (0%)	36	21

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	95	GLU
1	B	97	ARG
1	D	208	PRO
1	C	208	PRO

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	245/239 (102%)	244 (100%)	1 (0%)	84	80
1	B	243/239 (102%)	241 (99%)	2 (1%)	73	64
1	C	244/239 (102%)	241 (99%)	3 (1%)	63	49
1	D	251/239 (105%)	247 (98%)	4 (2%)	55	38
All	All	983/956 (103%)	973 (99%)	10 (1%)	68	56

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	175	THR
1	B	175	THR
1	B	233	LEU
1	C	175	THR
1	C	191	THR
1	C	233	LEU
1	D	118	PHE

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Mol	Chain	Res	Type
1	D	175	THR
1	D	191	THR
1	D	233	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	197	GLN
1	B	164	ASN
1	C	18	HIS
1	C	91	HIS
1	D	62	GLN
1	D	238	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 4 are monoatomic - leaving 14 for Mogul analysis.

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	ETX	C	503	-	5,5,5	0.76	0	4,4,4	0.20	0
4	ACT	B	504	-	3,3,3	0.93	0	3,3,3	1.44	0
5	IMD	D	502	-	5,5,5	0.68	0	5,5,5	0.35	0
3	ETX	B	502	-	5,5,5	0.48	0	4,4,4	0.67	0
5	IMD	D	503	-	5,5,5	0.58	0	5,5,5	0.58	0
3	ETX	B	501	-	5,5,5	0.70	0	4,4,4	0.17	0
5	IMD	C	505	-	5,5,5	0.56	0	5,5,5	0.57	0
5	IMD	B	505	-	5,5,5	0.70	0	5,5,5	0.57	0
4	ACT	C	504	-	3,3,3	0.86	0	3,3,3	1.51	0
3	ETX	C	501	-	5,5,5	0.54	0	4,4,4	0.51	0
3	ETX	B	503	-	5,5,5	0.80	0	4,4,4	0.64	0
5	IMD	D	501	-	5,5,5	0.67	0	5,5,5	0.45	0
3	ETX	A	501	-	5,5,5	0.65	0	4,4,4	0.62	0
3	ETX	C	502	-	5,5,5	0.65	0	4,4,4	0.62	0

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	ETX	C	503	-	-	1/3/3/3	-
5	IMD	D	502	-	-	-	0/1/1/1
3	ETX	B	502	-	-	3/3/3/3	-
5	IMD	D	503	-	-	-	0/1/1/1
3	ETX	B	501	-	-	2/3/3/3	-
5	IMD	C	505	-	-	-	0/1/1/1
5	IMD	B	505	-	-	-	0/1/1/1
5	IMD	D	501	-	-	-	0/1/1/1
3	ETX	C	501	-	-	2/3/3/3	-
3	ETX	B	503	-	-	1/3/3/3	-
3	ETX	A	501	-	-	2/3/3/3	-
3	ETX	C	502	-	-	2/3/3/3	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	501	ETX	O1-C1-C2-O2
3	B	503	ETX	O1-C1-C2-O2
3	B	502	ETX	O1-C1-C2-O2
3	C	501	ETX	O1-C1-C2-O2
3	C	502	ETX	O1-C1-C2-O2
3	B	502	ETX	C4-C3-O2-C2
3	C	502	ETX	C4-C3-O2-C2
3	B	501	ETX	C4-C3-O2-C2
3	C	501	ETX	C4-C3-O2-C2
3	B	502	ETX	C1-C2-O2-C3
3	C	503	ETX	C4-C3-O2-C2
3	B	501	ETX	C1-C2-O2-C3
3	A	501	ETX	C1-C2-O2-C3

There are no ring outliers.

7 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	503	ETX	2	0
4	B	504	ACT	1	0
5	D	502	IMD	1	0
3	B	502	ETX	3	0
5	C	505	IMD	1	0
4	C	504	ACT	1	0
5	D	501	IMD	4	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	299/311 (96%)	0.39	12 (4%) 42 48	8, 30, 51, 63	5 (1%)
1	B	299/311 (96%)	0.66	25 (8%) 17 20	7, 33, 51, 65	5 (1%)
1	C	299/311 (96%)	0.17	10 (3%) 49 55	12, 25, 46, 63	6 (2%)
1	D	299/311 (96%)	0.11	11 (3%) 45 51	11, 24, 42, 62	9 (3%)
All	All	1196/1244 (96%)	0.33	58 (4%) 35 41	7, 28, 48, 65	25 (2%)

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	90	PRO	5.7
1	A	94	VAL	5.6
1	A	90	PRO	5.2
1	C	94	VAL	5.0
1	B	95	GLU	4.9
1	D	90	PRO	4.9
1	D	94	VAL	4.9
1	B	147	PHE	4.8
1	B	90	PRO	4.8
1	A	147	PHE	4.7
1	B	94	VAL	4.7
1	C	93	THR	4.6
1	B	98	LEU	4.4
1	A	88	GLY	4.2
1	C	88	GLY	3.9
1	B	93	THR	3.8
1	A	98	LEU	3.6
1	A	91	HIS	3.5
1	D	96	GLU	3.5
1	D	95	GLU	3.4
1	B	97	ARG	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	93	THR	3.3
1	C	91	HIS	3.3
1	B	88	GLY	3.1
1	C	89	SER	3.1
1	B	91	HIS	2.9
1	B	68	ALA	2.8
1	C	95	GLU	2.8
1	D	88	GLY	2.8
1	B	134	TRP	2.7
1	D	91	HIS	2.6
1	B	87	GLY	2.6
1	B	28	VAL	2.6
1	B	66	ALA	2.6
1	A	89	SER	2.5
1	D	147	PHE	2.5
1	B	310	PHE	2.5
1	B	103	HIS	2.5
1	B	5	ARG	2.5
1	A	146	VAL	2.4
1	B	89	SER	2.4
1	B	305	PRO	2.4
1	B	67	PHE	2.4
1	B	308	VAL	2.3
1	C	87	GLY	2.3
1	C	202	LEU	2.2
1	B	96	GLU	2.2
1	D	118	PHE	2.2
1	A	97	ARG	2.2
1	A	202	LEU	2.2
1	C	127	PHE	2.2
1	D	93	THR	2.2
1	B	131	ALA	2.1
1	A	87	GLY	2.1
1	D	89	SER	2.1
1	B	56	ARG	2.1
1	B	102	THR	2.1
1	D	87	GLY	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 [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	ACT	B	504	4/4	0.58	0.17	59,60,61,62	0
5	IMD	D	503	5/5	0.58	0.19	58,58,61,62	0
3	ETX	B	503	6/6	0.66	0.28	53,56,61,64	0
4	ACT	C	504	4/4	0.67	0.16	59,59,60,60	0
3	ETX	C	502	6/6	0.67	0.21	62,67,68,69	0
3	ETX	C	503	6/6	0.76	0.21	52,55,58,62	0
3	ETX	B	501	6/6	0.77	0.18	51,56,58,58	0
5	IMD	D	501	5/5	0.78	0.17	57,58,58,60	0
5	IMD	C	505	5/5	0.78	0.19	50,52,53,57	0
3	ETX	C	501	6/6	0.79	0.18	48,52,54,63	0
3	ETX	A	501	6/6	0.81	0.16	37,47,51,53	0
5	IMD	B	505	5/5	0.85	0.14	49,51,56,58	0
5	IMD	D	502	5/5	0.86	0.15	47,50,54,58	0
3	ETX	B	502	6/6	0.94	0.11	29,42,43,52	0
2	ZN	A	500	1/1	0.98	0.05	32,32,32,32	0
2	ZN	D	500	1/1	0.99	0.04	27,27,27,27	0
2	ZN	B	500	1/1	0.99	0.08	33,33,33,33	0
2	ZN	C	500	1/1	0.99	0.03	29,29,29,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.