



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:56 PM UTC

PDB ID : 1U24 / pdb_00001u24
Title : Crystal structure of Selenomonas ruminantium phytase
Authors : Chu, H.M.; Guo, R.T.; Lin, T.W.; Chou, C.C.; Shr, H.L.; Lai, H.L.; Tang, T.Y.; Cheng, K.J.; Selinger, B.L.; Wang, A.H.-J.
Deposited on : 2004-07-16
Resolution : 2.00 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

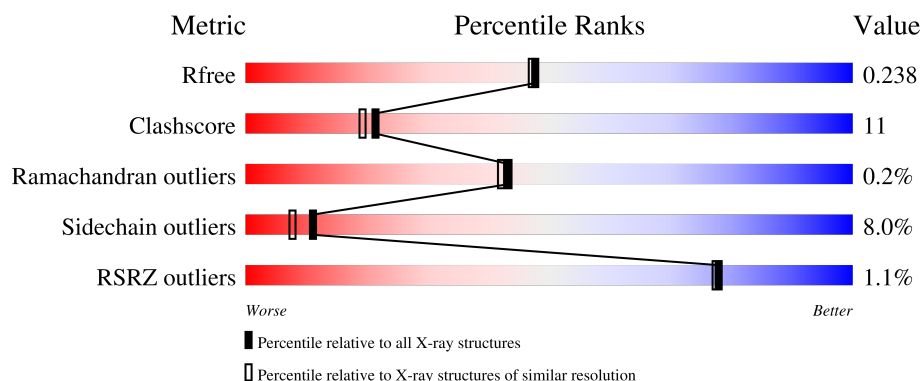
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	337	70% 19% • 7%
1	B	337	73% 18% • 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6380 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 myo-inositol hexaphosphate phosphohydrolase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	314	Total	C	N	O	S	Se	0	0	0
			2556	1629	457	463	1	6			
1	B	314	Total	C	N	O	S	Se	0	0	0
			2556	1629	457	463	1	6			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	cloning artifact	UNP Q7WUJ1
A	2	ALA	-	cloning artifact	UNP Q7WUJ1
A	3	SER	-	cloning artifact	UNP Q7WUJ1
A	4	MSE	-	cloning artifact	UNP Q7WUJ1
A	5	THR	-	cloning artifact	UNP Q7WUJ1
A	6	GLY	-	cloning artifact	UNP Q7WUJ1
A	7	GLY	-	cloning artifact	UNP Q7WUJ1
A	8	GLN	-	cloning artifact	UNP Q7WUJ1
A	9	GLN	-	cloning artifact	UNP Q7WUJ1
A	10	MSE	-	cloning artifact	UNP Q7WUJ1
A	11	GLY	-	cloning artifact	UNP Q7WUJ1
A	12	ARG	-	cloning artifact	UNP Q7WUJ1
A	13	GLY	-	cloning artifact	UNP Q7WUJ1
A	14	SER	-	cloning artifact	UNP Q7WUJ1
A	15	GLU	-	cloning artifact	UNP Q7WUJ1
A	16	PHE	-	cloning artifact	UNP Q7WUJ1
A	86	MSE	MET	modified residue	UNP Q7WUJ1
A	203	MSE	MET	modified residue	UNP Q7WUJ1
A	252	MSE	MET	modified residue	UNP Q7WUJ1
A	254	MSE	MET	modified residue	UNP Q7WUJ1
A	257	MSE	MET	modified residue	UNP Q7WUJ1
A	304	MSE	MET	modified residue	UNP Q7WUJ1
A	336	LEU	-	cloning artifact	UNP Q7WUJ1
A	337	GLU	-	cloning artifact	UNP Q7WUJ1
B	1	MSE	-	cloning artifact	UNP Q7WUJ1

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	2	ALA	-	cloning artifact	UNP Q7WUJ1
B	3	SER	-	cloning artifact	UNP Q7WUJ1
B	4	MSE	-	cloning artifact	UNP Q7WUJ1
B	5	THR	-	cloning artifact	UNP Q7WUJ1
B	6	GLY	-	cloning artifact	UNP Q7WUJ1
B	7	GLY	-	cloning artifact	UNP Q7WUJ1
B	8	GLN	-	cloning artifact	UNP Q7WUJ1
B	9	GLN	-	cloning artifact	UNP Q7WUJ1
B	10	MSE	-	cloning artifact	UNP Q7WUJ1
B	11	GLY	-	cloning artifact	UNP Q7WUJ1
B	12	ARG	-	cloning artifact	UNP Q7WUJ1
B	13	GLY	-	cloning artifact	UNP Q7WUJ1
B	14	SER	-	cloning artifact	UNP Q7WUJ1
B	15	GLU	-	cloning artifact	UNP Q7WUJ1
B	16	PHE	-	cloning artifact	UNP Q7WUJ1
B	86	MSE	MET	modified residue	UNP Q7WUJ1
B	203	MSE	MET	modified residue	UNP Q7WUJ1
B	252	MSE	MET	modified residue	UNP Q7WUJ1
B	254	MSE	MET	modified residue	UNP Q7WUJ1
B	257	MSE	MET	modified residue	UNP Q7WUJ1
B	304	MSE	MET	modified residue	UNP Q7WUJ1
B	336	LEU	-	cloning artifact	UNP Q7WUJ1
B	337	GLU	-	cloning artifact	UNP Q7WUJ1

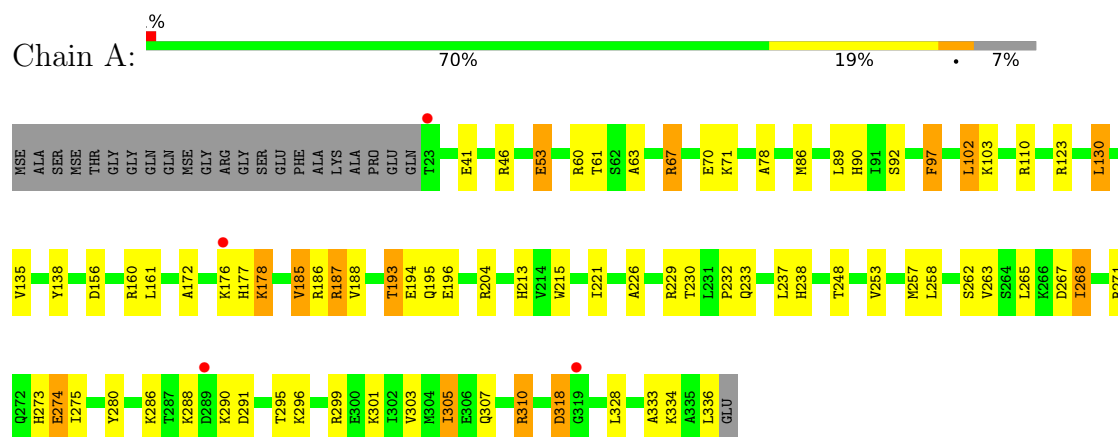
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	643	Total O 643 643	0	0
2	B	625	Total O 625 625	0	0

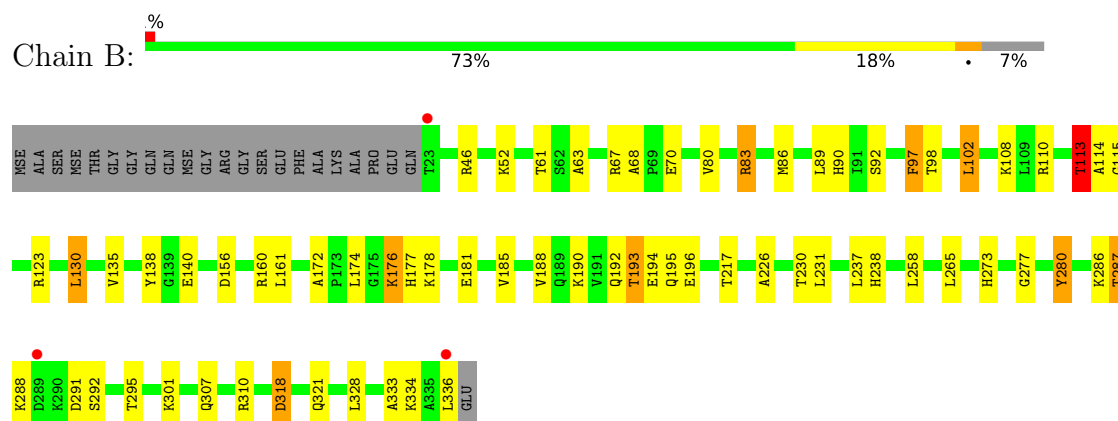
3 Residue-property plots [i](#)

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: myo-inositol hexaphosphate phosphohydrolase



- Molecule 1: myo-inositol hexaphosphate phosphohydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	46.47Å 140.38Å 80.68Å 90.00° 102.85° 90.00°	Depositor
Resolution (Å)	15.00 – 2.00 15.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (15.00-2.00) 90.7 (15.00-2.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.180 , 0.221 0.199 , 0.238	Depositor DCC
R_{free} test set	5166 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å ²)	16.0	Xtriage
Anisotropy	0.450	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 66.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6380	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.76% 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

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	0.82	1/2624 (0.0%)	1.09	17/3545 (0.5%)
1	B	0.88	0/2624	1.13	14/3545 (0.4%)
All	All	0.85	1/5248 (0.0%)	1.11	31/7090 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	185	VAL	CA-CB	5.18	1.61	1.54

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	318	ASP	N-CA-C	-8.82	98.18	110.35
1	B	188	VAL	N-CA-C	7.93	119.21	109.30
1	A	188	VAL	N-CA-C	7.25	118.46	108.89
1	B	231	LEU	N-CA-C	6.62	119.22	110.08
1	A	187	ARG	N-CA-C	-6.52	98.61	109.24
1	A	295	THR	N-CA-C	6.51	118.96	111.02
1	B	295	THR	N-CA-C	6.43	118.86	111.02
1	B	86	MSE	N-CA-C	6.37	118.22	111.28
1	B	138	TYR	N-CA-C	6.29	119.90	109.46
1	B	113	THR	CB-CA-C	-6.17	98.88	110.16
1	B	98	THR	N-CA-C	-6.14	103.07	110.13
1	A	86	MSE	N-CA-C	5.92	117.81	111.36
1	B	172	ALA	N-CA-C	5.87	118.73	108.82
1	A	213	HIS	N-CA-C	5.86	122.63	113.72
1	A	123	ARG	N-CA-C	5.80	118.19	108.96
1	B	123	ARG	N-CA-C	5.78	118.15	108.96
1	A	310	ARG	N-CA-C	-5.74	105.11	111.36
1	A	221	ILE	N-CA-C	5.69	116.47	110.72
1	A	138	TYR	N-CA-C	5.59	118.35	109.24
1	A	275	ILE	N-CA-C	-5.55	107.33	112.83
1	B	280	TYR	N-CA-C	5.54	119.13	112.93

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	217	THR	CA-C-N	-5.46	113.27	119.28
1	B	217	THR	C-N-CA	-5.46	113.27	119.28
1	A	172	ALA	N-CA-C	5.40	116.98	108.23
1	A	274	GLU	N-CA-C	5.37	119.45	113.01
1	A	204	ARG	N-CA-C	-5.35	103.33	110.55
1	A	232	PRO	N-CA-C	-5.29	102.27	111.32
1	A	97	PHE	N-CA-C	5.08	116.92	109.24
1	A	71	LYS	N-CA-C	5.04	116.78	111.28
1	B	97	PHE	N-CA-C	5.02	116.81	109.24
1	A	177	HIS	N-CA-C	-5.01	106.59	113.56

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	2556	0	2480	56	0
1	B	2556	0	2480	55	0
2	A	643	0	0	11	0
2	B	625	0	0	12	0
All	All	6380	0	4960	110	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 (110) 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:61:THR:HG22	1:A:63:ALA:H	1.21	1.05
1:B:61:THR:HG22	1:B:63:ALA:H	1.21	1.03
1:A:60:ARG:HH22	1:A:248:THR:HG23	1.33	0.93
1:B:113:THR:HG22	1:B:115:GLY:H	1.35	0.91
1:B:83:ARG:HG2	1:B:83:ARG:HH11	1.37	0.87
1:B:52:LYS:HE3	2:B:1168:HOH:O	1.76	0.85

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:THR:HG22	1:A:196:GLU:H	1.49	0.77
1:B:67:ARG:NH1	1:B:68:ALA:H	1.82	0.76
1:A:67:ARG:HH11	1:A:67:ARG:HB3	1.51	0.74
1:A:60:ARG:HH22	1:A:248:THR:CG2	1.99	0.73
1:A:307:GLN:HG3	1:A:333:ALA:HB2	1.71	0.71
1:B:193:THR:HG22	1:B:196:GLU:H	1.55	0.71
1:A:307:GLN:NE2	1:A:310:ARG:HH11	1.89	0.71
1:B:83:ARG:HG2	1:B:83:ARG:NH1	2.08	0.69
1:B:287:THR:HG21	1:B:292:SER:HA	1.75	0.68
1:A:299:ARG:HD3	2:A:1227:HOH:O	1.93	0.67
1:B:83:ARG:HH11	1:B:83:ARG:CG	2.08	0.66
1:A:90:HIS:HD2	2:A:1272:HOH:O	1.78	0.65
1:A:60:ARG:NH2	1:A:248:THR:HG23	2.08	0.65
1:B:130:LEU:HD22	1:B:135:VAL:HG21	1.78	0.64
1:A:288:LYS:H	1:A:288:LYS:CE	2.11	0.64
1:A:248:THR:HG21	2:A:1268:HOH:O	1.97	0.64
1:A:70:GLU:H	1:A:273:HIS:HE1	1.47	0.63
1:A:307:GLN:NE2	1:A:310:ARG:NH1	2.46	0.63
1:B:176:LYS:N	1:B:176:LYS:HE2	2.15	0.62
1:B:61:THR:HG22	1:B:63:ALA:N	2.05	0.61
1:B:328:LEU:HD13	1:B:328:LEU:O	2.01	0.61
1:A:280:TYR:HE2	1:A:301:LYS:HD3	1.64	0.61
1:A:61:THR:HG22	1:A:63:ALA:N	2.04	0.61
1:B:110:ARG:NH2	1:B:115:GLY:O	2.30	0.61
1:B:46:ARG:HD3	1:B:174:LEU:HG	1.83	0.60
1:B:83:ARG:NH1	1:B:83:ARG:CG	2.63	0.59
1:A:288:LYS:H	1:A:288:LYS:HE2	1.68	0.59
1:B:288:LYS:HG3	2:B:1297:HOH:O	2.03	0.58
1:B:92:SER:OG	1:B:238:HIS:HD2	1.85	0.58
1:B:70:GLU:H	1:B:273:HIS:HE1	1.52	0.57
1:A:67:ARG:HH11	1:A:67:ARG:CB	2.18	0.57
1:A:262:SER:HB2	2:A:705:HOH:O	2.03	0.57
1:A:67:ARG:HG2	1:A:274:GLU:O	2.04	0.57
1:B:307:GLN:HG3	1:B:333:ALA:HB2	1.87	0.56
1:A:226:ALA:O	1:A:230:THR:HG23	2.05	0.56
1:B:156:ASP:O	1:B:160:ARG:HG3	2.07	0.55
1:A:336:LEU:HD13	2:A:965:HOH:O	2.06	0.55
1:A:92:SER:OG	1:A:238:HIS:HD2	1.90	0.54
1:B:177:HIS:N	2:B:625:HOH:O	2.40	0.54
1:A:215:TRP:CZ3	1:A:305:ILE:HG22	2.43	0.53
1:B:113:THR:CG2	1:B:115:GLY:H	2.15	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:PHE:CE1	1:A:102:LEU:HG	2.43	0.53
1:B:193:THR:HG23	1:B:195:GLN:H	1.74	0.53
1:A:307:GLN:CG	1:A:333:ALA:HB2	2.39	0.53
1:A:288:LYS:N	1:A:288:LYS:HD3	2.25	0.52
1:A:263:VAL:CG2	1:A:268:ILE:HD11	2.39	0.52
1:B:176:LYS:HE2	1:B:176:LYS:CA	2.40	0.52
1:A:53:GLU:HG3	2:A:1658:HOH:O	2.10	0.51
1:A:267:ASP:O	1:A:271:ARG:HG2	2.11	0.51
1:B:287:THR:HG22	1:B:288:LYS:O	2.11	0.50
1:B:113:THR:HG22	1:B:115:GLY:N	2.16	0.50
1:B:328:LEU:HD13	1:B:328:LEU:C	2.37	0.50
1:B:334:LYS:N	1:B:334:LYS:HD2	2.26	0.50
1:B:156:ASP:OD1	1:B:160:ARG:HD2	2.12	0.50
1:A:307:GLN:HE21	1:A:310:ARG:NH1	2.08	0.50
1:B:113:THR:HG22	1:B:114:ALA:N	2.27	0.50
1:A:229:ARG:HD2	2:A:1413:HOH:O	2.13	0.49
1:A:156:ASP:O	1:A:160:ARG:HG3	2.13	0.49
1:A:288:LYS:N	1:A:288:LYS:CD	2.76	0.49
1:A:193:THR:CG2	1:A:196:GLU:H	2.22	0.49
1:B:307:GLN:NE2	1:B:310:ARG:HD2	2.27	0.49
1:B:113:THR:HG21	2:B:701:HOH:O	2.13	0.49
1:B:193:THR:HG22	1:B:196:GLU:HG3	1.93	0.49
1:B:90:HIS:HD2	2:B:637:HOH:O	1.95	0.49
1:B:226:ALA:O	1:B:230:THR:HG23	2.13	0.49
1:A:233:GLN:HB2	2:A:969:HOH:O	2.13	0.48
1:B:280:TYR:HE2	1:B:301:LYS:HD2	1.78	0.48
1:A:303:VAL:HG23	1:A:336:LEU:HG	1.96	0.48
1:A:193:THR:HG21	2:A:1624:HOH:O	2.14	0.48
1:B:273:HIS:HD2	1:B:277:GLY:O	1.97	0.48
1:B:97:PHE:CE1	1:B:102:LEU:HG	2.50	0.47
1:B:181:GLU:HG3	2:B:750:HOH:O	2.14	0.47
1:A:67:ARG:HB3	1:A:67:ARG:NH1	2.26	0.46
1:A:288:LYS:H	1:A:288:LYS:CD	2.27	0.46
1:A:70:GLU:H	1:A:273:HIS:CE1	2.28	0.46
1:B:70:GLU:H	1:B:273:HIS:CE1	2.33	0.46
1:B:336:LEU:HD23	2:B:1301:HOH:O	2.16	0.46
1:A:178:LYS:N	1:A:178:LYS:HD2	2.31	0.45
1:A:336:LEU:CD1	2:A:965:HOH:O	2.64	0.45
1:B:287:THR:HG21	1:B:291:ASP:C	2.41	0.45
1:B:334:LYS:HD3	2:B:1017:HOH:O	2.15	0.45
1:B:176:LYS:HE2	1:B:176:LYS:HA	1.99	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:113:THR:CG2	2:B:701:HOH:O	2.65	0.45
1:A:286:LYS:NZ	2:A:1739:HOH:O	2.51	0.44
1:A:288:LYS:CE	1:A:291:ASP:OD2	2.66	0.44
1:B:176:LYS:HG2	1:B:181:GLU:OE1	2.17	0.44
1:A:318:ASP:OD1	1:A:318:ASP:N	2.51	0.44
1:A:253:VAL:CG1	1:A:257:MSE:HE3	2.48	0.44
1:B:318:ASP:CG	1:B:321:GLN:HG2	2.44	0.43
1:A:288:LYS:HE3	1:A:291:ASP:OD2	2.19	0.43
1:A:263:VAL:HG23	1:A:268:ILE:HG12	1.99	0.43
1:B:192:GLN:OE1	2:B:1019:HOH:O	2.21	0.43
1:A:102:LEU:HD23	1:A:102:LEU:HA	1.89	0.42
1:B:92:SER:OG	1:B:238:HIS:CD2	2.70	0.42
1:B:307:GLN:HE21	1:B:310:ARG:NH1	2.17	0.42
1:A:130:LEU:HD22	1:A:135:VAL:HG21	2.02	0.42
1:A:187:ARG:HB3	1:A:187:ARG:NH1	2.35	0.42
1:B:286:LYS:HD2	1:B:286:LYS:HA	1.91	0.42
1:A:193:THR:HG23	1:A:195:GLN:N	2.35	0.41
1:A:78:ALA:HB1	1:B:286:LYS:HB2	2.03	0.41
1:A:178:LYS:N	1:A:178:LYS:CD	2.84	0.41
1:B:328:LEU:CD1	2:B:1283:HOH:O	2.69	0.41
1:A:176:LYS:HD3	1:A:176:LYS:HA	1.94	0.41
1:B:178:LYS:HE3	2:B:935:HOH:O	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	312/337 (93%)	301 (96%)	11 (4%)	0	100	100
1	B	312/337 (93%)	303 (97%)	8 (3%)	1 (0%)	36	35
All	All	624/674 (93%)	604 (97%)	19 (3%)	1 (0%)	43	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	176	LYS

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	262/269 (97%)	237 (90%)	25 (10%)	8	5
1	B	262/269 (97%)	245 (94%)	17 (6%)	15	12
All	All	524/538 (97%)	482 (92%)	42 (8%)	11	8

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	GLU
1	A	46	ARG
1	A	53	GLU
1	A	67	ARG
1	A	89	LEU
1	A	102	LEU
1	A	103	LYS
1	A	110	ARG
1	A	130	LEU
1	A	161	LEU
1	A	178	LYS
1	A	185	VAL
1	A	186	ARG
1	A	193	THR
1	A	194	GLU
1	A	237	LEU
1	A	258	LEU
1	A	265	LEU
1	A	268	ILE
1	A	290	LYS
1	A	296	LYS
1	A	305	ILE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	318	ASP
1	A	328	LEU
1	A	334	LYS
1	B	80	VAL
1	B	83	ARG
1	B	89	LEU
1	B	102	LEU
1	B	108	LYS
1	B	113	THR
1	B	130	LEU
1	B	140	GLU
1	B	161	LEU
1	B	185	VAL
1	B	190	LYS
1	B	193	THR
1	B	194	GLU
1	B	237	LEU
1	B	258	LEU
1	B	265	LEU
1	B	287	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	90	HIS
1	A	162	HIS
1	A	166	HIS
1	A	189	GLN
1	A	192	GLN
1	A	233	GLN
1	A	238	HIS
1	A	260	ASN
1	A	272	GLN
1	A	273	HIS
1	A	307	GLN
1	A	315	ASN
1	B	90	HIS
1	B	159	HIS
1	B	162	HIS
1	B	166	HIS
1	B	192	GLN
1	B	220	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	238	HIS
1	B	240	HIS
1	B	260	ASN
1	B	272	GLN
1	B	273	HIS
1	B	307	GLN
1	B	313	GLN
1	B	315	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 oligosaccharides 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 [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	308/337 (91%)	-0.39	4 (1%) 75 74	8, 19, 39, 48	0
1	B	308/337 (91%)	-0.48	3 (0%) 79 79	8, 17, 32, 46	0
All	All	616/674 (91%)	-0.44	7 (1%) 78 77	8, 18, 37, 48	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	319	GLY	3.7
1	B	336	LEU	2.9
1	B	23	THR	2.7
1	A	176	LYS	2.3
1	A	289	ASP	2.2
1	B	289	ASP	2.2
1	A	23	THR	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 oligosaccharides 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.