



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 31, 2021 – 02:25 PM EDT

PDB ID : 3M0X
Title : Crystal structure of Pseudomonas stutzeri L-rhamnose isomerase mutant S329L in complex with D-psicose
Authors : Yoshida, H.; Takeda, K.; Izumori, K.; Kamitori, S.
Deposited on : 2010-03-03
Resolution : 1.79 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.23.2
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) : 2.23.2

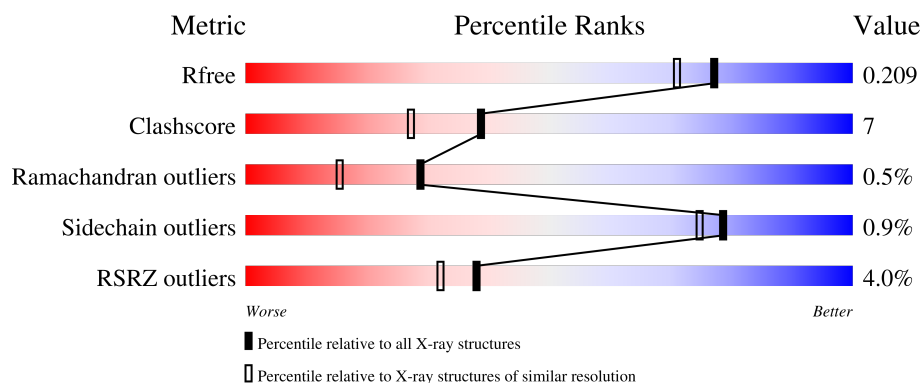
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.79 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	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	438	<div> <div>3%</div> <div>83%</div> <div>13%</div> <div>.</div> </div>
1	B	438	<div> <div>3%</div> <div>84%</div> <div>12%</div> <div>.</div> </div>
1	C	438	<div> <div>7%</div> <div>79%</div> <div>18%</div> <div>..</div> </div>
1	D	438	<div> <div>3%</div> <div>84%</div> <div>11%</div> <div>.</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14571 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 L-rhamnose isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	421	Total	C	N	O	S	0	0	0
			3262	2051	584	618	9			
1	B	421	Total	C	N	O	S	0	0	0
			3262	2051	584	618	9			
1	C	430	Total	C	N	O	S	0	0	0
			3313	2081	593	630	9			
1	D	419	Total	C	N	O	S	0	0	0
			3253	2045	582	617	9			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	150	ASN	ASP	engineered mutation	UNP Q75WH8
A	329	LEU	SER	engineered mutation	UNP Q75WH8
A	431	GLY	-	expression tag	UNP Q75WH8
A	432	SER	-	expression tag	UNP Q75WH8
A	433	HIS	-	expression tag	UNP Q75WH8
A	434	HIS	-	expression tag	UNP Q75WH8
A	435	HIS	-	expression tag	UNP Q75WH8
A	436	HIS	-	expression tag	UNP Q75WH8
A	437	HIS	-	expression tag	UNP Q75WH8
A	438	HIS	-	expression tag	UNP Q75WH8
B	150	ASN	ASP	engineered mutation	UNP Q75WH8
B	329	LEU	SER	engineered mutation	UNP Q75WH8
B	431	GLY	-	expression tag	UNP Q75WH8
B	432	SER	-	expression tag	UNP Q75WH8
B	433	HIS	-	expression tag	UNP Q75WH8
B	434	HIS	-	expression tag	UNP Q75WH8
B	435	HIS	-	expression tag	UNP Q75WH8
B	436	HIS	-	expression tag	UNP Q75WH8
B	437	HIS	-	expression tag	UNP Q75WH8
B	438	HIS	-	expression tag	UNP Q75WH8
C	150	ASN	ASP	engineered mutation	UNP Q75WH8

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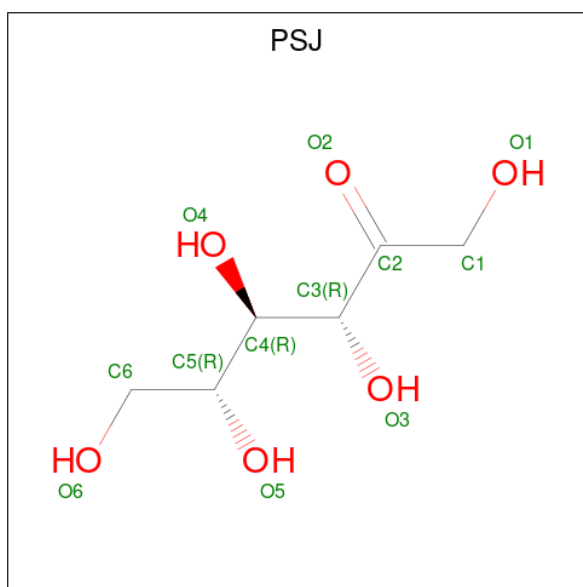
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Chain	Residue	Modelled	Actual	Comment	Reference
C	329	LEU	SER	engineered mutation	UNP Q75WH8
C	431	GLY	-	expression tag	UNP Q75WH8
C	432	SER	-	expression tag	UNP Q75WH8
C	433	HIS	-	expression tag	UNP Q75WH8
C	434	HIS	-	expression tag	UNP Q75WH8
C	435	HIS	-	expression tag	UNP Q75WH8
C	436	HIS	-	expression tag	UNP Q75WH8
C	437	HIS	-	expression tag	UNP Q75WH8
C	438	HIS	-	expression tag	UNP Q75WH8
D	150	ASN	ASP	engineered mutation	UNP Q75WH8
D	329	LEU	SER	engineered mutation	UNP Q75WH8
D	431	GLY	-	expression tag	UNP Q75WH8
D	432	SER	-	expression tag	UNP Q75WH8
D	433	HIS	-	expression tag	UNP Q75WH8
D	434	HIS	-	expression tag	UNP Q75WH8
D	435	HIS	-	expression tag	UNP Q75WH8
D	436	HIS	-	expression tag	UNP Q75WH8
D	437	HIS	-	expression tag	UNP Q75WH8
D	438	HIS	-	expression tag	UNP Q75WH8

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

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

- Molecule 3 is D-psicose (three-letter code: PSJ) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		
3	C	1	Total	C	O	0	0
			12	6	6		
3	D	1	Total	C	O	0	0
			12	6	6		

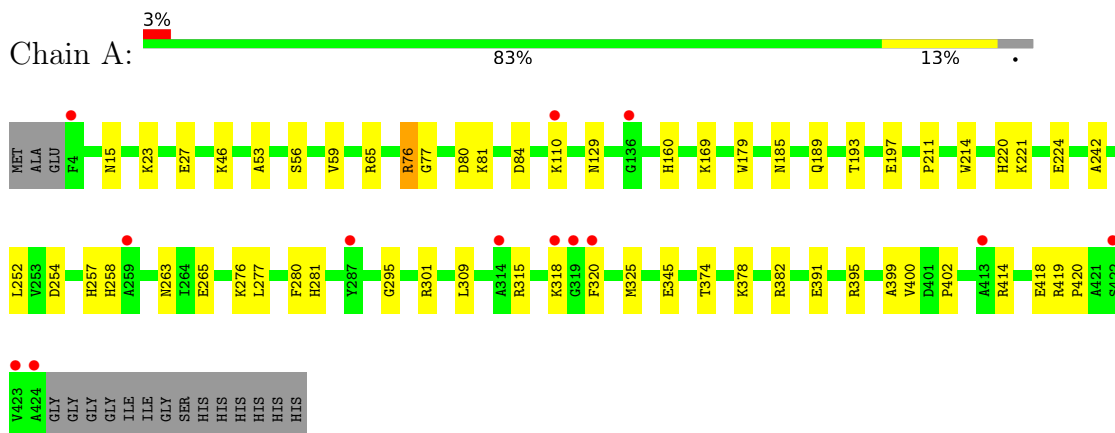
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	367	Total	O	0	0
			367	367		
4	B	361	Total	O	0	0
			361	361		
4	C	320	Total	O	0	0
			320	320		
4	D	377	Total	O	0	0
			377	377		

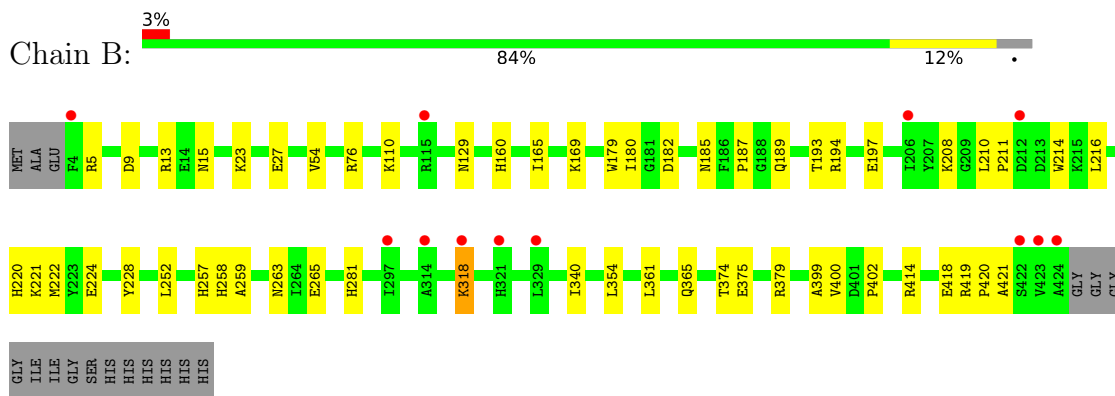
3 Residue-property plots [i](#)

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.

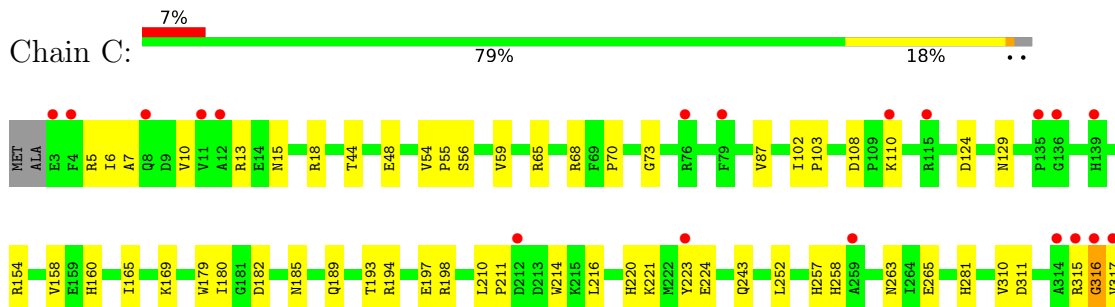
- Molecule 1: L-rhamnose isomerase



- Molecule 1: L-rhamnose isomerase

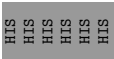
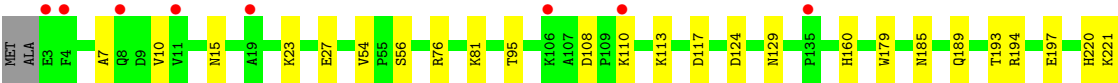
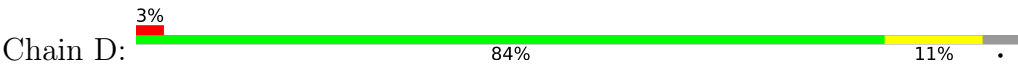


- Molecule 1: L-rhamnose isomerase





● Molecule 1: L-rhamnose isomerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.71Å 104.53Å 115.83Å 90.00° 112.03° 90.00°	Depositor
Resolution (Å)	47.76 – 1.79 47.76 – 1.79	Depositor EDS
% Data completeness (in resolution range)	91.8 (47.76-1.79) 91.9 (47.76-1.79)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.42 (at 1.79Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.196 , 0.218 0.187 , 0.209	Depositor DCC
R_{free} test set	14123 reflections (9.86%)	wwPDB-VP
Wilson B-factor (Å ²)	14.9	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 49.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	14571	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.30% 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: PSJ, MN

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	0.31	0/3334	0.56	0/4521
1	B	0.30	0/3334	0.55	0/4521
1	C	0.30	0/3385	0.54	0/4588
1	D	0.31	0/3325	0.56	0/4508
All	All	0.31	0/13378	0.55	0/18138

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3262	0	3171	44	0
1	B	3262	0	3171	42	0
1	C	3313	0	3219	71	0
1	D	3253	0	3158	35	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	12	0	10	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	12	0	10	3	0
3	C	12	0	10	2	0
3	D	12	0	10	3	0
4	A	367	0	0	5	0
4	B	361	0	0	10	0
4	C	320	0	0	4	0
4	D	377	0	0	2	0
All	All	14571	0	12759	194	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 (194) 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:129:ASN:H	1:A:160:HIS:HE1	1.05	0.98
1:A:56:SER:O	1:A:59:VAL:HG12	1.74	0.88
1:B:129:ASN:H	1:B:160:HIS:HE1	1.17	0.88
1:A:129:ASN:H	1:A:160:HIS:CE1	1.90	0.87
1:C:129:ASN:H	1:C:160:HIS:HE1	1.24	0.83
1:C:6:ILE:HD11	1:C:87:VAL:HG13	1.59	0.83
1:C:68:ARG:HH12	1:C:70:PRO:HB3	1.46	0.79
1:C:129:ASN:H	1:C:160:HIS:CE1	2.02	0.77
1:A:76:ARG:HH11	1:A:419:ARG:HG2	1.50	0.76
1:B:129:ASN:H	1:B:160:HIS:CE1	2.04	0.75
1:D:129:ASN:H	1:D:160:HIS:HE1	1.32	0.75
1:C:220:HIS:HE2	1:C:258:HIS:CE1	2.05	0.74
1:B:220:HIS:HE2	1:B:258:HIS:CE1	2.05	0.73
1:D:185:ASN:H	1:D:189:GLN:HE22	1.37	0.72
1:B:23:LYS:O	1:B:27:GLU:HG3	1.91	0.71
1:A:129:ASN:N	1:A:160:HIS:HE1	1.85	0.70
1:D:129:ASN:H	1:D:160:HIS:CE1	2.09	0.70
1:D:220:HIS:HE2	1:D:258:HIS:CE1	2.10	0.69
1:B:220:HIS:HE2	1:B:258:HIS:HE1	1.39	0.69
1:C:220:HIS:HE2	1:C:258:HIS:HE1	1.41	0.69
1:C:68:ARG:NH1	1:C:70:PRO:HB3	2.09	0.68
1:C:185:ASN:H	1:C:189:GLN:HE22	1.40	0.68
1:A:15:ASN:ND2	1:A:400:VAL:H	1.91	0.68
1:D:185:ASN:H	1:D:189:GLN:NE2	1.93	0.66
1:B:414:ARG:O	1:B:418:GLU:HG3	1.96	0.65
1:A:110:LYS:HD2	1:A:110:LYS:H	1.62	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:110:LYS:HG3	4:B:1084:HOH:O	1.97	0.65
1:C:185:ASN:H	1:C:189:GLN:NE2	1.95	0.64
1:B:15:ASN:ND2	1:B:400:VAL:H	1.97	0.63
1:C:6:ILE:HD11	1:C:87:VAL:HG22	1.80	0.63
1:B:361:LEU:O	1:B:365:GLN:HG3	1.99	0.62
1:C:210:LEU:HD11	1:C:216:LEU:HB2	1.81	0.62
1:C:44:THR:O	1:C:48:GLU:HG3	1.99	0.62
1:D:23:LYS:O	1:D:27:GLU:HG3	1.99	0.62
1:C:6:ILE:HD12	1:C:6:ILE:N	2.15	0.62
1:D:220:HIS:HE2	1:D:258:HIS:HE1	1.48	0.62
1:A:15:ASN:HD21	1:A:400:VAL:H	1.46	0.61
1:B:185:ASN:H	1:B:189:GLN:HE22	1.49	0.61
1:C:165:ILE:O	1:C:169:LYS:HG3	2.00	0.61
1:A:185:ASN:H	1:A:189:GLN:HE22	1.46	0.61
4:A:1210:HOH:O	1:C:243:GLN:HG2	2.02	0.60
1:C:429:ILE:HG23	1:C:430:ILE:HD12	1.83	0.60
1:C:15:ASN:ND2	1:C:400:VAL:H	2.00	0.60
1:A:220:HIS:HE2	1:A:258:HIS:CE1	2.20	0.59
1:B:15:ASN:HD21	1:B:400:VAL:H	1.52	0.57
1:B:13:ARG:HD2	4:B:1038:HOH:O	2.04	0.57
1:B:185:ASN:H	1:B:189:GLN:NE2	2.02	0.57
1:C:15:ASN:HD21	1:C:399:ALA:HA	1.69	0.57
1:D:402:PRO:HG2	4:D:563:HOH:O	2.05	0.57
1:C:340:ILE:HG21	1:C:402:PRO:HB2	1.87	0.56
1:D:378:LYS:HB3	1:D:382:ARG:NH2	2.21	0.56
1:C:179:TRP:CE2	3:C:603:PSJ:H1A	2.41	0.56
1:B:15:ASN:HD21	1:B:399:ALA:HA	1.71	0.56
1:B:165:ILE:O	1:B:169:LYS:HG3	2.06	0.55
1:B:375:GLU:O	1:B:379:ARG:HG3	2.06	0.55
1:C:18:ARG:HB3	4:C:544:HOH:O	2.06	0.55
1:C:311:ASP:O	1:C:315:ARG:HG2	2.07	0.55
1:B:402:PRO:HG2	4:B:454:HOH:O	2.06	0.54
1:B:179:TRP:CE2	3:B:602:PSJ:H1A	2.42	0.54
3:C:603:PSJ:O5	3:C:603:PSJ:C2	2.56	0.54
1:C:7:ALA:HB3	1:C:10:VAL:HG23	1.88	0.54
1:C:318:LYS:O	1:C:318:LYS:HG3	2.09	0.53
1:D:179:TRP:CE2	3:D:604:PSJ:H1A	2.43	0.53
1:C:430:ILE:H	1:C:430:ILE:HD13	1.73	0.53
1:B:54:VAL:HG11	4:B:966:HOH:O	2.08	0.53
1:D:15:ASN:ND2	1:D:400:VAL:H	2.07	0.53
1:A:185:ASN:H	1:A:189:GLN:NE2	2.06	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:108:ASP:OD2	1:C:110:LYS:HB2	2.10	0.52
1:D:221:LYS:HA	1:D:257:HIS:HB3	1.90	0.52
1:A:263:ASN:OD1	1:A:265:GLU:HG2	2.10	0.52
1:C:5:ARG:HB2	1:C:6:ILE:HD12	1.91	0.52
1:A:76:ARG:HD3	1:A:80:ASP:OD2	2.10	0.51
1:C:6:ILE:CD1	1:C:87:VAL:HG22	2.41	0.51
1:C:429:ILE:HG23	1:C:430:ILE:CD1	2.40	0.51
1:C:15:ASN:HD21	1:C:400:VAL:H	1.58	0.51
1:A:301:ARG:HD3	4:A:473:HOH:O	2.10	0.51
1:A:242:ALA:HB1	1:A:276:LYS:HD2	1.93	0.51
1:B:263:ASN:OD1	1:B:265:GLU:HG2	2.11	0.51
1:C:211:PRO:HD2	1:C:214:TRP:CG	2.46	0.50
1:B:210:LEU:HD11	1:B:216:LEU:HB2	1.92	0.50
1:A:59:VAL:HG21	1:A:84:ASP:HB2	1.94	0.50
1:A:277:LEU:HD23	1:A:320:PHE:CE1	2.47	0.50
1:A:402:PRO:HG2	4:A:485:HOH:O	2.11	0.50
1:D:263:ASN:OD1	1:D:265:GLU:HG2	2.12	0.50
1:C:318:LYS:O	1:C:320:PHE:N	2.45	0.49
1:A:221:LYS:HA	1:A:257:HIS:HB3	1.94	0.49
1:D:15:ASN:HD21	1:D:399:ALA:HA	1.76	0.49
1:C:6:ILE:HD11	1:C:87:VAL:CG1	2.36	0.49
1:B:76:ARG:HH11	1:B:76:ARG:HG2	1.78	0.48
1:C:180:ILE:HD12	1:C:182:ASP:CG	2.33	0.48
1:C:367:ASP:HB3	4:C:1097:HOH:O	2.12	0.48
1:A:23:LYS:O	1:A:27:GLU:HG3	2.13	0.48
1:C:318:LYS:C	1:C:320:PHE:N	2.65	0.48
1:C:73:GLY:CA	1:C:416:ALA:HA	2.44	0.48
3:A:601:PSJ:O5	3:A:601:PSJ:C2	2.62	0.48
1:C:48:GLU:HG2	1:C:392:ALA:HB1	1.96	0.48
1:D:193:THR:O	1:D:197:GLU:HG3	2.13	0.48
1:A:15:ASN:HD21	1:A:399:ALA:HA	1.79	0.47
1:A:169:LYS:HD3	1:A:211:PRO:CG	2.44	0.47
1:A:277:LEU:HD23	1:A:320:PHE:HE1	1.79	0.47
1:C:263:ASN:OD1	1:C:265:GLU:HG2	2.13	0.47
1:A:220:HIS:HE2	1:A:258:HIS:HE1	1.59	0.47
1:B:318:LYS:O	1:B:318:LYS:HD3	2.13	0.47
1:C:315:ARG:O	1:C:316:GLY:C	2.53	0.47
1:C:361:LEU:O	1:C:365:GLN:HG3	2.15	0.47
1:D:194:ARG:HG2	1:D:194:ARG:HH11	1.79	0.47
3:D:604:PSJ:O5	3:D:604:PSJ:C2	2.63	0.47
1:D:108:ASP:OD2	1:D:110:LYS:HB2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:242:ALA:HB1	1:D:276:LYS:HD2	1.96	0.46
1:B:76:ARG:HH22	1:B:421:ALA:HA	1.81	0.46
1:A:77:GLY:O	1:A:81:LYS:HG3	2.15	0.46
1:A:211:PRO:HD2	1:A:214:TRP:CG	2.50	0.46
1:C:315:ARG:O	1:C:315:ARG:HG3	2.16	0.46
1:D:56:SER:O	1:D:81:LYS:HD3	2.16	0.46
1:C:221:LYS:HA	1:C:257:HIS:HB3	1.98	0.46
1:B:211:PRO:HD2	1:B:214:TRP:CG	2.51	0.46
1:C:430:ILE:HD13	1:C:430:ILE:N	2.30	0.46
1:B:340:ILE:HG21	1:B:402:PRO:HB2	1.97	0.45
1:A:76:ARG:HD2	1:A:420:PRO:HD2	1.97	0.45
1:B:169:LYS:HD3	1:B:211:PRO:HG2	1.99	0.45
1:D:252:LEU:HD21	1:D:281:HIS:CG	2.52	0.45
1:C:310:VAL:HG21	1:C:353:ALA:CB	2.46	0.45
1:C:340:ILE:CG2	1:C:402:PRO:HB2	2.47	0.45
3:B:602:PSJ:O5	3:B:602:PSJ:C2	2.65	0.45
4:B:590:HOH:O	1:C:382:ARG:HD3	2.16	0.45
1:D:113:LYS:NZ	1:D:117:ASP:OD2	2.49	0.45
1:B:208:LYS:HG2	4:B:835:HOH:O	2.16	0.45
1:C:367:ASP:O	1:C:368:ASN:HB2	2.17	0.44
1:D:15:ASN:HD21	1:D:400:VAL:H	1.64	0.44
1:C:317:VAL:CG1	1:C:318:LYS:N	2.79	0.44
1:A:391:GLU:O	1:A:395:ARG:HG3	2.17	0.44
1:B:221:LYS:HG3	1:B:257:HIS:CG	2.52	0.44
1:C:318:LYS:O	1:C:319:GLY:C	2.54	0.44
1:A:59:VAL:HG11	1:A:81:LYS:HB3	1.98	0.44
1:C:310:VAL:HG21	1:C:353:ALA:HB3	2.00	0.44
1:C:375:GLU:O	1:C:379:ARG:HG3	2.17	0.44
1:A:280:PHE:CE2	1:A:309:LEU:HD21	2.53	0.44
1:B:221:LYS:HA	1:B:257:HIS:HB3	1.99	0.43
1:C:193:THR:O	1:C:197:GLU:HG3	2.18	0.43
1:B:194:ARG:NH2	4:B:707:HOH:O	2.50	0.43
1:C:13:ARG:HD2	4:C:915:HOH:O	2.16	0.43
1:A:193:THR:O	1:A:197:GLU:HG3	2.17	0.43
1:D:221:LYS:HE2	1:D:223:TYR:O	2.18	0.43
1:D:378:LYS:HB3	1:D:382:ARG:CZ	2.48	0.43
1:D:386:GLU:N	1:D:387:PRO:CD	2.82	0.43
1:A:414:ARG:O	1:A:418:GLU:HG3	2.19	0.43
1:C:252:LEU:HD21	1:C:281:HIS:CG	2.54	0.43
1:D:295:GLY:HA3	1:D:345:GLU:HG2	2.01	0.43
1:C:55:PRO:HD3	1:C:326:ILE:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:169:LYS:HA	4:B:720:HOH:O	2.19	0.43
1:D:76:ARG:HD3	1:D:420:PRO:HD2	2.00	0.43
1:A:179:TRP:CE2	3:A:601:PSJ:H1A	2.54	0.43
1:A:295:GLY:HA3	1:A:345:GLU:HG2	2.01	0.43
1:A:254:ASP:HA	1:A:281:HIS:HB2	2.01	0.42
1:A:65:ARG:HG2	4:A:615:HOH:O	2.18	0.42
1:C:65:ARG:HG2	4:D:441:HOH:O	2.19	0.42
1:D:108:ASP:OD2	1:D:110:LYS:HE2	2.20	0.42
1:C:56:SER:O	1:C:59:VAL:HG22	2.19	0.42
1:D:179:TRP:CD2	3:D:604:PSJ:H1A	2.55	0.42
1:A:252:LEU:HD21	1:A:281:HIS:CG	2.55	0.42
1:B:193:THR:O	1:B:197:GLU:HG3	2.20	0.42
1:C:194:ARG:O	1:C:198:ARG:HG3	2.20	0.41
1:A:280:PHE:CZ	1:A:309:LEU:HD21	2.55	0.41
1:B:76:ARG:HH11	1:B:76:ARG:CG	2.32	0.41
1:A:59:VAL:HG11	1:A:81:LYS:CB	2.51	0.41
1:B:252:LEU:HD21	1:B:281:HIS:CG	2.55	0.41
1:B:419:ARG:HA	1:B:420:PRO:HD3	1.81	0.41
1:C:386:GLU:N	1:C:387:PRO:CD	2.84	0.41
1:A:169:LYS:HD3	1:A:211:PRO:HG2	2.02	0.41
1:B:222:MET:CE	1:B:259:ALA:HB2	2.51	0.41
1:C:221:LYS:HE2	1:C:223:TYR:O	2.20	0.41
1:A:53:ALA:HB3	1:A:325:MET:HG2	2.02	0.41
1:C:6:ILE:HD11	1:C:87:VAL:CG2	2.50	0.41
1:C:54:VAL:HG13	1:C:55:PRO:HD2	2.01	0.41
1:C:102:ILE:CG2	1:C:103:PRO:HA	2.51	0.41
1:D:54:VAL:HG13	1:D:95:THR:HB	2.02	0.41
4:B:533:HOH:O	1:D:194:ARG:HD2	2.20	0.41
1:C:154:ARG:O	1:C:158:VAL:HG23	2.20	0.41
1:D:7:ALA:HB3	1:D:10:VAL:HG23	2.03	0.41
1:A:378:LYS:O	1:A:382:ARG:HG3	2.21	0.40
1:B:187:PRO:HG3	1:B:228:TYR:CE1	2.55	0.40
1:C:317:VAL:CG1	1:C:318:LYS:HG2	2.51	0.40
1:C:429:ILE:HG13	1:D:329:LEU:HD11	2.03	0.40
1:B:179:TRP:CD2	3:B:602:PSJ:H1A	2.55	0.40
1:A:46:LYS:HB3	1:A:46:LYS:HE2	1.88	0.40
1:C:414:ARG:O	1:C:418:GLU:HG3	2.21	0.40
1:A:315:ARG:NH2	4:A:531:HOH:O	2.54	0.40
1:B:5:ARG:HG2	4:B:1235:HOH:O	2.20	0.40
1:B:169:LYS:HD3	1:B:211:PRO:CG	2.51	0.40
1:C:329:LEU:HG	4:C:1301:HOH:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:429:ILE:HG13	1:D:329:LEU:CD1	2.51	0.40
1:B:180:ILE:HD12	1:B:182:ASP:CG	2.41	0.40
1:C:7:ALA:HB3	1:C:10:VAL:CG2	2.52	0.40
1:D:411:TYR:O	1:D:414:ARG:HB3	2.22	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	419/438 (96%)	410 (98%)	8 (2%)	1 (0%)	47	33
1	B	419/438 (96%)	409 (98%)	9 (2%)	1 (0%)	47	33
1	C	428/438 (98%)	412 (96%)	11 (3%)	5 (1%)	13	3
1	D	417/438 (95%)	407 (98%)	9 (2%)	1 (0%)	47	33
All	All	1683/1752 (96%)	1638 (97%)	37 (2%)	8 (0%)	29	15

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	316	GLY
1	C	318	LYS
1	C	320	PHE
1	A	224	GLU
1	B	224	GLU
1	C	224	GLU
1	D	224	GLU
1	C	319	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	330/341 (97%)	327 (99%)	3 (1%)	78	75
1	B	330/341 (97%)	326 (99%)	4 (1%)	71	65
1	C	334/341 (98%)	331 (99%)	3 (1%)	78	75
1	D	329/341 (96%)	327 (99%)	2 (1%)	86	84
All	All	1323/1364 (97%)	1311 (99%)	12 (1%)	78	75

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	ARG
1	A	318	LYS
1	A	374	THR
1	B	9	ASP
1	B	318	LYS
1	B	354	LEU
1	B	374	THR
1	C	124	ASP
1	C	374	THR
1	C	430	ILE
1	D	124	ASP
1	D	374	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	15	ASN
1	A	160	HIS
1	A	189	GLN
1	A	243	GLN
1	A	258	HIS
1	B	15	ASN

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Mol	Chain	Res	Type
1	B	160	HIS
1	B	189	GLN
1	B	258	HIS
1	C	15	ASN
1	C	160	HIS
1	C	189	GLN
1	C	258	HIS
1	C	344	ASN
1	D	8	GLN
1	D	15	ASN
1	D	160	HIS
1	D	189	GLN
1	D	258	HIS
1	D	344	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 8 are monoatomic - leaving 4 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	PSJ	C	603	2	10,11,11	0.42	0	9,14,14	0.70	0
3	PSJ	A	601	2	10,11,11	0.50	0	9,14,14	0.59	0
3	PSJ	B	602	2	10,11,11	0.57	0	9,14,14	0.65	0
3	PSJ	D	604	2	10,11,11	0.51	0	9,14,14	0.65	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	PSJ	C	603	2	-	0/16/16/16	-
3	PSJ	A	601	2	-	0/16/16/16	-
3	PSJ	B	602	2	-	0/16/16/16	-
3	PSJ	D	604	2	-	0/16/16/16	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	603	PSJ	2	0
3	A	601	PSJ	2	0
3	B	602	PSJ	3	0
3	D	604	PSJ	3	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	421/438 (96%)	0.43	13 (3%) 49 43	8, 14, 25, 39	0
1	B	421/438 (96%)	0.41	12 (2%) 51 46	8, 15, 26, 47	0
1	C	430/438 (98%)	0.74	30 (6%) 16 13	8, 17, 33, 53	0
1	D	419/438 (95%)	0.40	12 (2%) 51 46	8, 14, 25, 46	0
All	All	1691/1752 (96%)	0.50	67 (3%) 38 32	8, 15, 28, 53	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	423	VAL	10.2
1	B	424	ALA	9.2
1	C	319	GLY	8.4
1	C	3	GLU	6.5
1	A	4	PHE	6.1
1	C	4	PHE	4.8
1	C	320	PHE	4.7
1	D	4	PHE	4.7
1	C	315	ARG	4.6
1	C	321	HIS	4.6
1	D	3	GLU	4.5
1	C	314	ALA	4.4
1	C	317	VAL	4.3
1	C	316	GLY	4.1
1	C	318	LYS	4.1
1	A	424	ALA	3.8
1	D	421	ALA	3.8
1	B	314	ALA	3.7
1	C	79	PHE	3.6
1	D	11	VAL	3.5
1	A	319	GLY	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	422	SER	3.2
1	C	136	GLY	3.1
1	C	135	PRO	3.0
1	A	320	PHE	3.0
1	D	420	PRO	3.0
1	B	318	LYS	3.0
1	D	8	GLN	2.9
1	D	110	LYS	2.9
1	C	76	ARG	2.9
1	C	12	ALA	2.9
1	C	429	ILE	2.7
1	C	8	GLN	2.7
1	B	4	PHE	2.6
1	A	318	LYS	2.6
1	B	115	ARG	2.5
1	A	287	TYR	2.5
1	C	110	LYS	2.5
1	C	11	VAL	2.5
1	C	395	ARG	2.5
1	D	287	TYR	2.5
1	D	135	PRO	2.5
1	D	106	LYS	2.4
1	B	329	LEU	2.4
1	C	430	ILE	2.3
1	D	297	ILE	2.3
1	C	115	ARG	2.3
1	A	413	ALA	2.3
1	C	223	TYR	2.3
1	C	432	SER	2.2
1	D	19	ALA	2.2
1	C	392	ALA	2.2
1	C	212	ASP	2.2
1	B	297	ILE	2.2
1	A	136	GLY	2.2
1	A	422	SER	2.2
1	C	415	VAL	2.2
1	A	110	LYS	2.2
1	A	259	ALA	2.2
1	C	259	ALA	2.2
1	B	206	ILE	2.1
1	A	314	ALA	2.1
1	B	212	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	139	HIS	2.1
1	C	400	VAL	2.1
1	B	321	HIS	2.1
1	A	423	VAL	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 monosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	PSJ	D	604	12/12	0.90	0.14	10,14,18,20	0
3	PSJ	B	602	12/12	0.91	0.14	10,14,19,21	0
3	PSJ	C	603	12/12	0.91	0.11	10,16,19,19	0
3	PSJ	A	601	12/12	0.91	0.12	13,17,21,21	0
2	MN	C	506	1/1	0.99	0.06	12,12,12,12	0
2	MN	D	507	1/1	0.99	0.08	11,11,11,11	0
2	MN	A	502	1/1	1.00	0.04	10,10,10,10	0
2	MN	D	508	1/1	1.00	0.05	11,11,11,11	0
2	MN	B	503	1/1	1.00	0.07	13,13,13,13	0
2	MN	B	504	1/1	1.00	0.07	12,12,12,12	0
2	MN	C	505	1/1	1.00	0.06	15,15,15,15	0
2	MN	A	501	1/1	1.00	0.05	13,13,13,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.