



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 03:02 AM GMT

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 validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

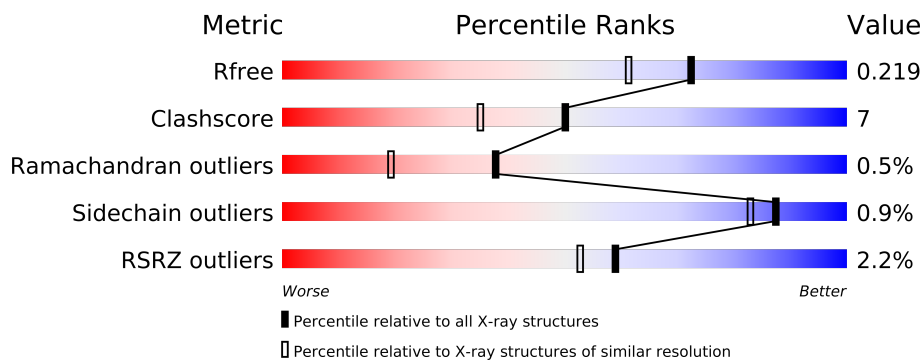
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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}	66092	3513 (1.80-1.80)
Clashscore	79885	4461 (1.80-1.80)
Ramachandran outliers	78287	4404 (1.80-1.80)
Sidechain outliers	78261	4403 (1.80-1.80)
RSRZ outliers	66119	3515 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	438	
1	B	438	
1	C	438	
1	D	438	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14571 atoms, of which 0 are hydrogen and 0 are deuterium.

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	UNP Q75WH8
A	329	LEU	SER	ENGINEERED	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	UNP Q75WH8
B	329	LEU	SER	ENGINEERED	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	UNP Q75WH8

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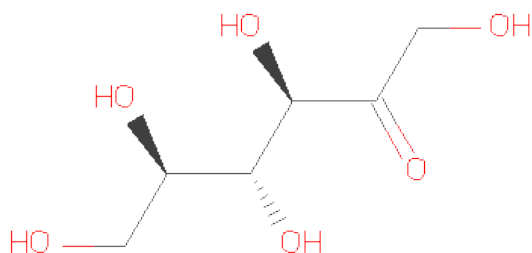
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Chain	Residue	Modelled	Actual	Comment	Reference
C	329	LEU	SER	ENGINEERED	UNP Q75WH8
C	431	GLY	-	EXPRESSION TAG	UNP Q75WH8
C	432	SER	-	EXPRESSION TAG	UNP Q75WH8
C	433	HIS	-	EXPRESSION TAG	UNP Q75WH8
C	434	HIS	-	EXPRESSION TAG	UNP Q75WH8
C	435	HIS	-	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	UNP Q75WH8
D	329	LEU	SER	ENGINEERED	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	B	2	Total Mn 2 2	0	0
2	A	2	Total Mn 2 2	0	0
2	D	2	Total Mn 2 2	0	0
2	C	2	Total Mn 2 2	0	0

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



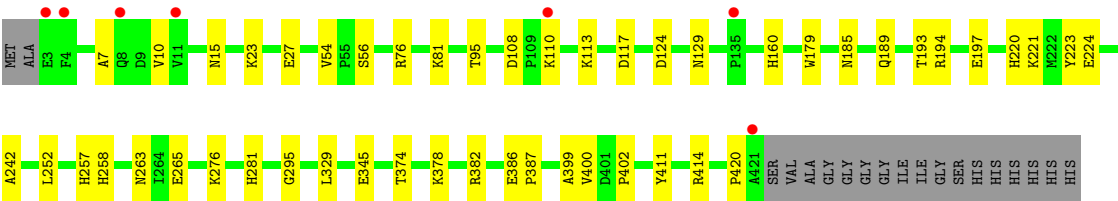
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		
3	C	1	Total	C	O	0	0
			12	6	6		
3	D	1	Total	C	O	0	0
			12	6	6		

- 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		

● Molecule 1: L-rhamnose isomerase

Chain D: 



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) 92.0 (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 0.197 , 0.219	Depositor DCC
R_{free} test set	14123 reflections (9.97%)	DCC
Wilson B-factor (Å ²)	14.9	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 36.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 143305 reflections	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.

5 Model quality

5.1 Standard geometry

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 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

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

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	12	0	10	2	0
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

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

All (194) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(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:H	1:A:110:LYS:HD2	1.62	0.65
1:B:110:LYS:HG3	4:B:1084:HOH:O	1.97	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:185:ASN:H	1:C:189:GLN:NE2	1.95	0.64
1:C:6:ILE:HD11	1:C:87:VAL:HG22	1.80	0.63
1:B:15:ASN:ND2	1:B:400:VAL:H	1.97	0.63
1:C:210:LEU:HD11	1:C:216:LEU:HB2	1.81	0.62
1:B:361:LEU:O	1:B:365:GLN:HG3	1.99	0.62
1:D:23:LYS:O	1:D:27:GLU:HG3	1.99	0.62
1:C:44:THR:O	1:C:48:GLU:HG3	1.99	0.62
1:C:6:ILE:N	1:C:6:ILE:HD12	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:D:378:LYS:HB3	1:D:382:ARG:NH2	2.21	0.56
1:C:340:ILE:HG21	1:C:402:PRO:HB2	1.87	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:HG3	1:C:318:LYS:O	2.09	0.53
1:D:179:TRP:CE2	3:D:604:PSJ:H1A	2.43	0.53
1:C:430:ILE:HD13	1:C:430:ILE:H	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
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

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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:402:PRO:HG2	4:A:485:HOH:O	2.11	0.50
1:A:277:LEU:HD23	1:A:320:PHE:CE1	2.47	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:D:15:ASN:HD21	1:D:399:ALA:HA	1.76	0.49
1:A:221:LYS:HA	1:A:257:HIS:HB3	1.94	0.49
1:C:6:ILE:HD11	1:C:87:VAL:CG1	2.36	0.49
1:B:76:ARG:HG2	1:B:76:ARG:HH11	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:C:318:LYS:C	1:C:320:PHE:N	2.65	0.48
1:A:23:LYS:O	1:A:27:GLU:HG3	2.13	0.48
1:C:73:GLY:CA	1:C:416:ALA:HA	2.44	0.48
3:A:601:PSJ:C2	3:A:601:PSJ:O5	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: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:B:318:LYS:O	1:B:318:LYS:HD3	2.13	0.47
1:D:194:ARG:HH11	1:D:194:ARG:HG2	1.79	0.47
3:D:604:PSJ:C2	3:D:604:PSJ:O5	2.63	0.47
1:D:108:ASP:OD2	1:D:110:LYS:HB2	2.15	0.47
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

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:315:ARG:O	1:C:315:ARG:HG3	2.16	0.46
1:A:211:PRO:HD2	1:A:214:TRP:CG	2.50	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:C2	3:B:602:PSJ:O5	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:C:318:LYS:O	1:C:319:GLY:C	2.54	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: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:D:378:LYS:HB3	1:D:382:ARG:CZ	2.48	0.43
1:D:221:LYS:HE2	1:D:223:TYR:O	2.18	0.43
1:D:386:GLU:N	1:D:387:PRO:CD	2.82	0.43
1:A:193:THR:O	1:A:197:GLU:HG3	2.17	0.43
1:D:295:GLY:HA3	1:D:345:GLU:HG2	2.01	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:C:55:PRO:HD3	1:C:326:ILE:O	2.19	0.43
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

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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:C:386:GLU:N	1:C:387:PRO:CD	2.84	0.41
1:B:419:ARG:HA	1:B:420:PRO:HD3	1.81	0.41
1:A:169:LYS:HD3	1:A:211:PRO:HG2	2.02	0.41
1:C:221:LYS:HE2	1:C:223:TYR:O	2.20	0.41
1:B:222:MET:CE	1:B:259:ALA:HB2	2.51	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
1:A:53:ALA:HB3	1:A:325:MET:HG2	2.02	0.41
4:B:533:HOH:O	1:D:194:ARG:HD2	2.20	0.41
1:D:7:ALA:HB3	1:D:10:VAL:HG23	2.03	0.41
1:C:154:ARG:O	1:C:158:VAL:HG23	2.20	0.41
1:C:429:ILE:HG13	1:D:329:LEU:HD11	2.03	0.40
1:C:317:VAL:CG1	1:C:318:LYS:HG2	2.51	0.40
1:B:187:PRO:HG3	1:B:228:TYR:CE1	2.55	0.40
1:A:378:LYS:O	1:A:382:ARG:HG3	2.21	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:C:429:ILE:HG13	1:D:329:LEU:CD1	2.51	0.40
1:B:169:LYS:HD3	1:B:211:PRO:CG	2.51	0.40
1:B:5:ARG:HG2	4:B:1235:HOH:O	2.20	0.40
1:A:315:ARG:NH2	4:A:531:HOH:O	2.54	0.40
1:C:329:LEU:HG	4:C:1301:HOH:O	2.21	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
1:B:180:ILE:HD12	1:B:182:ASP:CG	2.41	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%)	56	38
1	B	419/438 (96%)	409 (98%)	9 (2%)	1 (0%)	56	38
1	C	428/438 (98%)	412 (96%)	11 (3%)	5 (1%)	19	5
1	D	417/438 (95%)	407 (98%)	9 (2%)	1 (0%)	56	38
All	All	1683/1752 (96%)	1638 (97%)	37 (2%)	8 (0%)	38	19

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%)	87	83
1	B	330/341 (97%)	326 (99%)	4 (1%)	82	74
1	C	334/341 (98%)	331 (99%)	3 (1%)	87	83
1	D	329/341 (96%)	327 (99%)	2 (1%)	92	90
All	All	1323/1364 (97%)	1311 (99%)	12 (1%)	87	83

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

Some 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
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 ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	A	601	2	11,11,11	0.47	0	14,14,14	0.64	0
3	PSJ	B	602	2	11,11,11	0.54	0	14,14,14	0.69	0
3	PSJ	C	603	2	11,11,11	0.40	0	14,14,14	0.75	0
3	PSJ	D	604	2	11,11,11	0.52	0	14,14,14	0.72	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	A	601	2	-	0/16/16/16	0/0/0/0
3	PSJ	B	602	2	-	0/16/16/16	0/0/0/0
3	PSJ	C	603	2	-	0/16/16/16	0/0/0/0
3	PSJ	D	604	2	-	0/16/16/16	0/0/0/0

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.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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%)	0.14	6 (1%) 72 68	8, 14, 25, 39	0
1	B	421/438 (96%)	0.11	7 (1%) 67 62	8, 15, 26, 47	0
1	C	430/438 (98%)	0.43	18 (4%) 35 28	8, 17, 33, 53	0
1	D	419/438 (95%)	0.09	7 (1%) 67 62	8, 14, 25, 46	0
All	All	1691/1752 (96%)	0.19	38 (2%) 59 53	8, 15, 28, 53	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	424	ALA	7.4
1	B	423	VAL	6.8
1	C	319	GLY	5.1
1	A	4	PHE	4.8
1	C	3	GLU	4.5
1	C	320	PHE	4.4
1	D	3	GLU	4.3
1	C	317	VAL	3.9
1	C	314	ALA	3.8
1	D	4	PHE	3.6
1	D	421	ALA	3.6
1	B	422	SER	3.5
1	C	321	HIS	3.5
1	C	315	ARG	3.3
1	A	424	ALA	3.2
1	C	4	PHE	3.2
1	C	318	LYS	3.2
1	C	316	GLY	3.1
1	C	79	PHE	2.9
1	A	319	GLY	2.6
1	A	318	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	314	ALA	2.4
1	D	8	GLN	2.4
1	B	115	ARG	2.4
1	C	8	GLN	2.3
1	C	135	PRO	2.3
1	C	136	GLY	2.3
1	D	135	PRO	2.3
1	C	432	SER	2.2
1	B	421	ALA	2.2
1	A	320	PHE	2.2
1	D	11	VAL	2.2
1	C	429	ILE	2.2
1	B	318	LYS	2.2
1	C	76	ARG	2.1
1	A	423	VAL	2.1
1	D	110	LYS	2.1
1	C	428	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	PSJ	B	602	12/12	0.12	1.68	10,14,19,21	0
3	PSJ	D	604	12/12	0.12	0.65	10,14,18,20	0
3	PSJ	A	601	12/12	0.11	0.54	13,17,21,21	0
3	PSJ	C	603	12/12	0.09	-0.88	10,16,19,19	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	MN	D	507	1/1	0.07	-1.85	11,11,11,11	0
2	MN	C	506	1/1	0.06	-2.66	12,12,12,12	0
2	MN	B	504	1/1	0.04	-3.45	12,12,12,12	0
2	MN	B	503	1/1	0.05	-4.00	13,13,13,13	0
2	MN	C	505	1/1	0.05	-4.61	15,15,15,15	0
2	MN	D	508	1/1	0.03	-5.10	11,11,11,11	0
2	MN	A	502	1/1	0.02	-6.40	10,10,10,10	0
2	MN	A	501	1/1	0.03	-7.58	13,13,13,13	0

6.5 Other polymers

There are no such residues in this entry.