



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:36 PM GMT

PDB ID : 4FFH
Title : Crystal Structure of Levan Fructotransferase D54N mutant from *Arthrobacter ureafaciens* in complex with sucrose
Authors : Park, J.; Rhee, S.
Deposited on : 2012-06-01
Resolution : 2.20 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

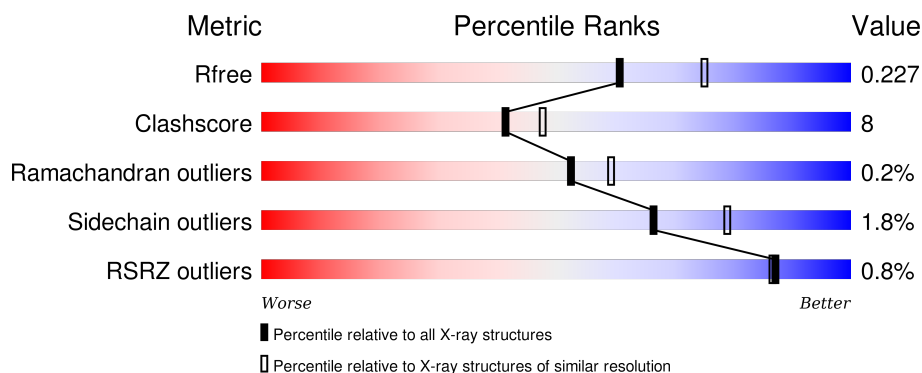
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	91344	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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. 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	492	<div> <div></div> <div>78%19%..</div> </div>
1	B	492	<div> <div></div> <div>77%21%.</div> </div>
1	C	492	<div> <div></div> <div>83%15%.</div> </div>
1	D	492	<div> <div></div> <div>80%16%..</div> </div>

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
2	SUC	D	601	-	-	-	X
2	SUC	D	602	-	-	-	X
2	SUC	D	605	-	-	-	X

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 15917 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 Levan fructotransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	480	Total	C	N	O	S	0	0	0
			3739	2375	646	710	8			
1	B	480	Total	C	N	O	S	0	0	0
			3739	2375	646	710	8			
1	C	480	Total	C	N	O	S	0	0	0
			3739	2375	646	710	8			
1	D	478	Total	C	N	O	S	0	0	0
			3725	2367	643	707	8			

There are 52 discrepancies between the modelled and reference sequences:

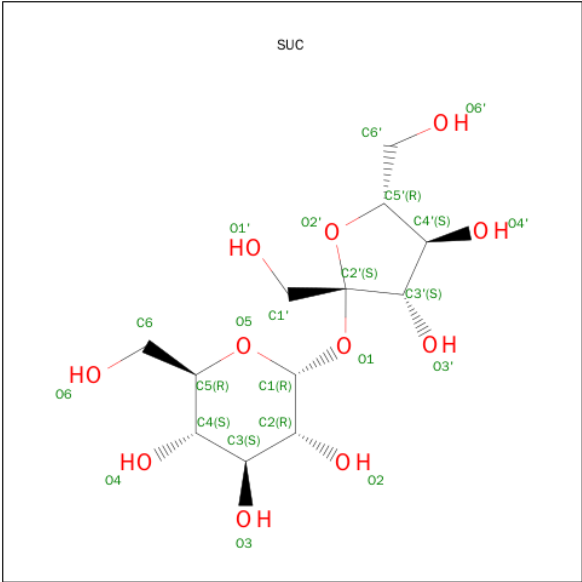
Chain	Residue	Modelled	Actual	Comment	Reference
A	40	MET	-	EXPRESSION TAG	UNP Q9KJD0
A	54	ASN	ASP	ENGINEERED MUTATION	UNP Q9KJD0
A	115	ASP	GLY	CONFLICT	UNP Q9KJD0
A	522	LEU	-	EXPRESSION TAG	UNP Q9KJD0
A	523	GLU	-	EXPRESSION TAG	UNP Q9KJD0
A	524	HIS	-	EXPRESSION TAG	UNP Q9KJD0
A	525	HIS	-	EXPRESSION TAG	UNP Q9KJD0
A	526	HIS	-	EXPRESSION TAG	UNP Q9KJD0
A	527	HIS	-	EXPRESSION TAG	UNP Q9KJD0
A	528	HIS	-	EXPRESSION TAG	UNP Q9KJD0
A	529	HIS	-	EXPRESSION TAG	UNP Q9KJD0
A	530	HIS	-	EXPRESSION TAG	UNP Q9KJD0
A	531	HIS	-	EXPRESSION TAG	UNP Q9KJD0
B	40	MET	-	EXPRESSION TAG	UNP Q9KJD0
B	54	ASN	ASP	ENGINEERED MUTATION	UNP Q9KJD0
B	115	ASP	GLY	CONFLICT	UNP Q9KJD0
B	522	LEU	-	EXPRESSION TAG	UNP Q9KJD0
B	523	GLU	-	EXPRESSION TAG	UNP Q9KJD0
B	524	HIS	-	EXPRESSION TAG	UNP Q9KJD0
B	525	HIS	-	EXPRESSION TAG	UNP Q9KJD0
B	526	HIS	-	EXPRESSION TAG	UNP Q9KJD0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	527	HIS	-	EXPRESSION TAG	UNP Q9KJD0
B	528	HIS	-	EXPRESSION TAG	UNP Q9KJD0
B	529	HIS	-	EXPRESSION TAG	UNP Q9KJD0
B	530	HIS	-	EXPRESSION TAG	UNP Q9KJD0
B	531	HIS	-	EXPRESSION TAG	UNP Q9KJD0
C	40	MET	-	EXPRESSION TAG	UNP Q9KJD0
C	54	ASN	ASP	ENGINEERED MUTATION	UNP Q9KJD0
C	115	ASP	GLY	CONFLICT	UNP Q9KJD0
C	522	LEU	-	EXPRESSION TAG	UNP Q9KJD0
C	523	GLU	-	EXPRESSION TAG	UNP Q9KJD0
C	524	HIS	-	EXPRESSION TAG	UNP Q9KJD0
C	525	HIS	-	EXPRESSION TAG	UNP Q9KJD0
C	526	HIS	-	EXPRESSION TAG	UNP Q9KJD0
C	527	HIS	-	EXPRESSION TAG	UNP Q9KJD0
C	528	HIS	-	EXPRESSION TAG	UNP Q9KJD0
C	529	HIS	-	EXPRESSION TAG	UNP Q9KJD0
C	530	HIS	-	EXPRESSION TAG	UNP Q9KJD0
C	531	HIS	-	EXPRESSION TAG	UNP Q9KJD0
D	40	MET	-	EXPRESSION TAG	UNP Q9KJD0
D	54	ASN	ASP	ENGINEERED MUTATION	UNP Q9KJD0
D	115	ASP	GLY	CONFLICT	UNP Q9KJD0
D	522	LEU	-	EXPRESSION TAG	UNP Q9KJD0
D	523	GLU	-	EXPRESSION TAG	UNP Q9KJD0
D	524	HIS	-	EXPRESSION TAG	UNP Q9KJD0
D	525	HIS	-	EXPRESSION TAG	UNP Q9KJD0
D	526	HIS	-	EXPRESSION TAG	UNP Q9KJD0
D	527	HIS	-	EXPRESSION TAG	UNP Q9KJD0
D	528	HIS	-	EXPRESSION TAG	UNP Q9KJD0
D	529	HIS	-	EXPRESSION TAG	UNP Q9KJD0
D	530	HIS	-	EXPRESSION TAG	UNP Q9KJD0
D	531	HIS	-	EXPRESSION TAG	UNP Q9KJD0

- Molecule 2 is SUCROSE (three-letter code: SUC, HOH) (formula: C₁₂H₂₂O₁₁, H₂O).

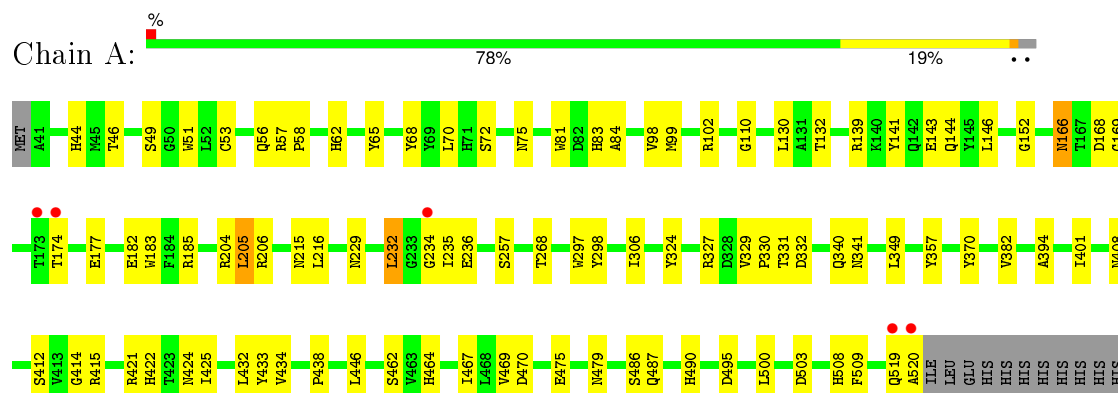


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	843	Total	C	O	0	0
			975	72	903		

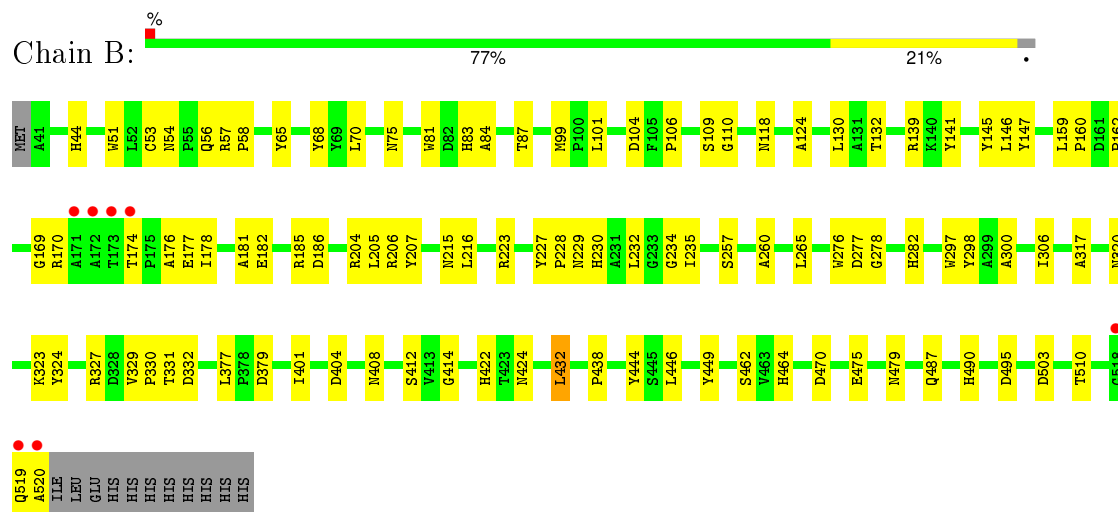
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

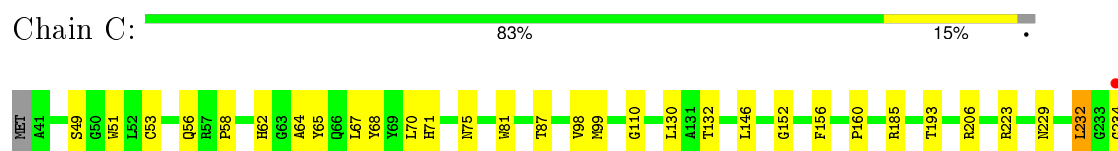
- Molecule 1: Levan fructotransferase

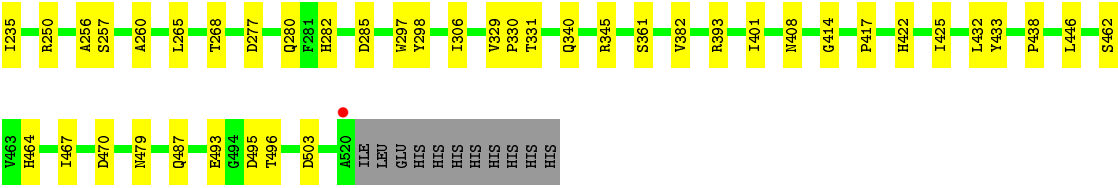


- Molecule 1: Levan fructotransferase

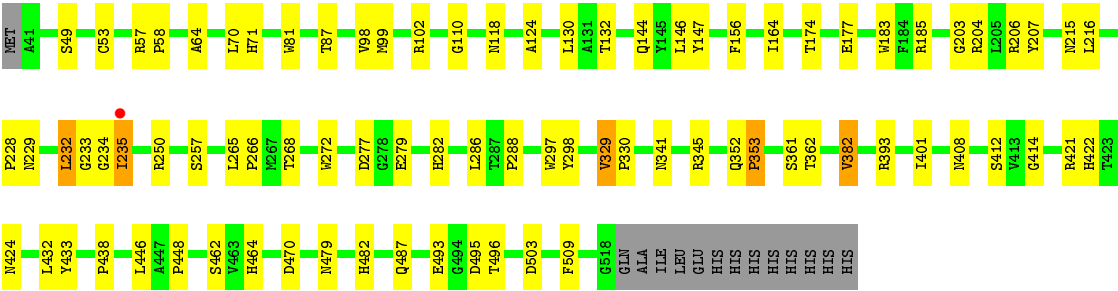
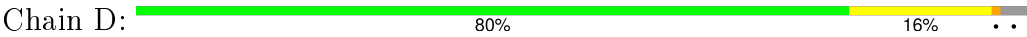


- Molecule 1: Levan fructotransferase





• Molecule 1: Levan fructotransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	82.28Å 167.00Å 261.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.20 44.38 – 2.20	Depositor EDS
% Data completeness (in resolution range)	83.0 (50.00-2.20) 90.2 (44.38-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.18 (at 2.20Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.190 , 0.218 0.200 , 0.227	Depositor DCC
R_{free} test set	16422 reflections (9.93%)	DCC
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.403	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 36.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 170519 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	15917	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.73% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: SUC

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	0.33	0/3863	0.63	1/5312 (0.0%)
1	B	0.33	0/3863	0.64	1/5312 (0.0%)
1	C	0.33	0/3863	0.63	1/5312 (0.0%)
1	D	0.34	0/3849	0.66	2/5293 (0.0%)
All	All	0.34	0/15438	0.64	5/21229 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	432	LEU	N-CA-C	-6.49	93.48	111.00
1	D	234	GLY	N-CA-C	6.40	129.10	113.10
1	B	432	LEU	N-CA-C	-6.26	94.11	111.00
1	A	432	LEU	N-CA-C	-6.19	94.28	111.00
1	C	432	LEU	N-CA-C	-5.68	95.67	111.00

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	3739	0	3491	73	0
1	B	3739	0	3491	71	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	3739	0	3491	51	0
1	D	3725	0	3478	52	0
2	D	975	0	120	29	0
All	All	15917	0	14071	247	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 8.

All (247) 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:139:ARG:HH22	2:D:601:SUC:H61	1.27	0.97
1:B:174:THR:HB	1:B:177:GLU:HG3	1.46	0.95
1:B:139:ARG:HH22	2:D:602:SUC:H61	1.33	0.90
1:A:329:VAL:HG13	1:A:330:PRO:HD2	1.55	0.85
1:C:425:ILE:HD13	1:C:467:ILE:HD13	1.61	0.82
1:A:434:VAL:HB	1:A:487:GLN:HE22	1.44	0.81
1:B:408:ASN:HB3	1:B:503:ASP:HB2	1.65	0.79
1:C:329:VAL:HG13	1:C:330:PRO:HD2	1.66	0.78
1:A:174:THR:OG1	1:A:177:GLU:HG3	1.85	0.77
1:B:329:VAL:HG13	1:B:330:PRO:HD2	1.66	0.76
1:B:206:ARG:HH11	1:B:234:GLY:HA3	1.52	0.73
1:D:203:GLY:HA2	1:D:235:ILE:HG22	1.68	0.73
1:C:417:PRO:HG2	2:D:965:HOH:O	1.88	0.73
1:A:139:ARG:NH2	2:D:601:SUC:H61	2.03	0.73
1:D:329:VAL:HG22	1:D:330:PRO:HD2	1.73	0.70
1:A:206:ARG:HH11	1:A:234:GLY:HA3	1.57	0.69
1:A:139:ARG:HH22	2:D:601:SUC:C6	2.03	0.69
1:A:434:VAL:CB	1:A:487:GLN:HE22	2.06	0.69
1:A:141:TYR:O	1:A:143:GLU:HG2	1.92	0.69
1:B:145:TYR:CZ	1:B:162:PRO:HG3	2.29	0.68
1:D:464:HIS:H	1:D:479:ASN:ND2	1.91	0.67
1:B:54:ASN:HD21	2:D:602:SUC:H1'1	1.60	0.67
1:C:464:HIS:H	1:C:479:ASN:ND2	1.92	0.67
1:C:408:ASN:HB3	1:C:503:ASP:HB2	1.76	0.66
1:D:207:TYR:O	1:D:235:ILE:HG21	1.96	0.66
1:D:174:THR:OG1	1:D:177:GLU:HG3	1.96	0.65
1:D:53:CYS:HB3	1:D:70:LEU:HB2	1.78	0.65
1:A:464:HIS:H	1:A:479:ASN:ND2	1.95	0.65
1:C:393:ARG:HD3	1:C:496:THR:HG22	1.80	0.64
1:B:206:ARG:HD3	1:B:234:GLY:HA2	1.77	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:393:ARG:HD3	1:D:496:THR:HG22	1.80	0.63
1:C:206:ARG:NH1	1:C:234:GLY:N	2.47	0.63
1:B:145:TYR:CE2	1:B:162:PRO:HG3	2.34	0.63
1:A:44:HIS:HD2	1:A:470:ASP:OD2	1.81	0.62
1:B:464:HIS:H	1:B:479:ASN:ND2	1.97	0.62
1:A:206:ARG:HD3	1:A:234:GLY:HA2	1.80	0.62
1:A:229:ASN:O	1:A:232:LEU:HB2	1.99	0.62
1:B:159:LEU:HD22	1:B:160:PRO:HD2	1.81	0.62
1:A:102:ARG:HD3	2:D:957:HOH:O	2.00	0.62
1:C:81:TRP:HB2	1:C:99:MET:HB2	1.82	0.61
1:C:53:CYS:HB3	1:C:70:LEU:HB2	1.81	0.61
1:C:206:ARG:HH11	1:C:234:GLY:CA	2.12	0.61
1:B:81:TRP:HB2	1:B:99:MET:HB2	1.83	0.61
1:D:464:HIS:H	1:D:479:ASN:HD21	1.48	0.61
1:A:44:HIS:HE1	1:A:475:GLU:OE1	1.84	0.61
1:A:324:TYR:O	1:A:327:ARG:HG2	2.02	0.60
1:B:87:THR:HG22	2:D:1042:HOH:O	2.01	0.60
1:B:329:VAL:HG12	1:B:331:THR:H	1.67	0.60
1:C:193:THR:HG22	2:D:1195:HOH:O	2.01	0.59
1:A:229:ASN:HB3	1:A:232:LEU:HD22	1.82	0.59
1:B:229:ASN:O	1:B:232:LEU:HB2	2.02	0.59
1:C:229:ASN:O	1:C:232:LEU:HB2	2.04	0.58
1:A:434:VAL:HB	1:A:487:GLN:NE2	2.17	0.58
1:A:81:TRP:HB2	1:A:99:MET:HB2	1.85	0.58
1:D:257:SER:HA	1:D:268:THR:O	2.03	0.58
1:C:422:HIS:O	1:C:438:PRO:HB2	2.04	0.57
1:D:203:GLY:CA	1:D:235:ILE:HG22	2.34	0.57
1:C:49:SER:O	1:C:71:HIS:HE1	1.87	0.57
1:B:206:ARG:NH1	1:B:234:GLY:HA3	2.20	0.57
1:D:49:SER:O	1:D:71:HIS:HE1	1.87	0.57
1:B:519:GLN:N	1:B:519:GLN:OE1	2.36	0.57
1:D:229:ASN:HB3	1:D:232:LEU:HD22	1.86	0.57
1:D:206:ARG:NE	1:D:233:GLY:HA2	2.20	0.56
1:C:56:GLN:HB2	1:C:68:TYR:HB2	1.87	0.56
1:A:329:VAL:HG13	1:A:330:PRO:CD	2.31	0.56
1:D:64:ALA:HB2	1:D:87:THR:HG22	1.87	0.56
1:A:215:ASN:O	1:A:216:LEU:HB2	2.05	0.56
1:B:44:HIS:HD2	1:B:470:ASP:OD2	1.88	0.56
1:D:414:GLY:O	1:D:495:ASP:HB3	2.05	0.56
1:B:204:ARG:O	1:B:234:GLY:O	2.24	0.56
1:A:408:ASN:HB3	1:A:503:ASP:HB2	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:146:LEU:HD23	1:B:147:TYR:N	2.21	0.56
1:C:64:ALA:HB2	1:C:87:THR:HG22	1.88	0.55
1:A:72:SER:HB3	2:D:1252:HOH:O	2.06	0.55
1:B:54:ASN:HD21	2:D:602:SUC:C1'	2.19	0.55
1:C:464:HIS:H	1:C:479:ASN:HD21	1.54	0.55
1:B:53:CYS:HB3	1:B:70:LEU:HB2	1.87	0.55
1:B:174:THR:HG22	1:B:176:ALA:H	1.70	0.55
1:C:206:ARG:NH1	1:C:234:GLY:H	2.03	0.55
1:B:404:ASP:HB3	2:D:1106:HOH:O	2.07	0.55
1:B:519:GLN:O	1:B:520:ALA:HB3	2.07	0.55
1:A:433:TYR:CE2	2:D:603:SUC:H1'2	2.41	0.55
1:A:53:CYS:HB3	1:A:70:LEU:HB2	1.89	0.54
1:B:118:ASN:HB2	1:B:124:ALA:HA	1.90	0.54
1:C:285:ASP:HB3	2:D:859:HOH:O	2.08	0.54
1:B:446:LEU:HD21	1:B:490:HIS:CD2	2.43	0.54
1:C:277:ASP:OD1	1:C:282:HIS:HE1	1.90	0.54
1:A:166:ASN:HD22	1:A:166:ASN:C	2.11	0.54
1:A:204:ARG:O	1:A:234:GLY:O	2.26	0.54
1:C:257:SER:HB3	1:C:298:TYR:CE2	2.43	0.54
1:D:228:PRO:HD3	1:D:286:LEU:HD13	1.90	0.53
1:D:81:TRP:HB2	1:D:99:MET:HB2	1.90	0.53
1:B:324:TYR:O	1:B:327:ARG:HG2	2.09	0.53
1:A:236:GLU:OE2	2:D:601:SUC:H1'2	2.07	0.53
1:C:206:ARG:HD3	1:C:234:GLY:HA2	1.89	0.52
1:D:412:SER:OG	1:D:424:ASN:ND2	2.41	0.52
1:D:329:VAL:HG22	1:D:330:PRO:CD	2.39	0.52
1:D:257:SER:HB3	1:D:298:TYR:CE2	2.43	0.52
1:A:139:ARG:HG2	1:A:182:GLU:HG2	1.91	0.52
1:A:519:GLN:O	1:A:520:ALA:C	2.48	0.52
1:D:422:HIS:O	1:D:438:PRO:HB2	2.09	0.51
1:B:51:TRP:HB2	1:B:75:ASN:HD22	1.74	0.51
1:A:486:SER:O	1:A:487:GLN:HG3	2.10	0.51
1:A:464:HIS:H	1:A:479:ASN:HD21	1.59	0.51
1:D:433:TYR:CE2	2:D:606:SUC:H1'2	2.46	0.51
1:C:250:ARG:NH1	2:D:1171:HOH:O	2.43	0.51
1:A:206:ARG:NH1	1:A:234:GLY:HA3	2.24	0.50
1:C:62:HIS:HE1	1:C:152:GLY:HA3	1.75	0.50
1:B:401:ILE:O	1:B:462:SER:HA	2.11	0.50
1:C:433:TYR:CE2	2:D:605:SUC:H1'2	2.47	0.50
1:B:146:LEU:HD23	1:B:146:LEU:C	2.32	0.50
1:D:132:THR:HG21	1:D:185:ARG:HB3	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:215:ASN:O	1:B:216:LEU:HB2	2.10	0.50
1:C:257:SER:HA	1:C:268:THR:O	2.12	0.50
1:B:110:GLY:HA3	1:B:130:LEU:O	2.12	0.49
1:D:393:ARG:HG3	1:D:393:ARG:HH11	1.78	0.49
1:C:110:GLY:HA3	1:C:130:LEU:O	2.13	0.49
1:A:62:HIS:HE1	1:A:152:GLY:HA3	1.76	0.49
1:D:118:ASN:HB2	1:D:124:ALA:HA	1.95	0.49
1:C:329:VAL:HG13	1:C:330:PRO:CD	2.40	0.49
1:B:206:ARG:HD3	1:B:234:GLY:CA	2.42	0.49
1:D:352:GLN:OE1	1:D:482:HIS:HE1	1.94	0.49
1:C:51:TRP:HB2	1:C:75:ASN:HD22	1.78	0.49
1:B:141:TYR:CE1	1:B:170:ARG:HD3	2.47	0.49
1:C:393:ARG:CD	1:C:496:THR:HG22	2.43	0.48
1:C:345:ARG:HD3	1:C:361:SER:HB3	1.95	0.48
1:C:414:GLY:O	1:C:495:ASP:HB3	2.13	0.48
1:A:206:ARG:HH11	1:A:234:GLY:CA	2.26	0.48
1:C:160:PRO:HG2	2:D:1218:HOH:O	2.14	0.48
1:B:276:TRP:NE1	1:B:278:GLY:HA2	2.29	0.48
1:C:65:TYR:OH	1:C:306:ILE:HG13	2.12	0.48
1:D:470:ASP:HB3	2:D:699:HOH:O	2.12	0.48
1:A:49:SER:HB2	2:D:1258:HOH:O	2.13	0.48
1:A:257:SER:HA	1:A:268:THR:O	2.14	0.48
1:C:146:LEU:HD23	1:C:146:LEU:C	2.34	0.48
1:C:329:VAL:HG12	1:C:331:THR:H	1.78	0.48
1:C:65:TYR:CZ	1:C:306:ILE:HG13	2.48	0.48
1:B:101:LEU:HD12	1:B:106:PRO:HA	1.96	0.47
1:B:412:SER:OG	1:B:424:ASN:ND2	2.47	0.47
1:B:414:GLY:O	1:B:495:ASP:HB3	2.14	0.47
1:A:433:TYR:CZ	2:D:603:SUC:H1'2	2.50	0.47
1:A:62:HIS:CE1	1:A:152:GLY:HA3	2.49	0.47
1:D:146:LEU:HD23	1:D:146:LEU:C	2.35	0.47
1:D:408:ASN:HB3	1:D:503:ASP:HB2	1.97	0.47
1:A:183:TRP:O	1:A:204:ARG:HD2	2.14	0.47
1:A:44:HIS:CE1	1:A:475:GLU:OE1	2.65	0.47
1:B:44:HIS:HE1	1:B:475:GLU:OE2	1.97	0.47
1:A:486:SER:C	1:A:487:GLN:HG3	2.34	0.47
1:D:393:ARG:CD	1:D:496:THR:HG22	2.44	0.47
1:A:370:TYR:HE2	1:A:520:ALA:HA	1.79	0.47
1:A:132:THR:HG21	1:A:185:ARG:HB3	1.95	0.47
1:B:323:LYS:HD3	1:B:449:TYR:CZ	2.50	0.47
1:C:98:VAL:HG22	1:C:156:PHE:HD1	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:178:ILE:O	1:B:182:GLU:HG3	2.15	0.47
1:B:118:ASN:HB2	1:B:124:ALA:CA	2.45	0.47
1:D:204:ARG:O	1:D:235:ILE:HB	2.15	0.46
1:C:223:ARG:HB2	1:C:280:GLN:HB3	1.97	0.46
1:D:99:MET:HG2	1:D:147:TYR:CD1	2.50	0.46
1:B:464:HIS:H	1:B:479:ASN:HD21	1.62	0.46
2:D:603:SUC:H1	2:D:603:SUC:H1'1	1.83	0.46
1:A:401:ILE:O	1:A:462:SER:HA	2.15	0.46
1:C:493:GLU:HG2	1:D:493:GLU:HG2	1.96	0.46
1:A:425:ILE:HD13	1:A:467:ILE:HD13	1.98	0.46
1:A:166:ASN:ND2	1:A:169:GLY:H	2.13	0.46
1:B:56:GLN:HB2	1:B:68:TYR:HB2	1.97	0.46
1:D:215:ASN:O	1:D:216:LEU:HB2	2.15	0.46
1:D:421:ARG:O	1:D:422:HIS:HB3	2.16	0.46
1:B:257:SER:HB3	1:B:298:TYR:CE2	2.51	0.46
1:C:206:ARG:NH1	1:C:234:GLY:CA	2.78	0.45
1:B:44:HIS:CE1	1:B:475:GLU:OE2	2.70	0.45
1:A:422:HIS:O	1:A:438:PRO:HB2	2.17	0.45
1:B:109:SER:HB2	1:B:186:ASP:OD1	2.15	0.45
1:A:146:LEU:C	1:A:146:LEU:HD23	2.37	0.45
1:B:277:ASP:OD1	1:B:282:HIS:HE1	1.99	0.45
1:B:379:ASP:HB3	1:B:510:THR:HG22	1.99	0.45
1:D:362:THR:HG22	2:D:995:HOH:O	2.17	0.45
1:B:206:ARG:HA	1:B:234:GLY:HA2	1.99	0.45
1:A:65:TYR:CE1	1:A:306:ILE:HG13	2.52	0.45
1:C:260:ALA:CB	1:C:265:LEU:HB2	2.47	0.45
1:D:345:ARG:HD3	1:D:361:SER:HB3	1.99	0.45
1:D:277:ASP:OD1	1:D:282:HIS:HE1	1.99	0.45
1:A:434:VAL:CG1	1:A:487:GLN:HE22	2.30	0.44
1:B:227:TYR:HA	1:B:228:PRO:HD3	1.83	0.44
1:A:46:THR:HG22	1:A:340:GLN:HB3	1.99	0.44
1:D:203:GLY:HA2	1:D:235:ILE:CG2	2.42	0.44
1:B:223:ARG:HG2	1:B:223:ARG:HH11	1.83	0.44
1:A:51:TRP:HB2	1:A:75:ASN:HA	1.99	0.44
1:B:206:ARG:HH11	1:B:234:GLY:CA	2.23	0.44
1:D:110:GLY:HA3	1:D:130:LEU:O	2.18	0.44
1:B:169:GLY:HA2	1:B:181:ALA:CB	2.48	0.44
1:A:508:HIS:HB2	2:D:644:HOH:O	2.16	0.44
1:D:204:ARG:HB2	1:D:207:TYR:O	2.18	0.44
1:B:65:TYR:CE1	1:B:306:ILE:HG13	2.53	0.44
1:A:414:GLY:O	1:A:495:ASP:HB3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:206:ARG:HH11	1:C:234:GLY:HA3	1.80	0.43
1:B:260:ALA:HB1	1:B:265:LEU:HB2	1.99	0.43
1:D:144:GLN:HG2	1:D:164:ILE:HB	2.00	0.43
1:B:422:HIS:O	1:B:438:PRO:HB2	2.18	0.43
1:A:56:GLN:HB2	1:A:68:TYR:HB2	2.01	0.43
1:A:500:LEU:HD13	1:A:509:PHE:CG	2.54	0.43
1:D:57:ARG:HA	1:D:58:PRO:HD3	1.89	0.43
1:B:329:VAL:HG13	1:B:330:PRO:CD	2.41	0.43
1:A:470:ASP:HB3	2:D:822:HOH:O	2.18	0.42
1:A:110:GLY:HA3	1:A:130:LEU:O	2.18	0.42
1:A:57:ARG:HA	1:A:58:PRO:HD3	1.91	0.42
1:B:330:PRO:HD3	1:B:444:TYR:CD1	2.54	0.42
1:A:166:ASN:HD21	1:A:168:ASP:HB2	1.84	0.42
1:A:83:HIS:CD2	1:A:84:ALA:N	2.87	0.42
1:C:401:ILE:O	1:C:462:SER:HA	2.19	0.42
1:D:183:TRP:O	1:D:204:ARG:HD2	2.19	0.42
1:D:118:ASN:HB2	1:D:124:ALA:CA	2.49	0.42
1:D:265:LEU:HB3	1:D:266:PRO:CD	2.50	0.42
1:C:206:ARG:HD3	1:C:234:GLY:CA	2.49	0.42
1:D:382:VAL:HG13	1:D:509:PHE:CE2	2.55	0.42
1:C:132:THR:HG21	1:C:185:ARG:HB3	2.01	0.42
1:A:257:SER:HB3	1:A:298:TYR:CE2	2.55	0.41
1:C:235:ILE:HG23	1:C:256:ALA:CB	2.49	0.41
1:B:377:LEU:HD12	1:B:377:LEU:N	2.35	0.41
1:A:415:ARG:HA	1:A:421:ARG:O	2.20	0.41
1:A:205:LEU:CD2	1:A:206:ARG:HG2	2.50	0.41
1:A:446:LEU:HD21	1:A:490:HIS:CD2	2.55	0.41
1:B:470:ASP:HB3	2:D:757:HOH:O	2.19	0.41
1:C:260:ALA:HB1	1:C:265:LEU:HB2	2.02	0.41
1:D:98:VAL:HG13	1:D:156:PHE:CD1	2.55	0.41
1:B:169:GLY:HA2	1:B:181:ALA:HB1	2.02	0.41
1:B:132:THR:HG21	1:B:185:ARG:HB3	2.00	0.41
1:B:205:LEU:O	1:B:207:TYR:HD2	2.03	0.41
1:A:349:LEU:HG	1:A:357:TYR:HB3	2.01	0.41
1:A:329:VAL:O	1:A:332:ASP:HB2	2.21	0.41
1:B:300:ALA:HB2	1:B:317:ALA:HB2	2.03	0.41
1:B:83:HIS:CD2	1:B:84:ALA:N	2.89	0.41
1:B:329:VAL:O	1:B:332:ASP:HB2	2.21	0.41
1:C:229:ASN:HB3	1:C:232:LEU:HD22	2.03	0.41
1:D:341:ASN:HD22	1:D:341:ASN:HA	1.76	0.41
1:C:470:ASP:HB3	2:D:684:HOH:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:102:ARG:HD3	2:D:1081:HOH:O	2.20	0.41
1:C:58:PRO:HB3	1:C:67:LEU:HA	2.01	0.41
1:A:329:VAL:HG12	1:A:331:THR:H	1.86	0.41
1:B:432:LEU:HA	1:B:432:LEU:HD22	1.87	0.41
1:A:412:SER:OG	1:A:424:ASN:ND2	2.54	0.41
1:B:230:HIS:C	1:B:232:LEU:H	2.25	0.40
1:A:394:ALA:HA	1:A:469:VAL:O	2.22	0.40
1:A:341:ASN:HA	1:A:341:ASN:HD22	1.69	0.40
1:A:206:ARG:HD3	1:A:234:GLY:CA	2.47	0.40
1:D:401:ILE:O	1:D:462:SER:HA	2.22	0.40
1:D:272:TRP:CE2	1:D:288:PRO:HB3	2.56	0.40
1:B:57:ARG:HA	1:B:58:PRO:HD3	1.89	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	478/492 (97%)	454 (95%)	23 (5%)	1 (0%)	52	59
1	B	478/492 (97%)	454 (95%)	23 (5%)	1 (0%)	52	59
1	C	478/492 (97%)	454 (95%)	24 (5%)	0	100	100
1	D	476/492 (97%)	454 (95%)	20 (4%)	2 (0%)	39	42
All	All	1910/1968 (97%)	1816 (95%)	90 (5%)	4 (0%)	52	59

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	235	ILE
1	D	353	PRO
1	A	235	ILE

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Mol	Chain	Res	Type
1	B	235	ILE

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	381/393 (97%)	374 (98%)	7 (2%)	66	79
1	B	381/393 (97%)	377 (99%)	4 (1%)	82	91
1	C	381/393 (97%)	375 (98%)	6 (2%)	70	82
1	D	380/393 (97%)	370 (97%)	10 (3%)	54	66
All	All	1523/1572 (97%)	1496 (98%)	27 (2%)	66	79

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	98	VAL
1	A	144	GLN
1	A	166	ASN
1	A	205	LEU
1	A	232	LEU
1	A	297	TRP
1	A	382	VAL
1	B	104	ASP
1	B	297	TRP
1	B	320	ASN
1	B	487	GLN
1	C	232	LEU
1	C	297	TRP
1	C	340	GLN
1	C	382	VAL
1	C	446	LEU
1	C	487	GLN
1	D	232	LEU
1	D	250	ARG
1	D	279	GLU

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Mol	Chain	Res	Type
1	D	297	TRP
1	D	329	VAL
1	D	353	PRO
1	D	382	VAL
1	D	446	LEU
1	D	448	PRO
1	D	487	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	HIS
1	A	54	ASN
1	A	62	HIS
1	A	133	GLN
1	A	144	GLN
1	A	166	ASN
1	A	289	GLN
1	A	341	ASN
1	A	424	ASN
1	A	479	ASN
1	A	487	GLN
1	A	508	HIS
1	B	44	HIS
1	B	54	ASN
1	B	62	HIS
1	B	66	GLN
1	B	75	ASN
1	B	144	GLN
1	B	190	HIS
1	B	280	GLN
1	B	282	HIS
1	B	320	ASN
1	B	341	ASN
1	B	391	ASN
1	B	424	ASN
1	B	479	ASN
1	B	487	GLN
1	C	62	HIS
1	C	71	HIS
1	C	75	ASN
1	C	133	GLN

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Mol	Chain	Res	Type
1	C	144	GLN
1	C	282	HIS
1	C	289	GLN
1	C	341	ASN
1	C	424	ASN
1	C	479	ASN
1	D	71	HIS
1	D	133	GLN
1	D	144	GLN
1	D	282	HIS
1	D	289	GLN
1	D	341	ASN
1	D	424	ASN
1	D	479	ASN
1	D	482	HIS
1	D	487	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

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
2	SUC	D	601	-	24,24,24	1.59	4 (16%)	36,36,36	1.11	4 (11%)
2	SUC	D	602	-	24,24,24	1.56	4 (16%)	36,36,36	1.12	2 (5%)
2	SUC	D	603	-	24,24,24	1.60	5 (20%)	36,36,36	1.18	4 (11%)
2	SUC	D	604	-	24,24,24	1.56	4 (16%)	36,36,36	1.25	5 (13%)
2	SUC	D	605	-	24,24,24	1.62	4 (16%)	36,36,36	1.18	3 (8%)
2	SUC	D	606	-	24,24,24	1.59	6 (25%)	36,36,36	1.07	2 (5%)

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
2	SUC	D	601	-	-	0/12/51/51	0/2/2/2
2	SUC	D	602	-	-	0/12/51/51	0/2/2/2
2	SUC	D	603	-	-	0/12/51/51	0/2/2/2
2	SUC	D	604	-	-	0/12/51/51	0/2/2/2
2	SUC	D	605	-	-	0/12/51/51	0/2/2/2
2	SUC	D	606	-	-	0/12/51/51	0/2/2/2

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	605	SUC	C4-C3	-4.30	1.41	1.52
2	D	601	SUC	C4-C3	-4.19	1.41	1.52
2	D	602	SUC	C4-C3	-4.10	1.41	1.52
2	D	606	SUC	C4-C3	-3.98	1.41	1.52
2	D	604	SUC	C4-C3	-3.96	1.42	1.52
2	D	603	SUC	C4-C3	-3.95	1.42	1.52
2	D	602	SUC	O2-C2	-3.34	1.35	1.43
2	D	603	SUC	O2-C2	-3.24	1.35	1.43
2	D	604	SUC	O2-C2	-3.20	1.35	1.43
2	D	605	SUC	O2-C2	-3.13	1.35	1.43
2	D	601	SUC	O2-C2	-3.11	1.35	1.43
2	D	606	SUC	O2-C2	-2.95	1.35	1.43
2	D	605	SUC	C3-C2	-2.32	1.46	1.52
2	D	606	SUC	O4'-C4'	-2.23	1.37	1.43
2	D	606	SUC	C3-C2	-2.20	1.46	1.52
2	D	603	SUC	O2'-C5'	-2.19	1.38	1.43
2	D	606	SUC	O2'-C5'	-2.19	1.38	1.43
2	D	602	SUC	C3-C2	-2.18	1.46	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	603	SUC	C3-C2	-2.16	1.46	1.52
2	D	605	SUC	O4'-C4'	-2.12	1.37	1.43
2	D	603	SUC	O4'-C4'	-2.07	1.38	1.43
2	D	602	SUC	O4'-C4'	-2.06	1.38	1.43
2	D	604	SUC	C3-C2	-2.02	1.47	1.52
2	D	601	SUC	C3-C2	-2.02	1.47	1.52
2	D	606	SUC	O5-C5	2.02	1.49	1.44
2	D	601	SUC	O5-C5	2.02	1.49	1.44
2	D	604	SUC	O5-C5	2.06	1.49	1.44

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	603	SUC	C2'-O1-C1	-3.49	108.34	117.53
2	D	606	SUC	O3'-C3'-C4'	-2.67	103.71	113.29
2	D	605	SUC	C2'-O1-C1	-2.58	110.72	117.53
2	D	604	SUC	O3-C3-C4	-2.28	105.19	110.34
2	D	603	SUC	O3'-C3'-C4'	-2.18	105.47	113.29
2	D	602	SUC	C2'-O1-C1	-2.15	111.86	117.53
2	D	601	SUC	C2'-O1-C1	-2.10	111.98	117.53
2	D	604	SUC	C2'-O1-C1	-2.10	112.00	117.53
2	D	601	SUC	O3'-C3'-C4'	-2.09	105.79	113.29
2	D	602	SUC	O3'-C3'-C4'	-2.03	105.98	113.29
2	D	605	SUC	O2'-C2'-C1'	2.01	113.45	107.98
2	D	603	SUC	O1-C1-C2	2.04	115.17	108.36
2	D	601	SUC	O5-C5-C6	2.14	111.76	106.36
2	D	601	SUC	O2'-C2'-C1'	2.15	113.83	107.98
2	D	604	SUC	O2'-C2'-C1'	2.16	113.85	107.98
2	D	606	SUC	O5-C5-C6	2.21	111.94	106.36
2	D	604	SUC	C3-C4-C5	2.22	114.07	110.20
2	D	605	SUC	C3-C4-C5	2.43	114.43	110.20
2	D	604	SUC	O5-C5-C6	2.49	112.64	106.36
2	D	603	SUC	O5-C5-C4	2.56	114.49	109.68

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	601	SUC	4	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	602	SUC	3	0
2	D	603	SUC	3	0
2	D	605	SUC	1	0
2	D	606	SUC	1	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	480/492 (97%)	-0.35	5 (1%) 84 83	18, 28, 42, 66	0
1	B	480/492 (97%)	-0.35	7 (1%) 76 75	17, 28, 42, 63	0
1	C	480/492 (97%)	-0.49	2 (0%) 93 93	17, 26, 39, 65	0
1	D	478/492 (97%)	-0.53	1 (0%) 95 95	17, 26, 36, 45	0
All	All	1918/1968 (97%)	-0.43	15 (0%) 87 87	17, 27, 39, 66	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	520	ALA	7.9
1	C	520	ALA	7.6
1	B	520	ALA	5.5
1	A	519	GLN	3.9
1	B	173	THR	3.7
1	A	174	THR	3.6
1	B	172	ALA	3.2
1	C	234	GLY	3.0
1	A	173	THR	2.9
1	B	174	THR	2.9
1	A	234	GLY	2.8
1	D	235	ILE	2.5
1	B	519	GLN	2.4
1	B	518	GLY	2.4
1	B	171	ALA	2.3

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 carbohydrates 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. 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	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	SUC	D	601	23/23	0.67	0.25	5.91	57,62,66,67	0
2	SUC	D	605	23/23	0.93	0.15	3.30	28,37,47,53	0
2	SUC	D	602	23/23	0.67	0.21	2.55	54,60,64,65	0
2	SUC	D	606	23/23	0.94	0.13	0.44	23,32,44,51	0
2	SUC	D	603	23/23	0.95	0.09	-0.27	23,33,43,47	0
2	SUC	D	604	23/23	0.96	0.08	-0.34	25,33,43,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.