



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 10:17 AM UTC

PDB ID : 2PWG / pdb_00002pwg
Title : Crystal Structure of the Trehalulose Synthase MutB From Pseudomonas Mesoacidophila MX-45 Complexed to the Inhibitor Castanospermine
Authors : Ravaud, S.; Robert, X.; Haser, R.; Aghajari, N.
Deposited on : 2007-05-11
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

<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

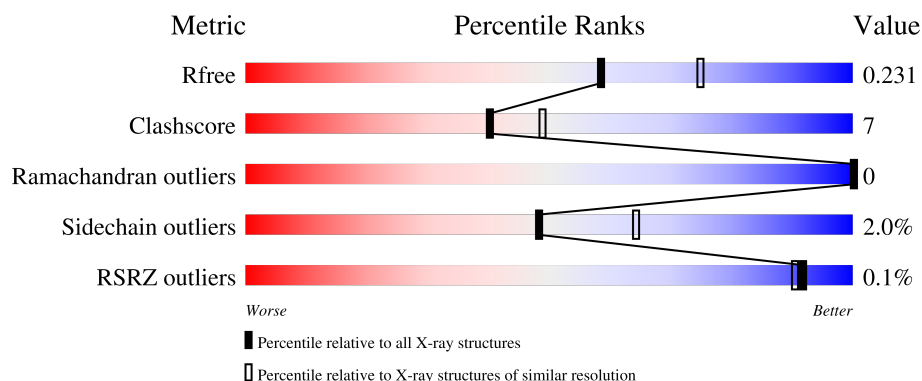
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (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 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	556	 83% 16% .
1	B	556	 84% 15% .

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 10449 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.

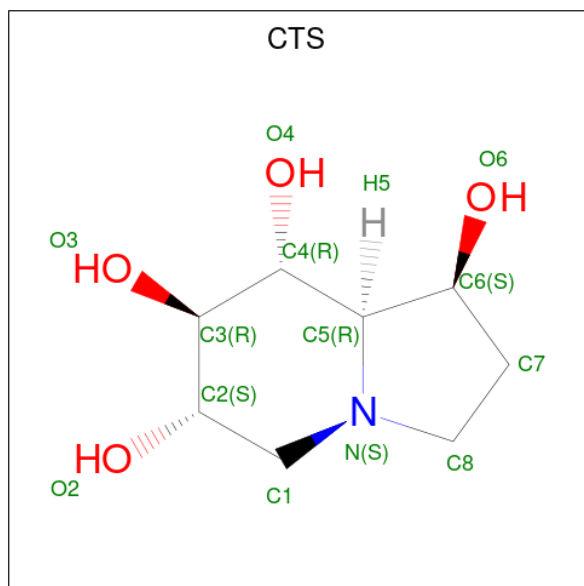
- Molecule 1 is a protein called Sucrose isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	556	Total	C	N	O	S	0	6	0
			4540	2900	768	860	12			
1	B	556	Total	C	N	O	S	0	3	0
			4514	2881	765	856	12			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		
2	B	1	Total	Ca	0	0
			1	1		

- Molecule 3 is CASTANOSPERMINE (CCD ID: CTS) (formula: C₈H₁₅NO₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			13	8	1	4		
3	B	1	Total	C	N	O	0	0
			13	8	1	4		

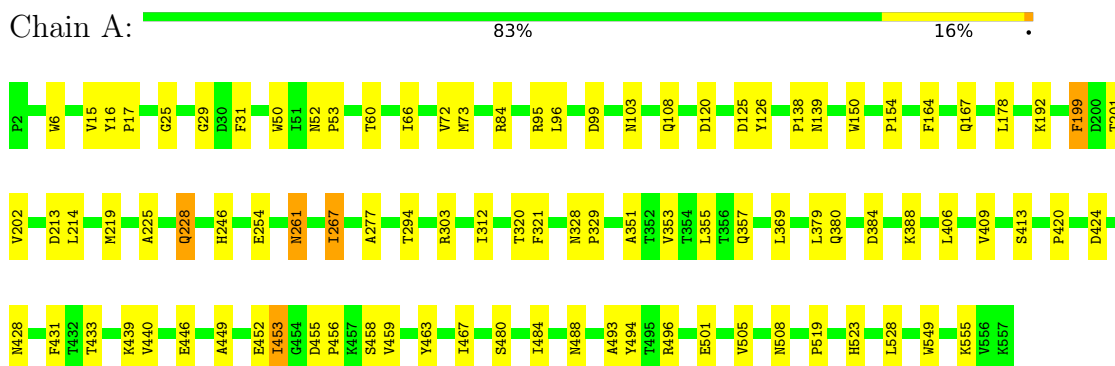
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	690	Total	O	0	0
			690	690		
4	B	677	Total	O	0	0
			677	677		

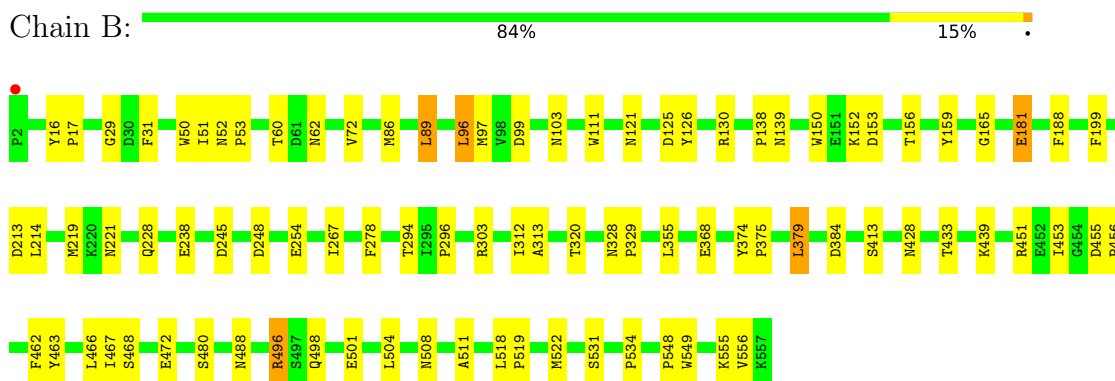
3 Residue-property plots

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: Sucrose isomerase



- Molecule 1: Sucrose isomerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.24Å 85.96Å 123.52Å 90.00° 99.20° 90.00°	Depositor
Resolution (Å)	19.91 – 2.20 19.91 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.9 (19.91-2.20) 99.8 (19.91-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.60 (at 2.19Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.186 , 0.230 0.188 , 0.231	Depositor DCC
R_{free} test set	6823 reflections (10.13%)	wwPDB-VP
Wilson B-factor (Å ²)	18.1	Xtriage
Anisotropy	0.444	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 56.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.55$, $\langle L^2 \rangle = 0.39$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10449	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 48.63 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.3685e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

Bond lengths and bond angles in the following residue types are not validated in this section: CA, CTS

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	0.64	5/4676 (0.1%)	1.02	18/6357 (0.3%)
1	B	0.60	4/4650 (0.1%)	1.01	16/6321 (0.3%)
All	All	0.62	9/9326 (0.1%)	1.01	34/12678 (0.3%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	428	ASN	CG-OD1	5.81	1.34	1.23
1	A	428	ASN	CG-OD1	5.79	1.34	1.23
1	B	121	ASN	CG-OD1	5.74	1.34	1.23
1	A	167	GLN	CD-OE1	5.69	1.34	1.23
1	A	108	GLN	CD-OE1	5.66	1.34	1.23
1	A	380	GLN	CD-OE1	5.34	1.33	1.23
1	B	221	ASN	CG-OD1	5.28	1.33	1.23
1	A	246	HIS	ND1-CE1	5.22	1.37	1.32
1	B	498	GLN	CD-OE1	5.10	1.33	1.23

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	320	THR	N-CA-C	-7.81	98.31	110.42
1	B	267	ILE	N-CA-C	7.70	121.82	113.43
1	B	126	TYR	N-CA-C	-7.61	103.98	113.18
1	B	320	THR	N-CA-C	-7.39	98.97	110.42
1	A	103	ASN	N-CA-C	-7.33	103.29	112.90
1	B	103	ASN	N-CA-C	-7.25	103.40	112.90
1	A	126	TYR	N-CA-C	-6.68	105.12	113.28
1	B	453	ILE	N-CA-C	6.58	117.33	110.62
1	A	246	HIS	CB-CG-CD2	-6.56	122.67	131.20
1	B	384	ASP	N-CA-C	6.53	120.09	110.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	267	ILE	N-CA-C	6.49	120.50	113.43
1	A	60	THR	N-CA-C	-6.16	105.02	112.54
1	B	199	PHE	N-CA-C	6.00	118.47	108.20
1	B	153	ASP	CA-C-N	5.95	126.10	119.32
1	B	153	ASP	C-N-CA	5.95	126.10	119.32
1	A	66	ILE	N-CA-C	5.82	116.96	108.58
1	A	246	HIS	CB-CG-ND1	5.70	131.25	122.70
1	A	384	ASP	N-CA-C	5.64	118.53	110.24
1	A	453	ILE	N-CA-C	5.62	116.26	110.36
1	B	16	TYR	N-CA-C	-5.61	97.41	109.81
1	B	433	THR	N-CA-C	-5.59	106.50	113.38
1	A	16	TYR	N-CA-C	-5.55	97.54	109.81
1	B	60	THR	N-CA-C	-5.49	105.84	112.54
1	A	433	THR	N-CA-C	-5.46	106.29	113.17
1	A	413	SER	N-CA-C	5.42	117.88	110.24
1	B	555	LYS	N-CA-C	-5.42	100.46	109.46
1	A	120	ASP	N-CA-C	5.36	119.12	112.59
1	A	199	PHE	N-CA-C	5.35	117.35	108.20
1	A	25	GLY	N-CA-C	5.32	122.97	115.30
1	B	413	SER	N-CA-C	5.32	118.03	110.10
1	A	555	LYS	N-CA-C	-5.17	100.81	109.24
1	A	388	LYS	N-CA-C	-5.08	105.75	111.28
1	B	496	ARG	N-CA-C	-5.05	100.22	108.76
1	B	165	GLY	N-CA-C	5.02	117.70	111.93

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	4540	0	4282	60	0
1	B	4514	0	4250	58	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	13	0	15	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	13	0	15	1	0
4	A	690	0	0	6	0
4	B	677	0	0	11	0
All	All	10449	0	8562	117	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 (117) 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:B:522:MET:HE3	1:B:556:VAL:HG11	1.39	0.99
1:A:31:PHE:HE2	1:A:72[A]:VAL:HG11	1.38	0.88
1:A:125:ASP:HB3	4:A:6018:HOH:O	1.77	0.85
1:B:125:ASP:HB2	4:B:5889:HOH:O	1.78	0.82
1:B:31:PHE:HE2	1:B:72:VAL:HG11	1.45	0.81
1:A:214:LEU:HB2	1:A:219:MET:HE2	1.64	0.80
1:B:130:ARG:NH2	4:B:4839:HOH:O	2.18	0.75
1:A:31:PHE:HE2	1:A:72[A]:VAL:CG1	2.00	0.75
1:B:31:PHE:HE2	1:B:72:VAL:CG1	2.01	0.74
1:A:199:PHE:CB	1:A:202[A]:VAL:HG12	2.17	0.73
1:A:449:ALA:O	1:A:453:ILE:HG12	1.89	0.72
1:B:214:LEU:HB2	1:B:219:MET:HE2	1.72	0.72
1:B:111:TRP:CZ2	1:B:181:GLU:HG2	2.25	0.70
1:B:238:GLU:HA	4:B:4865:HOH:O	1.91	0.69
1:A:199:PHE:HB2	1:A:202[A]:VAL:CG1	2.23	0.68
1:B:214:LEU:HB2	1:B:219:MET:CE	2.25	0.67
1:A:439:LYS:HE3	4:A:4118:HOH:O	1.96	0.66
1:A:214:LEU:HB2	1:A:219:MET:CE	2.26	0.66
1:A:199:PHE:O	1:A:202[A]:VAL:HG13	1.96	0.65
1:A:199:PHE:HB2	1:A:202[A]:VAL:HG12	1.77	0.64
1:B:355[B]:LEU:HD22	1:B:466:LEU:HB3	1.80	0.64
1:B:374:TYR:OH	1:B:439:LYS:HE2	1.98	0.63
1:A:199:PHE:HB3	1:A:202[A]:VAL:HG12	1.80	0.62
1:A:294:THR:HG23	4:A:4619:HOH:O	1.99	0.62
1:A:452:GLU:OE1	1:A:459[A]:VAL:HG22	2.01	0.60
1:B:355[B]:LEU:HD21	1:B:467:ILE:HG13	1.84	0.59
1:A:84:ARG:HB2	1:A:84:ARG:NH1	2.18	0.59
1:B:468:SER:O	1:B:472:GLU:HG3	2.04	0.58
1:B:451:ARG:HD2	4:B:4668:HOH:O	2.05	0.57
1:A:84:ARG:HB2	1:A:84:ARG:CZ	2.35	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:254:GLU:OE1	3:A:8000:CTS:H81	2.04	0.57
1:A:523:HIS:HD2	4:A:6055:HOH:O	1.89	0.56
1:B:355[B]:LEU:HD21	1:B:467:ILE:CG1	2.36	0.56
1:A:31:PHE:CE2	1:A:72[A]:VAL:HG11	2.30	0.55
1:A:452:GLU:HB3	1:A:458:SER:HB2	1.89	0.54
1:B:125:ASP:HB2	4:B:4824:HOH:O	2.07	0.53
1:A:406[A]:LEU:C	1:A:406[A]:LEU:HD13	2.33	0.53
1:A:369:LEU:C	1:A:459[A]:VAL:HG21	2.34	0.53
1:A:138:PRO:HD2	1:A:150:TRP:CG	2.45	0.52
1:A:15:VAL:O	1:A:17:PRO:HD3	2.10	0.51
1:B:355[B]:LEU:HD22	1:B:466:LEU:CB	2.41	0.51
1:B:504:LEU:HD13	1:B:522:MET:HE1	1.91	0.51
1:B:294[B]:THR:HG23	4:B:4879:HOH:O	2.12	0.50
1:B:31:PHE:CE2	1:B:72:VAL:HG11	2.36	0.50
1:B:504:LEU:HD22	1:B:522:MET:CE	2.42	0.50
1:B:294[A]:THR:HG22	4:B:4879:HOH:O	2.11	0.50
1:A:328:ASN:HB3	1:A:329:PRO:HD2	1.94	0.49
1:B:480:SER:O	1:B:496:ARG:HA	2.11	0.49
1:B:462:PHE:CZ	1:B:531:SER:HB3	2.47	0.49
1:B:508:ASN:O	1:B:549:TRP:HA	2.13	0.48
1:B:52:ASN:HB3	1:B:53:PRO:HD2	1.96	0.48
1:A:480:SER:O	1:A:496:ARG:HA	2.14	0.47
1:B:504:LEU:HD22	1:B:522:MET:HE2	1.96	0.47
1:B:522:MET:CE	1:B:556:VAL:HG11	2.28	0.47
1:B:312:ILE:C	1:B:312:ILE:HD12	2.40	0.47
1:A:484:ILE:HD12	1:A:519:PRO:HG2	1.97	0.47
1:B:152:LYS:HB2	1:B:159:TYR:CE2	2.49	0.47
1:B:511:ALA:C	1:B:548:PRO:HB3	2.40	0.46
1:A:261:ASN:HD22	1:A:261:ASN:C	2.23	0.46
1:B:138:PRO:HD2	1:B:150:TRP:CG	2.51	0.45
1:A:72[B]:VAL:HG22	4:A:4196:HOH:O	2.15	0.45
1:A:424:ASP:HB2	4:A:4325:HOH:O	2.17	0.45
1:B:139:ASN:HB2	1:B:213:ASP:OD1	2.15	0.45
1:B:303:ARG:NH1	1:B:488:ASN:HB3	2.32	0.45
1:B:86:MET:SD	1:B:96[B]:LEU:HD23	2.57	0.45
1:A:420:PRO:HB3	1:A:431:PHE:CG	2.52	0.44
1:B:455:ASP:HA	1:B:456:PRO:HD2	1.89	0.44
1:B:374:TYR:CD1	1:B:375:PRO:HD2	2.52	0.44
1:A:312:ILE:HD12	1:A:312:ILE:C	2.42	0.44
1:A:303:ARG:NH1	1:A:488:ASN:HB3	2.33	0.44
1:A:139:ASN:HB2	1:A:213:ASP:OD1	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:508:ASN:O	1:A:549:TRP:HA	2.18	0.44
1:A:321:PHE:CZ	1:A:357:GLN:HG3	2.53	0.44
1:B:254:GLU:OE1	3:B:8001:CTS:H81	2.18	0.44
1:B:96[A]:LEU:HD22	1:B:97:MET:N	2.33	0.43
1:B:534:PRO:HG2	4:B:5842:HOH:O	2.18	0.43
1:A:199:PHE:O	1:A:202[A]:VAL:CG1	2.66	0.43
1:A:31:PHE:CE2	1:A:72[A]:VAL:CG1	2.91	0.43
1:A:178:LEU:C	1:A:178:LEU:HD13	2.44	0.43
1:A:50:TRP:CD1	1:A:50:TRP:C	2.97	0.42
1:A:154:PRO:HB3	1:B:156:THR:HA	2.00	0.42
1:B:111:TRP:HZ2	1:B:181:GLU:HG2	1.79	0.42
1:B:379:LEU:HD12	1:B:379:LEU:HA	1.81	0.42
1:A:52:ASN:HB3	1:A:53:PRO:HD2	2.00	0.42
1:B:50:TRP:CD1	1:B:50:TRP:C	2.96	0.42
1:B:62:ASN:HB2	4:B:4090:HOH:O	2.20	0.42
1:B:296:PRO:HG2	4:B:4230:HOH:O	2.19	0.42
1:B:31:PHE:CE2	1:B:72:VAL:CG1	2.91	0.42
1:B:52:ASN:HB3	1:B:53:PRO:CD	2.50	0.42
1:B:188:PHE:HA	4:B:5060:HOH:O	2.19	0.42
1:A:199:PHE:HB2	1:A:202[A]:VAL:HG11	2.00	0.42
1:A:420:PRO:HB2	1:A:440:VAL:HG22	2.02	0.42
1:A:353:VAL:O	1:A:357:GLN:HG2	2.20	0.42
1:A:369:LEU:O	1:A:459[A]:VAL:HG21	2.20	0.42
1:A:452:GLU:HB3	1:A:458:SER:CB	2.50	0.42
1:A:455:ASP:HA	1:A:456:PRO:HD2	1.93	0.41
1:A:72[A]:VAL:HG12	1:A:73:MET:N	2.36	0.41
1:A:164:PHE:CE2	3:A:8000:CTS:H12	2.55	0.41
1:A:484:ILE:HG12	1:A:494:TYR:HA	2.01	0.41
1:A:267:ILE:HG21	1:A:277:ALA:HB2	2.03	0.41
1:B:51:ILE:O	1:B:51:ILE:HG13	2.21	0.41
1:B:89:LEU:HD23	1:B:96[B]:LEU:HB2	2.02	0.41
1:A:355[A]:LEU:HD11	1:A:467:ILE:HG13	2.02	0.41
1:B:152:LYS:HE3	1:B:159:TYR:CZ	2.56	0.41
1:A:225:ALA:HA	1:A:228:GLN:HG3	2.03	0.41
1:A:351:ALA:HB2	1:A:463:TYR:CE2	2.55	0.41
1:B:17:PRO:O	1:B:29:GLY:HA3	2.21	0.41
1:B:248:ASP:O	1:B:248:ASP:CG	2.64	0.41
1:B:312:ILE:HD12	1:B:313:ALA:N	2.36	0.41
1:A:528:LEU:HD12	1:A:528:LEU:N	2.36	0.41
1:A:17:PRO:O	1:A:29:GLY:HA3	2.20	0.40
1:A:201:THR:O	1:A:201:THR:HG23	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:328:ASN:HB3	1:B:329:PRO:HD2	2.03	0.40
1:A:6:TRP:CH2	1:A:95:ARG:HD2	2.56	0.40
1:B:254:GLU:HA	1:B:278:PHE:HB2	2.04	0.40
1:A:493:ALA:HA	1:A:505:VAL:O	2.21	0.40
1:B:368:GLU:HA	1:B:463:TYR:CD1	2.57	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	560/556 (101%)	536 (96%)	24 (4%)	0	100	100
1	B	557/556 (100%)	535 (96%)	22 (4%)	0	100	100
All	All	1117/1112 (100%)	1071 (96%)	46 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	475/474 (100%)	465 (98%)	10 (2%)	47	63
1	B	471/474 (99%)	460 (98%)	11 (2%)	44	59
All	All	946/948 (100%)	925 (98%)	21 (2%)	48	61

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	96[A]	LEU
1	A	96[B]	LEU
1	A	99	ASP
1	A	192	LYS
1	A	228	GLN
1	A	261	ASN
1	A	379	LEU
1	A	409	VAL
1	A	446	GLU
1	A	501	GLU
1	B	89	LEU
1	B	96[A]	LEU
1	B	96[B]	LEU
1	B	99	ASP
1	B	181	GLU
1	B	228	GLN
1	B	245	ASP
1	B	379	LEU
1	B	501	GLU
1	B	518	LEU
1	B	519	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	108	GLN
1	A	221	ASN
1	A	261	ASN
1	A	523	HIS
1	A	550	GLN
1	B	14	GLN
1	B	52	ASN
1	B	108	GLN
1	B	262	GLN
1	B	380	GLN
1	B	547	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	CTS	A	8000	-	14,14,14	1.70	3 (21%)	17,21,21	2.24	5 (29%)
3	CTS	B	8001	-	14,14,14	1.70	3 (21%)	17,21,21	1.96	3 (17%)

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	CTS	A	8000	-	-	-	0/2/2/2
3	CTS	B	8001	-	-	-	0/2/2/2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	8000	CTS	C2-C3	4.89	1.60	1.52
3	B	8001	CTS	C6-C5	3.34	1.57	1.54
3	B	8001	CTS	C2-C3	3.32	1.57	1.52
3	B	8001	CTS	C1-C2	-2.70	1.49	1.52
3	A	8000	CTS	C6-C5	2.21	1.56	1.54
3	A	8000	CTS	C1-C2	-2.19	1.49	1.52

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	8000	CTS	C1-N-C5	5.46	120.26	110.62
3	A	8000	CTS	C8-C7-C6	-4.72	96.42	103.95
3	B	8001	CTS	C1-N-C5	4.64	118.81	110.62
3	B	8001	CTS	C8-C7-C6	-4.24	97.18	103.95
3	B	8001	CTS	C2-C3-C4	3.01	116.15	110.86
3	A	8000	CTS	C2-C3-C4	3.00	116.14	110.86
3	A	8000	CTS	O3-C3-C4	-2.10	105.42	110.38
3	A	8000	CTS	C2-C1-N	2.04	114.51	110.72

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	8000	CTS	2	0
3	B	8001	CTS	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 [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	556/556 (100%)	-0.75	0 100 100	8, 18, 27, 37	6 (1%)
1	B	556/556 (100%)	-0.75	1 (0%) 91 90	8, 18, 27, 37	3 (0%)
All	All	1112/1112 (100%)	-0.75	1 (0%) 92 90	8, 18, 27, 37	9 (0%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	2	PRO	2.2

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](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	CTS	B	8001	13/13	0.92	0.09	15,18,20,21	0
3	CTS	A	8000	13/13	0.93	0.08	13,17,20,20	0
2	CA	A	7000	1/1	0.98	0.03	21,21,21,21	0
2	CA	B	7001	1/1	0.99	0.01	20,20,20,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.