



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 3WQE / pdb_00003wqe
Title : D-threo-3-hydroxyaspartate dehydratase from Delftia sp. HT23 complexed with D-allothreonine
Authors : Yasutake, Y.; Matsumoto, Y.; Wada, M.
Deposited on : 2014-01-25
Resolution : 1.60 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

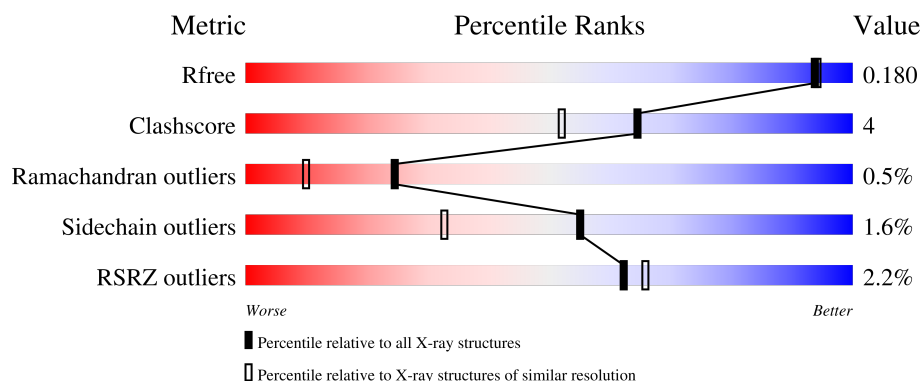
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (1.60-1.60)

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	390	83% 12% ..
1	B	390	4% 80% 17% .

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
4	2TL	A	403	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6505 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 D-threo-3-hydroxyaspartate dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	379	Total	C	N	O	S	0	6	0
			2862	1778	535	534	15			
1	B	389	Total	C	N	O	S	0	5	0
			2940	1825	554	544	17			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	GLY	-	expression tag	UNP B2DFG5
A	-8	HIS	-	expression tag	UNP B2DFG5
A	-7	HIS	-	expression tag	UNP B2DFG5
A	-6	HIS	-	expression tag	UNP B2DFG5
A	-5	HIS	-	expression tag	UNP B2DFG5
A	-4	HIS	-	expression tag	UNP B2DFG5
A	-3	HIS	-	expression tag	UNP B2DFG5
A	-2	ALA	-	expression tag	UNP B2DFG5
A	-1	MET	-	expression tag	UNP B2DFG5
A	0	SER	-	expression tag	UNP B2DFG5
B	-9	GLY	-	expression tag	UNP B2DFG5
B	-8	HIS	-	expression tag	UNP B2DFG5
B	-7	HIS	-	expression tag	UNP B2DFG5
B	-6	HIS	-	expression tag	UNP B2DFG5
B	-5	HIS	-	expression tag	UNP B2DFG5
B	-4	HIS	-	expression tag	UNP B2DFG5
B	-3	HIS	-	expression tag	UNP B2DFG5
B	-2	ALA	-	expression tag	UNP B2DFG5
B	-1	MET	-	expression tag	UNP B2DFG5
B	0	SER	-	expression tag	UNP B2DFG5

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: C₈H₁₀NO₆P).

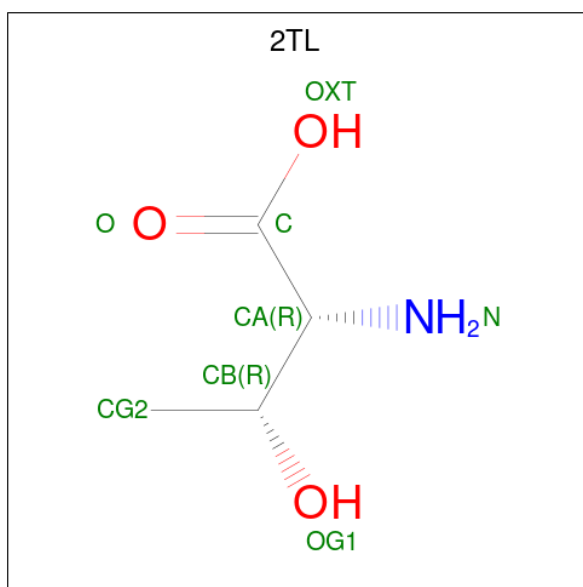


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 4 is D-allothreonine (CCD ID: 2TL) (formula: C₄H₉NO₃).



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

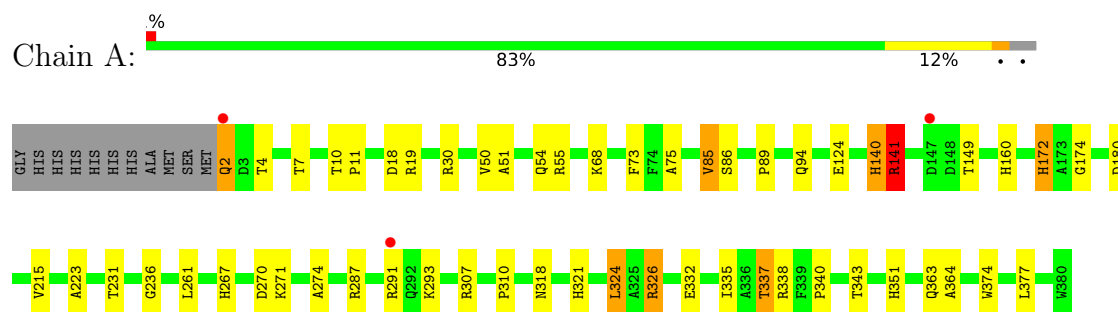
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	334	Total	O	0	0
			334	334		
5	B	320	Total	O	0	0
			320	320		

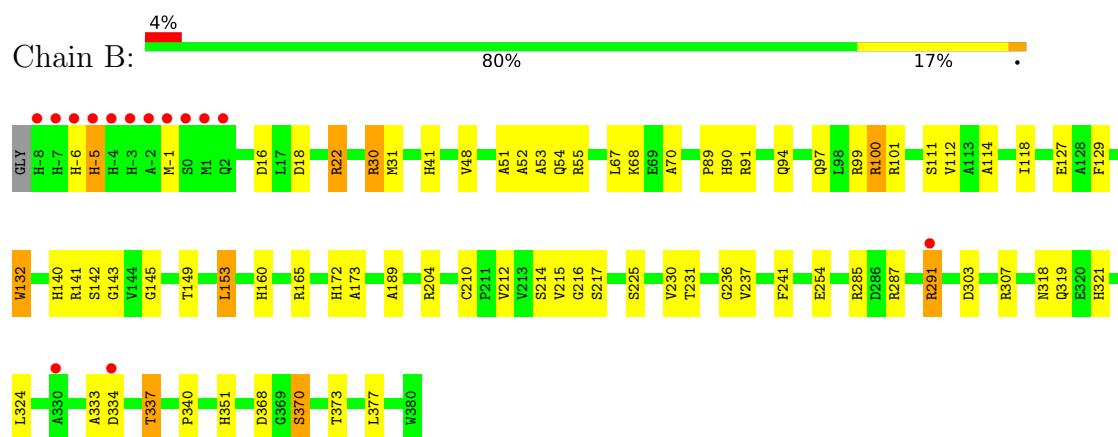
3 Residue-property plots [i](#)

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: D-threo-3-hydroxyaspartate dehydratase



- Molecule 1: D-threo-3-hydroxyaspartate dehydratase



4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	157.98Å 157.98Å 158.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	26.33 – 1.60 26.33 – 1.60	Depositor EDS
% Data completeness (in resolution range)	99.8 (26.33-1.60) 99.8 (26.33-1.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.10 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.166 , 0.182 0.163 , 0.180	Depositor DCC
R_{free} test set	6531 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	16.7	Xtriage
Anisotropy	0.019	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 45.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for -l,-k,-h 0.005 for -h,l,k	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6505	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.73% 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, 0.333 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: PLP, MG, 2TL

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	1.80	34/2927 (1.2%)	1.57	19/3970 (0.5%)
1	B	1.83	46/3008 (1.5%)	1.63	45/4078 (1.1%)
All	All	1.81	80/5935 (1.3%)	1.60	64/8048 (0.8%)

All (80) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	143	GLY	N-CA	12.48	1.56	1.45
1	B	172	HIS	ND1-CE1	11.30	1.43	1.32
1	B	141	ARG	CZ-NH2	10.64	1.47	1.33
1	A	351	HIS	ND1-CE1	9.96	1.42	1.32
1	B	351	HIS	CE1-NE2	9.40	1.42	1.32
1	A	307	ARG	N-CA	9.20	1.57	1.46
1	A	318	ASN	CG-OD1	-8.62	1.07	1.23
1	A	267	HIS	ND1-CE1	7.96	1.40	1.32
1	B	118	ILE	CA-CB	-7.86	1.45	1.54
1	A	351	HIS	CD2-NE2	7.84	1.46	1.37
1	B	285	ARG	CZ-NH2	7.65	1.43	1.33
1	A	172	HIS	CG-CD2	7.64	1.44	1.35
1	B	141	ARG	C-O	7.55	1.33	1.24
1	B	216	GLY	N-CA	7.54	1.52	1.45
1	B	351	HIS	CD2-NE2	7.50	1.46	1.37
1	A	338	ARG	CZ-NH2	7.43	1.43	1.33
1	A	172	HIS	ND1-CE1	7.14	1.39	1.32
1	B	145	GLY	N-CA	6.95	1.52	1.45
1	B	142	SER	CA-C	6.90	1.62	1.52
1	A	364	ALA	N-CA	6.87	1.54	1.46
1	A	141	ARG	CZ-NH1	6.82	1.42	1.32
1	A	321	HIS	CG-CD2	6.79	1.43	1.35
1	B	55	ARG	CZ-NH2	6.71	1.42	1.33
1	B	172	HIS	CG-CD2	6.65	1.43	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	141	ARG	CZ-NH2	6.46	1.41	1.33
1	B	30	ARG	CZ-NH2	6.38	1.41	1.33
1	B	204	ARG	CD-NE	-6.35	1.37	1.46
1	A	140	HIS	ND1-CE1	6.26	1.38	1.32
1	A	141	ARG	N-CA	6.25	1.54	1.46
1	B	-6	HIS	CG-CD2	6.24	1.42	1.35
1	A	19	ARG	CZ-NH1	6.18	1.41	1.32
1	A	287	ARG	CZ-NH2	6.14	1.41	1.33
1	B	-6	HIS	CE1-NE2	6.11	1.38	1.32
1	B	41	HIS	CG-CD2	6.10	1.42	1.35
1	A	343	THR	N-CA	5.99	1.53	1.46
1	A	335	ILE	CA-CB	5.98	1.63	1.54
1	A	174	GLY	C-O	5.92	1.30	1.24
1	B	97	GLN	N-CA	5.83	1.53	1.46
1	A	73	PHE	N-CA	-5.79	1.39	1.46
1	B	214	SER	N-CA	5.79	1.53	1.46
1	A	363	GLN	CD-NE2	5.74	1.45	1.33
1	B	141	ARG	NE-CZ	5.70	1.39	1.33
1	B	141	ARG	CB-CG	5.67	1.69	1.52
1	B	373	THR	N-CA	5.67	1.53	1.46
1	B	99	ARG	N-CA	5.57	1.53	1.46
1	B	70	ALA	N-CA	5.55	1.53	1.46
1	B	160	HIS	ND1-CE1	5.55	1.38	1.32
1	A	374	TRP	N-CA	5.47	1.53	1.46
1	B	142	SER	C-O	5.46	1.30	1.24
1	B	215	VAL	N-CA	5.45	1.52	1.46
1	B	48	VAL	N-CA	5.43	1.53	1.46
1	B	101	ARG	CA-CB	5.43	1.63	1.53
1	A	324	LEU	N-CA	5.43	1.53	1.46
1	A	85	VAL	N-CA	5.38	1.52	1.46
1	B	132	TRP	CA-C	-5.38	1.46	1.52
1	A	215	VAL	N-CA	5.35	1.52	1.46
1	B	-5	HIS	CG-CD2	5.32	1.41	1.35
1	B	129	PHE	C-O	5.31	1.30	1.23
1	B	230	VAL	C-O	5.30	1.31	1.24
1	B	337	THR	CA-CB	5.30	1.61	1.53
1	B	112	VAL	CA-C	5.27	1.59	1.52
1	A	160	HIS	ND1-CE1	5.25	1.37	1.32
1	B	231	THR	CB-OG1	-5.25	1.35	1.43
1	B	142	SER	N-CA	-5.23	1.39	1.46
1	B	318	ASN	CG-OD1	-5.20	1.13	1.23
1	A	19	ARG	C-O	-5.18	1.18	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	338	ARG	N-CA	-5.15	1.39	1.46
1	B	321	HIS	ND1-CE1	5.14	1.37	1.32
1	A	223	ALA	C-O	5.13	1.30	1.24
1	B	114	ALA	N-CA	5.13	1.52	1.46
1	A	307	ARG	CD-NE	5.11	1.53	1.46
1	B	340	PRO	CA-C	5.10	1.58	1.52
1	B	173	ALA	CA-CB	5.08	1.59	1.52
1	B	90	HIS	CE1-NE2	5.07	1.37	1.32
1	B	41	HIS	ND1-CE1	5.06	1.37	1.32
1	A	326	ARG	N-CA	5.05	1.52	1.46
1	A	340	PRO	N-CD	5.05	1.54	1.47
1	A	7	THR	N-CA	5.04	1.52	1.46
1	A	337	THR	N-CA	5.04	1.52	1.46
1	B	141	ARG	CZ-NH1	5.00	1.39	1.32

All (64) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	143	GLY	N-CA-C	-15.10	96.83	112.08
1	A	338	ARG	NE-CZ-NH1	-9.42	112.08	121.50
1	B	100[A]	ARG	NE-CZ-NH2	9.26	127.53	119.20
1	B	100[B]	ARG	NE-CZ-NH2	9.26	127.53	119.20
1	A	310	PRO	CA-C-N	-8.49	110.06	122.46
1	A	310	PRO	C-N-CA	-8.49	110.06	122.46
1	A	19	ARG	NE-CZ-NH2	-8.09	111.92	119.20
1	B	30	ARG	NE-CZ-NH1	-8.00	113.50	121.50
1	B	141	ARG	N-CA-C	-7.80	94.19	110.80
1	B	143	GLY	CA-C-O	7.60	127.92	122.37
1	B	55	ARG	NE-CZ-NH1	-7.45	114.05	121.50
1	B	217	SER	N-CA-C	-7.43	98.02	109.24
1	B	100[A]	ARG	NE-CZ-NH1	-6.77	114.73	121.50
1	B	100[B]	ARG	NE-CZ-NH1	-6.77	114.73	121.50
1	B	55	ARG	NE-CZ-NH2	6.76	125.28	119.20
1	B	231	THR	OG1-CB-CG2	-6.67	95.95	109.30
1	A	75	ALA	N-CA-C	-6.34	104.81	112.54
1	B	370	SER	N-CA-CB	-6.11	100.98	109.97
1	B	337	THR	CB-CA-C	6.00	121.68	110.70
1	B	165	ARG	CG-CD-NE	-5.98	98.85	112.00
1	B	287	ARG	NE-CZ-NH1	-5.97	115.53	121.50
1	B	52	ALA	N-CA-C	-5.94	104.89	111.36
1	B	153	LEU	CD1-CG-CD2	5.93	123.85	110.80
1	B	101	ARG	CG-CD-NE	5.92	125.03	112.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	149	THR	CA-C-N	-5.91	112.41	118.85
1	A	149	THR	C-N-CA	-5.91	112.41	118.85
1	A	271	LYS	CA-CB-CG	-5.89	102.32	114.10
1	B	215	VAL	CA-C-N	-5.80	114.72	122.54
1	B	215	VAL	C-N-CA	-5.80	114.72	122.54
1	A	338	ARG	CA-C-O	5.76	126.63	120.24
1	A	307	ARG	CD-NE-CZ	-5.74	116.36	124.40
1	A	10	THR	O-C-N	5.74	128.16	121.39
1	B	141	ARG	CB-CA-C	5.72	121.81	110.42
1	B	237	VAL	O-C-N	5.62	129.41	122.05
1	B	319	GLN	O-C-N	5.57	127.81	122.07
1	B	53	ALA	N-CA-C	-5.52	105.34	111.36
1	A	55	ARG	NE-CZ-NH1	-5.51	115.99	121.50
1	B	100[A]	ARG	CB-CG-CD	5.50	123.94	111.30
1	B	100[B]	ARG	CB-CG-CD	5.50	123.94	111.30
1	B	30	ARG	NE-CZ-NH2	5.48	124.14	119.20
1	B	48	VAL	O-C-N	5.45	123.91	120.42
1	A	4	THR	O-C-N	-5.43	116.56	122.92
1	B	165	ARG	NE-CZ-NH1	-5.43	116.07	121.50
1	B	91	ARG	N-CA-C	-5.41	106.45	113.16
1	A	307	ARG	N-CA-C	-5.38	101.89	109.96
1	B	149	THR	CA-C-N	-5.37	113.00	118.85
1	B	149	THR	C-N-CA	-5.37	113.00	118.85
1	B	94	GLN	CA-CB-CG	-5.32	103.47	114.10
1	A	287	ARG	O-C-N	5.31	128.90	121.83
1	B	141	ARG	CA-CB-CG	5.30	124.70	114.10
1	B	111	SER	O-C-N	5.29	129.31	123.33
1	B	189	ALA	N-CA-C	-5.18	105.75	111.71
1	A	274	ALA	O-C-N	-5.18	117.13	123.19
1	A	50	VAL	N-CA-C	-5.16	105.68	110.53
1	B	41	HIS	CA-CB-CG	-5.13	108.67	113.80
1	A	340	PRO	CB-CA-C	-5.12	104.68	111.23
1	B	210	CYS	CA-C-N	-5.10	113.47	119.84
1	B	210	CYS	C-N-CA	-5.10	113.47	119.84
1	A	270	ASP	CA-C-O	-5.09	114.12	119.97
1	B	16	ASP	CB-CG-OD1	5.08	130.08	118.40
1	B	225	SER	CB-CA-C	-5.07	102.89	110.90
1	B	204	ARG	CB-CG-CD	-5.06	99.65	111.30
1	B	142	SER	N-CA-CB	-5.02	102.01	110.49
1	B	22	ARG	CG-CD-NE	-5.01	100.98	112.00

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	2862	0	2869	16	0
1	B	2940	0	2932	28	0
2	A	15	0	6	1	0
2	B	15	0	6	0	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0
4	A	8	0	6	4	0
4	B	8	0	6	3	0
5	A	334	0	0	4	0
5	B	320	0	0	2	0
All	All	6505	0	5825	45	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 4.

All (45) 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:291:ARG:HD2	1:B:291:ARG:H	1.32	0.94
1:B:18:ASP:HB3	1:B:22:ARG:NH2	1.84	0.93
1:B:-5:HIS:NE2	1:B:-1:MET:HE2	2.02	0.74
1:A:231[B]:THR:HG22	5:A:788:HOH:O	1.88	0.72
1:B:368:ASP:OD1	1:B:370:SER:HB2	1.92	0.69
1:A:231[B]:THR:CG2	5:A:788:HOH:O	2.43	0.66
1:B:303:ASP:OD2	1:B:307:ARG:NH1	2.28	0.66
1:A:51:ALA:HA	1:A:54:GLN:HE21	1.61	0.66
1:B:68[A]:LYS:HG2	1:B:377:LEU:HD21	1.79	0.65
1:A:172:HIS:NE2	4:A:403:2TL:HG22	2.11	0.64
1:B:140:HIS:CE1	4:B:403:2TL:HG21	2.32	0.64
1:B:291:ARG:NH1	5:B:678:HOH:O	2.31	0.64
1:A:180:ASP:OD2	1:A:293:LYS:HE3	1.98	0.63
1:B:140:HIS:HE1	4:B:403:2TL:CG2	2.12	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94[A]:GLN:HG3	5:A:647:HOH:O	1.99	0.62
1:B:291:ARG:HD2	1:B:291:ARG:N	2.09	0.62
1:A:337:THR:HG23	5:A:691:HOH:O	1.98	0.62
1:B:18:ASP:HB3	1:B:22:ARG:HH22	1.62	0.61
1:B:140:HIS:CE1	4:B:403:2TL:CG2	2.83	0.61
1:B:51:ALA:HA	1:B:54:GLN:HE21	1.65	0.61
1:A:68:LYS:HG2	1:A:377:LEU:HD21	1.89	0.55
1:B:100[B]:ARG:HG2	1:B:100[B]:ARG:HH11	1.71	0.55
1:B:254:GLU:HG3	5:B:758:HOH:O	2.09	0.53
1:B:324:LEU:C	1:B:324:LEU:HD23	2.35	0.52
1:B:334:ASP:HB3	1:B:337:THR:OG1	2.10	0.51
1:B:68[A]:LYS:CD	1:B:377:LEU:HD11	2.43	0.48
1:B:100[B]:ARG:HG2	1:B:100[B]:ARG:NH1	2.29	0.48
1:A:141:ARG:NH1	4:A:403:2TL:HG23	2.29	0.47
1:B:68[A]:LYS:HD3	1:B:377:LEU:HD11	1.96	0.47
1:A:326:ARG:HD2	1:A:332:GLU:OE1	2.15	0.47
1:B:100[B]:ARG:HH11	1:B:100[B]:ARG:CG	2.27	0.47
1:A:2:GLN:HE21	1:A:2:GLN:C	2.23	0.46
1:B:368:ASP:OD1	1:B:370:SER:CB	2.62	0.46
1:B:333:ALA:O	1:B:334:ASP:C	2.58	0.46
1:A:2:GLN:NE2	1:A:2:GLN:CA	2.79	0.44
1:A:324:LEU:HD23	1:A:324:LEU:C	2.43	0.43
1:B:68[A]:LYS:HE2	1:B:377:LEU:HD11	2.00	0.43
1:B:100[B]:ARG:NH1	1:B:127:GLU:OE2	2.48	0.43
1:B:132:TRP:CH2	1:B:212:VAL:HG11	2.56	0.41
1:A:11:PRO:HA	1:A:261:LEU:O	2.21	0.41
1:A:85:VAL:O	1:A:86:SER:C	2.62	0.41
2:A:401:PLP:O3	4:A:403:2TL:N	2.52	0.41
1:B:30:ARG:HG2	1:B:31:MET:HE2	2.03	0.41
1:B:67:LEU:HD23	1:B:67:LEU:HA	1.94	0.41
1:A:140:HIS:NE2	4:A:403:2TL:HG21	2.37	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	383/390 (98%)	377 (98%)	4 (1%)	2 (0%)	24	10
1	B	392/390 (100%)	384 (98%)	6 (2%)	2 (0%)	24	10
All	All	775/780 (99%)	761 (98%)	10 (1%)	4 (0%)	24	10

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	141	ARG
1	B	241	PHE
1	A	236	GLY
1	B	236	GLY

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	286/289 (99%)	280 (98%)	6 (2%)	47	23
1	B	294/289 (102%)	291 (99%)	3 (1%)	68	50
All	All	580/578 (100%)	571 (98%)	9 (2%)	55	33

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	GLN
1	A	18	ASP
1	A	30	ARG
1	A	89	PRO
1	A	124	GLU
1	A	291	ARG
1	B	89	PRO
1	B	153	LEU
1	B	291	ARG

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	2	GLN
1	A	28	GLN
1	A	54	GLN
1	A	363	GLN
1	B	-4	HIS
1	B	2	GLN
1	B	21	GLN
1	B	28	GLN
1	B	54	GLN
1	B	94	GLN
1	B	125	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 3 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	$\# Z > 2$
2	PLP	B	401	4	15,15,16	2.07	5 (33%)	21,22,23	1.72	5 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	2TL	A	403	2	7,7,7	2.10	3 (42%)	8,9,9	3.07	4 (50%)
2	PLP	A	401	4	15,15,16	3.18	4 (26%)	21,22,23	1.74	8 (38%)
4	2TL	B	403	2	7,7,7	1.72	2 (28%)	8,9,9	2.11	3 (37%)

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	PLP	B	401	4	-	0/6/6/8	0/1/1/1
4	2TL	A	403	2	-	0/8/8/8	-
2	PLP	A	401	4	-	0/6/6/8	0/1/1/1
4	2TL	B	403	2	-	0/8/8/8	-

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	PLP	C5-C4	10.28	1.52	1.40
2	B	401	PLP	C5-C4	4.57	1.45	1.40
2	A	401	PLP	C3-C2	4.35	1.45	1.41
4	B	403	2TL	OXT-C	-3.44	1.19	1.30
2	A	401	PLP	C6-C5	3.33	1.44	1.37
4	A	403	2TL	OXT-C	-3.31	1.20	1.30
4	A	403	2TL	CA-N	3.24	1.56	1.47
2	B	401	PLP	C3-C2	3.23	1.44	1.41
2	B	401	PLP	C4A-C4	-3.06	1.45	1.51
2	A	401	PLP	C3-C4	2.89	1.45	1.40
2	B	401	PLP	C6-C5	2.67	1.43	1.37
2	B	401	PLP	C6-N1	-2.50	1.29	1.34
4	A	403	2TL	CG2-CB	-2.21	1.45	1.51
4	B	403	2TL	CG2-CB	-2.12	1.45	1.51

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	403	2TL	OG1-CB-CA	5.40	120.51	109.00
4	A	403	2TL	OG1-CB-CG2	-4.87	95.11	109.68
4	A	403	2TL	C-CA-N	-3.81	101.49	109.49
4	B	403	2TL	CG2-CB-CA	3.77	120.81	112.18
2	A	401	PLP	C5-C6-N1	-3.31	118.44	123.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	PLP	O2P-P-O4P	-3.20	98.33	106.67
2	A	401	PLP	C6-N1-C2	2.97	124.59	119.20
4	B	403	2TL	C-CA-N	-2.96	103.26	109.49
2	B	401	PLP	O3P-P-O2P	2.68	117.84	107.80
2	A	401	PLP	C6-C5-C4	2.52	120.16	118.10
2	A	401	PLP	C2A-C2-C3	-2.46	117.92	120.80
2	B	401	PLP	C4A-C4-C5	-2.43	118.43	120.94
2	B	401	PLP	C2A-C2-C3	-2.39	118.00	120.80
4	B	403	2TL	OXT-C-O	2.37	129.45	124.08
4	A	403	2TL	CG2-CB-CA	2.27	117.36	112.18
2	A	401	PLP	O2P-P-O4P	-2.16	101.04	106.67
2	A	401	PLP	O3P-P-O2P	2.05	115.48	107.80
2	B	401	PLP	C4-C3-C2	-2.04	116.83	119.89
2	A	401	PLP	C5A-C5-C4	-2.03	118.64	122.64
2	A	401	PLP	O3-C3-C4	2.03	123.39	118.10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	403	2TL	4	0
2	A	401	PLP	1	0
4	B	403	2TL	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	379/390 (97%)	-0.38	3 (0%) 82 86	6, 19, 36, 58	6 (1%)
1	B	389/390 (99%)	-0.30	14 (3%) 46 48	7, 19, 43, 82	5 (1%)
All	All	768/780 (98%)	-0.34	17 (2%) 62 66	6, 19, 39, 82	11 (1%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	-7	HIS	4.7
1	B	-8	HIS	3.8
1	B	-5	HIS	3.4
1	A	2	GLN	3.4
1	B	291	ARG	3.3
1	B	-2	ALA	3.3
1	B	-1	MET	3.3
1	B	-6	HIS	3.1
1	B	-4	HIS	2.6
1	A	147	ASP	2.4
1	B	-3	HIS	2.4
1	A	291	ARG	2.4
1	B	1	MET	2.4
1	B	330	ALA	2.4
1	B	2	GLN	2.3
1	B	0	SER	2.2
1	B	334	ASP	2.2

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 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($\text{\AA}^2$)	Q<0.9
4	2TL	B	403	8/8	0.96	0.08	14,17,18,27	0
3	MG	B	402	1/1	0.97	0.07	19,19,19,19	0
4	2TL	A	403	8/8	0.97	0.06	17,18,23,33	0
3	MG	A	402	1/1	0.97	0.08	11,11,11,11	1
2	PLP	A	401	15/16	0.98	0.05	14,16,18,19	0
3	MG	B	404	1/1	0.98	0.15	17,17,17,17	1
2	PLP	B	401	15/16	0.99	0.04	13,13,18,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.