



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 27, 2026 – 10:15 PM UTC

PDB ID : 4PB5 / pdb_00004pb5
Title : D-threo-3-hydroxyaspartate dehydratase H351A mutant complexed with L-erythro-3-hydroxyaspartate
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Deposited on : 2014-04-11
Resolution : 1.90 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

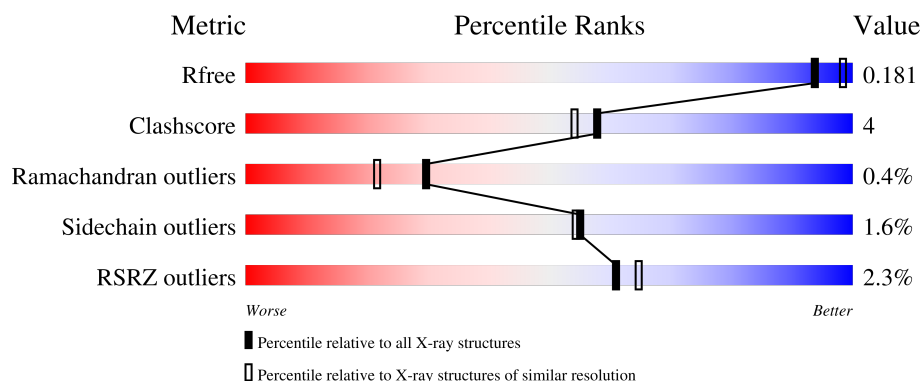
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	% 86% 11% ...
1	B	390	3% 86% 12% ...

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6314 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
			2857	1775	533	534	15			
1	B	388	Total	C	N	O	S	0	6	0
			2930	1820	549	543	18			

There are 22 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
A	351	ALA	HIS	engineered mutation	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
B	351	ALA	HIS	engineered mutation	UNP B2DFG5

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

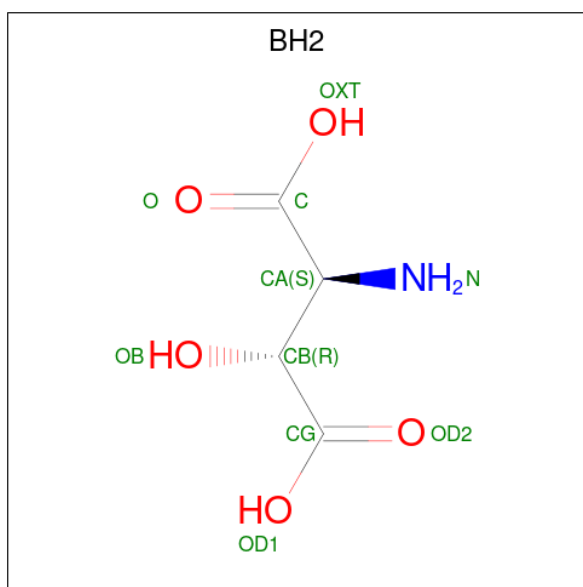


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- 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	1	Total	Mg	0	0
			1	1		

- Molecule 4 is (3R)-3-hydroxy-L-aspartic acid (CCD ID: BH2) (formula: C₄H₇NO₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			10	4	1	5		
4	B	1	Total	C	N	O	0	0
			10	4	1	5		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			6	3	3		

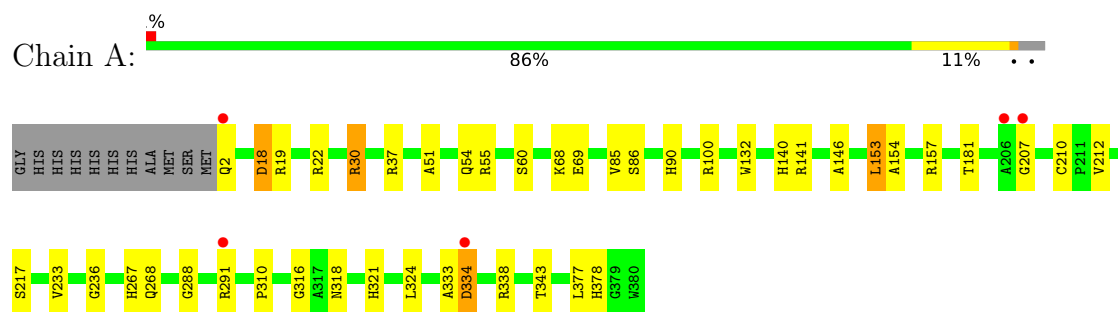
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	240	Total 240	O 240	0	0
6	B	229	Total 229	O 229	0	0

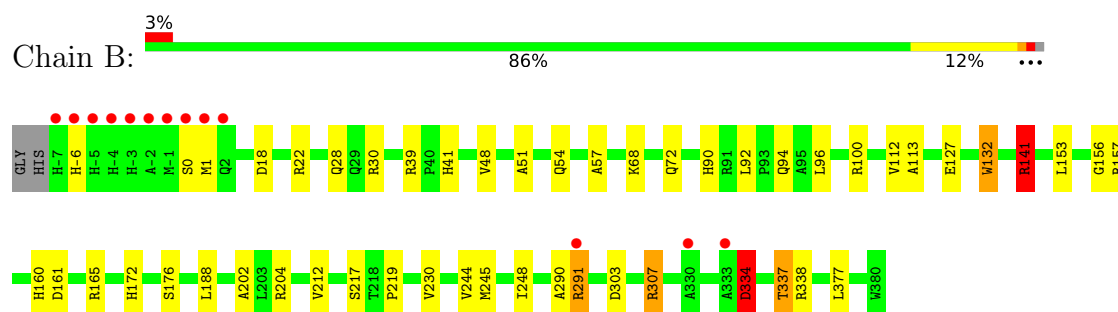
3 Residue-property plots [i](#)

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	157.63Å 157.63Å 157.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.46 – 1.90 40.46 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.8 (40.46-1.90) 99.8 (40.46-1.90)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.10 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.157 , 0.183 0.155 , 0.181	Depositor DCC
R_{free} test set	3835 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	18.7	Xtriage
Anisotropy	0.010	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 37.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.001 for l,-k,h 0.004 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6314	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 3.14% 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: BH2, MG, PLP, GOL

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	1.53	19/2921 (0.7%)	1.26	10/3962 (0.3%)
1	B	1.64	21/2999 (0.7%)	1.25	8/4065 (0.2%)
All	All	1.59	40/5920 (0.7%)	1.25	18/8027 (0.2%)

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	-6	HIS	CE1-NE2	21.65	1.54	1.32
1	B	-6	HIS	CG-ND1	11.03	1.50	1.38
1	A	267	HIS	ND1-CE1	9.16	1.41	1.32
1	A	267	HIS	CG-CD2	8.16	1.44	1.35
1	B	-6	HIS	CG-CD2	7.39	1.44	1.35
1	A	90	HIS	CG-CD2	7.28	1.43	1.35
1	B	202	ALA	N-CA	6.87	1.54	1.46
1	A	316	GLY	N-CA	6.22	1.51	1.45
1	B	172	HIS	ND1-CE1	6.15	1.38	1.32
1	A	140	HIS	ND1-CE1	6.07	1.38	1.32
1	A	19	ARG	CZ-NH1	6.05	1.41	1.32
1	B	39	ARG	CA-C	5.99	1.58	1.52
1	A	343	THR	N-CA	5.84	1.53	1.46
1	B	-6	HIS	ND1-CE1	5.80	1.38	1.32
1	A	207	GLY	N-CA	5.74	1.53	1.45
1	B	204	ARG	CD-NE	-5.73	1.38	1.46
1	A	378	HIS	CG-ND1	-5.68	1.32	1.38
1	B	113	ALA	N-CA	5.67	1.53	1.46
1	B	160	HIS	ND1-CE1	5.55	1.38	1.32
1	B	172	HIS	CG-CD2	5.50	1.42	1.35
1	B	112	VAL	CA-C	5.46	1.59	1.52
1	A	154	ALA	CA-C	-5.45	1.45	1.52
1	A	207	GLY	CA-C	5.44	1.58	1.51
1	A	153	LEU	CA-CB	5.41	1.61	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	140	HIS	CG-CD2	5.41	1.41	1.35
1	B	156	GLY	CA-C	5.38	1.58	1.52
1	B	92	LEU	CA-C	5.31	1.59	1.52
1	B	141	ARG	CA-C	5.31	1.59	1.52
1	B	244	VAL	N-CA	5.29	1.52	1.46
1	B	90	HIS	CE1-NE2	5.22	1.37	1.32
1	B	230	VAL	N-CA	5.16	1.52	1.46
1	A	18	ASP	N-CA	5.16	1.52	1.46
1	B	41	HIS	CE1-NE2	5.15	1.37	1.32
1	A	69	GLU	N-CA	5.11	1.52	1.46
1	B	132	TRP	CD2-CE2	5.10	1.50	1.41
1	A	181	THR	CA-C	5.08	1.57	1.52
1	A	233	VAL	CA-C	-5.06	1.46	1.52
1	A	321	HIS	ND1-CE1	5.04	1.37	1.32
1	B	94	GLN	CA-C	5.02	1.59	1.52
1	A	288	GLY	N-CA	5.02	1.52	1.45

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	-6	HIS	ND1-CG-CD2	7.45	113.55	106.10
1	A	217	SER	N-CA-C	-6.82	98.97	109.14
1	A	210	CYS	CA-C-N	-6.74	112.83	119.64
1	A	210	CYS	C-N-CA	-6.74	112.83	119.64
1	A	55	ARG	NE-CZ-NH2	6.58	125.13	119.20
1	A	310	PRO	CA-C-N	-6.39	113.14	122.46
1	A	310	PRO	C-N-CA	-6.39	113.14	122.46
1	A	334	ASP	N-CA-C	6.25	119.80	111.30
1	B	217	SER	N-CA-C	-6.04	100.14	109.14
1	A	146	ALA	CA-C-N	-5.94	113.14	122.65
1	A	146	ALA	C-N-CA	-5.94	113.14	122.65
1	B	290	ALA	CA-C-N	-5.81	113.43	122.67
1	B	290	ALA	C-N-CA	-5.81	113.43	122.67
1	B	48	VAL	O-C-N	5.57	123.98	120.42
1	B	334	ASP	N-CA-C	5.55	119.05	111.39
1	B	204	ARG	NE-CZ-NH2	-5.17	114.55	119.20
1	B	30	ARG	NE-CZ-NH1	-5.03	116.47	121.50
1	A	100	ARG	NE-CZ-NH1	-5.02	116.48	121.50

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	2857	0	2868	15	0
1	B	2930	0	2933	36	0
2	A	15	0	6	0	0
2	B	15	0	6	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	10	0	3	1	0
4	B	10	0	2	1	0
5	B	6	0	8	2	0
6	A	240	0	0	1	0
6	B	229	0	0	2	0
All	All	6314	0	5826	52	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 (52) 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:245[B]:MET:CE	1:B:245[B]:MET:HA	1.62	1.30
1:B:100[B]:ARG:CG	1:B:100[B]:ARG:HH11	1.55	1.16
1:B:100[B]:ARG:HH11	1:B:100[B]:ARG:HG2	1.03	1.14
1:B:245[B]:MET:HA	1:B:245[B]:MET:HE3	1.21	1.10
1:A:30:ARG:HH11	1:A:30:ARG:HG3	0.90	1.04
1:B:245[B]:MET:HA	1:B:245[B]:MET:HE2	1.44	0.99
1:A:30:ARG:HH11	1:A:30:ARG:CG	1.73	0.99
1:A:30:ARG:HG3	1:A:30:ARG:NH1	1.56	0.96
1:B:245[B]:MET:CE	1:B:245[B]:MET:CA	2.43	0.93
1:B:100[B]:ARG:HG2	1:B:100[B]:ARG:NH1	1.81	0.91
1:B:245[B]:MET:HE3	1:B:245[B]:MET:CA	2.07	0.84
1:B:245[B]:MET:HE2	1:B:245[B]:MET:CA	2.09	0.77
1:B:245[B]:MET:HE3	1:B:248:ILE:HD12	1.66	0.76
1:B:68[A]:LYS:HG2	1:B:377:LEU:HD21	1.74	0.70
1:B:100[B]:ARG:CG	1:B:100[B]:ARG:NH1	2.32	0.69
1:B:303:ASP:OD2	1:B:307:ARG:NH1	2.24	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:ALA:HA	1:A:54:GLN:HE21	1.59	0.68
1:B:334:ASP:OD2	1:B:337:THR:HG23	1.94	0.68
4:A:403:BH2:HA	6:A:739:HOH:O	1.95	0.66
1:B:100[B]:ARG:NH1	1:B:127:GLU:OE2	2.29	0.66
1:B:291:ARG:HD2	1:B:291:ARG:H	1.63	0.64
1:B:100[B]:ARG:HH11	1:B:100[B]:ARG:HG3	1.59	0.63
1:B:307:ARG:HH11	1:B:307:ARG:HB2	1.63	0.63
1:B:245[B]:MET:CE	1:B:248:ILE:HD12	2.31	0.61
1:B:51:ALA:HA	1:B:54:GLN:HE21	1.66	0.60
4:B:403:BH2:HA	6:B:662:HOH:O	2.05	0.56
1:B:18:ASP:HB3	1:B:22:ARG:NH2	2.21	0.56
1:B:291:ARG:HD2	1:B:291:ARG:N	2.23	0.53
1:A:153:LEU:O	1:A:157[B]:ARG:HG3	2.09	0.53
1:B:157:ARG:HB3	5:B:404:GOL:H32	1.92	0.52
1:B:219:PRO:HG3	1:B:245[B]:MET:HE1	1.93	0.51
1:A:333:ALA:O	1:A:334:ASP:C	2.56	0.47
1:A:132:TRP:CH2	1:A:212:VAL:HG11	2.50	0.46
1:B:132:TRP:CH2	1:B:212:VAL:HG11	2.51	0.46
1:A:338:ARG:O	1:A:338:ARG:HG3	2.15	0.46
1:B:0:SER:O	1:B:1:MET:HE2	2.16	0.45
1:A:324:LEU:HD23	1:A:324:LEU:C	2.41	0.45
1:A:18:ASP:HB3	1:A:22:ARG:NH2	2.32	0.44
1:A:37[B]:ARG:HD3	1:A:60:SER:OG	2.17	0.44
1:A:85:VAL:O	1:A:86:SER:C	2.60	0.44
1:A:318:ASN:OD1	1:B:141:ARG:HD3	2.17	0.43
1:B:307:ARG:HH11	1:B:307:ARG:CB	2.30	0.43
1:B:28:GLN:HG2	1:B:57:ALA:O	2.18	0.42
1:B:176:SER:HB3	1:B:188:LEU:HD23	2.01	0.42
1:B:161:ASP:HB2	5:B:404:GOL:H12	2.01	0.42
1:B:96:LEU:O	1:B:100[A]:ARG:HG3	2.19	0.41
1:B:165:ARG:HD3	6:B:542:HOH:O	2.20	0.41
1:A:68:LYS:HG2	1:A:377:LEU:HD21	2.01	0.41
1:B:96:LEU:HD21	1:B:100[A]:ARG:CZ	2.51	0.41
1:B:72:GLN:OE1	1:B:377:LEU:HD13	2.20	0.41
1:A:30:ARG:CG	1:A:30:ARG:NH1	2.40	0.40
1:B:338:ARG:HH11	1:B:338:ARG:HD2	1.74	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%)	375 (98%)	6 (2%)	2 (0%)	24	16
1	B	392/390 (100%)	382 (97%)	9 (2%)	1 (0%)	36	29
All	All	775/780 (99%)	757 (98%)	15 (2%)	3 (0%)	30	22

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	141	ARG
1	B	141	ARG
1	A	236	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	285/288 (99%)	281 (99%)	4 (1%)	59	59
1	B	293/288 (102%)	288 (98%)	5 (2%)	53	52
All	All	578/576 (100%)	569 (98%)	9 (2%)	55	54

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

Mol	Chain	Res	Type
1	A	2	GLN
1	A	30	ARG
1	A	268	GLN

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Mol	Chain	Res	Type
1	A	291	ARG
1	B	153	LEU
1	B	291	ARG
1	B	307	ARG
1	B	334	ASP
1	B	337	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	GLN
1	A	28	GLN
1	A	54	GLN
1	A	305	GLN
1	B	-4	HIS
1	B	28	GLN
1	B	54	GLN
1	B	94	GLN
1	B	363	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, 2 are monoatomic - leaving 5 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	A	401	4	15,15,16	2.48	6 (40%)	21,22,23	2.28	11 (52%)
4	BH2	B	403	3,2	8,9,9	1.17	1 (12%)	11,12,12	1.38	2 (18%)
2	PLP	B	401	4	15,15,16	2.38	4 (26%)	21,22,23	1.93	8 (38%)
5	GOL	B	404	-	5,5,5	1.21	1 (20%)	5,5,5	1.05	0
4	BH2	A	403	3,2	8,9,9	1.72	3 (37%)	11,12,12	0.71	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
2	PLP	A	401	4	-	0/6/6/8	0/1/1/1
4	BH2	B	403	3,2	-	3/12/12/12	-
2	PLP	B	401	4	-	0/6/6/8	0/1/1/1
5	GOL	B	404	-	-	2/4/4/4	-
4	BH2	A	403	3,2	-	5/12/12/12	-

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	PLP	C3-C2	6.29	1.47	1.41
2	A	401	PLP	C3-C2	5.32	1.46	1.41
2	B	401	PLP	C3-C4	4.81	1.49	1.40
2	A	401	PLP	C3-C4	4.52	1.48	1.40
2	A	401	PLP	C5-C4	4.19	1.45	1.40
2	B	401	PLP	C5-C4	3.39	1.44	1.40
2	A	401	PLP	C4A-C4	-3.07	1.45	1.51
4	A	403	BH2	OD2-CG	2.80	1.30	1.22
4	B	403	BH2	OXT-C	-2.72	1.22	1.30
5	B	404	GOL	O3-C3	2.42	1.52	1.42
4	A	403	BH2	OXT-C	-2.40	1.23	1.30
4	A	403	BH2	OD1-CG	-2.29	1.23	1.30
2	B	401	PLP	C6-C5	2.18	1.42	1.37
2	A	401	PLP	C6-C5	2.16	1.42	1.37
2	A	401	PLP	C2-N1	2.06	1.37	1.33

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	PLP	C4A-C4-C5	-4.41	116.39	120.94
2	A	401	PLP	C6-N1-C2	3.92	126.31	119.20
2	A	401	PLP	C2A-C2-N1	3.68	124.57	117.64
2	B	401	PLP	C2A-C2-N1	3.25	123.77	117.64
2	B	401	PLP	C6-N1-C2	3.15	124.91	119.20
2	A	401	PLP	O3-C3-C4	3.03	126.02	118.10
2	B	401	PLP	C4A-C4-C5	-2.89	117.96	120.94
2	B	401	PLP	O2P-P-O4P	-2.81	99.34	106.67
2	A	401	PLP	C2A-C2-C3	-2.78	117.54	120.80
2	A	401	PLP	C3-C4-C5	2.72	121.86	118.59
2	B	401	PLP	C3-C2-N1	-2.56	117.73	120.96
4	B	403	BH2	OD1-CG-CB	2.56	120.43	113.31
2	A	401	PLP	C3-C2-N1	-2.47	117.85	120.96
2	A	401	PLP	C4-C3-C2	-2.42	116.27	119.89
2	B	401	PLP	O2P-P-O1P	2.35	120.01	110.83
2	A	401	PLP	C6-C5-C4	-2.30	116.21	118.10
4	B	403	BH2	O-C-CA	-2.21	114.23	121.72
2	A	401	PLP	O3P-P-O2P	2.16	115.90	107.80
2	B	401	PLP	C5A-C5-C6	2.13	122.84	119.36
2	A	401	PLP	O2P-P-O1P	2.09	118.98	110.83
2	B	401	PLP	O3P-P-O2P	2.06	115.52	107.80

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	403	BH2	O-C-CA-CB
4	A	403	BH2	OXT-C-CA-CB
4	B	403	BH2	O-C-CA-CB
4	B	403	BH2	OXT-C-CA-CB
5	B	404	GOL	O1-C1-C2-C3
5	B	404	GOL	O1-C1-C2-O2
4	A	403	BH2	C-CA-CB-CG
4	B	403	BH2	C-CA-CB-CG
4	A	403	BH2	OXT-C-CA-N
4	A	403	BH2	O-C-CA-N

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	403	BH2	1	0
5	B	404	GOL	2	0
4	A	403	BH2	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	379/390 (97%)	-0.50	5 (1%) 75 78	7, 19, 36, 57	6 (1%)
1	B	388/390 (99%)	-0.39	13 (3%) 48 51	7, 19, 42, 88	6 (1%)
All	All	767/780 (98%)	-0.45	18 (2%) 61 65	7, 19, 39, 88	12 (1%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	-7	HIS	5.4
1	B	-5	HIS	4.6
1	B	-6	HIS	4.2
1	B	-1	MET	4.2
1	B	-4	HIS	3.9
1	B	1	MET	3.9
1	A	206	ALA	3.7
1	B	0	SER	3.7
1	A	2	GLN	3.6
1	B	-2	ALA	3.4
1	B	330	ALA	3.0
1	B	291	ARG	2.9
1	B	-3	HIS	2.8
1	B	333	ALA	2.6
1	B	2	GLN	2.6
1	A	334	ASP	2.5
1	A	207	GLY	2.4
1	A	291	ARG	2.1

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($\text{\AA}^2$)	Q<0.9
5	GOL	B	404	6/6	0.88	0.16	29,32,34,40	0
3	MG	B	402	1/1	0.91	0.14	29,29,29,29	0
4	BH2	A	403	10/10	0.94	0.07	22,27,30,30	0
4	BH2	B	403	10/10	0.95	0.07	20,23,31,33	0
3	MG	A	402	1/1	0.95	0.10	31,31,31,31	0
2	PLP	B	401	15/16	0.99	0.04	13,15,17,18	0
2	PLP	A	401	15/16	0.99	0.04	15,16,18,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.