



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 05:29 AM GMT

PDB ID : 1I3M
Title : MOLECULAR BASIS FOR SEVERE EPIMERASE-DEFICIENCY GALACTOSEMIA: X-RAY STRUCTURE OF THE HUMAN V94M-SUBSTITUTED UDP-GALACTOSE 4-EPIMERASE
Authors : Thoden, J.B.; Wohlers, T.M.; Fridovich-Keil, J.L.; Holden, H.M.
Deposited on : 2001-02-15
Resolution : 1.50 Å(reported)

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

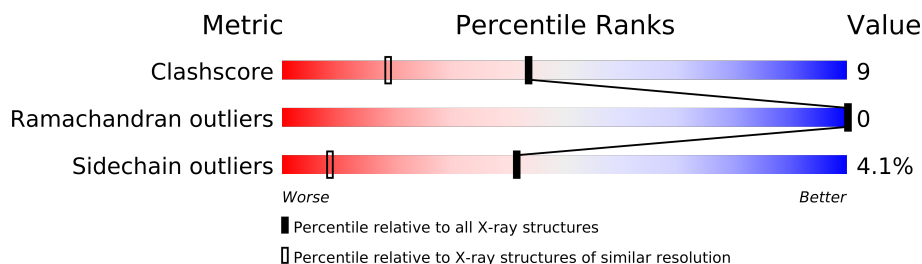
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.50 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))
Clashscore	79885	1768 (1.50-1.50)
Ramachandran outliers	78287	1720 (1.50-1.50)
Sidechain outliers	78261	1718 (1.50-1.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	348	
1	B	348	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6402 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called UDP-GLUCOSE 4-EPIMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	347	Total	C	N	O	S	0	3	0
			2699	1707	468	508	16			
1	B	345	Total	C	N	O	S	0	5	0
			2689	1705	464	504	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	94	MET	VAL	ENGINEERED	UNP Q14376
B	94	MET	VAL	ENGINEERED	UNP Q14376

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

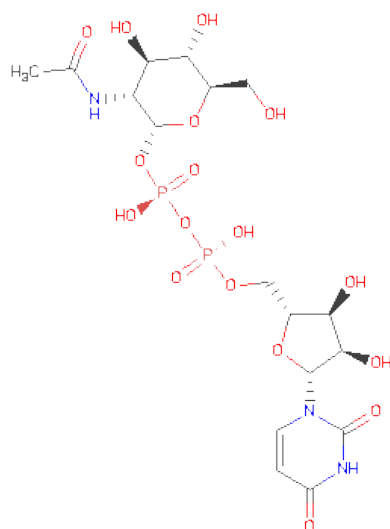
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

- Molecule 4 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
4	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 5 is URIDINE-DIPHOSPHATE-N-ACETYLGLUCOSAMINE (three-letter code: UD1) (formula: $C_{17}H_{27}N_3O_{17}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	1
			40	17	3	18	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	P	0	0
			39	17	3	17	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	418	Total	O	0	0
			418	418		
6	B	424	Total	O	0	0
			424	424		

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.

Note EDS was not executed.

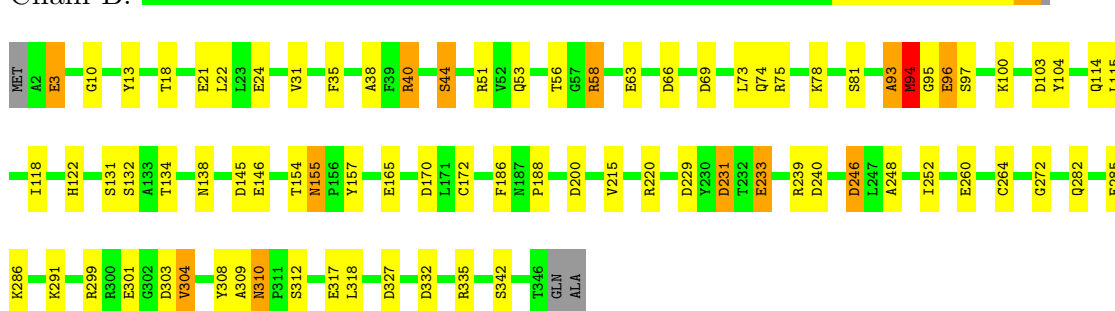
• Molecule 1: UDP-GLUCOSE 4-EPIMERASE

Chain A:



• Molecule 1: UDP-GLUCOSE 4-EPIMERASE

Chain B:



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.90Å 89.90Å 96.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.50	Depositor
% Data completeness (in resolution range)	98.4 (50.00-1.50)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT	Depositor
R, R_{free}	0.184 , 0.214	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6402	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: MG, UD1, NAD, CL

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	0.94	16/2768 (0.6%)	1.26	35/3744 (0.9%)
1	B	0.95	12/2770 (0.4%)	1.25	30/3748 (0.8%)
All	All	0.95	28/5538 (0.5%)	1.26	65/7492 (0.9%)

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	317	GLU	CD-OE2	7.54	1.33	1.25
1	B	146	GLU	CD-OE2	7.33	1.33	1.25
1	B	233	GLU	CD-OE2	7.00	1.33	1.25
1	A	316	GLU	CD-OE2	6.68	1.32	1.25
1	A	21	GLU	CD-OE2	6.60	1.32	1.25
1	A	317	GLU	CD-OE2	6.53	1.32	1.25
1	A	165	GLU	CD-OE2	6.32	1.32	1.25
1	B	24	GLU	CD-OE2	6.30	1.32	1.25
1	B	260	GLU	CD-OE2	6.27	1.32	1.25
1	A	47	GLU	CD-OE2	6.16	1.32	1.25
1	A	63[A]	GLU	CD-OE2	6.15	1.32	1.25
1	A	63[B]	GLU	CD-OE2	6.15	1.32	1.25
1	B	165	GLU	CD-OE2	6.04	1.32	1.25
1	B	301	GLU	CD-OE2	5.93	1.32	1.25
1	A	24[A]	GLU	CD-OE2	5.82	1.32	1.25
1	A	24[B]	GLU	CD-OE2	5.82	1.32	1.25
1	B	96	GLU	CD-OE2	5.74	1.31	1.25
1	B	63	GLU	CD-OE2	5.72	1.31	1.25
1	B	3	GLU	CD-OE2	5.70	1.31	1.25
1	A	96	GLU	CD-OE2	5.69	1.31	1.25
1	A	146	GLU	CD-OE2	5.66	1.31	1.25
1	A	301	GLU	CD-OE2	5.62	1.31	1.25
1	A	61	GLU	CD-OE2	5.55	1.31	1.25
1	A	221	GLU	CD-OE2	5.44	1.31	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	285	GLU	CD-OE2	5.35	1.31	1.25
1	A	260	GLU	CD-OE2	5.33	1.31	1.25
1	B	21	GLU	CD-OE2	5.01	1.31	1.25
1	A	117	GLU	CD-OE2	5.01	1.31	1.25

All (65) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	145	ASP	CB-CG-OD2	-9.92	109.37	118.30
1	A	220	ARG	NE-CZ-NH1	9.14	124.87	120.30
1	A	234	ASP	CB-CG-OD2	-8.71	110.46	118.30
1	B	145	ASP	CB-CG-OD1	8.02	125.52	118.30
1	B	335	ARG	NE-CZ-NH1	7.80	124.20	120.30
1	A	335	ARG	NE-CZ-NH1	7.69	124.14	120.30
1	A	80	TYR	CB-CG-CD1	-7.55	116.47	121.00
1	A	145	ASP	CB-CG-OD2	-7.33	111.70	118.30
1	A	69	ASP	CB-CG-OD2	-7.15	111.87	118.30
1	B	69	ASP	CB-CG-OD2	-7.12	111.89	118.30
1	B	75	ARG	NE-CZ-NH2	-7.12	116.74	120.30
1	B	231	ASP	CB-CG-OD1	7.08	124.67	118.30
1	A	66	ASP	CB-CG-OD1	6.79	124.42	118.30
1	A	220	ARG	CB-CA-C	6.71	123.82	110.40
1	B	93	ALA	N-CA-CB	6.63	119.38	110.10
1	B	69	ASP	CB-CG-OD1	6.54	124.19	118.30
1	A	69	ASP	CB-CG-OD1	6.53	124.17	118.30
1	A	170	ASP	CB-CG-OD2	-6.50	112.45	118.30
1	A	66	ASP	CB-CG-OD2	-6.49	112.46	118.30
1	B	38	ALA	N-CA-CB	6.47	119.16	110.10
1	B	200	ASP	CB-CG-OD1	6.46	124.11	118.30
1	B	246	ASP	CB-CG-OD2	-6.46	112.49	118.30
1	A	75	ARG	NE-CZ-NH1	6.45	123.53	120.30
1	B	246	ASP	CB-CG-OD1	6.43	124.08	118.30
1	A	231	ASP	CB-CG-OD2	-6.36	112.58	118.30
1	B	303	ASP	CB-CG-OD2	-6.36	112.58	118.30
1	A	145	ASP	CB-CG-OD1	6.18	123.86	118.30
1	B	170	ASP	CB-CG-OD1	6.18	123.86	118.30
1	A	300	ARG	NE-CZ-NH1	6.13	123.37	120.30
1	A	327	ASP	CB-CG-OD2	-6.12	112.79	118.30
1	B	66	ASP	CB-CG-OD2	-6.12	112.79	118.30
1	A	239	ARG	NE-CZ-NH1	6.09	123.34	120.30
1	A	332	ASP	CB-CG-OD1	6.02	123.72	118.30
1	B	229	ASP	CB-CG-OD2	-5.97	112.93	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	13	TYR	CB-CG-CD2	-5.96	117.42	121.00
1	A	184	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	A	299	ARG	NE-CZ-NH2	-5.92	117.34	120.30
1	A	106	ARG	NE-CZ-NH2	-5.88	117.36	120.30
1	A	246	ASP	CB-CG-OD1	5.82	123.54	118.30
1	B	66	ASP	CB-CG-OD1	5.76	123.49	118.30
1	B	291	LYS	CA-CB-CG	-5.76	100.73	113.40
1	B	332	ASP	CB-CG-OD1	5.70	123.43	118.30
1	B	51	ARG	NE-CZ-NH2	-5.69	117.46	120.30
1	A	300	ARG	NE-CZ-NH2	-5.62	117.49	120.30
1	B	200	ASP	CB-CG-OD2	-5.54	113.31	118.30
1	B	93	ALA	CB-CA-C	5.53	118.39	110.10
1	A	220	ARG	NE-CZ-NH2	-5.50	117.55	120.30
1	A	200	ASP	CB-CG-OD2	-5.48	113.36	118.30
1	B	335	ARG	NE-CZ-NH2	-5.47	117.56	120.30
1	B	231	ASP	CB-CG-OD2	-5.41	113.43	118.30
1	B	327	ASP	CB-CG-OD2	-5.39	113.45	118.30
1	A	200	ASP	CB-CG-OD1	5.39	123.15	118.30
1	A	234	ASP	CB-CG-OD1	5.35	123.12	118.30
1	A	332	ASP	CB-CG-OD2	-5.34	113.49	118.30
1	A	256	ARG	NE-CZ-NH1	5.31	122.96	120.30
1	A	246	ASP	CB-CG-OD2	-5.31	113.52	118.30
1	A	229	ASP	CB-CG-OD2	-5.29	113.53	118.30
1	A	231	ASP	CB-CG-OD1	5.26	123.03	118.30
1	A	169	ARG	CD-NE-CZ	-5.14	116.41	123.60
1	A	153	CYS	N-CA-CB	5.13	119.84	110.60
1	B	94	MET	N-CA-C	-5.11	97.21	111.00
1	B	318	LEU	CB-CA-C	-5.09	100.53	110.20
1	B	239	ARG	NE-CZ-NH1	5.05	122.82	120.30
1	A	155	ASN	N-CA-CB	-5.03	101.55	110.60
1	B	299	ARG	NE-CZ-NH1	5.01	122.81	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the

chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2699	0	2649	44	0
1	B	2689	0	2644	46	0
2	A	1	0	0	0	0
2	B	3	0	0	1	0
3	A	1	0	0	0	0
4	A	44	0	26	4	0
4	B	44	0	26	3	0
5	A	40	0	6	0	0
5	B	39	0	25	4	0
6	A	418	0	0	3	1
6	B	424	0	0	8	1
All	All	6402	0	5376	95	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 9.

All (95) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:132:SER:HB2	4:A:400:NAD:H5N	1.49	0.95
1:B:73:LEU:HB3	1:B:118[A]:ILE:HD13	1.54	0.89
1:B:58:ARG:HG3	1:B:58:ARG:HH11	1.53	0.74
1:A:100:LYS:HD3	1:A:103:ASP:HB2	1.70	0.73
1:A:155:ASN:OD1	1:A:157:TYR:HB3	1.92	0.70
1:B:44:SER:HB2	6:B:1669:HOH:O	1.91	0.69
1:A:94:MET:CB	1:A:96:GLU:HG3	2.23	0.69
1:A:132:SER:HB2	4:A:400:NAD:C5N	2.22	0.68
1:B:73:LEU:HD13	1:B:118[A]:ILE:CD1	2.22	0.68
1:B:73:LEU:CB	1:B:118[A]:ILE:HD13	2.23	0.66
1:B:155:ASN:H	1:B:155:ASN:HD22	1.44	0.65
1:B:56:THR:HB	1:B:58:ARG:NH1	2.12	0.65
1:B:94:MET:CB	1:B:96:GLU:HG3	2.27	0.64
1:B:122:HIS:ND1	6:B:1683:HOH:O	2.30	0.64
1:A:97:SER:HA	1:A:104:TYR:CE1	2.34	0.63
1:A:312:SER:HA	1:A:315:GLN:CG	2.28	0.62
1:B:58:ARG:HG3	1:B:58:ARG:NH1	2.13	0.61
1:A:113:ILE:HG12	1:A:167:MET:HE1	1.83	0.60
1:B:114:GLN:O	1:B:118[A]:ILE:HD12	2.01	0.60
1:B:310:ASN:ND2	1:B:312:SER:H	2.00	0.60
1:A:94:MET:HG3	1:A:206:ASN:OD1	2.01	0.59
1:B:310:ASN:HD22	1:B:312:SER:H	1.49	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:97:SER:HA	1:B:104:TYR:CE1	2.36	0.59
1:B:155:ASN:HD22	1:B:155:ASN:N	2.02	0.57
1:A:312:SER:HA	1:A:315:GLN:HG2	1.87	0.57
1:B:96:GLU:HB3	1:B:100:LYS:HD3	1.87	0.57
1:B:310:ASN:C	1:B:310:ASN:HD22	2.08	0.57
1:B:74:GLN:O	1:B:78:LYS:HG3	2.05	0.57
1:B:56:THR:HB	1:B:58:ARG:HH11	1.70	0.56
1:B:94:MET:HB3	1:B:96:GLU:HG3	1.88	0.55
1:B:93:ALA:C	1:B:95:GLY:H	2.10	0.54
1:B:35:PHE:CE2	1:B:40:ARG:HG3	2.42	0.54
1:B:286:LYS:HE2	6:B:1726:HOH:O	2.07	0.54
1:B:282:GLN:NE2	6:B:1908:HOH:O	2.28	0.54
4:B:900:NAD:H4N	5:B:901:UD1:H6'2	1.88	0.54
1:B:155:ASN:ND2	1:B:155:ASN:H	2.05	0.54
1:B:186:PHE:O	1:B:188:PRO:HD3	2.09	0.53
1:A:94:MET:HB2	1:A:96:GLU:HG3	1.91	0.52
1:B:286:LYS:HD3	6:B:1908:HOH:O	2.09	0.52
1:A:155:ASN:ND2	1:A:155:ASN:H	2.07	0.52
1:B:53:GLN:HG3	1:B:58:ARG:O	2.11	0.51
1:A:37:ASN:OD1	1:A:202:GLN:NE2	2.43	0.51
1:B:246:ASP:OD2	2:B:990:CL:CL	2.66	0.51
1:B:138:ASN:HD22	1:B:304:VAL:HG22	1.76	0.50
1:A:94:MET:HB3	1:A:96:GLU:HG3	1.91	0.50
1:A:100:LYS:HD3	1:A:103:ASP:CB	2.41	0.50
1:A:202:GLN:HG2	6:A:1475:HOH:O	2.11	0.49
1:B:115:LEU:O	1:B:118[B]:ILE:HG22	2.12	0.49
1:A:310:ASN:OD1	1:A:312:SER:OG	2.30	0.49
1:A:174:ALA:HA	6:B:1778:HOH:O	2.12	0.49
1:A:267:TYR:CE1	1:A:318:LEU:HD13	2.48	0.49
1:A:207:ASN:OD1	1:A:210:PRO:HG2	2.13	0.48
1:B:10:GLY:HA3	1:B:31:VAL:HG13	1.95	0.48
1:A:73:LEU:HD23	1:A:118:ILE:HD12	1.96	0.47
1:B:240:ASP:HB2	1:B:308:TYR:HA	1.96	0.47
1:B:131:SER:O	4:B:900:NAD:H6N	2.15	0.47
5:B:901:UD1:H4'	6:B:1901:HOH:O	2.14	0.47
1:B:22:LEU:HD23	1:B:252:ILE:HD12	1.96	0.47
1:A:312:SER:O	1:A:315:GLN:HG3	2.15	0.47
1:B:155:ASN:OD1	1:B:157:TYR:HB3	2.15	0.46
1:A:240:ASP:HB2	1:A:308:TYR:HA	1.98	0.46
1:A:214:GLN:OE1	1:A:220:ARG:HD2	2.16	0.45
1:B:215:VAL:HG22	1:B:220:ARG:HB2	1.97	0.45
1:A:328:ARG:HA	1:A:328:ARG:HD2	1.79	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:315:GLN:HB2	6:A:1646:HOH:O	2.18	0.44
1:B:73:LEU:HD13	1:B:118[A]:ILE:HD13	1.97	0.44
4:B:900:NAD:O7N	5:B:901:UD1:C6'	2.65	0.44
1:A:10:GLY:HA3	1:A:31:VAL:HG13	1.98	0.44
1:B:18:THR:HA	1:B:248:ALA:HB1	2.00	0.44
1:B:138:ASN:ND2	1:B:304:VAL:HG22	2.32	0.44
1:B:94:MET:C	1:B:96:GLU:HG3	2.39	0.43
1:A:311:PRO:O	1:A:315:GLN:HG2	2.18	0.43
1:A:131:SER:O	4:A:400:NAD:H6N	2.18	0.43
1:A:94:MET:HB3	1:A:96:GLU:CG	2.49	0.43
1:B:100:LYS:HE3	1:B:103:ASP:OD1	2.18	0.43
1:B:132:SER:HB3	1:B:134[A]:THR:HG22	2.01	0.42
1:B:272:GLY:HA2	1:B:309:ALA:O	2.19	0.42
1:A:172:CYS:SG	1:A:264:CYS:HB2	2.59	0.42
1:A:167:MET:HB3	1:A:167:MET:HE3	1.91	0.42
1:A:176:LYS:N	1:A:176:LYS:HD2	2.34	0.42
1:A:186:PHE:CE2	1:A:309:ALA:HB2	2.55	0.42
1:A:155:ASN:H	1:A:155:ASN:HD22	1.68	0.42
1:A:186:PHE:O	1:A:188:PRO:HD3	2.20	0.41
1:A:318:LEU:HD12	1:A:318:LEU:HA	1.83	0.41
1:A:132:SER:CB	4:A:400:NAD:H5N	2.33	0.41
5:B:901:UD1:H1'	6:B:1828:HOH:O	2.19	0.41
1:B:186:PHE:CE2	1:B:309:ALA:HB2	2.56	0.41
1:A:169:ARG:HD3	1:A:169:ARG:HH11	1.50	0.41
1:A:53:GLN:HG3	1:A:58:ARG:O	2.20	0.41
1:A:300:ARG:O	1:A:303:ASP:HB2	2.21	0.40
1:B:172:CYS:SG	1:B:264:CYS:HB2	2.61	0.40
1:A:77:PHE:HB3	1:A:122:HIS:CG	2.56	0.40
1:A:8:THR:O	1:A:88:PHE:HB2	2.20	0.40
1:A:125:LYS:HD3	1:A:178:TRP:NE1	2.37	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
6:A:1189:HOH:O	6:B:1638:HOH:O[4_456]	2.18	0.02

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	348/348 (100%)	340 (98%)	8 (2%)	0	100	100
1	B	349/348 (100%)	342 (98%)	7 (2%)	0	100	100
All	All	697/696 (100%)	682 (98%)	15 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	283/282 (100%)	273 (96%)	10 (4%)	48	13
1	B	284/282 (101%)	271 (95%)	13 (5%)	37	7
All	All	567/564 (100%)	544 (96%)	23 (4%)	41	8

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	94	MET
1	A	100	LYS
1	A	132	SER
1	A	153	CYS
1	A	169	ARG
1	A	176	LYS
1	A	220	ARG
1	A	247	LEU
1	A	295	LYS
1	A	303	ASP
1	B	3	GLU

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Mol	Chain	Res	Type
1	B	40	ARG
1	B	44	SER
1	B	58	ARG
1	B	81	SER
1	B	94	MET
1	B	154	THR
1	B	155	ASN
1	B	231	ASP
1	B	233	GLU
1	B	304	VAL
1	B	310	ASN
1	B	342	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	GLN
1	A	173	GLN
1	A	202	GLN
1	A	224	ASN
1	A	315	GLN
1	A	337	GLN
1	A	339	GLN
1	B	36	HIS
1	B	114	GLN
1	B	138	ASN
1	B	224	ASN
1	B	310	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 10 ligands modelled in this entry, 5 are monoatomic and 2 are modelled with single atom - leaving 3 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$
4	NAD	A	400	-	48,48,48	1.08	3 (6%)	73,73,73	1.31	6 (8%)
4	NAD	B	900	-	48,48,48	1.09	2 (4%)	73,73,73	1.60	8 (10%)
5	UD1	B	901	-	41,41,41	1.02	3 (7%)	58,62,62	1.15	4 (6%)

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
4	NAD	A	400	-	-	0/30/62/62	0/3/5/5
4	NAD	B	900	-	-	0/30/62/62	0/3/5/5
5	UD1	B	901	-	-	0/25/63/63	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	400	NAD	C2N-N1N	3.93	1.40	1.35
4	B	900	NAD	C4N-C3N	3.14	1.44	1.39
4	A	400	NAD	C4N-C3N	2.69	1.43	1.39
5	B	901	UD1	C6-N1	2.60	1.40	1.35
4	B	900	NAD	PA-O2A	-2.55	1.43	1.55
5	B	901	UD1	C2'-N2'	2.43	1.50	1.45
5	B	901	UD1	C6-C5	-2.33	1.32	1.36
4	A	400	NAD	PA-O2A	-2.24	1.45	1.55

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	900	NAD	C5N-C6N-N1N	-5.49	111.17	120.43
4	B	900	NAD	C3N-C7N-N7N	5.12	123.59	117.77
4	A	400	NAD	C5N-C4N-C3N	-4.90	113.95	120.32
4	A	400	NAD	C5N-C6N-N1N	-4.89	112.20	120.43
4	A	400	NAD	C6N-C5N-C4N	4.71	126.92	119.44
4	B	900	NAD	C5N-C4N-C3N	-4.50	114.47	120.32
4	B	900	NAD	C6N-C5N-C4N	4.49	126.56	119.44
4	B	900	NAD	O7N-C7N-N7N	-3.89	116.97	122.59
4	B	900	NAD	C6N-N1N-C2N	3.66	126.17	122.04
5	B	901	UD1	C4'-C3'-C2'	3.40	115.27	110.44
5	B	901	UD1	O7'-C7'-C8'	-2.98	116.22	122.04
4	B	900	NAD	O4D-C1D-N1N	-2.51	105.38	107.95
5	B	901	UD1	C6'-C5'-C4'	-2.45	107.08	113.00
4	A	400	NAD	C3N-C7N-N7N	2.31	120.40	117.77
4	B	900	NAD	O4D-C1D-C2D	-2.29	103.27	106.77
5	B	901	UD1	O7'-C7'-N2'	2.09	126.27	121.90
4	A	400	NAD	C2N-C3N-C4N	2.07	120.66	118.31
4	A	400	NAD	O4B-C1B-C2B	-2.03	103.66	106.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.