



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:06 PM UTC

PDB ID : 1BGV / pdb_00001bgv
Title : GLUTAMATE DEHYDROGENASE
Authors : Stillman, T.J.; Baker, P.J.; Britton, K.L.; Rice, D.W.
Deposited on : 1998-06-01
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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

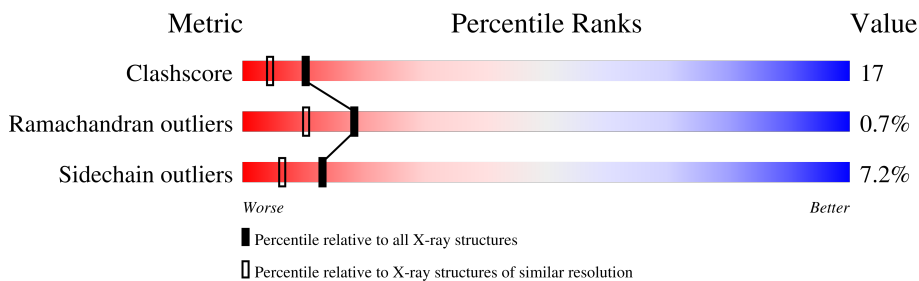
1 Overall quality at a glance

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))
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	449	

2 Entry composition [i](#)

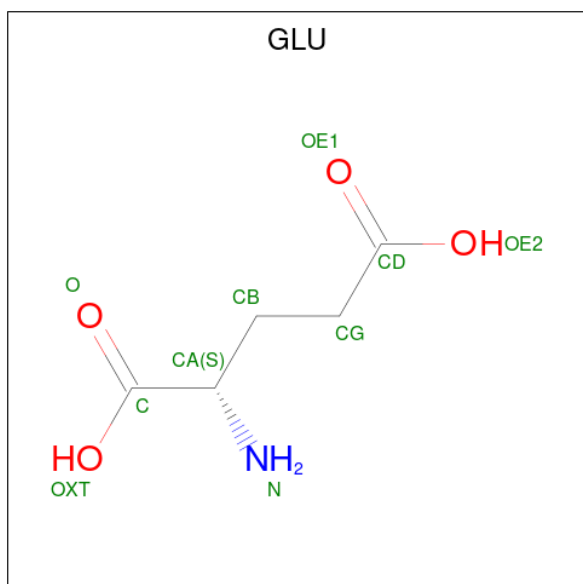
There are 3 unique types of molecules in this entry. The entry contains 3687 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 GLUTAMATE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	449	3460	2198	583	660	19	0	0	0

- Molecule 2 is GLUTAMIC ACID (CCD ID: GLU) (formula: C₅H₉NO₄).



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	217	Total	O	0	0
			217	217		

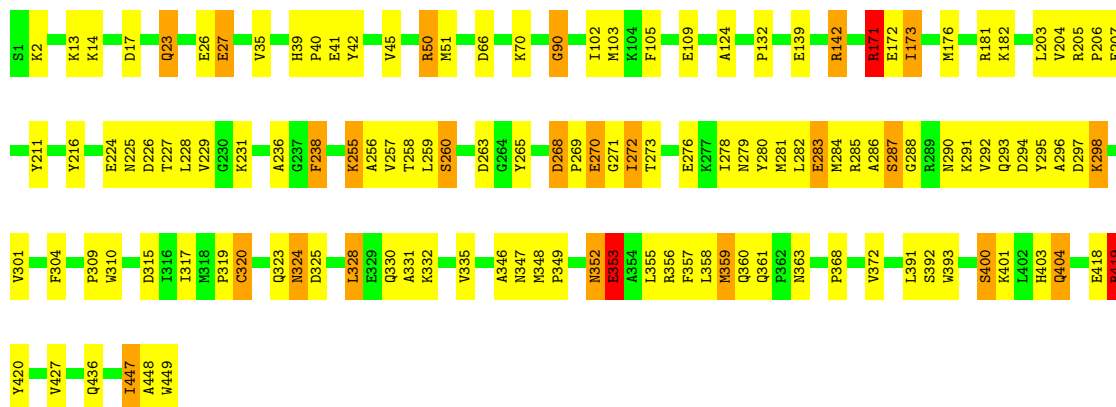
3 Residue-property plots

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. 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. 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: GLUTAMATE DEHYDROGENASE

Chain A: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	162.80Å 162.80Å 102.80Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 1.90	Depositor
% Data completeness (in resolution range)	99.0 (10.00-1.90)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT 5E	Depositor
R, R_{free}	0.173 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3687	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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	1.38	12/3535 (0.3%)	1.45	27/4781 (0.6%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	173	ILE	CG1-CD1	7.86	1.82	1.51
1	A	359	MET	SD-CE	7.18	1.97	1.79
1	A	51	MET	SD-CE	-6.25	1.64	1.79
1	A	173	ILE	CB-CG1	5.87	1.65	1.53
1	A	142	ARG	CD-NE	-5.55	1.38	1.46
1	A	176	MET	SD-CE	-5.53	1.65	1.79
1	A	102	ILE	CA-C	5.40	1.59	1.52
1	A	109	GLU	C-O	-5.31	1.18	1.24
1	A	229	VAL	CA-CB	-5.27	1.47	1.54
1	A	90	GLY	N-CA	5.16	1.50	1.45
1	A	392	SER	CA-C	-5.16	1.46	1.52
1	A	448	ALA	CA-CB	5.06	1.62	1.53

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	320	CYS	CB-CA-C	-8.83	96.27	111.45
1	A	353	GLU	N-CA-C	-8.82	101.67	111.28
1	A	142	ARG	NE-CZ-NH2	-7.43	112.51	119.20
1	A	173	ILE	CB-CG1-CD1	-7.37	98.32	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	447	ILE	CB-CA-C	-6.23	104.72	112.45
1	A	268	ASP	CA-C-N	6.04	127.38	119.84
1	A	268	ASP	C-N-CA	6.04	127.38	119.84
1	A	448	ALA	N-CA-C	-6.01	97.99	110.80
1	A	285	ARG	N-CA-C	-5.99	104.68	111.14
1	A	40	PRO	CB-CA-C	-5.98	102.88	112.21
1	A	372	VAL	CA-C-N	-5.88	113.25	122.65
1	A	372	VAL	C-N-CA	-5.88	113.25	122.65
1	A	211	TYR	CB-CA-C	-5.87	101.04	110.79
1	A	26	GLU	CA-C-N	-5.79	112.52	120.28
1	A	26	GLU	C-N-CA	-5.79	112.52	120.28
1	A	203	LEU	N-CA-C	-5.73	102.92	110.43
1	A	142	ARG	NE-CZ-NH1	5.58	127.08	121.50
1	A	419	ARG	CA-C-N	-5.51	113.86	122.73
1	A	419	ARG	C-N-CA	-5.51	113.86	122.73
1	A	171	ARG	CG-CD-NE	-5.35	100.22	112.00
1	A	27	GLU	CB-CA-C	5.35	119.67	110.79
1	A	50	ARG	NE-CZ-NH1	5.26	126.76	121.50
1	A	23	GLN	CB-CA-C	-5.26	102.06	110.79
1	A	363	ASN	N-CA-C	5.24	119.52	113.18
1	A	238	PHE	N-CA-C	-5.23	104.88	113.50
1	A	393	TRP	N-CA-C	-5.21	102.75	110.46
1	A	297	ASP	N-CA-C	-5.16	105.55	111.07

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	142	ARG	Sidechain

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	3460	0	3390	116	1
2	A	10	0	5	1	0
3	A	217	0	0	9	5

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3687	0	3395	117	5

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 17.

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:A:173:ILE:CD1	1:A:173:ILE:CG1	1.82	1.50
1:A:352:ASN:H	1:A:352:ASN:ND2	1.42	1.10
1:A:352:ASN:HD22	1:A:352:ASN:N	1.58	1.01
1:A:352:ASN:ND2	1:A:352:ASN:N	2.03	0.95
1:A:280:TYR:CZ	1:A:298:LYS:HD3	2.07	0.90
1:A:352:ASN:ND2	3:A:672:HOH:O	2.06	0.88
1:A:284:MET:HE1	1:A:295:TYR:HB2	1.57	0.86
1:A:227:THR:O	1:A:231:LYS:HE3	1.76	0.85
1:A:284:MET:CE	1:A:295:TYR:HB2	2.14	0.77
1:A:173:ILE:CD1	1:A:173:ILE:CB	2.63	0.76
1:A:328:LEU:HD13	1:A:332:LYS:HE3	1.68	0.75
1:A:259:LEU:CD1	1:A:272:ILE:HD13	2.18	0.73
1:A:352:ASN:H	1:A:352:ASN:HD22	0.76	0.72
1:A:335:VAL:HG21	1:A:357:PHE:CE2	2.24	0.72
1:A:259:LEU:HD11	1:A:272:ILE:HD13	1.68	0.72
1:A:182:LYS:CE	3:A:710:HOH:O	2.36	0.71
1:A:132:PRO:HB3	3:A:636:HOH:O	1.90	0.69
1:A:296:ALA:HA	1:A:301:VAL:CG2	2.23	0.68
1:A:259:LEU:HG	1:A:272:ILE:CD1	2.23	0.68
1:A:284:MET:HE1	1:A:295:TYR:CD2	2.32	0.65
1:A:182:LYS:HG3	3:A:710:HOH:O	1.96	0.64
1:A:348:MET:O	1:A:348:MET:HG2	1.97	0.64
1:A:226:ASP:CG	1:A:231:LYS:HZ1	2.06	0.63
1:A:41:GLU:HG2	1:A:42:TYR:N	2.12	0.63
1:A:355:LEU:HG	1:A:359:MET:HE2	1.80	0.63
1:A:400:SER:O	1:A:404:GLN:CG	2.47	0.62
1:A:265:TYR:CD1	1:A:309:PRO:HB3	2.35	0.61
1:A:35:VAL:CG1	1:A:436:GLN:HE21	2.14	0.61
1:A:283:GLU:O	1:A:286:ALA:HB3	2.01	0.60
1:A:171:ARG:CD	1:A:172:GLU:OE2	2.50	0.60
1:A:257:VAL:HG23	1:A:258:THR:HG23	1.83	0.60
1:A:352:ASN:O	1:A:353:GLU:C	2.45	0.59
1:A:348:MET:O	1:A:348:MET:CG	2.50	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:ASN:O	1:A:356:ARG:HG2	2.02	0.59
1:A:280:TYR:CD2	1:A:284:MET:HE2	2.37	0.59
1:A:296:ALA:HA	1:A:301:VAL:HG22	1.83	0.59
1:A:296:ALA:HB1	1:A:301:VAL:O	2.04	0.58
1:A:171:ARG:HD3	1:A:172:GLU:OE2	2.03	0.57
1:A:226:ASP:OD2	1:A:231:LYS:NZ	2.36	0.57
1:A:286:ALA:O	1:A:288:GLY:N	2.38	0.57
1:A:268:ASP:OD1	1:A:270:GLU:HG3	2.05	0.56
1:A:284:MET:HE1	1:A:295:TYR:HD2	1.69	0.56
1:A:284:MET:CE	1:A:295:TYR:HD2	2.19	0.55
1:A:205:ARG:N	1:A:206:PRO:HD2	2.22	0.55
1:A:323:GLN:HA	1:A:348:MET:O	2.07	0.55
1:A:255:LYS:HB2	1:A:273:THR:CG2	2.36	0.55
1:A:400:SER:O	1:A:404:GLN:HG2	2.07	0.55
1:A:50:ARG:HD2	1:A:449:TRP:OXT	2.06	0.54
1:A:263:ASP:OD1	1:A:292:VAL:HG12	2.08	0.54
1:A:182:LYS:HE3	3:A:710:HOH:O	2.06	0.54
1:A:259:LEU:CD1	1:A:272:ILE:CD1	2.85	0.53
1:A:280:TYR:O	1:A:284:MET:HG3	2.09	0.52
2:A:501:GLU:C	3:A:502:HOH:O	2.53	0.52
1:A:323:GLN:O	1:A:324:ASN:C	2.53	0.52
1:A:400:SER:O	1:A:404:GLN:HG3	2.11	0.51
1:A:401:LYS:HA	1:A:404:GLN:HG3	1.93	0.51
1:A:231:LYS:HA	1:A:315:ASP:OD2	2.10	0.50
1:A:205:ARG:HB3	1:A:206:PRO:HD3	1.93	0.50
1:A:304:PHE:CD1	1:A:304:PHE:N	2.79	0.50
1:A:236:ALA:HB2	1:A:310:TRP:CZ2	2.47	0.50
1:A:280:TYR:OH	1:A:298:LYS:HD3	2.12	0.49
1:A:103:MET:HA	1:A:103:MET:HE2	1.95	0.49
1:A:335:VAL:HG21	1:A:357:PHE:HE2	1.74	0.49
1:A:323:GLN:O	1:A:325:ASP:N	2.46	0.49
1:A:182:LYS:HE2	3:A:710:HOH:O	2.06	0.48
1:A:356:ARG:O	1:A:360:GLN:HG3	2.12	0.48
1:A:70:LYS:HE2	3:A:695:HOH:O	2.13	0.48
1:A:255:LYS:HB2	1:A:273:THR:HG21	1.94	0.48
1:A:293:GLN:O	1:A:296:ALA:N	2.46	0.48
1:A:228:LEU:HD23	1:A:228:LEU:HA	1.66	0.48
1:A:357:PHE:O	1:A:361:GLN:HG2	2.13	0.48
1:A:268:ASP:O	1:A:271:GLY:N	2.46	0.48
1:A:280:TYR:HD2	1:A:284:MET:HE2	1.77	0.48
1:A:35:VAL:HG11	1:A:436:GLN:HE21	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:317:ILE:HG22	1:A:319:PRO:HD3	1.96	0.48
1:A:173:ILE:HD13	1:A:173:ILE:HA	1.96	0.47
1:A:35:VAL:O	1:A:39:HIS:HD2	1.97	0.47
1:A:279:ASN:O	1:A:282:LEU:HB2	2.14	0.47
1:A:205:ARG:CB	1:A:206:PRO:HD3	2.45	0.47
1:A:256:ALA:O	1:A:271:GLY:HA2	2.14	0.47
1:A:259:LEU:HD11	1:A:272:ILE:CD1	2.41	0.46
1:A:278:ILE:O	1:A:281:MET:HG3	2.15	0.46
1:A:348:MET:HA	3:A:560:HOH:O	2.15	0.46
1:A:205:ARG:CB	1:A:206:PRO:CD	2.94	0.46
1:A:284:MET:HE1	1:A:295:TYR:CB	2.36	0.46
1:A:310:TRP:HA	1:A:310:TRP:CE3	2.51	0.46
1:A:328:LEU:O	1:A:332:LYS:HG3	2.15	0.46
1:A:403:HIS:CD2	1:A:403:HIS:C	2.94	0.45
1:A:207:GLU:OE1	1:A:207:GLU:N	2.45	0.45
1:A:310:TRP:HB2	1:A:330:GLN:HE21	1.81	0.45
1:A:331:ALA:O	1:A:335:VAL:HG23	2.16	0.45
1:A:294:ASP:O	1:A:298:LYS:HB2	2.17	0.45
1:A:296:ALA:HA	1:A:301:VAL:HG23	1.94	0.45
1:A:356:ARG:HG2	1:A:356:ARG:H	1.30	0.44
1:A:224:GLU:O	1:A:225:ASN:C	2.59	0.44
1:A:259:LEU:HG	1:A:272:ILE:HD12	2.00	0.44
1:A:295:TYR:C	1:A:295:TYR:CD1	2.96	0.44
1:A:348:MET:N	1:A:349:PRO:HD3	2.32	0.43
1:A:66:ASP:HB2	1:A:139:GLU:OE2	2.18	0.43
1:A:171:ARG:HD2	1:A:172:GLU:OE2	2.17	0.43
1:A:310:TRP:HA	1:A:310:TRP:HE3	1.84	0.43
1:A:403:HIS:HD2	1:A:403:HIS:O	2.02	0.43
1:A:259:LEU:CG	1:A:272:ILE:CD1	2.95	0.43
1:A:358:LEU:HA	1:A:358:LEU:HD23	1.77	0.42
1:A:236:ALA:HA	1:A:260:SER:HB3	2.02	0.42
1:A:368:PRO:HG3	1:A:427:VAL:HA	2.01	0.42
1:A:45:VAL:HB	1:A:447:ILE:HD11	2.02	0.42
1:A:238:PHE:HA	1:A:281:MET:HE3	2.01	0.42
1:A:268:ASP:HA	1:A:269:PRO:HD2	1.55	0.42
1:A:268:ASP:OD1	1:A:268:ASP:C	2.61	0.41
1:A:419:ARG:HG2	1:A:420:TYR:CD1	2.55	0.41
1:A:320:CYS:C	1:A:346:ALA:HB2	2.46	0.41
1:A:90:GLY:HA3	1:A:124:ALA:O	2.20	0.41
1:A:205:ARG:C	1:A:205:ARG:HD3	2.46	0.41
1:A:216:TYR:CD1	1:A:216:TYR:C	2.98	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:ARG:N	1:A:206:PRO:CD	2.83	0.41
1:A:23:GLN:HE21	1:A:23:GLN:HB3	1.60	0.41

All (5) 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	Interatomic distance (Å)	Clash overlap (Å)
3:A:571:HOH:O	3:A:571:HOH:O[6_555]	0.78	1.42
3:A:652:HOH:O	3:A:652:HOH:O[6_556]	1.51	0.69
3:A:643:HOH:O	3:A:643:HOH:O[6_555]	1.54	0.66
3:A:586:HOH:O	3:A:586:HOH:O[5_555]	1.77	0.43
1:A:17:ASP:CB	3:A:692:HOH:O[6_555]	2.16	0.04

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	447/449 (100%)	429 (96%)	15 (3%)	3 (1%)	18 10

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	287	SER
1	A	204	VAL
1	A	324	ASN

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	358/358 (100%)	332 (93%)	26 (7%)	13 6

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	13	LYS
1	A	14	LYS
1	A	27	GLU
1	A	105	PHE
1	A	171	ARG
1	A	181	ARG
1	A	255	LYS
1	A	260	SER
1	A	270	GLU
1	A	272	ILE
1	A	276	GLU
1	A	283	GLU
1	A	287	SER
1	A	290	ASN
1	A	291	LYS
1	A	298	LYS
1	A	328	LEU
1	A	347	ASN
1	A	352	ASN
1	A	353	GLU
1	A	391	LEU
1	A	400	SER
1	A	404	GLN
1	A	418	GLU
1	A	419	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	23	GLN
1	A	68	ASN
1	A	279	ASN
1	A	293	GLN

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Mol	Chain	Res	Type
1	A	312	GLN
1	A	330	GLN
1	A	352	ASN
1	A	360	GLN
1	A	363	ASN
1	A	403	HIS
1	A	436	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

1 ligand is 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	GLU	A	501	-	8,9,9	1.29	2 (25%)	8,11,11	1.13	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	GLU	A	501	-	-	3/9/9/9	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	GLU	OE1-CD	2.23	1.29	1.22
2	A	501	GLU	OXT-C	-2.15	1.23	1.30

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	GLU	O-C-CA-CB
2	A	501	GLU	OE2-CD-CG-CB
2	A	501	GLU	OE1-CD-CG-CB

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	GLU	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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.