



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 06:18 am BST

PDB ID : 3IH9
Title : Crystal Structure Analysis of Mglu in its tris form
Authors : Yoshimune, K.; Shirakihara, Y.
Deposited on : 2009-07-29
Resolution : 2.50 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

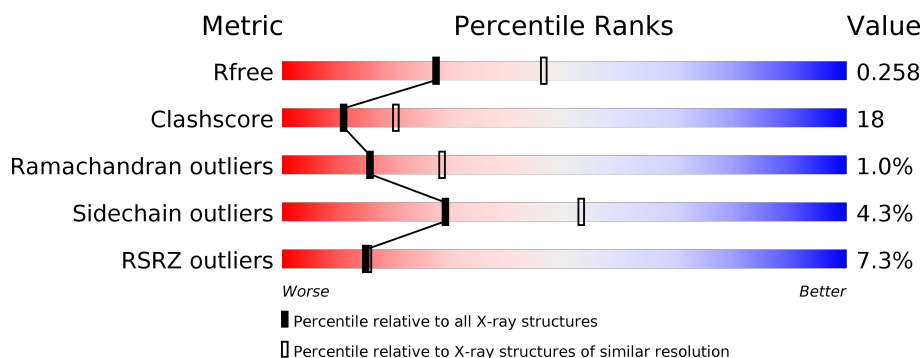
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
R_{free}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 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	456	<div> <div>6%</div> <div> <div></div> <div>64%</div> <div>32%</div> <div>• •</div> </div> </div>
1	B	456	<div> <div>8%</div> <div> <div></div> <div>65%</div> <div>32%</div> <div>• •</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6950 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.

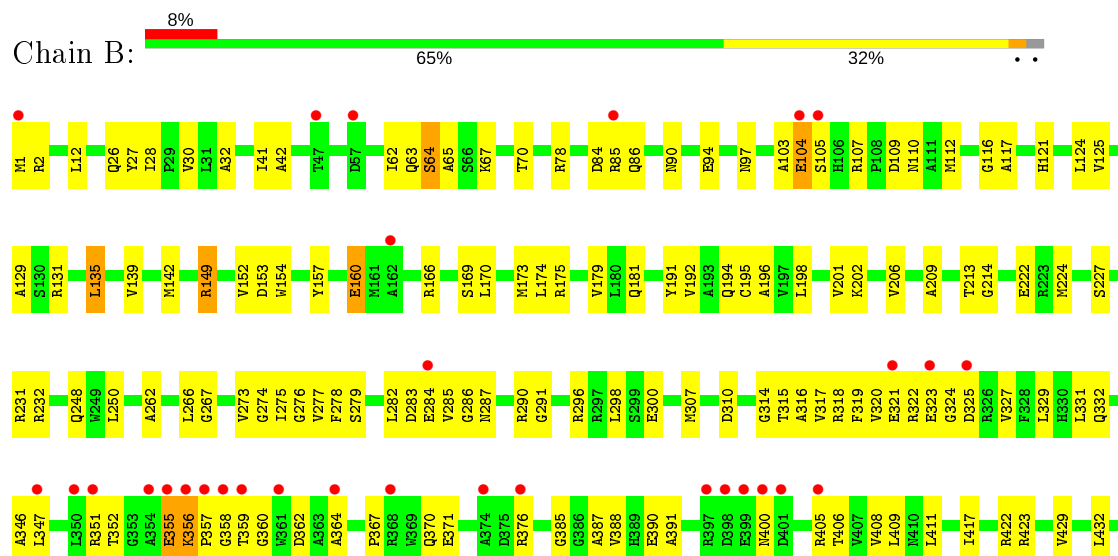
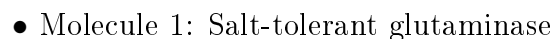
- Molecule 1 is a protein called Salt-tolerant glutaminase.

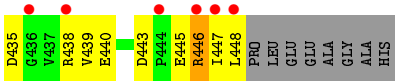
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	448	Total	C	N	O	S	0	0	0
			3331	2067	611	638	15			
1	B	448	Total	C	N	O	S	0	0	0
			3331	2067	611	638	15			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	150	Total	O	0	0
			150	150		
2	B	138	Total	O	0	0
			138	138		

- Molecule 1: Salt-tolerant glutaminase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	117.70 Å 141.39 Å 75.23 Å 90.00° 104.48° 90.00°	Depositor
Resolution (Å)	19.90 – 2.50 19.89 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.3 (19.90-2.50) 98.3 (19.89-2.40)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.08 (at 2.41 Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.217 , 0.258 0.217 , 0.258	Depositor DCC
R_{free} test set	2263 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	45.7	Xtriage
Anisotropy	0.192	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6950	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.69% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.39	0/3384	0.66	2/4594 (0.0%)
1	B	0.40	0/3384	0.65	1/4594 (0.0%)
All	All	0.39	0/6768	0.66	3/9188 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	266	LEU	CA-CB-CG	5.75	128.53	115.30
1	A	124	LEU	N-CA-C	5.26	125.19	111.00
1	B	124	LEU	N-CA-C	5.07	124.70	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3331	0	3312	123	0
1	B	3331	0	3312	122	0
2	A	150	0	0	15	0
2	B	138	0	0	7	0
All	All	6950	0	6624	242	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 18.

All (242) 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:104:GLU:N	1:B:104:GLU:OE2	1.72	1.19
1:A:62:ILE:HG21	1:A:142:MET:HE1	1.38	1.05
1:B:400:ASN:ND2	2:B:519:HOH:O	2.01	0.92
1:A:277:VAL:HG11	1:A:291:GLY:HA2	1.55	0.89
1:A:201:VAL:HG23	2:A:467:HOH:O	1.74	0.87
1:B:62:ILE:HD13	1:B:142:MET:HE3	1.55	0.86
1:B:411:LEU:HD11	1:B:448:LEU:HD12	1.57	0.85
1:B:62:ILE:HD13	1:B:142:MET:CE	2.13	0.77
1:A:397:ARG:NH1	1:A:397:ARG:HB3	2.00	0.76
1:B:320:VAL:HG22	1:B:329:LEU:HD12	1.69	0.73
1:A:376:ARG:HG2	1:A:389:HIS:HB2	1.69	0.73
1:A:322:ARG:HD2	1:A:327:VAL:HG22	1.69	0.73
1:B:149:ARG:HH11	1:B:149:ARG:HB3	1.54	0.72
1:A:105:SER:O	1:A:106:HIS:HB2	1.89	0.71
1:A:446:ARG:HG3	2:A:565:HOH:O	1.91	0.71
1:A:112:MET:HE2	1:A:258:LYS:HD3	1.73	0.70
1:B:63:GLN:HA	1:B:195:CYS:HB3	1.73	0.69
1:A:134:ARG:HH11	1:A:134:ARG:HG2	1.57	0.69
1:A:52:CYS:SG	1:A:202:LYS:HE2	2.33	0.69
1:A:284:GLU:OE1	1:A:285:VAL:HG13	1.92	0.69
1:A:279:SER:OG	1:A:290:ARG:HD2	1.93	0.69
1:B:279:SER:OG	1:B:290:ARG:HD2	1.93	0.68
1:A:445:GLU:O	1:A:447:ILE:HG13	1.94	0.68
1:A:252:ASP:OD1	2:A:492:HOH:O	2.11	0.68
1:B:110:ASN:ND2	1:B:112:MET:H	1.92	0.68
1:A:259:SER:HB2	1:A:265:VAL:HG22	1.76	0.67
1:B:160:GLU:HG2	1:B:192:VAL:HG13	1.75	0.67
1:B:356:LYS:HD3	1:B:356:LYS:O	1.94	0.67
1:B:376:ARG:HB2	1:B:376:ARG:NH1	2.10	0.67
1:B:408:VAL:HG22	1:B:440:GLU:HG3	1.75	0.67
1:A:356:LYS:O	1:A:356:LYS:HD3	1.95	0.67
1:A:397:ARG:HH11	1:A:397:ARG:CB	2.08	0.67
1:A:324:GLY:N	2:A:465:HOH:O	2.27	0.66
1:B:376:ARG:HB2	1:B:376:ARG:HH11	1.59	0.66
1:B:231:ARG:NH2	2:B:515:HOH:O	2.27	0.66
1:B:423:ARG:HD3	2:B:544:HOH:O	1.95	0.66
1:B:105:SER:HB2	2:B:501:HOH:O	1.96	0.65
1:A:397:ARG:HH11	1:A:397:ARG:HB3	1.61	0.65
1:A:99:LEU:HD23	1:A:102:GLU:OE1	1.97	0.65
1:A:160:GLU:HG2	1:A:192:VAL:HG13	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:160:GLU:HG2	1:A:192:VAL:CG1	2.26	0.65
1:B:125:VAL:HB	1:B:129:ALA:CB	2.27	0.64
1:B:411:LEU:CD1	1:B:448:LEU:HD12	2.25	0.64
1:B:320:VAL:HB	1:B:390:GLU:HB3	1.79	0.64
1:A:296:ARG:O	1:A:300:GLU:HG3	1.97	0.64
1:A:322:ARG:CD	1:A:327:VAL:HG22	2.27	0.64
1:A:107:ARG:NH2	2:A:606:HOH:O	2.31	0.63
1:A:110:ASN:HD22	1:A:112:MET:H	1.46	0.63
1:A:277:VAL:CG1	1:A:291:GLY:HA2	2.28	0.63
1:B:160:GLU:HG2	1:B:192:VAL:CG1	2.29	0.63
1:A:62:ILE:HD13	1:A:142:MET:HE1	1.80	0.63
1:B:104:GLU:N	1:B:104:GLU:CD	2.53	0.62
1:A:397:ARG:HH11	1:A:397:ARG:CA	2.13	0.62
1:A:62:ILE:HD13	1:A:142:MET:CE	2.30	0.62
1:B:201:VAL:HG13	1:B:276:GLY:C	2.19	0.61
1:B:320:VAL:HG11	1:B:391:ALA:HA	1.83	0.61
1:A:63:GLN:HA	1:A:195:CYS:HB3	1.83	0.61
1:B:209:ALA:O	1:B:213:THR:HG23	2.01	0.61
1:A:173:MET:O	1:A:176:SER:HB3	2.01	0.60
1:A:153:ASP:HB2	1:A:198:LEU:HD11	1.83	0.60
1:B:135:LEU:O	1:B:139:VAL:HG23	2.00	0.60
1:B:296:ARG:O	1:B:300:GLU:HG3	2.01	0.60
1:A:29:PRO:O	1:A:33:GLU:HG2	2.02	0.60
1:A:408:VAL:HG22	1:A:440:GLU:HG3	1.83	0.59
1:B:406:THR:HG23	1:B:438:ARG:HB3	1.84	0.59
1:B:282:LEU:HA	1:B:287:ASN:O	2.02	0.59
1:A:317:VAL:HG21	1:A:388:VAL:HG22	1.85	0.59
1:B:27:TYR:CD1	1:B:28:ILE:HG23	2.38	0.59
1:B:28:ILE:HD12	1:B:30:VAL:HG12	1.85	0.59
1:A:121:HIS:HD2	1:A:194:GLN:OE1	1.86	0.58
1:B:351:ARG:HG3	1:B:351:ARG:HH21	1.69	0.58
1:B:103:ALA:C	1:B:105:SER:H	2.06	0.58
1:B:325:ASP:OD2	1:B:405:ARG:HG3	2.04	0.58
1:B:282:LEU:HD22	1:B:286:GLY:O	2.04	0.57
1:B:351:ARG:HD2	1:B:355:GLU:HB3	1.86	0.57
1:B:67:LYS:HE2	1:B:117:ALA:HB2	1.86	0.57
1:A:27:TYR:CD1	1:A:28:ILE:HG23	2.40	0.57
1:A:5:ILE:HB	1:A:6:PRO:HD3	1.86	0.56
1:B:352:THR:O	1:B:435:ASP:HB3	2.05	0.56
1:B:70:THR:HG21	1:B:194:GLN:CD	2.25	0.56
1:B:121:HIS:HD2	1:B:194:GLN:OE1	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:429:VAL:HG13	1:B:439:VAL:HG11	1.88	0.56
1:A:429:VAL:HG13	1:A:439:VAL:HG11	1.86	0.55
1:A:110:ASN:ND2	1:A:112:MET:H	2.03	0.55
1:A:347:LEU:HD22	1:A:432:LEU:HD11	1.88	0.55
1:B:64:SER:HB2	2:B:457:HOH:O	2.06	0.55
1:A:361:TRP:CZ3	1:A:370:GLN:HG2	2.40	0.55
1:A:41:ILE:HA	1:A:201:VAL:HG11	1.89	0.55
1:B:370:GLN:HE21	1:B:370:GLN:HA	1.71	0.55
1:A:400:ASN:ND2	2:A:481:HOH:O	2.39	0.55
1:B:103:ALA:C	1:B:105:SER:N	2.59	0.54
1:B:201:VAL:HG13	1:B:276:GLY:O	2.07	0.54
1:B:62:ILE:HG21	1:B:142:MET:CE	2.37	0.54
1:B:355:GLU:HG3	1:B:355:GLU:O	2.07	0.54
1:A:112:MET:CE	1:A:258:LYS:HD3	2.37	0.53
1:B:376:ARG:CB	1:B:376:ARG:HH11	2.21	0.53
1:B:110:ASN:HD22	1:B:112:MET:H	1.55	0.53
1:A:324:GLY:CA	2:A:465:HOH:O	2.56	0.53
1:A:310:ASP:HB3	2:A:567:HOH:O	2.09	0.53
1:B:125:VAL:HB	1:B:129:ALA:HB2	1.91	0.53
1:B:320:VAL:HG22	1:B:329:LEU:CD1	2.36	0.53
1:A:320:VAL:HB	1:A:390:GLU:HB3	1.90	0.53
1:A:166:ARG:O	1:A:169:SER:HB3	2.09	0.52
1:B:110:ASN:HD21	1:B:112:MET:HB2	1.74	0.52
1:A:326:ARG:HG3	2:A:465:HOH:O	2.09	0.52
1:A:85:ARG:HD3	2:A:528:HOH:O	2.10	0.52
1:B:62:ILE:HG13	1:B:65:ALA:HB3	1.91	0.52
1:B:317:VAL:HG21	1:B:388:VAL:HG22	1.92	0.52
1:B:408:VAL:HG22	1:B:440:GLU:CG	2.40	0.52
1:B:315:THR:HG22	1:B:315:THR:O	2.10	0.51
1:A:172:HIS:NE2	2:A:457:HOH:O	2.34	0.51
1:A:422:ARG:CD	1:A:448:LEU:O	2.59	0.51
1:B:135:LEU:O	1:B:135:LEU:HD22	2.09	0.51
1:B:443:ASP:OD1	1:B:446:ARG:HA	2.11	0.51
1:A:355:GLU:HG3	1:A:355:GLU:O	2.11	0.51
1:B:62:ILE:HG21	1:B:142:MET:HE3	1.93	0.50
1:B:351:ARG:HD2	1:B:355:GLU:CB	2.42	0.50
1:B:445:GLU:O	1:B:447:ILE:HG13	2.10	0.50
1:A:315:THR:O	1:A:315:THR:HG22	2.11	0.50
1:A:422:ARG:HG3	1:A:448:LEU:O	2.12	0.50
1:A:397:ARG:NH1	1:A:397:ARG:CB	2.70	0.50
1:A:401:ASP:N	1:A:401:ASP:OD2	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:370:GLN:NE2	1:B:370:GLN:HA	2.25	0.50
1:B:202:LYS:O	1:B:206:VAL:HG23	2.12	0.50
1:B:367:PRO:O	1:B:371:GLU:HG3	2.12	0.50
1:A:321:GLU:HB2	1:A:328:PHE:HB2	1.93	0.49
1:A:322:ARG:HG2	1:A:322:ARG:HH21	1.77	0.49
1:A:361:TRP:HZ3	1:A:370:GLN:HG2	1.78	0.49
1:B:322:ARG:HD3	1:B:327:VAL:HG22	1.93	0.49
1:B:86:GLN:NE2	1:B:179:VAL:HA	2.28	0.49
1:B:214:GLY:O	1:B:227:SER:HA	2.13	0.49
1:A:84:ASP:O	1:A:107:ARG:NH1	2.46	0.49
1:A:104:GLU:OE2	1:A:104:GLU:N	2.33	0.49
1:B:325:ASP:OD1	1:B:405:ARG:HD2	2.12	0.48
1:B:411:LEU:HD11	1:B:448:LEU:CD1	2.37	0.48
1:A:339:GLY:O	1:A:343:VAL:HG23	2.13	0.48
1:A:393:ALA:O	1:A:396:ALA:HB3	2.13	0.48
1:B:78:ARG:NH2	2:B:532:HOH:O	2.45	0.48
1:B:166:ARG:O	1:B:169:SER:HB3	2.14	0.48
1:A:273:VAL:HG22	1:A:274:GLY:N	2.29	0.48
1:A:101:LEU:O	1:A:106:HIS:HA	2.14	0.47
1:A:134:ARG:HH11	1:A:134:ARG:CG	2.24	0.47
1:B:322:ARG:CD	1:B:327:VAL:HG22	2.44	0.47
1:A:191:TYR:O	1:A:194:GLN:HB3	2.14	0.47
1:A:201:VAL:CG1	1:A:276:GLY:O	2.62	0.47
1:B:262:ALA:HB3	2:B:465:HOH:O	2.14	0.47
1:A:184:ALA:O	1:A:188:VAL:HG23	2.14	0.47
1:A:35:ASP:OD2	1:A:38:ARG:HG3	2.15	0.47
1:A:181:GLN:N	1:A:181:GLN:CD	2.68	0.47
1:B:319:PHE:CD2	1:B:321:GLU:HG3	2.50	0.47
1:A:336:ARG:HA	1:A:418:ASP:OD1	2.15	0.47
1:B:277:VAL:HG11	1:B:291:GLY:HA2	1.97	0.47
1:A:318:ARG:NH1	1:A:332:GLN:OE1	2.47	0.47
1:B:283:ASP:OD1	1:B:285:VAL:HG22	2.15	0.47
1:B:445:GLU:O	1:B:446:ARG:C	2.52	0.47
1:B:42:ALA:HB3	1:B:201:VAL:HG12	1.96	0.46
1:A:9:LEU:O	1:A:13:VAL:HG23	2.15	0.46
1:B:222:GLU:HB3	1:B:224:MET:HE3	1.97	0.46
1:B:42:ALA:CB	1:B:201:VAL:HG12	2.45	0.46
1:A:284:GLU:HG2	2:A:495:HOH:O	2.15	0.46
1:A:319:PHE:CD2	1:A:321:GLU:HG2	2.51	0.46
1:A:3:HIS:O	1:A:6:PRO:HD2	2.16	0.46
1:B:152:VAL:HG13	1:B:196:ALA:HB3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:70:THR:HG21	1:B:194:GLN:OE1	2.16	0.46
1:B:358:GLY:O	1:B:360:GLY:N	2.50	0.45
1:B:417:ILE:O	1:B:422:ARG:NH1	2.49	0.45
1:B:85:ARG:HH11	1:B:85:ARG:HG3	1.81	0.45
1:A:209:ALA:O	1:A:213:THR:HG23	2.17	0.45
1:B:110:ASN:ND2	1:B:112:MET:HB2	2.31	0.45
1:A:175:ARG:HD2	1:A:182:ASP:O	2.16	0.45
1:A:358:GLY:O	1:A:360:GLY:N	2.50	0.45
1:A:165:ASP:OD2	1:B:175:ARG:NH1	2.49	0.45
1:A:56:ALA:HA	1:A:201:VAL:CG2	2.47	0.45
1:A:432:LEU:O	1:A:435:ASP:HB2	2.16	0.45
1:B:26:GLN:HA	1:B:32:ALA:HB2	1.99	0.45
1:B:273:VAL:HG22	1:B:274:GLY:N	2.31	0.44
1:A:90:ASN:O	1:A:109:ASP:HB3	2.16	0.44
1:A:362:ASP:OD1	1:A:364:ALA:N	2.48	0.44
1:B:170:LEU:O	1:B:174:LEU:HG	2.17	0.44
1:B:323:GLU:HA	1:B:323:GLU:OE1	2.17	0.44
1:B:346:ALA:HB2	1:B:388:VAL:HG13	2.00	0.44
1:A:350:LEU:HA	1:A:395:ALA:HB2	1.99	0.44
1:B:41:ILE:HD11	1:B:298:LEU:HD11	1.99	0.44
1:B:191:TYR:O	1:B:194:GLN:HB3	2.17	0.44
1:A:445:GLU:O	1:A:446:ARG:C	2.56	0.44
1:A:282:LEU:HA	1:A:287:ASN:O	2.18	0.44
1:A:405:ARG:HH11	1:A:405:ARG:HG3	1.81	0.44
1:B:27:TYR:HD1	1:B:28:ILE:HG23	1.80	0.44
1:B:2:ARG:HG3	1:B:2:ARG:O	2.18	0.44
1:B:84:ASP:O	1:B:107:ARG:NH1	2.51	0.44
1:B:154:TRP:HA	1:B:157:TYR:HB3	1.99	0.43
1:A:134:ARG:HG2	1:A:134:ARG:NH1	2.29	0.43
1:A:284:GLU:CG	2:A:495:HOH:O	2.65	0.43
1:A:20:ASN:HB2	1:A:290:ARG:NH2	2.33	0.43
1:A:240:SER:HB3	2:A:463:HOH:O	2.18	0.43
1:A:201:VAL:HG12	1:A:276:GLY:O	2.19	0.43
1:A:351:ARG:HD2	1:A:355:GLU:HB3	2.01	0.43
1:A:169:SER:OG	1:B:173:MET:HB2	2.19	0.43
1:A:110:ASN:O	1:A:116:GLY:HA3	2.18	0.43
1:A:93:GLY:HA2	1:A:243:TYR:HB3	2.01	0.43
1:A:27:TYR:CE1	1:A:28:ILE:HG12	2.53	0.43
1:B:267:GLY:HA3	1:B:275:ILE:HB	2.01	0.43
1:A:134:ARG:CG	1:A:134:ARG:NH1	2.81	0.42
1:B:316:ALA:O	1:B:331:LEU:HA	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:214:GLY:O	1:A:227:SER:HA	2.19	0.42
1:A:283:ASP:OD1	1:A:285:VAL:HG22	2.19	0.42
1:B:387:ALA:O	1:B:390:GLU:HB2	2.19	0.42
1:A:316:ALA:HB1	1:A:335:ILE:CD1	2.49	0.42
1:A:419:ASP:HA	1:A:422:ARG:NH1	2.35	0.42
1:B:201:VAL:HG23	1:B:278:PHE:HB2	2.02	0.42
1:A:138:ALA:O	1:A:141:ILE:HG22	2.19	0.42
1:A:130:SER:OG	1:A:133:GLU:HG3	2.20	0.42
1:A:423:ARG:HG2	1:A:423:ARG:HH11	1.85	0.42
1:B:110:ASN:O	1:B:116:GLY:HA3	2.20	0.42
1:B:125:VAL:HB	1:B:129:ALA:HB1	2.01	0.42
1:B:85:ARG:HB2	1:B:85:ARG:CZ	2.50	0.42
1:B:90:ASN:O	1:B:109:ASP:HB3	2.19	0.42
1:A:104:GLU:CD	1:A:104:GLU:H	2.18	0.42
1:A:422:ARG:NE	1:A:448:LEU:O	2.53	0.42
1:B:222:GLU:HG2	1:B:224:MET:CE	2.50	0.41
1:B:27:TYR:CE1	1:B:28:ILE:HG12	2.55	0.41
1:B:347:LEU:HD22	1:B:432:LEU:HD11	2.01	0.41
1:A:317:VAL:O	1:A:387:ALA:HB3	2.21	0.41
1:B:318:ARG:NH1	1:B:332:GLN:OE1	2.53	0.41
1:B:62:ILE:HG21	1:B:142:MET:HE1	2.01	0.41
1:A:77:ASP:OD2	1:A:124:LEU:O	2.39	0.41
1:B:318:ARG:O	1:B:319:PHE:HB3	2.21	0.41
1:B:362:ASP:OD1	1:B:364:ALA:N	2.45	0.41
1:A:258:LYS:HA	1:A:258:LYS:HD2	1.95	0.41
1:A:292:VAL:HG12	1:A:293:LEU:N	2.36	0.41
1:A:367:PRO:O	1:A:371:GLU:HG3	2.20	0.41
1:B:232:ARG:NH1	1:B:232:ARG:HG2	2.36	0.41
1:A:157:TYR:HD1	1:A:192:VAL:HG12	1.85	0.41
1:A:296:ARG:NH1	2:A:543:HOH:O	2.53	0.41
1:B:67:LYS:HE3	1:B:191:TYR:OH	2.21	0.41
1:B:153:ASP:HB2	1:B:198:LEU:HD11	2.03	0.41
1:B:355:GLU:C	1:B:357:PRO:HD3	2.40	0.40
1:B:322:ARG:HH21	1:B:322:ARG:HG2	1.85	0.40
1:A:6:PRO:HG3	1:A:51:HIS:HB3	2.03	0.40
1:A:201:VAL:CG1	1:A:276:GLY:C	2.89	0.40
1:B:314:GLY:HA3	1:B:385:GLY:C	2.42	0.40
1:A:97:ASN:ND2	1:B:97:ASN:ND2	2.69	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	446/456 (98%)	413 (93%)	29 (6%)	4 (1%)	17	31
1	B	446/456 (98%)	409 (92%)	32 (7%)	5 (1%)	14	26
All	All	892/912 (98%)	822 (92%)	61 (7%)	9 (1%)	15	28

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	446	ARG
1	A	359	THR
1	A	446	ARG
1	B	355	GLU
1	B	359	THR
1	B	310	ASP
1	A	444	PRO
1	B	324	GLY
1	A	324	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	336/341 (98%)	324 (96%)	12 (4%)	35	61
1	B	336/341 (98%)	319 (95%)	17 (5%)	24	45
All	All	672/682 (98%)	643 (96%)	29 (4%)	29	53

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	92	SER
1	A	135	LEU
1	A	228	ILE
1	A	248	GLN
1	A	250	LEU
1	A	266	LEU
1	A	284	GLU
1	A	326	ARG
1	A	329	LEU
1	A	351	ARG
1	A	356	LYS
1	A	397	ARG
1	B	1	MET
1	B	12	LEU
1	B	64	SER
1	B	94	GLU
1	B	104	GLU
1	B	131	ARG
1	B	135	LEU
1	B	149	ARG
1	B	160	GLU
1	B	181	GLN
1	B	248	GLN
1	B	250	LEU
1	B	266	LEU
1	B	284	GLU
1	B	307	MET
1	B	356	LYS
1	B	409	LEU

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	26	GLN
1	A	86	GLN
1	A	110	ASN
1	A	121	HIS
1	A	248	GLN
1	B	86	GLN
1	B	90	ASN
1	B	110	ASN

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Mol	Chain	Res	Type
1	B	121	HIS
1	B	181	GLN
1	B	248	GLN
1	B	370	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	448/456 (98%)	0.34	28 (6%)	20 21	24, 43, 81, 108	0
1	B	448/456 (98%)	0.39	37 (8%)	11 11	26, 45, 81, 107	0
All	All	896/912 (98%)	0.37	65 (7%)	15 15	24, 44, 81, 108	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	357	PRO	7.1
1	A	400	ASN	6.9
1	B	357	PRO	6.7
1	A	359	THR	5.7
1	B	358	GLY	5.3
1	B	359	THR	5.0
1	B	356	LYS	4.8
1	A	397	ARG	4.8
1	A	448	LEU	4.7
1	A	356	LYS	4.7
1	A	446	ARG	4.6
1	A	355	GLU	4.6
1	B	355	GLU	4.3
1	A	399	GLU	4.3
1	B	104	GLU	4.2
1	A	401	ASP	4.2
1	B	400	ASN	4.1
1	B	399	GLU	4.0
1	B	446	ARG	4.0
1	B	398	ASP	3.9
1	A	351	ARG	3.9
1	A	374	ALA	3.8
1	B	447	ILE	3.8
1	A	104	GLU	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	401	ASP	3.5
1	A	358	GLY	3.4
1	B	444	PRO	3.3
1	A	438	ARG	3.2
1	B	374	ALA	3.1
1	A	354	ALA	2.9
1	B	354	ALA	2.8
1	B	105	SER	2.7
1	A	347	LEU	2.6
1	B	350	LEU	2.6
1	A	284	GLU	2.6
1	A	103	ALA	2.6
1	B	368	ARG	2.6
1	B	323	GLU	2.5
1	B	448	LEU	2.5
1	A	375	ASP	2.4
1	B	284	GLU	2.4
1	A	398	ASP	2.3
1	B	376	ARG	2.3
1	B	364	ALA	2.3
1	B	436	GLY	2.3
1	B	1	MET	2.2
1	B	85	ARG	2.2
1	B	438	ARG	2.2
1	B	57	ASP	2.2
1	B	351	ARG	2.2
1	B	321	GLU	2.2
1	A	436	GLY	2.2
1	A	262	ALA	2.1
1	A	402	GLY	2.1
1	B	47	THR	2.1
1	B	361	TRP	2.1
1	B	405	ARG	2.1
1	A	447	ILE	2.1
1	B	162	ALA	2.1
1	A	102	GLU	2.0
1	B	397	ARG	2.0
1	A	364	ALA	2.0
1	B	325	ASP	2.0
1	B	347	LEU	2.0
1	A	321	GLU	2.0

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 carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.