



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:12 AM GMT

PDB ID : 2ZY9
Title : Improved crystal structure of magnesium transporter MgtE
Authors : Nureki, O.; Ishitani, R.; Hattori, M.
Deposited on : 2009-01-19
Resolution : 2.94 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
<http://wwpdb.org/validation/2016/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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

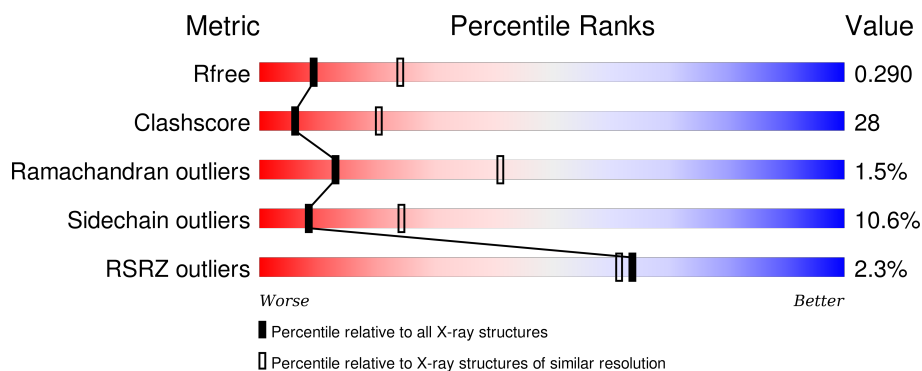
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.94 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}	91344	2057 (2.98-2.90)
Clashscore	102246	2308 (2.98-2.90)
Ramachandran outliers	100387	2245 (2.98-2.90)
Sidechain outliers	100360	2247 (2.98-2.90)
RSRZ outliers	91569	2065 (2.98-2.90)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	473	<div> <div>47%</div> <div>38%</div> <div>5%</div> <div>10%</div> </div>
1	B	473	<div> <div>3%</div> <div>46%</div> <div>39%</div> <div>5%</div> <div>9%</div> </div>

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
2	MG	A	451	-	-	-	X
2	MG	A	454	-	-	-	X
2	MG	A	455	-	-	-	X
2	MG	A	456	-	-	-	X
2	MG	B	451	-	-	-	X
2	MG	B	454	-	-	-	X

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6468 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 Mg2+ transporter MgtE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	427	Total	C	N	O	S	0	0	0
			3242	2105	525	605	7			
1	B	429	Total	C	N	O	S	0	0	0
			3213	2094	507	605	7			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	EXPRESSION TAG	UNP Q5SMG8
A	-21	GLY	-	EXPRESSION TAG	UNP Q5SMG8
A	-20	SER	-	EXPRESSION TAG	UNP Q5SMG8
A	-19	SER	-	EXPRESSION TAG	UNP Q5SMG8
A	-18	HIS	-	EXPRESSION TAG	UNP Q5SMG8
A	-17	HIS	-	EXPRESSION TAG	UNP Q5SMG8
A	-16	HIS	-	EXPRESSION TAG	UNP Q5SMG8
A	-15	HIS	-	EXPRESSION TAG	UNP Q5SMG8
A	-14	HIS	-	EXPRESSION TAG	UNP Q5SMG8
A	-13	HIS	-	EXPRESSION TAG	UNP Q5SMG8
A	-12	SER	-	EXPRESSION TAG	UNP Q5SMG8
A	-11	SER	-	EXPRESSION TAG	UNP Q5SMG8
A	-10	GLY	-	EXPRESSION TAG	UNP Q5SMG8
A	-9	LEU	-	EXPRESSION TAG	UNP Q5SMG8
A	-8	GLY	-	EXPRESSION TAG	UNP Q5SMG8
A	-7	VAL	-	EXPRESSION TAG	UNP Q5SMG8
A	-6	LEU	-	EXPRESSION TAG	UNP Q5SMG8
A	-5	PRO	-	EXPRESSION TAG	UNP Q5SMG8
A	-4	GLY	-	EXPRESSION TAG	UNP Q5SMG8
A	-3	GLY	-	EXPRESSION TAG	UNP Q5SMG8
A	-2	PRO	-	EXPRESSION TAG	UNP Q5SMG8
A	-1	LEU	-	EXPRESSION TAG	UNP Q5SMG8
A	0	HIS	-	EXPRESSION TAG	UNP Q5SMG8
B	-22	MET	-	EXPRESSION TAG	UNP Q5SMG8
B	-21	GLY	-	EXPRESSION TAG	UNP Q5SMG8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	SER	-	EXPRESSION TAG	UNP Q5SMG8
B	-19	SER	-	EXPRESSION TAG	UNP Q5SMG8
B	-18	HIS	-	EXPRESSION TAG	UNP Q5SMG8
B	-17	HIS	-	EXPRESSION TAG	UNP Q5SMG8
B	-16	HIS	-	EXPRESSION TAG	UNP Q5SMG8
B	-15	HIS	-	EXPRESSION TAG	UNP Q5SMG8
B	-14	HIS	-	EXPRESSION TAG	UNP Q5SMG8
B	-13	HIS	-	EXPRESSION TAG	UNP Q5SMG8
B	-12	SER	-	EXPRESSION TAG	UNP Q5SMG8
B	-11	SER	-	EXPRESSION TAG	UNP Q5SMG8
B	-10	GLY	-	EXPRESSION TAG	UNP Q5SMG8
B	-9	LEU	-	EXPRESSION TAG	UNP Q5SMG8
B	-8	GLY	-	EXPRESSION TAG	UNP Q5SMG8
B	-7	VAL	-	EXPRESSION TAG	UNP Q5SMG8
B	-6	LEU	-	EXPRESSION TAG	UNP Q5SMG8
B	-5	PRO	-	EXPRESSION TAG	UNP Q5SMG8
B	-4	GLY	-	EXPRESSION TAG	UNP Q5SMG8
B	-3	GLY	-	EXPRESSION TAG	UNP Q5SMG8
B	-2	PRO	-	EXPRESSION TAG	UNP Q5SMG8
B	-1	LEU	-	EXPRESSION TAG	UNP Q5SMG8
B	0	HIS	-	EXPRESSION TAG	UNP Q5SMG8

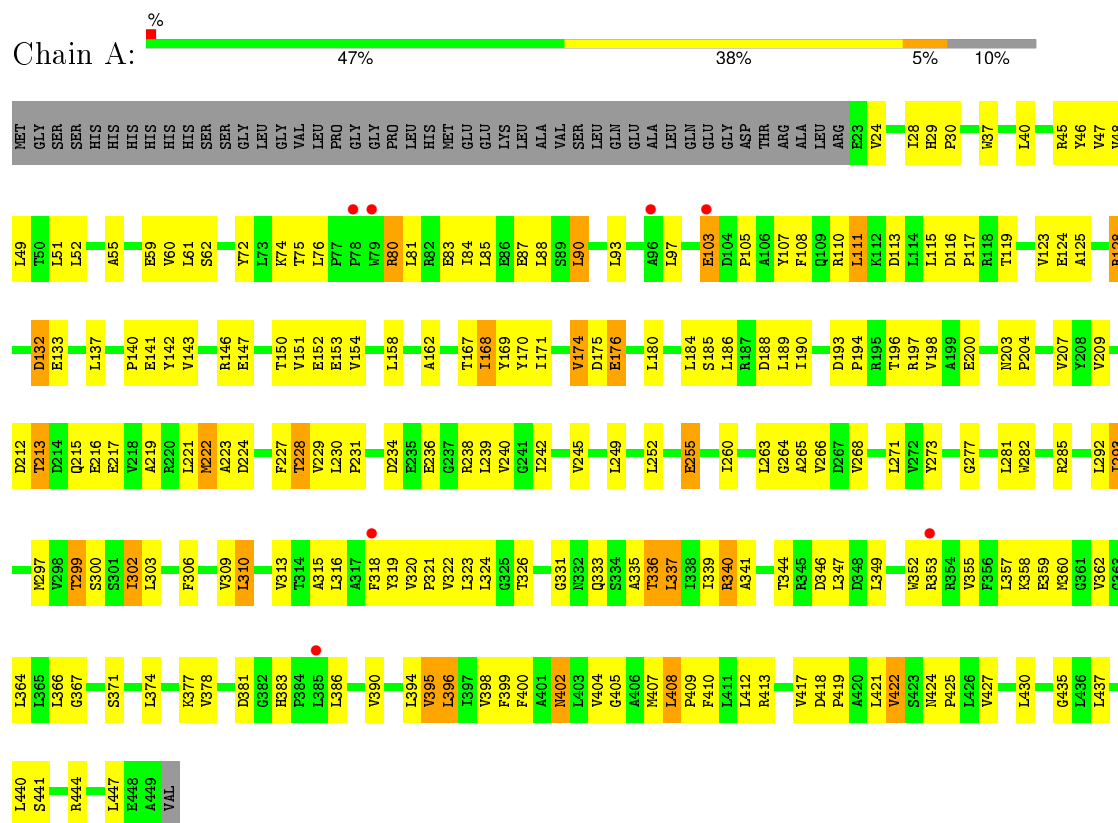
- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	6	Total Mg 6 6	0	0
2	A	7	Total Mg 7 7	0	0

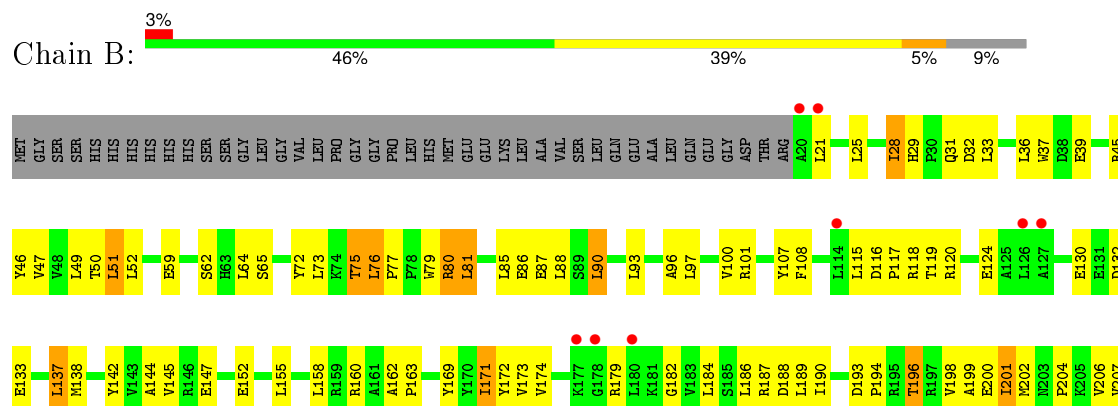
3 Residue-property plots [i](#)

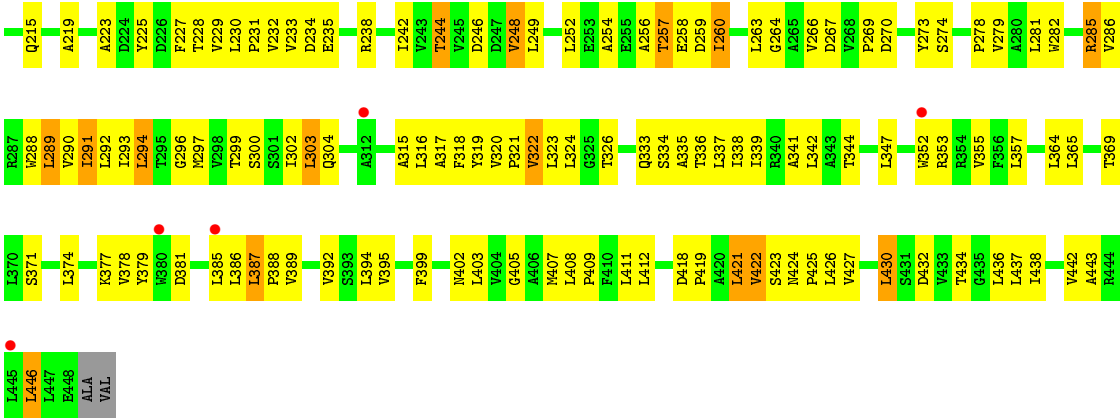
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

• Molecule 1: Mg²⁺ transporter MgtE



• Molecule 1: Mg²⁺ transporter MgtE





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	135.36Å 76.53Å 160.04Å 90.00° 102.17° 90.00°	Depositor
Resolution (Å)	26.71 – 2.94 48.11 – 2.86	Depositor EDS
% Data completeness (in resolution range)	95.0 (26.71-2.94) 92.8 (48.11-2.86)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.19 (at 2.86Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.255 , 0.295 0.252 , 0.290	Depositor DCC
R_{free} test set	1620 reflections (4.98%)	DCC
Wilson B-factor (Å ²)	74.8	Xtriage
Anisotropy	0.585	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 61.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	3 of 34584 reflections (0.009%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6468	wwPDB-VP
Average B, all atoms (Å ²)	119.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.85% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: MG

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	0.25	0/3302	0.44	0/4527
1	B	0.24	0/3273	0.46	0/4498
All	All	0.25	0/6575	0.45	0/9025

There are no bond length outliers.

There are no bond angle outliers.

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	3242	0	3262	222	0
1	B	3213	0	3201	198	0
2	A	7	0	0	0	0
2	B	6	0	0	0	0
All	All	6468	0	6463	366	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 28.

All (366) 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:299:THR:HG21	1:B:321:PRO:HB2	1.46	0.97
1:A:333:GLN:OE1	1:B:285:ARG:HB3	1.71	0.90
1:A:132:ASP:O	1:A:215:GLN:HB2	1.73	0.89
1:A:315:ALA:HA	1:A:318:PHE:CD1	2.11	0.85
1:A:169:TYR:HB3	1:B:169:TYR:HE2	1.41	0.84
1:A:90:LEU:HB2	1:A:141:GLU:OE1	1.77	0.83
1:B:37:TRP:CE2	1:B:45:ARG:HD2	2.14	0.82
1:B:120:ARG:O	1:B:124:GLU:HG2	1.79	0.82
1:B:201:ILE:H	1:B:201:ILE:HD13	1.46	0.81
1:A:336:THR:HG23	1:A:340:ARG:NH1	1.96	0.81
1:B:179:ARG:HA	1:B:238:ARG:HA	1.62	0.80
1:A:103:GLU:O	1:A:105:PRO:HD3	1.82	0.79
1:B:371:SER:HA	1:B:394:LEU:HD23	1.64	0.79
1:A:282:TRP:HZ2	1:B:369:THR:HG21	1.48	0.79
1:A:167:THR:HG21	1:A:229:VAL:HG21	1.66	0.78
1:A:340:ARG:CZ	1:B:260:ILE:HG22	2.13	0.78
1:A:255:GLU:OE1	1:A:255:GLU:HA	1.83	0.78
1:B:155:LEU:HD21	1:B:189:LEU:HB3	1.66	0.77
1:B:315:ALA:HA	1:B:318:PHE:CD1	2.20	0.77
1:B:244:THR:HB	1:B:246:ASP:OD1	1.83	0.77
1:A:268:VAL:HB	1:A:271:LEU:HD12	1.67	0.77
1:A:230:LEU:HD12	1:A:231:PRO:HD2	1.67	0.76
1:B:315:ALA:HA	1:B:318:PHE:HD1	1.50	0.75
1:A:336:THR:HG23	1:A:340:ARG:HH12	1.51	0.75
1:A:316:LEU:HD23	1:A:386:LEU:HD12	1.69	0.75
1:A:383:HIS:HE1	1:B:304:GLN:HE22	1.35	0.74
1:A:282:TRP:CZ2	1:B:369:THR:HG21	2.25	0.72
1:A:383:HIS:CE1	1:B:304:GLN:HE22	2.08	0.71
1:B:256:ALA:O	1:B:260:ILE:HG12	1.91	0.71
1:B:219:ALA:HB2	1:B:248:VAL:HG22	1.71	0.71
1:B:193:ASP:O	1:B:196:THR:HG23	1.90	0.71
1:A:197:ARG:O	1:A:200:GLU:HG2	1.91	0.71
1:A:158:LEU:HD22	1:A:168:ILE:HB	1.72	0.70
1:A:340:ARG:NH2	1:B:260:ILE:HG22	2.07	0.70
1:A:137:LEU:HD11	1:A:215:GLN:HG3	1.71	0.69
1:A:299:THR:HG22	1:B:322:VAL:HG22	1.74	0.69
1:B:193:ASP:OD2	1:B:194:PRO:HD2	1.93	0.69
1:A:318:PHE:HA	1:B:303:LEU:HD12	1.75	0.69
1:B:387:LEU:N	1:B:388:PRO:HD2	2.07	0.68
1:B:101:ARG:HG3	1:B:108:PHE:CD1	2.28	0.68
1:A:263:LEU:HD21	1:A:340:ARG:HH21	1.58	0.68
1:A:169:TYR:HB3	1:B:169:TYR:CE2	2.26	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:LEU:HD23	1:A:123:VAL:HG21	1.74	0.68
1:A:46:TYR:HE2	1:A:75:THR:HB	1.59	0.68
1:A:336:THR:HG23	1:A:340:ARG:CZ	2.25	0.67
1:A:383:HIS:HB3	1:A:386:LEU:CD2	2.24	0.67
1:A:340:ARG:HD3	1:A:340:ARG:N	2.10	0.67
1:B:158:LEU:HD21	1:B:171:ILE:HD12	1.77	0.66
1:A:273:TYR:OH	1:B:355:VAL:HG13	1.94	0.66
1:B:145:VAL:HG23	1:B:173:VAL:HG12	1.77	0.66
1:B:45:ARG:O	1:B:49:LEU:HB2	1.96	0.65
1:B:388:PRO:O	1:B:392:VAL:HB	1.96	0.65
1:A:180:LEU:HD11	1:A:231:PRO:HB2	1.78	0.65
1:A:337:LEU:HD23	1:B:266:VAL:HG21	1.77	0.65
1:A:24:VAL:HG12	1:A:28:ILE:HG23	1.79	0.65
1:B:33:LEU:CD2	1:B:52:LEU:HD11	2.27	0.64
1:B:25:LEU:HA	1:B:28:ILE:HG23	1.78	0.64
1:A:216:GLU:HG3	1:A:255:GLU:HG3	1.79	0.64
1:A:383:HIS:CE1	1:B:304:GLN:NE2	2.65	0.64
1:B:50:THR:HA	1:B:80:ARG:HD2	1.80	0.64
1:B:324:LEU:HD12	1:B:432:ASP:HA	1.79	0.64
1:A:336:THR:HG23	1:A:340:ARG:NH2	2.12	0.63
1:A:125:ALA:HA	1:A:128:ARG:NH1	2.14	0.63
1:A:124:GLU:O	1:A:128:ARG:HD3	1.98	0.62
1:B:142:TYR:H	1:B:142:TYR:HD2	1.46	0.62
1:B:132:ASP:O	1:B:215:GLN:HB2	1.99	0.62
1:B:399:PHE:O	1:B:403:LEU:HD23	2.00	0.62
1:A:395:VAL:O	1:A:395:VAL:HG12	2.00	0.61
1:B:377:LYS:O	1:B:381:ASP:HB2	1.99	0.61
1:A:29:HIS:ND1	1:A:30:PRO:HD2	2.15	0.61
1:A:383:HIS:HE1	1:B:304:GLN:NE2	1.97	0.61
1:A:84:ILE:O	1:A:88:LEU:HB2	2.01	0.61
1:B:334:SER:HB3	1:B:402:ASN:OD1	2.01	0.61
1:B:418:ASP:O	1:B:421:LEU:HB2	2.01	0.60
1:A:352:TRP:HB3	1:A:410:PHE:CD2	2.36	0.60
1:A:418:ASP:HB3	1:A:421:LEU:CD2	2.30	0.60
1:A:263:LEU:CD2	1:A:340:ARG:HH21	2.15	0.60
1:A:170:TYR:HD2	1:A:229:VAL:HB	1.66	0.60
1:A:303:LEU:HD23	1:A:440:LEU:HD11	1.83	0.60
1:A:146:ARG:HG3	1:A:147:GLU:H	1.66	0.60
1:B:72:TYR:HB3	1:B:76:LEU:HD23	1.82	0.60
1:A:360:MET:HG3	1:A:399:PHE:CZ	2.37	0.60
1:A:74:LYS:HA	1:A:107:TYR:HE1	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:378:VAL:HG21	1:A:390:VAL:HG11	1.84	0.60
1:A:152:GLU:HB2	1:A:194:PRO:HB2	1.84	0.60
1:A:293:ILE:HD11	1:B:326:THR:OG1	2.01	0.59
1:A:319:TYR:OH	1:A:378:VAL:HA	2.03	0.59
1:A:140:PRO:HD2	1:A:141:GLU:OE1	2.02	0.59
1:B:299:THR:O	1:B:302:ILE:HG22	2.02	0.59
1:B:302:ILE:HD11	1:B:437:LEU:HD21	1.85	0.59
1:A:107:TYR:CZ	1:A:111:LEU:HD23	2.38	0.59
1:B:282:TRP:CE2	1:B:286:VAL:HG21	2.37	0.59
1:A:133:GLU:OE1	1:A:212:ASP:HA	2.02	0.59
1:B:47:VAL:HA	1:B:50:THR:HG22	1.84	0.59
1:A:190:ILE:CD1	1:B:186:LEU:HD11	2.33	0.59
1:B:379:TYR:HB2	1:B:387:LEU:HD21	1.85	0.58
1:A:61:LEU:HD23	1:A:84:ILE:HG21	1.85	0.58
1:A:90:LEU:HD23	1:A:123:VAL:HG22	1.85	0.58
1:B:101:ARG:HG3	1:B:108:PHE:CG	2.38	0.58
1:B:323:LEU:HD21	1:B:374:LEU:HD21	1.84	0.58
1:A:186:LEU:HD21	1:B:190:ILE:HD12	1.84	0.58
1:A:93:LEU:HD23	1:A:123:VAL:CG2	2.33	0.58
1:A:142:TYR:HB3	1:A:242:ILE:HG12	1.84	0.58
1:A:358:LYS:O	1:A:362:VAL:HG23	2.02	0.58
1:B:282:TRP:CZ2	1:B:286:VAL:HG21	2.38	0.58
1:B:423:SER:O	1:B:427:VAL:HG23	2.04	0.58
1:A:299:THR:CG2	1:B:321:PRO:HB2	2.28	0.58
1:B:290:VAL:O	1:B:294:LEU:HD22	2.03	0.58
1:A:323:LEU:HD12	1:A:394:LEU:HD13	1.85	0.57
1:A:409:PRO:HD3	1:A:422:VAL:HG11	1.86	0.57
1:A:340:ARG:CD	1:A:340:ARG:N	2.67	0.57
1:A:74:LYS:HA	1:A:107:TYR:CE1	2.39	0.57
1:A:315:ALA:HA	1:A:318:PHE:HD1	1.67	0.57
1:A:333:GLN:HG2	1:B:264:GLY:O	2.04	0.57
1:A:418:ASP:HB3	1:A:421:LEU:HD23	1.86	0.57
1:A:190:ILE:HD11	1:B:186:LEU:HD11	1.85	0.56
1:B:344:THR:O	1:B:344:THR:HG22	2.04	0.56
1:B:199:ALA:HA	1:B:202:MET:HG2	1.85	0.56
1:A:320:VAL:N	1:A:321:PRO:HD2	2.18	0.56
1:A:377:LYS:O	1:A:381:ASP:HB2	2.05	0.56
1:A:52:LEU:O	1:A:80:ARG:NH2	2.38	0.56
1:B:37:TRP:CZ2	1:B:45:ARG:HB3	2.40	0.56
1:A:306:PHE:HB2	1:A:440:LEU:HD13	1.88	0.56
1:B:145:VAL:HG23	1:B:173:VAL:CG1	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:422:VAL:HG12	1:B:422:VAL:O	2.06	0.56
1:A:143:VAL:HG11	1:A:158:LEU:HD23	1.88	0.56
1:A:336:THR:HG23	1:A:340:ARG:HH22	1.71	0.55
1:A:234:ASP:HB3	1:A:240:VAL:HG11	1.89	0.55
1:B:137:LEU:H	1:B:137:LEU:HD23	1.72	0.55
1:B:389:VAL:HG12	1:B:446:LEU:HG	1.89	0.55
1:B:408:LEU:N	1:B:409:PRO:HD2	2.22	0.55
1:A:169:TYR:CG	1:B:169:TYR:CD2	2.95	0.54
1:B:316:LEU:HD21	1:B:443:ALA:HB1	1.88	0.54
1:A:169:TYR:CG	1:B:169:TYR:HD2	2.25	0.54
1:A:154:VAL:HG21	1:A:198:VAL:HG11	1.89	0.54
1:A:150:THR:OG1	1:A:153:GLU:HG3	2.08	0.54
1:B:147:GLU:HB2	1:B:199:ALA:HB2	1.89	0.54
1:B:142:TYR:HB3	1:B:242:ILE:HG12	1.90	0.53
1:A:97:LEU:HD21	1:A:111:LEU:HD12	1.90	0.53
1:A:55:ALA:HA	1:A:87:GLU:HG2	1.89	0.53
1:A:326:THR:OG1	1:B:293:ILE:HG13	2.07	0.53
1:B:233:VAL:HG23	1:B:234:ASP:O	2.09	0.53
1:A:169:TYR:CD2	1:B:169:TYR:CD2	2.96	0.53
1:A:335:ALA:HA	1:A:405:GLY:O	2.08	0.53
1:B:33:LEU:HD23	1:B:52:LEU:HD11	1.91	0.53
1:A:150:THR:O	1:A:154:VAL:HG23	2.09	0.53
1:A:337:LEU:HD12	1:B:273:TYR:CD1	2.44	0.53
1:A:219:ALA:HB1	1:A:252:LEU:CD1	2.39	0.53
1:B:424:ASN:HB2	1:B:425:PRO:HD3	1.91	0.53
1:B:162:ALA:N	1:B:163:PRO:HD2	2.24	0.53
1:A:170:TYR:CD2	1:A:229:VAL:HB	2.44	0.52
1:B:81:LEU:O	1:B:85:LEU:HB2	2.08	0.52
1:B:172:TYR:CE2	1:B:231:PRO:HG3	2.45	0.52
1:A:371:SER:OG	1:A:395:VAL:HG23	2.09	0.52
1:B:32:ASP:O	1:B:36:LEU:HG	2.09	0.52
1:A:193:ASP:HB3	1:A:196:THR:HG23	1.91	0.52
1:A:209:VAL:HB	1:A:213:THR:HG21	1.90	0.52
1:B:73:LEU:HD11	1:B:96:ALA:HB1	1.92	0.52
1:A:355:VAL:HG13	1:B:273:TYR:OH	2.10	0.52
1:A:105:PRO:O	1:A:108:PHE:HB3	2.09	0.52
1:B:207:VAL:HG22	1:B:227:PHE:CE2	2.45	0.52
1:B:130:GLU:HB2	1:B:133:GLU:HG3	1.92	0.52
1:A:47:VAL:HG12	1:A:51:LEU:HG	1.91	0.52
1:A:151:VAL:HB	1:A:194:PRO:HA	1.92	0.51
1:B:37:TRP:CH2	1:B:45:ARG:HB3	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:394:LEU:C	1:A:396:LEU:H	2.14	0.51
1:B:204:PRO:C	1:B:206:VAL:H	2.14	0.51
1:B:173:VAL:HG21	1:B:202:MET:SD	2.51	0.51
1:A:303:LEU:HD22	1:A:310:LEU:HD21	1.91	0.51
1:B:225:TYR:HB3	1:B:227:PHE:CE1	2.46	0.51
1:A:174:VAL:HG23	1:A:175:ASP:O	2.10	0.51
1:A:344:THR:O	1:A:344:THR:HG22	2.11	0.51
1:B:152:GLU:HG3	1:B:194:PRO:HB2	1.93	0.51
1:A:184:LEU:HD21	1:A:189:LEU:HD21	1.92	0.51
1:B:225:TYR:HB3	1:B:227:PHE:CD1	2.46	0.50
1:B:182:GLY:HA3	1:B:202:MET:HE1	1.92	0.50
1:A:119:THR:O	1:A:123:VAL:HG23	2.12	0.50
1:A:371:SER:OG	1:A:394:LEU:HB3	2.11	0.50
1:A:306:PHE:CD2	1:A:440:LEU:HB3	2.46	0.50
1:A:116:ASP:HB2	1:A:117:PRO:HD2	1.92	0.50
1:A:260:ILE:HG22	1:B:336:THR:HG23	1.94	0.50
1:B:77:PRO:O	1:B:81:LEU:HB2	2.12	0.50
1:B:352:TRP:CE2	1:B:353:ARG:HG2	2.47	0.49
1:A:383:HIS:CE1	1:B:304:GLN:OE1	2.65	0.49
1:A:371:SER:HA	1:A:394:LEU:HD23	1.94	0.49
1:A:80:ARG:O	1:A:84:ILE:HG12	2.12	0.49
1:A:333:GLN:HB3	1:B:285:ARG:HH11	1.77	0.49
1:B:182:GLY:HA3	1:B:202:MET:CE	2.41	0.49
1:B:353:ARG:O	1:B:357:LEU:HG	2.12	0.49
1:A:28:ILE:HD12	1:A:28:ILE:O	2.13	0.49
1:A:366:LEU:HD22	1:B:289:LEU:HD23	1.95	0.49
1:B:293:ILE:CG2	1:B:294:LEU:N	2.75	0.49
1:B:352:TRP:CZ2	1:B:353:ARG:CZ	2.96	0.49
1:B:386:LEU:O	1:B:389:VAL:HG22	2.12	0.48
1:B:207:VAL:HG22	1:B:227:PHE:HE2	1.79	0.48
1:A:40:LEU:HB3	1:A:45:ARG:HD2	1.95	0.48
1:A:222:MET:HA	1:A:222:MET:HE2	1.95	0.48
1:A:437:LEU:O	1:A:441:SER:HB3	2.14	0.48
1:B:86:GLU:C	1:B:88:LEU:H	2.17	0.48
1:A:264:GLY:O	1:A:265:ALA:HB3	2.14	0.48
1:A:222:MET:HE2	1:A:227:PHE:HB2	1.96	0.48
1:A:223:ALA:HA	1:B:249:LEU:HD23	1.95	0.48
1:B:317:ALA:HA	1:B:320:VAL:HG23	1.96	0.48
1:A:383:HIS:HB3	1:A:386:LEU:HD21	1.96	0.47
1:B:426:LEU:HD23	1:B:426:LEU:C	2.34	0.47
1:B:72:TYR:O	1:B:76:LEU:HB2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:238:ARG:O	1:A:240:VAL:HG13	2.15	0.47
1:B:25:LEU:HA	1:B:28:ILE:CG2	2.44	0.47
1:B:324:LEU:HD11	1:B:436:LEU:HG	1.96	0.47
1:A:302:ILE:HA	1:A:302:ILE:HD13	1.61	0.47
1:B:80:ARG:O	1:B:80:ARG:HG3	2.14	0.47
1:B:73:LEU:O	1:B:107:TYR:HE1	1.97	0.47
1:B:352:TRP:HE3	1:B:407:MET:HE1	1.79	0.47
1:A:340:ARG:NE	1:B:260:ILE:HG22	2.30	0.47
1:B:412:LEU:HD23	1:B:422:VAL:HG23	1.97	0.47
1:B:438:ILE:O	1:B:442:VAL:HG23	2.14	0.47
1:A:277:GLY:O	1:A:281:LEU:HD13	2.14	0.47
1:A:359:GLU:O	1:A:402:ASN:ND2	2.48	0.47
1:B:323:LEU:CD2	1:B:374:LEU:HD21	2.45	0.46
1:A:299:THR:CG2	1:B:322:VAL:HG22	2.45	0.46
1:A:331:GLY:C	1:A:427:VAL:HG11	2.36	0.46
1:A:137:LEU:HD11	1:A:215:GLN:CG	2.43	0.46
1:A:24:VAL:CG1	1:A:28:ILE:HG23	2.44	0.46
1:A:352:TRP:CG	1:A:353:ARG:N	2.84	0.46
1:B:188:ASP:C	1:B:190:ILE:H	2.19	0.46
1:A:320:VAL:O	1:A:324:LEU:HB2	2.15	0.46
1:A:324:LEU:HD12	1:A:435:GLY:HA3	1.97	0.46
1:A:60:VAL:C	1:A:62:SER:H	2.18	0.46
1:A:162:ALA:HB1	1:B:187:ARG:HH12	1.80	0.46
1:A:352:TRP:HE3	1:A:407:MET:HE1	1.81	0.46
1:A:374:LEU:HA	1:B:297:MET:HE1	1.98	0.46
1:B:281:LEU:O	1:B:285:ARG:HG3	2.16	0.46
1:A:171:ILE:HD12	1:A:171:ILE:N	2.32	0.45
1:B:319:TYR:CE2	1:B:378:VAL:HG22	2.51	0.45
1:A:266:VAL:HG21	1:B:337:LEU:HD21	1.98	0.45
1:A:341:ALA:CB	1:A:347:LEU:HD12	2.46	0.45
1:A:344:THR:O	1:A:344:THR:CG2	2.64	0.45
1:A:55:ALA:CA	1:A:87:GLU:HG2	2.47	0.45
1:A:340:ARG:CZ	1:B:260:ILE:CG2	2.91	0.45
1:A:400:PHE:O	1:A:404:VAL:HG23	2.16	0.45
1:B:138:MET:SD	1:B:232:VAL:HG11	2.57	0.45
1:A:158:LEU:HD21	1:A:171:ILE:CG1	2.47	0.45
1:B:434:THR:O	1:B:438:ILE:HG12	2.17	0.45
1:B:374:LEU:O	1:B:378:VAL:HG23	2.17	0.45
1:A:249:LEU:HD12	1:B:223:ALA:HA	1.98	0.45
1:B:46:TYR:HE1	1:B:75:THR:HB	1.81	0.45
1:A:349:LEU:HD22	1:A:413:ARG:HG2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:257:THR:HA	1:B:260:ILE:HG13	1.99	0.45
1:A:418:ASP:HA	1:A:419:PRO:HD2	1.80	0.45
1:B:198:VAL:C	1:B:200:GLU:H	2.20	0.45
1:A:340:ARG:NE	1:B:260:ILE:CG2	2.81	0.44
1:A:93:LEU:CD2	1:A:123:VAL:HG21	2.45	0.44
1:A:48:VAL:O	1:A:48:VAL:HG12	2.16	0.44
1:A:346:ASP:OD1	1:B:274:SER:HB2	2.16	0.44
1:A:158:LEU:O	1:A:162:ALA:N	2.48	0.44
1:A:285:ARG:HB3	1:B:333:GLN:OE1	2.16	0.44
1:A:336:THR:HG22	1:A:337:LEU:N	2.32	0.44
1:A:421:LEU:HA	1:A:421:LEU:HD13	1.82	0.44
1:A:228:THR:HG22	1:A:245:VAL:HG11	1.99	0.44
1:B:341:ALA:O	1:B:347:LEU:N	2.50	0.44
1:A:333:GLN:O	1:A:337:LEU:HB2	2.17	0.44
1:B:169:TYR:N	1:B:169:TYR:CD1	2.86	0.44
1:A:128:ARG:HB2	1:A:128:ARG:HH11	1.82	0.44
1:B:59:GLU:O	1:B:62:SER:HB3	2.17	0.44
1:A:303:LEU:HB3	1:B:318:PHE:HD2	1.83	0.44
1:B:352:TRP:HZ2	1:B:353:ARG:CZ	2.30	0.44
1:B:288:TRP:CZ3	1:B:426:LEU:HD12	2.52	0.44
1:B:184:LEU:CD2	1:B:201:ILE:HG12	2.48	0.44
1:A:61:LEU:CD2	1:A:84:ILE:HG21	2.48	0.44
1:A:140:PRO:O	1:A:142:TYR:CD2	2.70	0.44
1:A:340:ARG:HD3	1:A:340:ARG:H	1.80	0.43
1:A:347:LEU:HD11	1:B:273:TYR:CD2	2.53	0.43
1:A:171:ILE:HB	1:A:184:LEU:HB3	2.00	0.43
1:A:408:LEU:N	1:A:409:PRO:HD2	2.33	0.43
1:A:249:LEU:HD21	1:B:249:LEU:HD13	1.99	0.43
1:A:224:ASP:O	1:B:31:GLN:NE2	2.51	0.43
1:A:37:TRP:CZ3	1:A:45:ARG:HG2	2.53	0.43
1:A:80:ARG:NH1	1:A:83:GLU:HG2	2.34	0.43
1:B:100:VAL:HG12	1:B:108:PHE:HB2	2.00	0.43
1:B:21:LEU:O	1:B:25:LEU:HD12	2.18	0.43
1:A:422:VAL:HG12	1:A:422:VAL:O	2.18	0.43
1:A:234:ASP:C	1:A:236:GLU:H	2.22	0.43
1:A:72:TYR:HB3	1:A:76:LEU:HD23	2.00	0.43
1:B:21:LEU:O	1:B:25:LEU:HB2	2.19	0.43
1:A:360:MET:HG3	1:A:399:PHE:HZ	1.80	0.43
1:A:297:MET:C	1:A:299:THR:N	2.71	0.43
1:A:46:TYR:CE2	1:A:75:THR:HB	2.46	0.43
1:B:79:TRP:CE3	1:B:80:ARG:N	2.87	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:353:ARG:HG2	1:A:353:ARG:H	1.59	0.43
1:B:207:VAL:HG11	1:B:225:TYR:CE2	2.54	0.43
1:B:338:ILE:O	1:B:342:LEU:HG	2.18	0.43
1:A:374:LEU:HA	1:B:297:MET:CE	2.48	0.43
1:B:116:ASP:O	1:B:118:ARG:N	2.52	0.43
1:A:97:LEU:CD2	1:A:111:LEU:HD12	2.48	0.43
1:B:387:LEU:N	1:B:388:PRO:CD	2.78	0.43
1:B:47:VAL:O	1:B:51:LEU:HB2	2.19	0.43
1:A:249:LEU:CD2	1:B:249:LEU:HD13	2.48	0.43
1:A:424:ASN:HB2	1:A:425:PRO:HD3	2.00	0.43
1:B:137:LEU:N	1:B:137:LEU:HD23	2.34	0.43
1:B:426:LEU:O	1:B:430:LEU:HB2	2.19	0.43
1:B:335:ALA:HA	1:B:405:GLY:O	2.19	0.43
1:A:142:TYR:CE1	1:A:239:LEU:HG	2.54	0.43
1:A:322:VAL:HG22	1:B:296:GLY:O	2.19	0.43
1:B:201:ILE:HD13	1:B:201:ILE:N	2.25	0.42
1:B:144:ALA:HA	1:B:172:TYR:O	2.19	0.42
1:A:132:ASP:O	1:A:215:GLN:N	2.46	0.42
1:A:85:LEU:HB3	1:A:115:LEU:HD21	2.00	0.42
1:A:309:VAL:HG21	1:A:444:ARG:HB2	2.00	0.42
1:B:145:VAL:O	1:B:173:VAL:HA	2.19	0.42
1:B:418:ASP:HA	1:B:419:PRO:HD2	1.84	0.42
1:A:40:LEU:HG	1:A:45:ARG:HG3	2.01	0.42
1:B:442:VAL:O	1:B:442:VAL:HG12	2.19	0.42
1:A:335:ALA:O	1:A:339:ILE:HG13	2.20	0.42
1:A:110:ARG:O	1:A:113:ASP:HB2	2.18	0.42
1:B:90:LEU:HA	1:B:90:LEU:HD23	1.69	0.42
1:A:326:THR:O	1:A:326:THR:HG22	2.20	0.42
1:B:130:GLU:HB2	1:B:133:GLU:CG	2.50	0.42
1:A:302:ILE:CD1	1:A:437:LEU:HD21	2.49	0.42
1:A:176:GLU:HG3	1:A:176:GLU:H	1.59	0.42
1:A:367:GLY:N	1:A:398:VAL:HG11	2.35	0.42
1:A:417:VAL:O	1:A:419:PRO:HD3	2.20	0.42
1:A:260:ILE:HG12	1:B:260:ILE:HD13	2.02	0.42
1:B:230:LEU:HA	1:B:231:PRO:HD3	1.80	0.42
1:B:442:VAL:O	1:B:446:LEU:HB2	2.20	0.42
1:A:203:ASN:HA	1:A:204:PRO:HD2	1.66	0.42
1:A:437:LEU:O	1:A:441:SER:CB	2.68	0.41
1:A:268:VAL:CB	1:A:271:LEU:HD12	2.45	0.41
1:B:28:ILE:HD12	1:B:29:HIS:O	2.20	0.41
1:A:207:VAL:HB	1:A:227:PHE:HE2	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:333:GLN:CB	1:B:285:ARG:HH11	2.32	0.41
1:B:231:PRO:HA	1:B:242:ILE:HG22	2.01	0.41
1:A:29:HIS:CG	1:A:30:PRO:HD2	2.56	0.41
1:A:309:VAL:HG12	1:A:316:LEU:HD12	2.02	0.41
1:B:302:ILE:HA	1:B:302:ILE:HD12	1.92	0.41
1:A:49:LEU:HD12	1:A:76:LEU:HD21	2.01	0.41
1:B:152:GLU:HG3	1:B:194:PRO:O	2.21	0.41
1:B:187:ARG:NH1	1:B:187:ARG:HG3	2.34	0.41
1:A:352:TRP:CE3	1:A:407:MET:HE1	2.56	0.41
1:A:59:GLU:O	1:A:62:SER:HB3	2.20	0.41
1:B:335:ALA:O	1:B:339:ILE:HG13	2.20	0.41
1:B:47:VAL:O	1:B:50:THR:HG22	2.21	0.41
1:B:51:LEU:HD12	1:B:51:LEU:HA	1.86	0.41
1:B:229:VAL:HG12	1:B:230:LEU:N	2.35	0.41
1:B:263:LEU:HA	1:B:263:LEU:HD23	1.85	0.41
1:A:97:LEU:HD12	1:A:123:VAL:HG11	2.02	0.41
1:B:152:GLU:CG	1:B:194:PRO:HB2	2.50	0.41
1:A:395:VAL:O	1:A:396:LEU:HD23	2.21	0.41
1:A:404:VAL:O	1:A:408:LEU:HD23	2.20	0.41
1:B:385:LEU:H	1:B:385:LEU:HD12	1.86	0.41
1:A:221:LEU:HD23	1:A:221:LEU:HA	1.86	0.41
1:A:273:TYR:HE2	1:B:355:VAL:HG22	1.85	0.40
1:A:378:VAL:CG2	1:A:390:VAL:HG11	2.50	0.40
1:B:85:LEU:O	1:B:93:LEU:HD13	2.21	0.40
1:A:37:TRP:CE3	1:A:45:ARG:HG2	2.56	0.40
1:B:364:LEU:HD13	1:B:364:LEU:C	2.41	0.40
1:B:371:SER:OG	1:B:395:VAL:HG23	2.21	0.40
1:A:383:HIS:CE1	1:B:304:GLN:CD	2.94	0.40
1:A:412:LEU:HB3	1:A:417:VAL:HB	2.03	0.40
1:B:316:LEU:HD21	1:B:443:ALA:CB	2.51	0.40
1:A:185:SER:HB3	1:A:188:ASP:CG	2.41	0.40
1:B:256:ALA:O	1:B:259:ASP:HB2	2.22	0.40
1:A:417:VAL:HG12	1:A:418:ASP:N	2.36	0.40
1:B:76:LEU:HD13	1:B:76:LEU:HA	1.93	0.40
1:B:290:VAL:O	1:B:291:ILE:C	2.59	0.40
1:B:187:ARG:HH11	1:B:187:ARG:HG3	1.87	0.40
1:A:61:LEU:O	1:A:61:LEU:HG	2.21	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	425/473 (90%)	372 (88%)	48 (11%)	5 (1%)	16	47
1	B	427/473 (90%)	367 (86%)	52 (12%)	8 (2%)	10	34
All	All	852/946 (90%)	739 (87%)	100 (12%)	13 (2%)	13	41

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	269	PRO
1	A	103	GLU
1	A	395	VAL
1	A	402	ASN
1	A	213	THR
1	B	64	LEU
1	B	87	GLU
1	B	254	ALA
1	A	422	VAL
1	B	422	VAL
1	B	278	PRO
1	B	291	ILE
1	B	117	PRO

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	339/400 (85%)	310 (91%)	29 (9%)	13	35

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	333/400 (83%)	291 (87%)	42 (13%)	5	16
All	All	672/800 (84%)	601 (89%)	71 (11%)	8	24

All (71) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	ARG
1	A	81	LEU
1	A	90	LEU
1	A	111	LEU
1	A	128	ARG
1	A	132	ASP
1	A	168	ILE
1	A	174	VAL
1	A	176	GLU
1	A	217	GLU
1	A	222	MET
1	A	228	THR
1	A	255	GLU
1	A	292	LEU
1	A	293	ILE
1	A	299	THR
1	A	300	SER
1	A	302	ILE
1	A	310	LEU
1	A	313	VAL
1	A	336	THR
1	A	337	LEU
1	A	340	ARG
1	A	357	LEU
1	A	364	LEU
1	A	396	LEU
1	A	408	LEU
1	A	430	LEU
1	A	447	LEU
1	B	28	ILE
1	B	39	GLU
1	B	51	LEU
1	B	65	SER
1	B	75	THR
1	B	76	LEU
1	B	80	ARG

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Mol	Chain	Res	Type
1	B	81	LEU
1	B	90	LEU
1	B	97	LEU
1	B	115	LEU
1	B	119	THR
1	B	137	LEU
1	B	160	ARG
1	B	171	ILE
1	B	174	VAL
1	B	196	THR
1	B	201	ILE
1	B	228	THR
1	B	235	GLU
1	B	244	THR
1	B	248	VAL
1	B	252	LEU
1	B	257	THR
1	B	258	GLU
1	B	260	ILE
1	B	267	ASP
1	B	270	ASP
1	B	279	VAL
1	B	285	ARG
1	B	289	LEU
1	B	292	LEU
1	B	294	LEU
1	B	300	SER
1	B	303	LEU
1	B	322	VAL
1	B	365	LEU
1	B	387	LEU
1	B	411	LEU
1	B	421	LEU
1	B	430	LEU
1	B	446	LEU

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

Mol	Chain	Res	Type
1	A	31	GLN
1	A	109	GLN
1	A	215	GLN

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Mol	Chain	Res	Type
1	A	383	HIS
1	A	424	ASN
1	B	63	HIS
1	B	215	GLN
1	B	329	ASN
1	B	424	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 13 ligands modelled in this entry, 13 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 ⓘ

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	427/473 (90%)	-0.18	7 (1%) 74 74	68, 106, 150, 184	0
1	B	429/473 (90%)	-0.06	13 (3%) 54 50	80, 123, 175, 235	0
All	All	856/946 (90%)	-0.12	20 (2%) 64 61	68, 115, 164, 235	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	20	ALA	7.8
1	B	21	LEU	5.4
1	A	78	PRO	4.9
1	B	177	LYS	4.0
1	A	385	LEU	3.8
1	B	352	TRP	3.3
1	B	126	LEU	3.0
1	A	79	TRP	2.9
1	B	127	ALA	2.7
1	A	353	ARG	2.7
1	B	114	LEU	2.5
1	B	178	GLY	2.4
1	B	180	LEU	2.3
1	B	445	LEU	2.3
1	A	103	GLU	2.2
1	B	385	LEU	2.2
1	B	380	TRP	2.2
1	B	312	ALA	2.2
1	A	96	ALA	2.1
1	A	318	PHE	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	MG	A	455	1/1	0.94	0.46	8.24	62,62,62,62	0
2	MG	B	454	1/1	0.87	0.48	7.67	81,81,81,81	0
2	MG	A	451	1/1	0.99	0.62	7.15	72,72,72,72	0
2	MG	B	451	1/1	0.82	0.32	5.06	90,90,90,90	0
2	MG	A	454	1/1	0.98	0.36	3.51	60,60,60,60	0
2	MG	A	456	1/1	0.67	0.32	2.96	79,79,79,79	0
2	MG	B	452	1/1	0.97	0.47	-	78,78,78,78	0
2	MG	B	453	1/1	0.84	0.50	-	74,74,74,74	0
2	MG	B	456	1/1	0.93	0.82	-	98,98,98,98	0
2	MG	A	452	1/1	0.99	0.32	-	70,70,70,70	0
2	MG	A	453	1/1	0.84	0.30	-	75,75,75,75	0
2	MG	B	455	1/1	0.95	0.24	-	61,61,61,61	0
2	MG	A	457	1/1	0.92	0.41	-	75,75,75,75	0

6.5 Other polymers [i](#)

There are no such residues in this entry.