



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 01:41 AM GMT

PDB ID : 2IMO
Title : Crystal structure of allantoate amidohydrolase from Escherichia coli at pH 4.6
Authors : Agarwal, R.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center
for Structural Genomics (NYSGXRC)
Deposited on : 2006-10-04
Resolution : 2.80 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>

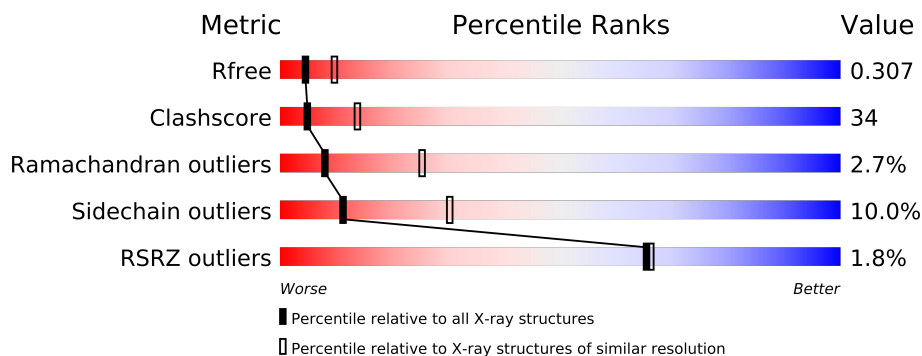
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
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 2.80 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}	66092	1799 (2.80-2.80)
Clashscore	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)
RSRZ outliers	66119	1802 (2.80-2.80)

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.

Mol	Chain	Length	Quality of chain
1	A	423	
1	B	423	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6265 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 Allantoate amidohydrolase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	399	Total	C	N	O	S	Se	0	0	0
			3128	1975	541	592	7	13			
1	B	394	Total	C	N	O	S	Se	0	0	0
			3080	1946	535	579	7	13			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	CLONING ARTIFACT	UNP P77425
A	2	SER	-	CLONING ARTIFACT	UNP P77425
A	3	LEU	-	CLONING ARTIFACT	UNP P77425
A	29	MSE	MET	MODIFIED RESIDUE	UNP P77425
A	49	MSE	MET	MODIFIED RESIDUE	UNP P77425
A	125	MSE	MET	MODIFIED RESIDUE	UNP P77425
A	167	MSE	MET	MODIFIED RESIDUE	UNP P77425
A	234	MSE	MET	MODIFIED RESIDUE	UNP P77425
A	258	MSE	MET	MODIFIED RESIDUE	UNP P77425
A	308	MSE	MET	MODIFIED RESIDUE	UNP P77425
A	315	MSE	MET	MODIFIED RESIDUE	UNP P77425
A	325	MSE	MET	MODIFIED RESIDUE	UNP P77425
A	332	MSE	MET	MODIFIED RESIDUE	UNP P77425
A	354	MSE	MET	MODIFIED RESIDUE	UNP P77425
A	373	MSE	MET	MODIFIED RESIDUE	UNP P77425
A	405	MSE	MET	MODIFIED RESIDUE	UNP P77425
A	414	GLU	-	EXPRESSION TAG	UNP P77425
A	415	GLY	-	EXPRESSION TAG	UNP P77425
A	416	GLY	-	EXPRESSION TAG	UNP P77425
A	417	SER	-	EXPRESSION TAG	UNP P77425
A	418	HIS	-	EXPRESSION TAG	UNP P77425
A	419	HIS	-	EXPRESSION TAG	UNP P77425
A	420	HIS	-	EXPRESSION TAG	UNP P77425
A	421	HIS	-	EXPRESSION TAG	UNP P77425
A	422	HIS	-	EXPRESSION TAG	UNP P77425

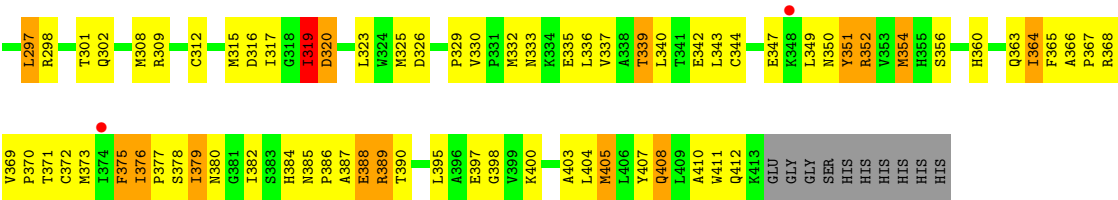
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Chain	Residue	Modelled	Actual	Comment	Reference
A	423	HIS	-	EXPRESSION TAG	UNP P77425
B	1	MSE	-	CLONING ARTIFACT	UNP P77425
B	2	SER	-	CLONING ARTIFACT	UNP P77425
B	3	LEU	-	CLONING ARTIFACT	UNP P77425
B	29	MSE	MET	MODIFIED RESIDUE	UNP P77425
B	49	MSE	MET	MODIFIED RESIDUE	UNP P77425
B	125	MSE	MET	MODIFIED RESIDUE	UNP P77425
B	167	MSE	MET	MODIFIED RESIDUE	UNP P77425
B	234	MSE	MET	MODIFIED RESIDUE	UNP P77425
B	258	MSE	MET	MODIFIED RESIDUE	UNP P77425
B	308	MSE	MET	MODIFIED RESIDUE	UNP P77425
B	315	MSE	MET	MODIFIED RESIDUE	UNP P77425
B	325	MSE	MET	MODIFIED RESIDUE	UNP P77425
B	332	MSE	MET	MODIFIED RESIDUE	UNP P77425
B	354	MSE	MET	MODIFIED RESIDUE	UNP P77425
B	373	MSE	MET	MODIFIED RESIDUE	UNP P77425
B	405	MSE	MET	MODIFIED RESIDUE	UNP P77425
B	414	GLU	-	EXPRESSION TAG	UNP P77425
B	415	GLY	-	EXPRESSION TAG	UNP P77425
B	416	GLY	-	EXPRESSION TAG	UNP P77425
B	417	SER	-	EXPRESSION TAG	UNP P77425
B	418	HIS	-	EXPRESSION TAG	UNP P77425
B	419	HIS	-	EXPRESSION TAG	UNP P77425
B	420	HIS	-	EXPRESSION TAG	UNP P77425
B	421	HIS	-	EXPRESSION TAG	UNP P77425
B	422	HIS	-	EXPRESSION TAG	UNP P77425
B	423	HIS	-	EXPRESSION TAG	UNP P77425

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	37	Total O 37 37	0	0
2	B	20	Total O 20 20	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	94.75Å 183.92Å 48.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.78 – 2.80 48.78 – 2.77	Depositor EDS
% Data completeness (in resolution range)	90.0 (48.78-2.80) 88.5 (48.78-2.77)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 2.77Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.245 , 0.302 0.264 , 0.307	Depositor DCC
R_{free} test set	578 reflections (2.94%)	DCC
Wilson B-factor (Å ²)	50.1	Xtriage
Anisotropy	0.469	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 22.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 21933 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	6265	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.31% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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	0.55	2/3183 (0.1%)	0.74	3/4302 (0.1%)
1	B	0.54	4/3136 (0.1%)	0.71	1/4239 (0.0%)
All	All	0.55	6/6319 (0.1%)	0.72	4/8541 (0.0%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	258	MSE	CG-SE	-5.50	1.76	1.95
1	B	29	MSE	CG-SE	-5.45	1.76	1.95
1	B	354	MSE	CG-SE	-5.08	1.78	1.95
1	B	125	MSE	CG-SE	-5.07	1.78	1.95
1	A	49	MSE	CG-SE	-5.05	1.78	1.95
1	B	308	MSE	CG-SE	-5.05	1.78	1.95

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	117	LEU	CA-CB-CG	6.11	129.36	115.30
1	B	66	GLY	N-CA-C	-5.41	99.57	113.10
1	A	343	LEU	CA-CB-CG	5.36	127.62	115.30
1	A	68	LEU	N-CA-C	-5.30	96.69	111.00

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	3128	0	3064	180	0
1	B	3080	0	3028	247	0
2	A	37	0	0	6	0
2	B	20	0	0	3	0
All	All	6265	0	6092	419	0

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 34.

All (419) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:344:CYS:HA	1:B:349:LEU:HD12	1.42	1.02
1:A:365:PHE:HB3	1:A:371:THR:HG21	1.49	0.94
1:A:379:ILE:HD11	1:A:389:ARG:HB2	1.49	0.94
1:B:209:VAL:HG11	1:B:373:MSE:HE3	1.49	0.93
1:A:208:VAL:HG13	1:A:340:LEU:HD23	1.51	0.92
1:B:240:THR:HG22	1:B:280:PRO:O	1.75	0.87
1:B:228:HIS:O	1:B:232:THR:HG22	1.76	0.86
1:B:74:PRO:HG2	1:B:75:GLN:HE21	1.39	0.86
1:A:240:THR:HG22	1:A:280:PRO:O	1.76	0.85
1:A:149:PRO:HB3	1:A:174:LEU:HD23	1.57	0.84
1:B:340:LEU:HB2	1:B:405:MSE:HE2	1.58	0.84
1:B:228:HIS:HD2	1:B:230:GLY:H	1.23	0.83
1:B:379:ILE:HA	1:B:389:ARG:CZ	2.12	0.80
1:A:4:ILE:HG13	1:A:400:LYS:HD3	1.65	0.79
1:B:379:ILE:HD13	1:B:382:ILE:O	1.85	0.77
1:A:204:GLN:HB2	2:A:439:HOH:O	1.83	0.77
1:B:246:ARG:HH21	1:B:315:MSE:SE	2.17	0.76
1:B:142:ASN:ND2	1:B:147:ALA:HB3	2.01	0.75
1:B:277:ASN:HD22	1:B:277:ASN:H	1.33	0.75
1:A:199:LEU:HA	2:A:439:HOH:O	1.85	0.75
1:B:376:ILE:HG23	1:B:390:THR:HG21	1.67	0.74
1:B:379:ILE:HA	1:B:389:ARG:NH1	2.02	0.73
1:B:68:LEU:HD21	1:B:110:LYS:HE2	1.69	0.73
1:A:68:LEU:HD22	1:A:116:PRO:HD2	1.71	0.73
1:B:63:ASN:HD22	1:B:125:MSE:HA	1.54	0.72
1:B:39:LEU:HD13	1:B:170:CYS:HB3	1.72	0.71
1:B:376:ILE:HD11	1:B:398:GLY:HA3	1.71	0.71
1:B:116:PRO:HG3	1:B:120:VAL:HG23	1.72	0.70
1:B:379:ILE:H	1:B:379:ILE:HD12	1.57	0.70

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:332:MSE:SE	2:A:451:HOH:O	2.60	0.70
1:A:29:MSE:HE3	1:A:92:ASN:HB3	1.74	0.69
1:B:46:LYS:HG3	1:B:56:THR:OG1	1.91	0.69
1:A:14:THR:CG2	1:A:100:LEU:HD11	2.22	0.69
1:A:286:THR:HG23	2:B:427:HOH:O	1.92	0.69
1:B:61:VAL:HG11	1:B:142:ASN:OD1	1.93	0.69
1:A:274:ASN:HD21	1:B:269:VAL:H	1.40	0.69
1:A:272:ARG:HD3	1:A:282:LYS:HE2	1.74	0.68
1:A:369:VAL:O	1:A:371:THR:HG23	1.94	0.68
1:B:31:ARG:NH1	1:B:41:THR:HG21	2.07	0.68
1:A:202:ASN:HB3	1:A:204:GLN:HE21	1.59	0.68
1:A:204:GLN:NE2	1:A:352:ARG:HE	1.91	0.67
1:A:217:ARG:HB2	1:A:325:MSE:HB3	1.76	0.67
1:A:333:ASN:O	1:A:337:VAL:HG23	1.94	0.67
1:A:267:GLY:O	1:B:274:ASN:HA	1.95	0.67
1:B:83:HIS:CG	1:B:128:GLU:HG2	2.30	0.67
1:B:365:PHE:HB3	1:B:371:THR:HG21	1.77	0.67
1:B:360:HIS:O	1:B:363:GLN:HG2	1.95	0.66
1:A:185:ILE:HB	1:A:369:VAL:HG11	1.77	0.66
1:B:369:VAL:O	1:B:371:THR:HG23	1.95	0.66
1:B:34:TYR:OH	1:B:167:MSE:HE3	1.95	0.66
1:A:246:ARG:NH2	1:A:315:MSE:HE2	2.11	0.66
1:B:90:GLY:HA3	1:B:384:HIS:O	1.96	0.66
1:B:379:ILE:HD12	1:B:379:ILE:N	2.11	0.65
1:B:228:HIS:CD2	1:B:230:GLY:H	2.10	0.65
1:A:14:THR:HG22	1:A:100:LEU:HD11	1.79	0.65
1:B:385:ASN:HD21	1:B:387:ALA:HB3	1.62	0.64
1:A:246:ARG:HH21	1:A:315:MSE:HE2	1.61	0.64
1:B:116:PRO:HB3	1:B:411:TRP:CE2	2.33	0.64
1:A:368:ARG:HH11	1:A:368:ARG:HG3	1.63	0.64
1:A:49:MSE:HE1	1:A:122:VAL:HB	1.80	0.64
1:B:167:MSE:SE	1:B:174:LEU:HB3	2.48	0.63
1:B:312:CYS:HB3	1:B:317:ILE:O	1.99	0.63
1:B:277:ASN:H	1:B:277:ASN:ND2	1.95	0.63
1:B:8:ARG:HG2	1:B:8:ARG:HH11	1.64	0.63
1:B:63:ASN:ND2	1:B:126:ALA:H	1.96	0.62
1:A:248:CYS:O	1:A:252:VAL:HG23	2.00	0.61
1:B:49:MSE:O	1:B:54:LEU:HB2	2.00	0.61
1:B:309:ARG:HD3	1:B:319:ILE:HD13	1.82	0.61
1:B:60:GLU:CD	1:B:60:GLU:H	2.04	0.61
1:B:36:PRO:O	1:B:37:GLU:HB2	1.99	0.61
1:A:194:GLU:O	1:A:195:GLN:HB2	2.01	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:39:LEU:HD12	1:A:170:CYS:HA	1.82	0.61
1:A:343:LEU:HB3	1:A:346:ARG:HH22	1.65	0.61
1:B:185:ILE:HB	1:B:369:VAL:HG11	1.83	0.61
1:B:286:THR:HG23	2:B:425:HOH:O	2.02	0.60
1:B:174:LEU:HD12	1:B:174:LEU:O	2.01	0.60
1:A:343:LEU:HB2	1:A:346:ARG:HH12	1.66	0.60
1:B:226:SER:HA	1:B:279:VAL:O	2.00	0.60
1:B:408:GLN:HA	1:B:412:GLN:HG2	1.83	0.60
1:B:46:LYS:NZ	1:B:56:THR:HG23	2.16	0.60
1:B:319:ILE:HD12	1:B:319:ILE:H	1.66	0.60
1:B:268:LYS:H	1:B:286:THR:HG22	1.67	0.59
1:B:205:SER:O	1:B:377:PRO:HD3	2.01	0.59
1:B:109:LEU:O	1:B:113:TYR:HB2	2.02	0.59
1:A:61:VAL:HG21	1:A:142:ASN:CG	2.23	0.59
1:B:141:LYS:HB3	1:B:147:ALA:HB2	1.83	0.59
1:A:168:LYS:HB2	1:A:168:LYS:HZ3	1.67	0.59
1:B:277:ASN:HD22	1:B:277:ASN:N	1.95	0.59
1:B:271:PRO:HG2	1:B:274:ASN:HB2	1.84	0.59
1:B:379:ILE:HG12	1:B:388:GLU:HA	1.85	0.59
1:A:8:ARG:NH1	1:A:8:ARG:HG2	2.18	0.58
1:A:31:ARG:HH12	1:A:96:GLN:HG3	1.68	0.58
1:A:273:PRO:HG3	1:A:278:VAL:HG11	1.86	0.58
1:B:271:PRO:HG2	1:B:274:ASN:HD22	1.69	0.58
1:B:333:ASN:O	1:B:337:VAL:HG23	2.02	0.58
1:A:205:SER:O	1:A:377:PRO:HD3	2.03	0.58
1:B:175:PRO:HG2	1:B:179:LEU:HD11	1.85	0.58
1:A:136:VAL:O	1:A:137:PHE:HB2	2.04	0.58
1:A:204:GLN:OE1	1:A:352:ARG:HB2	2.03	0.58
1:B:33:LEU:HG	1:B:127:GLU:HB2	1.84	0.58
1:A:46:LYS:HG3	1:A:56:THR:HB	1.84	0.58
1:B:213:VAL:HG21	1:B:290:ARG:HB3	1.85	0.58
1:B:214:GLY:HA3	1:B:329:PRO:HG3	1.86	0.58
1:A:174:LEU:HD22	1:A:174:LEU:O	2.04	0.57
1:B:120:VAL:HG21	1:B:411:TRP:CZ3	2.40	0.57
1:A:218:TYR:CZ	1:A:323:LEU:HG	2.40	0.57
1:A:274:ASN:ND2	1:B:268:LYS:HA	2.20	0.57
1:A:147:ALA:O	1:A:148:ASN:HB3	2.05	0.57
1:A:209:VAL:HA	1:A:354:MSE:O	2.03	0.57
1:A:365:PHE:CB	1:A:371:THR:HG21	2.29	0.57
1:A:268:LYS:HB3	1:A:286:THR:HG22	1.87	0.57
1:A:8:ARG:HH11	1:A:8:ARG:HG2	1.69	0.57
1:B:333:ASN:ND2	1:B:370:PRO:HB3	2.20	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:218:TYR:CE2	1:A:323:LEU:HG	2.39	0.57
1:A:196:GLY:O	1:A:381:GLY:HA2	2.04	0.57
1:B:132:ARG:HD3	1:B:133:PHE:CE1	2.40	0.57
1:B:367:PRO:HB2	1:B:368:ARG:NH1	2.19	0.57
1:A:188:PHE:HB3	1:A:371:THR:HG22	1.86	0.56
1:B:93:LEU:HB3	1:B:193:ILE:HG21	1.85	0.56
1:B:105:ALA:O	1:B:109:LEU:HB2	2.05	0.56
1:B:268:LYS:HB3	1:B:286:THR:CG2	2.34	0.56
1:B:8:ARG:HG2	1:B:8:ARG:NH1	2.18	0.56
1:B:239:ASP:HB3	1:B:242:TYR:HB3	1.86	0.56
1:A:264:LEU:H	1:B:234:MSE:HE1	1.71	0.56
1:B:116:PRO:HB3	1:B:411:TRP:CD2	2.40	0.56
1:B:179:LEU:H	1:B:179:LEU:HD12	1.70	0.56
1:B:379:ILE:CG1	1:B:389:ARG:H	2.19	0.56
1:B:71:THR:O	1:B:72:GLU:HB3	2.06	0.56
1:A:168:LYS:HZ1	1:A:173:THR:HG23	1.69	0.56
1:A:128:GLU:O	1:A:129:GLU:HG3	2.06	0.56
1:B:347:GLU:HB2	1:B:349:LEU:HG	1.88	0.56
1:A:33:LEU:HD22	1:A:34:TYR:CD1	2.40	0.56
1:B:228:HIS:CD2	1:B:231:THR:HG23	2.41	0.56
1:B:222:LEU:HD13	1:B:285:PHE:HE2	1.69	0.56
1:B:228:HIS:CD2	1:B:231:THR:H	2.24	0.56
1:B:191:LEU:HD21	1:B:376:ILE:HD13	1.86	0.56
1:A:272:ARG:HD3	1:A:282:LYS:CE	2.35	0.55
1:A:204:GLN:NE2	1:A:352:ARG:HH21	2.04	0.55
1:A:15:LEU:O	1:A:19:SER:HB2	2.06	0.55
1:A:116:PRO:HG3	1:A:120:VAL:HG22	1.87	0.55
1:A:14:THR:O	1:A:17:TRP:HB3	2.07	0.55
1:B:29:MSE:HB2	1:B:90:GLY:O	2.06	0.55
1:A:216:ARG:HA	1:A:325:MSE:O	2.07	0.55
1:B:83:HIS:CD2	1:B:128:GLU:HG2	2.42	0.55
1:A:137:PHE:HB3	1:A:141:LYS:HG3	1.89	0.55
1:B:250:GLN:O	1:B:254:LYS:HG3	2.06	0.55
1:B:52:SER:C	1:B:54:LEU:H	2.09	0.55
1:A:209:VAL:HG13	1:A:373:MSE:HB2	1.89	0.55
1:A:268:LYS:HB3	1:A:286:THR:CG2	2.37	0.55
1:A:345:GLU:OE2	1:A:345:GLU:HA	2.06	0.55
1:B:352:ARG:HG2	1:B:352:ARG:HH11	1.72	0.55
1:A:37:GLU:CD	1:A:37:GLU:H	2.10	0.55
1:B:273:PRO:HG3	1:B:278:VAL:HG11	1.88	0.55
1:A:31:ARG:HD2	1:A:84:ILE:O	2.06	0.54
1:A:226:SER:HA	1:A:279:VAL:O	2.07	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:204:GLN:HA	1:B:350:ASN:ND2	2.22	0.54
1:A:68:LEU:HD22	1:A:116:PRO:CD	2.35	0.54
1:A:67:ARG:HA	1:A:120:VAL:O	2.08	0.54
1:B:63:ASN:HD22	1:B:125:MSE:CA	2.20	0.54
1:A:103:TRP:HA	1:A:122:VAL:HG21	1.89	0.54
1:A:130:GLY:H	1:A:136:VAL:HG12	1.72	0.54
1:B:216:ARG:HG2	1:B:323:LEU:HD21	1.88	0.54
1:B:188:PHE:HB3	1:B:371:THR:HG22	1.89	0.54
1:B:193:ILE:O	1:B:195:GLN:N	2.34	0.54
1:B:368:ARG:HG3	1:B:368:ARG:HH11	1.73	0.53
1:B:379:ILE:HG13	1:B:389:ARG:O	2.08	0.53
1:B:243:ALA:HB2	1:B:315:MSE:HE2	1.90	0.53
1:A:294:ALA:HB2	1:A:329:PRO:HD2	1.91	0.53
1:A:383:SER:HA	1:A:388:GLU:OE1	2.09	0.53
1:B:240:THR:HG21	1:B:282:LYS:N	2.22	0.53
1:B:91:GLY:O	1:B:388:GLU:HG3	2.09	0.53
1:B:379:ILE:HD11	1:B:389:ARG:H	1.73	0.53
1:B:366:ALA:HB3	1:B:367:PRO:HD3	1.90	0.53
1:B:319:ILE:HG22	1:B:320:ASP:H	1.72	0.53
1:B:68:LEU:HD12	1:B:116:PRO:HD2	1.91	0.53
1:B:213:VAL:CG2	1:B:290:ARG:HB3	2.39	0.53
1:A:130:GLY:CA	1:A:136:VAL:HG12	2.39	0.52
1:B:385:ASN:ND2	1:B:387:ALA:H	2.08	0.52
1:A:204:GLN:HE22	1:A:352:ARG:NH2	2.07	0.52
1:B:54:LEU:HD22	1:B:68:LEU:HD23	1.90	0.52
1:A:130:GLY:N	1:A:136:VAL:HG12	2.25	0.52
1:B:218:TYR:OH	1:B:301:THR:HB	2.10	0.52
1:B:70:GLY:O	1:B:74:PRO:HB3	2.10	0.52
1:B:28:GLY:HA3	1:B:386:PRO:HG3	1.92	0.52
1:A:91:GLY:HA3	1:A:388:GLU:CG	2.39	0.52
1:A:255:ALA:O	1:A:258:MSE:HB2	2.10	0.52
1:B:68:LEU:CD2	1:B:110:LYS:HE2	2.39	0.52
1:B:379:ILE:HA	1:B:389:ARG:NH2	2.26	0.51
1:B:128:GLU:O	1:B:128:GLU:HG3	2.07	0.51
1:B:60:GLU:HB2	1:B:179:LEU:HD21	1.92	0.51
1:B:315:MSE:HE2	1:B:317:ILE:HD12	1.91	0.51
1:B:53:GLY:H	1:B:110:LYS:HZ1	1.57	0.51
1:A:368:ARG:HD2	1:A:368:ARG:N	2.25	0.51
1:A:8:ARG:O	1:A:12:GLU:HB2	2.10	0.51
1:A:39:LEU:CD1	1:A:170:CYS:HA	2.39	0.51
1:B:216:ARG:HH21	1:B:298:ARG:HG3	1.76	0.51
1:B:227:ASN:CB	1:B:232:THR:HG21	2.41	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:268:LYS:HE2	1:A:270:GLU:OE1	2.10	0.51
1:B:40:GLU:O	1:B:44:GLN:HG2	2.10	0.51
1:B:291:HIS:ND1	1:B:292:THR:N	2.59	0.51
1:A:106:ILE:HD12	1:A:122:VAL:CG2	2.41	0.51
1:A:200:GLU:OE2	1:A:379:ILE:HA	2.11	0.50
1:B:149:PRO:HB3	1:B:174:LEU:HD13	1.92	0.50
1:B:368:ARG:NH1	1:B:368:ARG:HG3	2.26	0.50
1:B:319:ILE:O	1:B:320:ASP:HB2	2.09	0.50
1:B:323:LEU:HD11	1:B:326:ASP:HB2	1.94	0.50
1:A:55:GLU:HG2	1:A:55:GLU:O	2.11	0.50
1:A:367:PRO:HG2	1:A:368:ARG:HD2	1.91	0.50
1:B:194:GLU:HG3	1:B:375:PHE:CD2	2.46	0.50
1:B:194:GLU:HA	1:B:375:PHE:CE2	2.46	0.50
1:A:79:LEU:HD23	1:A:121:GLU:O	2.11	0.50
1:B:33:LEU:HD23	1:B:34:TYR:H	1.77	0.50
1:A:298:ARG:O	1:A:302:GLN:HG3	2.10	0.50
1:A:219:THR:HG1	1:A:324:TRP:HD1	1.58	0.50
1:A:128:GLU:C	1:A:129:GLU:HG3	2.32	0.50
1:B:85:ASP:OD1	1:B:127:GLU:HA	2.12	0.50
1:A:45:PHE:O	1:A:49:MSE:HG3	2.12	0.50
1:A:206:ILE:HD11	1:A:398:GLY:HA2	1.94	0.50
1:B:39:LEU:HG	1:B:43:GLN:HE21	1.77	0.50
1:B:212:ILE:HG12	1:B:356:SER:HB3	1.92	0.50
1:B:232:THR:HG23	1:B:232:THR:O	2.11	0.50
1:A:42:GLN:NE2	1:A:84:ILE:HG12	2.27	0.50
1:A:347:GLU:HB2	2:A:427:HOH:O	2.11	0.49
1:B:233:PRO:O	1:B:237:ARG:HG3	2.13	0.49
1:B:340:LEU:O	1:B:343:LEU:HB3	2.12	0.49
1:B:340:LEU:HD13	1:B:405:MSE:HG3	1.94	0.49
1:A:168:LYS:HB2	1:A:168:LYS:NZ	2.27	0.49
1:A:15:LEU:N	1:A:16:PRO:HD2	2.27	0.49
1:B:21:PHE:CE2	1:B:44:GLN:HG3	2.47	0.49
1:B:227:ASN:HB3	1:B:232:THR:HG21	1.93	0.49
1:B:379:ILE:HG22	1:B:380:ASN:N	2.26	0.49
1:A:204:GLN:NE2	1:A:352:ARG:NE	2.60	0.49
1:B:68:LEU:HD12	1:B:116:PRO:CD	2.42	0.49
1:B:71:THR:O	1:B:72:GLU:CB	2.61	0.49
1:B:200:GLU:HG3	1:B:201:SER:N	2.28	0.49
1:B:382:ILE:HG21	1:B:387:ALA:HB3	1.93	0.49
1:B:199:LEU:HD21	1:B:352:ARG:HB3	1.94	0.49
1:A:206:ILE:HG22	1:A:351:TYR:HB2	1.93	0.49
1:B:53:GLY:H	1:B:110:LYS:NZ	2.10	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:216:ARG:HA	1:B:325:MSE:O	2.13	0.49
1:A:130:GLY:HA3	1:A:136:VAL:HG12	1.94	0.49
1:B:50:ALA:C	1:B:52:SER:H	2.15	0.48
1:B:269:VAL:O	1:B:270:GLU:HG3	2.12	0.48
1:A:148:ASN:HD21	1:A:150:ASP:HB2	1.77	0.48
1:A:182:ARG:NE	1:A:185:ILE:HD11	2.28	0.48
1:B:15:LEU:N	1:B:16:PRO:HD2	2.27	0.48
1:B:52:SER:HB3	1:B:110:LYS:HZ1	1.78	0.48
1:B:60:GLU:OE1	1:B:179:LEU:HD23	2.14	0.48
1:B:214:GLY:O	1:B:297:LEU:HD12	2.14	0.48
1:B:57:ARG:O	1:B:64:LEU:HD12	2.12	0.48
1:A:391:ASN:O	1:A:394:ASP:HB2	2.14	0.48
1:B:404:LEU:O	1:B:407:TYR:HB3	2.13	0.48
1:B:90:GLY:HA2	1:B:386:PRO:HA	1.95	0.48
1:A:168:LYS:HA	1:A:172:PHE:O	2.14	0.48
1:A:193:ILE:HA	1:A:376:ILE:HG13	1.96	0.48
1:B:379:ILE:CD1	1:B:389:ARG:H	2.27	0.47
1:A:182:ARG:CZ	1:A:185:ILE:HD11	2.44	0.47
1:B:379:ILE:HG23	1:B:389:ARG:NH1	2.29	0.47
1:B:85:ASP:CG	1:B:127:GLU:HA	2.34	0.47
1:A:305:GLU:O	1:A:309:ARG:HD2	2.13	0.47
1:B:271:PRO:O	1:B:272:ARG:HD2	2.15	0.47
1:A:353:VAL:HG13	1:A:353:VAL:O	2.13	0.47
1:B:198:VAL:HB	1:B:354:MSE:HE1	1.96	0.47
1:A:174:LEU:H	1:A:174:LEU:HD13	1.79	0.47
1:B:330:VAL:HG21	1:B:363:GLN:HB2	1.97	0.47
1:A:246:ARG:O	1:A:250:GLN:HB2	2.14	0.47
1:A:39:LEU:HD12	1:A:170:CYS:CA	2.45	0.47
1:A:21:PHE:CE2	1:A:44:GLN:HG3	2.49	0.47
1:A:372:CYS:HB2	2:A:446:HOH:O	2.15	0.47
1:B:57:ARG:HH11	1:B:57:ARG:HG2	1.78	0.47
1:A:52:SER:O	1:A:110:LYS:HE3	2.15	0.47
1:B:14:THR:O	1:B:17:TRP:CD1	2.68	0.47
1:B:140:SER:C	1:B:364:ILE:HD11	2.35	0.46
1:B:120:VAL:HG22	1:B:410:ALA:CB	2.45	0.46
1:B:333:ASN:HB3	1:B:336:LEU:HB2	1.97	0.46
1:B:335:GLU:O	1:B:339:THR:HG22	2.15	0.46
1:B:379:ILE:HG13	1:B:389:ARG:H	1.81	0.46
1:B:298:ARG:HG3	1:B:298:ARG:HH11	1.80	0.46
1:B:5:THR:HA	1:B:8:ARG:HB3	1.98	0.46
1:A:343:LEU:HB2	1:A:346:ARG:NH1	2.30	0.46
1:A:343:LEU:HD12	1:A:343:LEU:C	2.36	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:298:ARG:O	1:B:302:GLN:HB2	2.16	0.46
1:B:249:HIS:O	1:B:250:GLN:HB2	2.16	0.46
1:A:63:ASN:HD22	1:A:125:MSE:HA	1.81	0.46
1:B:112:GLN:HE21	1:B:112:GLN:HB3	1.53	0.46
1:B:141:LYS:N	1:B:364:ILE:HD11	2.31	0.46
1:B:179:LEU:CD1	1:B:179:LEU:H	2.28	0.46
1:A:8:ARG:HD2	1:A:393:THR:HG23	1.97	0.46
1:B:42:GLN:HE21	1:B:84:ILE:HG12	1.80	0.46
1:A:174:LEU:CD1	1:A:174:LEU:H	2.28	0.46
1:B:268:LYS:HB3	1:B:286:THR:HG22	1.96	0.46
1:A:67:ARG:H	1:A:121:GLU:HA	1.80	0.45
1:A:209:VAL:CG1	1:A:373:MSE:HB2	2.46	0.45
1:B:92:ASN:HB3	1:B:388:GLU:CD	2.36	0.45
1:A:35:SER:H	1:A:38:TRP:HB3	1.80	0.45
1:A:206:ILE:O	1:A:351:TYR:HA	2.17	0.45
1:A:293:ASP:OD2	1:A:296:VAL:HG23	2.15	0.45
1:A:162:SER:HB3	1:A:165:ASP:HB2	1.98	0.45
1:B:376:ILE:CG2	1:B:390:THR:HG21	2.42	0.45
1:B:14:THR:HG21	1:B:103:TRP:HZ3	1.80	0.45
1:A:239:ASP:HB3	1:A:242:TYR:HB3	1.99	0.45
1:A:206:ILE:HG22	1:A:351:TYR:CB	2.46	0.45
1:B:199:LEU:CD2	1:B:352:ARG:HB3	2.46	0.45
1:B:79:LEU:HD21	1:B:123:VAL:HB	1.98	0.45
1:A:176:ASN:N	1:A:176:ASN:OD1	2.50	0.45
1:A:324:TRP:CZ3	1:A:325:MSE:HB2	2.51	0.45
1:B:52:SER:O	1:B:54:LEU:N	2.50	0.45
1:B:94:ASP:OD1	1:B:192:HIS:HE1	2.00	0.45
1:B:274:ASN:CG	1:B:274:ASN:O	2.55	0.44
1:B:179:LEU:HD12	1:B:179:LEU:N	2.31	0.44
1:B:35:SER:OG	1:B:36:PRO:HD2	2.17	0.44
1:A:204:GLN:HE22	1:A:352:ARG:NE	2.15	0.44
1:B:63:ASN:HD22	1:B:126:ALA:H	1.64	0.44
1:B:309:ARG:CD	1:B:319:ILE:HD13	2.48	0.44
1:A:39:LEU:O	1:A:43:GLN:HG2	2.17	0.44
1:B:332:MSE:HE3	1:B:372:CYS:HA	1.98	0.44
1:A:148:ASN:HA	1:A:178:PRO:HG3	1.99	0.44
1:A:91:GLY:HA3	1:A:388:GLU:HG2	1.98	0.44
1:B:36:PRO:O	1:B:37:GLU:CB	2.63	0.44
1:B:332:MSE:HE3	1:B:372:CYS:CA	2.47	0.44
1:B:48:ARG:HB3	1:B:103:TRP:CZ2	2.52	0.44
1:A:364:ILE:O	1:A:368:ARG:HD3	2.18	0.44
1:B:216:ARG:NH2	1:B:298:ARG:HG3	2.33	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:390:THR:HB	1:B:395:LEU:HD21	2.00	0.44
1:A:368:ARG:HG3	1:A:368:ARG:NH1	2.32	0.44
1:B:221:THR:HB	1:B:320:ASP:HB3	2.00	0.44
1:B:66:GLY:O	1:B:122:VAL:HG12	2.18	0.44
1:A:120:VAL:HG11	1:A:411:TRP:HZ3	1.83	0.43
1:B:105:ALA:HB1	1:B:403:ALA:HA	2.00	0.43
1:B:251:SER:O	1:B:254:LYS:HB2	2.18	0.43
1:B:32:LEU:O	1:B:38:TRP:HB2	2.18	0.43
1:A:199:LEU:HB3	1:A:377:PRO:HG3	1.99	0.43
1:A:216:ARG:HG2	1:A:323:LEU:HD21	2.00	0.43
1:A:218:TYR:OH	1:A:301:THR:HB	2.18	0.43
1:A:343:LEU:O	1:A:347:GLU:HB3	2.17	0.43
1:A:105:ALA:HB1	1:A:403:ALA:HA	2.00	0.43
1:A:149:PRO:CB	1:A:174:LEU:HD23	2.39	0.43
1:A:39:LEU:HD12	1:A:170:CYS:CB	2.48	0.43
1:B:35:SER:HB2	1:B:132:ARG:NH2	2.33	0.43
1:A:14:THR:HG21	1:A:100:LEU:HD11	2.00	0.43
1:B:90:GLY:HA2	1:B:386:PRO:CA	2.49	0.43
1:A:33:LEU:HG	1:A:127:GLU:HB3	2.00	0.43
1:B:258:MSE:HB2	1:B:262:LEU:HD23	2.00	0.43
1:B:83:HIS:HB2	1:B:85:ASP:OD1	2.19	0.43
1:A:33:LEU:HD22	1:A:34:TYR:CG	2.53	0.43
1:A:147:ALA:O	1:A:148:ASN:CB	2.67	0.43
1:B:351:TYR:HD2	1:B:351:TYR:H	1.65	0.43
1:B:238:ARG:HD3	1:B:316:ASP:HB3	2.01	0.43
1:B:64:LEU:HB3	1:B:124:ALA:HB3	1.99	0.43
1:A:276:VAL:HA	1:B:286:THR:HG21	2.01	0.43
1:A:345:GLU:C	1:A:347:GLU:N	2.71	0.43
1:B:74:PRO:CG	1:B:75:GLN:HE21	2.21	0.42
1:A:253:GLU:OE1	1:A:253:GLU:HA	2.19	0.42
1:B:50:ALA:HB2	1:B:56:THR:HG21	2.01	0.42
1:B:46:LYS:HZ1	1:B:56:THR:HG23	1.84	0.42
1:A:106:ILE:HD12	1:A:122:VAL:HG22	2.01	0.42
1:B:3:LEU:HD12	1:B:400:LYS:NZ	2.34	0.42
1:B:3:LEU:O	1:B:6:HIS:HB3	2.18	0.42
1:B:240:THR:HG21	1:B:282:LYS:H	1.84	0.42
1:A:241:VAL:HB	1:A:280:PRO:HG3	2.01	0.42
1:B:379:ILE:HG13	1:B:389:ARG:HG3	2.00	0.42
1:A:315:MSE:HG3	1:A:317:ILE:CD1	2.49	0.42
1:B:64:LEU:O	1:B:123:VAL:HA	2.19	0.42
1:B:68:LEU:CG	1:B:110:LYS:HG3	2.50	0.42
1:B:35:SER:HB2	1:B:132:ARG:HH22	1.84	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:61:VAL:HG21	1:A:142:ASN:ND2	2.34	0.42
1:B:68:LEU:HD11	1:B:110:LYS:HG3	2.00	0.42
1:B:253:GLU:OE2	1:B:256:LYS:HD3	2.19	0.42
1:B:343:LEU:O	1:B:343:LEU:HD13	2.19	0.42
1:A:81:GLY:HA3	1:A:123:VAL:O	2.20	0.42
1:A:64:LEU:HB3	1:A:124:ALA:HB3	2.02	0.42
1:B:388:GLU:HG3	2:B:424:HOH:O	2.19	0.42
1:A:204:GLN:HE22	1:A:352:ARG:CZ	2.33	0.42
1:B:174:LEU:HA	1:B:175:PRO:HD3	1.88	0.42
1:B:108:TRP:O	1:B:112:GLN:HB2	2.20	0.42
1:A:15:LEU:O	1:A:19:SER:CB	2.68	0.42
1:B:194:GLU:O	1:B:195:GLN:HB2	2.20	0.42
1:B:240:THR:HG23	1:B:283:THR:OG1	2.19	0.41
1:A:194:GLU:O	1:A:195:GLN:CB	2.66	0.41
1:A:400:LYS:O	1:A:404:LEU:HG	2.20	0.41
1:B:46:LYS:HD2	1:B:58:PHE:CE1	2.55	0.41
1:B:29:MSE:H	1:B:90:GLY:H	1.68	0.41
1:A:89:ASN:O	1:A:386:PRO:HD3	2.20	0.41
1:A:222:LEU:HD13	1:A:285:PHE:HE2	1.85	0.41
1:A:61:VAL:HG22	1:A:138:TRP:NE1	2.35	0.41
1:B:129:GLU:O	1:B:129:GLU:HG2	2.20	0.41
1:A:102:ALA:HB3	1:A:122:VAL:HG13	2.02	0.41
1:B:323:LEU:O	1:B:323:LEU:HD13	2.20	0.41
1:A:189:VAL:O	1:A:189:VAL:HG23	2.19	0.41
1:A:68:LEU:HB3	1:A:120:VAL:HG23	2.02	0.41
1:A:217:ARG:HD3	1:B:276:VAL:HG21	2.03	0.41
1:B:60:GLU:CD	1:B:60:GLU:N	2.72	0.41
1:A:312:CYS:HB3	1:A:317:ILE:O	2.20	0.41
1:B:339:THR:HA	1:B:342:GLU:OE1	2.21	0.41
1:A:88:VAL:HG23	1:A:89:ASN:H	1.86	0.41
1:B:136:VAL:O	1:B:137:PHE:HB2	2.21	0.41
1:B:110:LYS:O	1:B:114:GLY:N	2.54	0.41
1:B:167:MSE:HG2	1:B:172:PHE:O	2.21	0.41
1:B:175:PRO:HG2	1:B:179:LEU:CD1	2.50	0.41
1:B:35:SER:H	1:B:132:ARG:HH12	1.68	0.41
1:B:336:LEU:HB3	1:B:372:CYS:HB3	2.03	0.41
1:A:108:TRP:O	1:A:112:GLN:HB2	2.21	0.41
1:A:174:LEU:N	1:A:174:LEU:HD13	2.36	0.41
1:B:30:THR:HG23	1:B:88:VAL:HA	2.03	0.41
1:A:106:ILE:HD12	1:A:122:VAL:HG23	2.03	0.40
1:A:264:LEU:N	1:B:234:MSE:HE1	2.34	0.40
1:B:64:LEU:O	1:B:123:VAL:HG23	2.21	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:365:PHE:HB3	1:A:371:THR:CG2	2.36	0.40
1:B:240:THR:HG22	1:B:280:PRO:C	2.39	0.40
1:B:364:ILE:O	1:B:364:ILE:HG13	2.21	0.40
1:A:262:LEU:HB2	1:A:291:HIS:HD2	1.86	0.40
1:A:240:THR:HG22	1:A:280:PRO:C	2.39	0.40
1:B:379:ILE:N	1:B:379:ILE:CD1	2.81	0.40
1:A:273:PRO:CG	1:A:278:VAL:HG11	2.51	0.40
1:B:57:ARG:HG2	1:B:57:ARG:NH1	2.37	0.40
1:B:73:TYR:N	1:B:74:PRO:CD	2.84	0.40
1:B:67:ARG:O	1:B:67:ARG:HG3	2.21	0.40
1:A:286:THR:HG21	1:B:276:VAL:HA	2.04	0.40
1:A:183:GLN:HA	2:A:443:HOH:O	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	393/423 (93%)	350 (89%)	35 (9%)	8 (2%)	11	35
1	B	390/423 (92%)	328 (84%)	49 (13%)	13 (3%)	6	19
All	All	783/846 (93%)	678 (87%)	84 (11%)	21 (3%)	8	25

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	34	TYR
1	B	36	PRO
1	B	320	ASP
1	B	379	ILE
1	A	129	GLU
1	B	32	LEU
1	B	250	GLN
1	B	37	GLU

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Mol	Chain	Res	Type
1	B	175	PRO
1	A	31	ARG
1	A	67	ARG
1	A	348	LYS
1	A	148	ASN
1	A	112	GLN
1	B	33	LEU
1	B	73	TYR
1	B	113	TYR
1	A	88	VAL
1	B	319	ILE
1	B	53	GLY
1	A	130	GLY

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	338/341 (99%)	306 (90%)	32 (10%)	12	33
1	B	331/341 (97%)	296 (89%)	35 (11%)	10	27
All	All	669/682 (98%)	602 (90%)	67 (10%)	11	30

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	LEU
1	A	33	LEU
1	A	45	PHE
1	A	48	ARG
1	A	56	THR
1	A	61	VAL
1	A	79	LEU
1	A	86	THR
1	A	88	VAL
1	A	93	LEU
1	A	120	VAL
1	A	128	GLU

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Mol	Chain	Res	Type
1	A	129	GLU
1	A	168	LYS
1	A	174	LEU
1	A	176	ASN
1	A	191	LEU
1	A	206	ILE
1	A	209	VAL
1	A	215	GLN
1	A	216	ARG
1	A	246	ARG
1	A	264	LEU
1	A	265	THR
1	A	274	ASN
1	A	304	LEU
1	A	323	LEU
1	A	327	GLU
1	A	343	LEU
1	A	347	GLU
1	A	368	ARG
1	A	393	THR
1	B	3	LEU
1	B	13	GLU
1	B	14	THR
1	B	35	SER
1	B	36	PRO
1	B	44	GLN
1	B	54	LEU
1	B	68	LEU
1	B	69	ASN
1	B	79	LEU
1	B	112	GLN
1	B	128	GLU
1	B	192	HIS
1	B	234	MSE
1	B	236	TYR
1	B	246	ARG
1	B	250	GLN
1	B	264	LEU
1	B	272	ARG
1	B	274	ASN
1	B	277	ASN
1	B	297	LEU

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Mol	Chain	Res	Type
1	B	319	ILE
1	B	339	THR
1	B	351	TYR
1	B	352	ARG
1	B	364	ILE
1	B	375	PHE
1	B	376	ILE
1	B	378	SER
1	B	388	GLU
1	B	389	ARG
1	B	397	GLU
1	B	405	MSE
1	B	408	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	GLN
1	A	63	ASN
1	A	148	ASN
1	A	195	GLN
1	A	202	ASN
1	A	204	GLN
1	A	274	ASN
1	A	303	GLN
1	A	363	GLN
1	B	42	GLN
1	B	43	GLN
1	B	44	GLN
1	B	63	ASN
1	B	75	GLN
1	B	89	ASN
1	B	92	ASN
1	B	96	GLN
1	B	112	GLN
1	B	195	GLN
1	B	228	HIS
1	B	277	ASN
1	B	303	GLN
1	B	360	HIS
1	B	363	GLN
1	B	384	HIS

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Mol	Chain	Res	Type
1	B	385	ASN
1	B	391	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 ⓘ

There are no ligands in this entry.

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	399/423 (94%)	-0.13	2 (0%) 88 90	26, 45, 66, 84	0
1	B	394/423 (93%)	0.15	12 (3%) 48 49	32, 58, 91, 99	0
All	All	793/846 (93%)	0.01	14 (1%) 65 66	26, 51, 84, 99	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	26	ALA	6.3
1	B	53	GLY	5.4
1	B	175	PRO	3.5
1	B	131	SER	3.5
1	B	27	GLY	3.4
1	B	129	GLU	3.0
1	A	86	THR	2.5
1	B	61	VAL	2.5
1	B	106	ILE	2.2
1	A	75	GLN	2.2
1	B	348	LYS	2.1
1	B	215	GLN	2.1
1	B	45	PHE	2.1
1	B	374	ILE	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.