



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2023 – 10:37 AM EDT

PDB ID : 2R89  
Title : Crystal structure of the long-chain fatty acid transporter FadL mutant delta N3  
Authors : Hearn, E.M.; Patel, D.R.; Lepore, B.W.; Indic, M.; van den Berg, B.  
Deposited on : 2007-09-10  
Resolution : 3.40 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](mailto: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>.

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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 : 2.35  
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) : 2.35

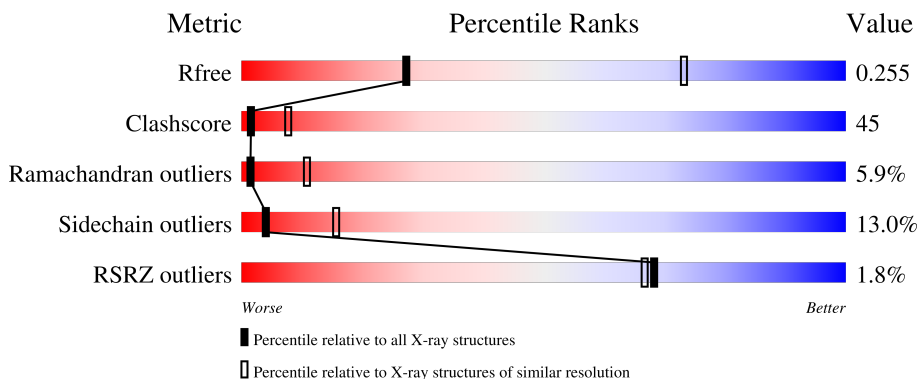
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.40 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	1026 (3.48-3.32)
Clashscore	141614	1055 (3.48-3.32)
Ramachandran outliers	138981	1038 (3.48-3.32)
Sidechain outliers	138945	1038 (3.48-3.32)
RSRZ outliers	127900	2173 (3.50-3.30)

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  $\geq 3$ , 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	424	
1	B	424	

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 5608 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 Long-chain fatty acid transport protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	363	2804	1774	475	549	6	0	0	0
1	B	363	2804	1774	475	549	6	0	0	0

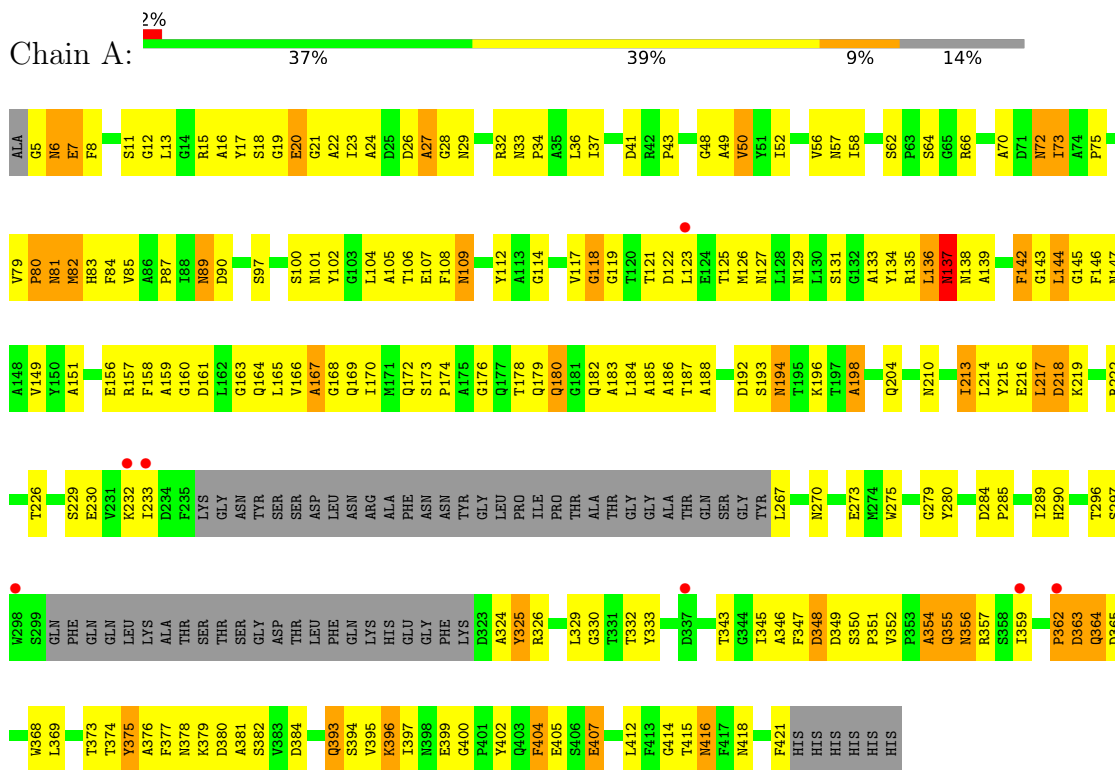
There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	PHE	deletion	UNP P10384
A	?	-	GLN	deletion	UNP P10384
A	?	-	LEU	deletion	UNP P10384
A	197	THR	ILE	conflict	UNP P10384
A	422	HIS	-	expression tag	UNP P10384
A	423	HIS	-	expression tag	UNP P10384
A	424	HIS	-	expression tag	UNP P10384
A	425	HIS	-	expression tag	UNP P10384
A	426	HIS	-	expression tag	UNP P10384
A	427	HIS	-	expression tag	UNP P10384
B	?	-	PHE	deletion	UNP P10384
B	?	-	GLN	deletion	UNP P10384
B	?	-	LEU	deletion	UNP P10384
B	197	THR	ILE	conflict	UNP P10384
B	422	HIS	-	expression tag	UNP P10384
B	423	HIS	-	expression tag	UNP P10384
B	424	HIS	-	expression tag	UNP P10384
B	425	HIS	-	expression tag	UNP P10384
B	426	HIS	-	expression tag	UNP P10384
B	427	HIS	-	expression tag	UNP P10384

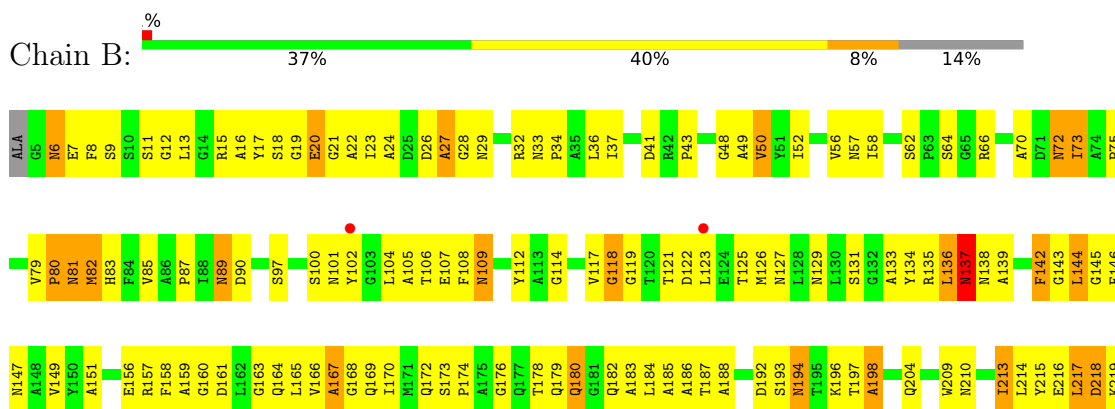
### 3 Residue-property plots

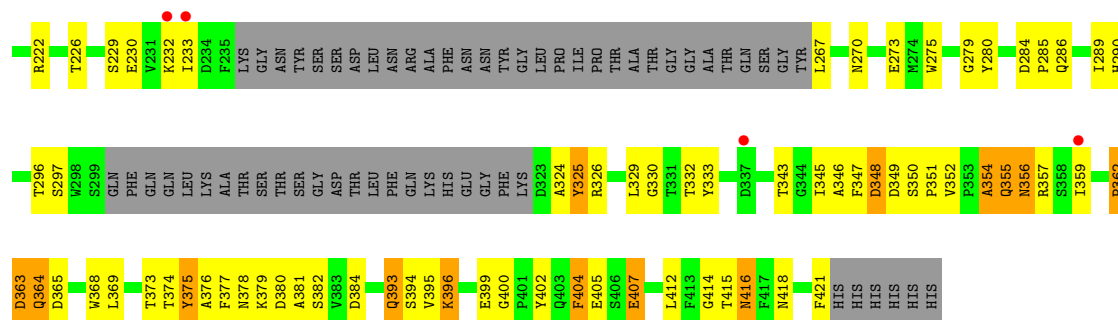
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Long-chain fatty acid transport protein



- Molecule 1: Long-chain fatty acid transport protein





## 4 Data and refinement statistics i

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	92.02Å 91.91Å 289.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 3.40 38.81 – 3.40	Depositor EDS
% Data completeness (in resolution range)	97.7 (10.00-3.40) 98.6 (38.81-3.40)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.48 (at 3.40Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.252 , 0.302 0.257 , 0.255	Depositor DCC
$R_{free}$ test set	1320 reflections (7.74%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	78.5	Xtrriage
Anisotropy	0.444	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , -1.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.459 for -k,-h,-l	Xtrriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	5608	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.40% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.50	0/2880	0.73	0/3924
1	B	0.50	0/2880	0.73	0/3924
All	All	0.50	0/5760	0.73	0/7848

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 [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	2804	0	2598	243	0
1	B	2804	0	2598	252	0
All	All	5608	0	5196	491	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 45.

All (491) 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:137:ASN:HD22	1:A:138:ASN:N	1.55	1.04
1:B:137:ASN:HD22	1:B:138:ASN:N	1.54	1.04

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:VAL:HG11	1:A:357:ARG:HA	1.47	0.97
1:B:52:ILE:HG12	1:B:412:LEU:HD22	1.48	0.96
1:B:352:VAL:HG11	1:B:357:ARG:HA	1.47	0.96
1:A:52:ILE:HG12	1:A:412:LEU:HD22	1.48	0.95
1:B:109:ASN:HD22	1:B:109:ASN:H	1.00	0.94
1:A:48:GLY:HA3	1:A:416:ASN:HD22	1.32	0.93
1:B:57:ASN:ND2	1:B:72:ASN:H	1.68	0.92
1:A:57:ASN:ND2	1:A:72:ASN:H	1.67	0.92
1:A:109:ASN:H	1:A:109:ASN:HD22	0.99	0.92
1:B:48:GLY:HA3	1:B:416:ASN:HD22	1.32	0.91
1:A:109:ASN:H	1:A:109:ASN:ND2	1.69	0.89
1:A:378:ASN:O	1:A:380:ASP:N	2.06	0.89
1:B:378:ASN:O	1:B:380:ASP:N	2.06	0.88
1:B:89:ASN:ND2	1:B:90:ASP:H	1.71	0.88
1:A:89:ASN:ND2	1:A:90:ASP:H	1.72	0.87
1:B:66:ARG:HH11	1:B:164:GLN:NE2	1.72	0.87
1:B:66:ARG:HH11	1:B:164:GLN:HE22	0.89	0.87
1:B:109:ASN:H	1:B:109:ASN:ND2	1.70	0.87
1:A:66:ARG:HH11	1:A:164:GLN:NE2	1.72	0.87
1:A:66:ARG:NH1	1:A:164:GLN:HE22	1.73	0.87
1:A:33:ASN:HB3	1:A:36:LEU:HD13	1.55	0.86
1:A:66:ARG:HH11	1:A:164:GLN:HE22	0.90	0.86
1:B:33:ASN:HB3	1:B:36:LEU:HD13	1.55	0.86
1:B:66:ARG:NH1	1:B:164:GLN:HE22	1.73	0.83
1:B:7:GLU:HB3	1:B:50:VAL:HG21	1.60	0.83
1:B:137:ASN:HD22	1:B:138:ASN:H	1.27	0.83
1:A:402:TYR:HB2	1:A:404:PHE:CE1	2.14	0.82
1:B:402:TYR:HB2	1:B:404:PHE:CE1	2.14	0.82
1:A:137:ASN:HD22	1:A:138:ASN:H	1.28	0.82
1:B:48:GLY:HA3	1:B:416:ASN:ND2	1.94	0.81
1:B:186:ALA:C	1:B:188:ALA:H	1.82	0.81
1:A:186:ALA:C	1:A:188:ALA:H	1.81	0.81
1:A:48:GLY:HA3	1:A:416:ASN:ND2	1.94	0.81
1:A:213:ILE:O	1:A:213:ILE:HG22	1.81	0.80
1:B:213:ILE:HG22	1:B:213:ILE:O	1.79	0.80
1:B:136:LEU:O	1:B:136:LEU:HG	1.83	0.78
1:B:178:THR:HG22	1:B:180:GLN:H	1.48	0.78
1:A:136:LEU:O	1:A:136:LEU:HG	1.82	0.77
1:B:73:ILE:HD12	1:B:108:PHE:CE2	2.19	0.77
1:B:158:PHE:HE2	1:B:196:LYS:HG3	1.50	0.77
1:A:73:ILE:HD12	1:A:108:PHE:CE2	2.19	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:THR:HG22	1:A:180:GLN:H	1.48	0.76
1:B:89:ASN:CG	1:B:90:ASP:H	1.88	0.76
1:B:49:ALA:HB2	1:B:80:PRO:HA	1.67	0.76
1:A:49:ALA:HB2	1:A:80:PRO:HA	1.67	0.76
1:A:158:PHE:HE2	1:A:196:LYS:HG3	1.50	0.76
1:A:109:ASN:HD22	1:A:109:ASN:N	1.80	0.75
1:A:137:ASN:ND2	1:A:138:ASN:N	2.34	0.75
1:A:363:ASP:C	1:A:364:GLN:HG3	2.06	0.75
1:A:52:ILE:HG12	1:A:412:LEU:CD2	2.17	0.75
1:A:89:ASN:CG	1:A:90:ASP:H	1.89	0.75
1:A:101:ASN:HB3	1:A:102:TYR:CE1	2.22	0.75
1:B:101:ASN:HB3	1:B:102:TYR:CE1	2.22	0.74
1:A:382:SER:OG	1:A:418:ASN:HB2	1.87	0.74
1:B:52:ILE:HG12	1:B:412:LEU:CD2	2.17	0.74
1:B:363:ASP:C	1:B:364:GLN:HG3	2.06	0.74
1:B:382:SER:OG	1:B:418:ASN:HB2	1.88	0.74
1:B:137:ASN:ND2	1:B:138:ASN:N	2.33	0.73
1:B:56:VAL:HG12	1:B:73:ILE:HG23	1.71	0.73
1:A:41:ASP:O	1:A:87:PRO:HG3	1.88	0.73
1:B:41:ASP:O	1:B:87:PRO:HG3	1.88	0.72
1:B:7:GLU:CB	1:B:50:VAL:HG21	2.19	0.72
1:B:326:ARG:HG3	1:B:348:ASP:HB3	1.71	0.72
1:A:56:VAL:HG12	1:A:73:ILE:HG23	1.70	0.72
1:A:186:ALA:O	1:A:188:ALA:N	2.21	0.72
1:B:109:ASN:HD22	1:B:109:ASN:N	1.81	0.71
1:A:109:ASN:ND2	1:A:109:ASN:N	2.36	0.71
1:A:326:ARG:HG3	1:A:348:ASP:HB3	1.72	0.71
1:B:186:ALA:O	1:B:188:ALA:N	2.22	0.70
1:A:89:ASN:ND2	1:A:90:ASP:N	2.40	0.70
1:A:125:THR:HG23	1:A:149:VAL:HG13	1.73	0.70
1:B:394:SER:HB2	1:B:407:GLU:HB2	1.74	0.70
1:B:117:VAL:HG23	1:B:359:ILE:HD11	1.73	0.70
1:A:106:THR:HG21	1:A:362:PRO:CD	2.22	0.70
1:B:158:PHE:CE2	1:B:196:LYS:HG3	2.27	0.70
1:A:34:PRO:HB2	1:A:143:GLY:HA3	1.74	0.69
1:A:117:VAL:HG23	1:A:359:ILE:HD11	1.73	0.69
1:B:89:ASN:ND2	1:B:90:ASP:N	2.39	0.69
1:B:125:THR:HG23	1:B:149:VAL:HG13	1.73	0.69
1:A:158:PHE:CE2	1:A:196:LYS:HG3	2.27	0.69
1:B:34:PRO:HB2	1:B:143:GLY:HA3	1.75	0.69
1:A:394:SER:HB2	1:A:407:GLU:HB2	1.74	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:106:THR:HG21	1:B:362:PRO:CD	2.22	0.68
1:A:24:ALA:HB2	1:A:36:LEU:HD23	1.75	0.68
1:B:109:ASN:ND2	1:B:109:ASN:N	2.37	0.68
1:A:50:VAL:HG12	1:A:414:GLY:HA3	1.77	0.67
1:B:381:ALA:HA	1:B:418:ASN:O	1.93	0.67
1:B:24:ALA:HB2	1:B:36:LEU:HD23	1.75	0.67
1:A:381:ALA:HA	1:A:418:ASN:O	1.94	0.67
1:B:131:SER:HB3	1:B:145:GLY:HA3	1.77	0.67
1:A:131:SER:HB3	1:A:145:GLY:HA3	1.77	0.67
1:B:50:VAL:HG12	1:B:414:GLY:HA3	1.77	0.66
1:A:8:PHE:CE2	1:A:32:ARG:HD2	2.30	0.66
1:A:82:MET:O	1:A:97:SER:HA	1.96	0.66
1:A:43:PRO:HG2	1:A:421:PHE:HB2	1.78	0.66
1:B:82:MET:O	1:B:97:SER:HA	1.96	0.66
1:B:7:GLU:HA	1:B:17:TYR:OH	1.96	0.65
1:B:49:ALA:CB	1:B:80:PRO:HA	2.26	0.65
1:B:157:ARG:HB2	1:B:198:ALA:HB3	1.78	0.65
1:B:20:GLU:HB2	1:B:32:ARG:HD3	1.79	0.65
1:A:20:GLU:HB2	1:A:32:ARG:HD3	1.79	0.64
1:A:157:ARG:HB2	1:A:198:ALA:HB3	1.78	0.64
1:A:407:GLU:HG3	1:A:407:GLU:O	1.98	0.64
1:B:43:PRO:HG2	1:B:421:PHE:HB2	1.79	0.64
1:B:186:ALA:C	1:B:188:ALA:N	2.51	0.64
1:A:32:ARG:HG3	1:A:32:ARG:HH11	1.63	0.64
1:A:49:ALA:CB	1:A:80:PRO:HA	2.27	0.64
1:A:378:ASN:C	1:A:380:ASP:H	2.02	0.64
1:B:378:ASN:C	1:B:380:ASP:H	2.01	0.64
1:B:73:ILE:HG12	1:B:73:ILE:O	1.97	0.64
1:B:32:ARG:HG3	1:B:32:ARG:HH11	1.63	0.63
1:A:73:ILE:HG12	1:A:73:ILE:O	1.97	0.63
1:A:347:PHE:CG	1:A:348:ASP:N	2.66	0.63
1:B:347:PHE:CG	1:B:348:ASP:N	2.67	0.63
1:A:232:LYS:HG2	1:A:270:ASN:HD21	1.63	0.63
1:B:232:LYS:HG2	1:B:270:ASN:HD21	1.63	0.63
1:B:64:SER:HB3	1:B:165:LEU:HD21	1.81	0.63
1:B:108:PHE:HB2	1:B:118:GLY:O	1.99	0.63
1:A:108:PHE:HB2	1:A:118:GLY:O	1.99	0.63
1:B:407:GLU:O	1:B:407:GLU:HG3	1.97	0.62
1:A:144:LEU:HD21	1:A:146:PHE:CE1	2.35	0.62
1:A:106:THR:HG21	1:A:362:PRO:HD3	1.80	0.62
1:B:144:LEU:HD21	1:B:146:PHE:CE1	2.35	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:ASN:HD22	1:A:81:ASN:C	2.01	0.62
1:B:106:THR:HG21	1:B:362:PRO:HD3	1.80	0.62
1:A:284:ASP:CG	1:A:285:PRO:HD2	2.21	0.61
1:B:6:ASN:ND2	1:B:101:ASN:O	2.33	0.61
1:B:284:ASP:CG	1:B:285:PRO:HD2	2.20	0.61
1:A:52:ILE:HG23	1:A:412:LEU:HD21	1.84	0.60
1:A:64:SER:HB3	1:A:165:LEU:HD21	1.81	0.60
1:B:52:ILE:HG23	1:B:412:LEU:HD21	1.83	0.60
1:A:48:GLY:CA	1:A:416:ASN:ND2	2.64	0.60
1:B:19:GLY:HA3	1:B:290:HIS:HB2	1.84	0.60
1:B:136:LEU:O	1:B:136:LEU:CG	2.50	0.60
1:A:19:GLY:HA3	1:A:290:HIS:HB2	1.84	0.59
1:A:5:GLY:HA2	1:A:17:TYR:CB	2.32	0.59
1:B:8:PHE:CE2	1:B:32:ARG:HD2	2.38	0.59
1:B:48:GLY:CA	1:B:416:ASN:ND2	2.64	0.59
1:B:213:ILE:O	1:B:213:ILE:CG2	2.51	0.59
1:A:75:PRO:HD2	1:A:105:ALA:O	2.03	0.59
1:A:57:ASN:ND2	1:A:72:ASN:N	2.46	0.59
1:B:89:ASN:CG	1:B:90:ASP:N	2.57	0.58
1:A:16:ALA:O	1:A:326:ARG:NH2	2.36	0.58
1:A:15:ARG:HD2	1:A:20:GLU:OE1	2.02	0.58
1:A:136:LEU:O	1:A:137:ASN:HB2	2.03	0.58
1:B:136:LEU:O	1:B:137:ASN:HB2	2.04	0.58
1:B:163:GLY:HA3	1:B:188:ALA:O	2.04	0.58
1:A:142:PHE:CD1	1:A:142:PHE:N	2.70	0.58
1:B:15:ARG:HD2	1:B:20:GLU:OE1	2.02	0.58
1:B:75:PRO:HD2	1:B:105:ALA:O	2.03	0.58
1:B:16:ALA:O	1:B:326:ARG:NH2	2.37	0.58
1:A:217:LEU:O	1:A:218:ASP:HB3	2.03	0.58
1:A:186:ALA:C	1:A:188:ALA:N	2.51	0.57
1:A:213:ILE:O	1:A:213:ILE:CG2	2.52	0.57
1:B:217:LEU:O	1:B:218:ASP:HB3	2.03	0.57
1:B:421:PHE:N	1:B:421:PHE:CD2	2.71	0.57
1:B:142:PHE:N	1:B:142:PHE:CD1	2.70	0.57
1:A:89:ASN:CG	1:A:90:ASP:N	2.57	0.57
1:A:136:LEU:O	1:A:136:LEU:CG	2.49	0.57
1:A:105:ALA:HA	1:A:121:THR:O	2.04	0.57
1:B:81:ASN:HD22	1:B:81:ASN:C	2.01	0.57
1:B:105:ALA:HA	1:B:121:THR:O	2.04	0.57
1:A:119:GLY:HA3	1:A:156:GLU:O	2.05	0.57
1:A:163:GLY:HA3	1:A:188:ALA:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:VAL:O	1:A:170:ILE:HG13	2.05	0.57
1:A:6:ASN:OD1	1:A:101:ASN:O	2.22	0.57
1:B:166:VAL:O	1:B:170:ILE:HG13	2.05	0.57
1:A:421:PHE:N	1:A:421:PHE:CD2	2.71	0.56
1:B:119:GLY:HA3	1:B:156:GLU:O	2.05	0.56
1:B:57:ASN:ND2	1:B:72:ASN:N	2.47	0.56
1:A:19:GLY:O	1:A:20:GLU:C	2.44	0.56
1:A:57:ASN:HD22	1:A:72:ASN:H	1.50	0.56
1:A:26:ASP:O	1:A:28:GLY:N	2.40	0.55
1:A:73:ILE:HD12	1:A:108:PHE:HE2	1.71	0.55
1:B:49:ALA:HB2	1:B:80:PRO:CA	2.35	0.55
1:B:81:ASN:HD21	1:B:416:ASN:HD21	1.55	0.55
1:B:26:ASP:O	1:B:28:GLY:N	2.40	0.55
1:B:375:TYR:HD2	1:B:375:TYR:C	2.10	0.55
1:B:19:GLY:O	1:B:20:GLU:C	2.44	0.54
1:A:49:ALA:HB2	1:A:80:PRO:CA	2.36	0.54
1:B:73:ILE:HD12	1:B:108:PHE:HE2	1.71	0.54
1:B:329:LEU:HD23	1:B:330:GLY:N	2.22	0.54
1:B:182:GLN:O	1:B:185:ALA:HB3	2.07	0.54
1:B:357:ARG:NH1	1:B:395:VAL:HB	2.23	0.54
1:A:84:PHE:HZ	1:B:325:TYR:CE1	2.25	0.54
1:A:182:GLN:O	1:A:185:ALA:HB3	2.07	0.54
1:A:375:TYR:HD2	1:A:375:TYR:C	2.11	0.54
1:B:144:LEU:C	1:B:144:LEU:HD23	2.28	0.54
1:A:352:VAL:HG11	1:A:357:ARG:CA	2.30	0.54
1:A:357:ARG:NH1	1:A:395:VAL:HB	2.23	0.54
1:B:33:ASN:CB	1:B:36:LEU:HD13	2.34	0.54
1:A:375:TYR:C	1:A:375:TYR:CD2	2.82	0.54
1:A:8:PHE:CD1	1:A:8:PHE:C	2.82	0.53
1:A:81:ASN:HD21	1:A:416:ASN:HD21	1.57	0.53
1:B:27:ALA:HB1	1:B:85:VAL:HG23	1.90	0.53
1:B:178:THR:HG22	1:B:180:GLN:N	2.21	0.53
1:A:144:LEU:C	1:A:144:LEU:HD23	2.28	0.53
1:B:125:THR:HG23	1:B:149:VAL:CG1	2.39	0.53
1:A:20:GLU:HA	1:A:29:ASN:ND2	2.24	0.53
1:A:125:THR:HG23	1:A:149:VAL:CG1	2.39	0.53
1:A:62:SER:C	1:A:64:SER:H	2.13	0.52
1:B:7:GLU:H	1:B:7:GLU:CD	2.11	0.52
1:B:17:TYR:HA	1:B:20:GLU:OE1	2.10	0.52
1:A:178:THR:HG22	1:A:180:GLN:N	2.20	0.52
1:A:329:LEU:HD23	1:A:330:GLY:N	2.23	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:374:THR:HG23	1:B:384:ASP:OD2	2.09	0.52
1:A:374:THR:HG23	1:A:384:ASP:OD2	2.09	0.52
1:B:144:LEU:HD23	1:B:144:LEU:O	2.10	0.52
1:A:214:LEU:HD11	1:A:222:ARG:HB2	1.92	0.52
1:A:27:ALA:HB1	1:A:85:VAL:HG23	1.90	0.52
1:A:174:PRO:C	1:A:176:GLY:H	2.13	0.52
1:B:214:LEU:HD11	1:B:222:ARG:HB2	1.92	0.52
1:A:33:ASN:CB	1:A:36:LEU:HD13	2.34	0.52
1:A:144:LEU:HD23	1:A:144:LEU:O	2.10	0.52
1:A:158:PHE:HD2	1:A:196:LYS:HA	1.75	0.52
1:B:57:ASN:HD22	1:B:72:ASN:H	1.51	0.52
1:B:73:ILE:HD12	1:B:108:PHE:CZ	2.44	0.52
1:A:421:PHE:N	1:A:421:PHE:HD2	2.07	0.51
1:B:20:GLU:HA	1:B:29:ASN:ND2	2.24	0.51
1:A:81:ASN:O	1:A:81:ASN:ND2	2.32	0.51
1:A:104:LEU:HD12	1:A:105:ALA:N	2.25	0.51
1:B:158:PHE:HD2	1:B:196:LYS:HA	1.75	0.51
1:A:404:PHE:N	1:A:404:PHE:CD1	2.79	0.51
1:B:18:SER:OG	1:B:290:HIS:HD2	1.94	0.51
1:B:174:PRO:C	1:B:176:GLY:H	2.13	0.51
1:A:73:ILE:HD12	1:A:108:PHE:CZ	2.45	0.51
1:A:131:SER:HB3	1:A:145:GLY:CA	2.41	0.51
1:B:8:PHE:C	1:B:8:PHE:CD1	2.84	0.51
1:B:34:PRO:HB2	1:B:143:GLY:CA	2.41	0.51
1:A:5:GLY:N	1:A:326:ARG:NE	2.58	0.51
1:B:421:PHE:N	1:B:421:PHE:HD2	2.08	0.51
1:B:21:GLY:HA2	1:B:33:ASN:HB2	1.93	0.51
1:B:52:ILE:HG23	1:B:412:LEU:CD2	2.41	0.51
1:B:79:VAL:CG1	1:B:100:SER:HB3	2.41	0.51
1:B:404:PHE:N	1:B:404:PHE:CD1	2.79	0.51
1:A:17:TYR:HA	1:A:20:GLU:OE1	2.10	0.50
1:A:79:VAL:CG1	1:A:100:SER:HB3	2.42	0.50
1:A:332:THR:HG22	1:A:333:TYR:N	2.26	0.50
1:B:62:SER:C	1:B:64:SER:H	2.13	0.50
1:B:81:ASN:O	1:B:81:ASN:ND2	2.32	0.50
1:A:284:ASP:CG	1:A:285:PRO:CD	2.80	0.50
1:B:104:LEU:HD12	1:B:105:ALA:N	2.26	0.50
1:B:332:THR:HG22	1:B:333:TYR:N	2.25	0.50
1:B:375:TYR:C	1:B:375:TYR:CD2	2.81	0.50
1:A:18:SER:OG	1:A:290:HIS:HD2	1.95	0.50
1:A:52:ILE:HG23	1:A:412:LEU:CD2	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:PRO:HB2	1:A:143:GLY:CA	2.40	0.50
1:A:280:TYR:HA	1:A:289:ILE:O	2.12	0.50
1:B:280:TYR:HA	1:B:289:ILE:O	2.12	0.50
1:B:402:TYR:HB2	1:B:404:PHE:CZ	2.47	0.50
1:A:101:ASN:HB3	1:A:102:TYR:CD1	2.46	0.50
1:A:12:GLY:O	1:A:13:LEU:C	2.50	0.49
1:A:21:GLY:HA2	1:A:33:ASN:HB2	1.94	0.49
1:A:84:PHE:HZ	1:B:325:TYR:HE1	1.60	0.49
1:B:12:GLY:O	1:B:13:LEU:C	2.51	0.49
1:B:131:SER:HB3	1:B:145:GLY:CA	2.41	0.49
1:B:284:ASP:CG	1:B:285:PRO:CD	2.80	0.49
1:A:402:TYR:HB2	1:A:404:PHE:CZ	2.47	0.49
1:B:167:ALA:O	1:B:169:GLN:N	2.46	0.49
1:B:7:GLU:CD	1:B:7:GLU:N	2.65	0.49
1:A:396:LYS:HA	1:A:405:GLU:HA	1.94	0.49
1:A:133:ALA:HA	1:A:143:GLY:HA2	1.94	0.49
1:A:350:SER:HB2	1:A:365:ASP:N	2.28	0.49
1:B:101:ASN:HB3	1:B:102:TYR:CD1	2.47	0.49
1:B:133:ALA:HA	1:B:143:GLY:HA2	1.95	0.49
1:B:26:ASP:C	1:B:28:GLY:N	2.66	0.49
1:B:167:ALA:O	1:B:170:ILE:N	2.46	0.49
1:B:350:SER:HB2	1:B:365:ASP:N	2.28	0.49
1:A:167:ALA:O	1:A:170:ILE:N	2.46	0.48
1:B:396:LYS:HA	1:B:405:GLU:HA	1.94	0.48
1:A:279:GLY:O	1:A:290:HIS:HA	2.13	0.48
1:A:382:SER:HG	1:A:418:ASN:HB2	1.78	0.48
1:A:194:ASN:OD1	1:A:194:ASN:C	2.52	0.48
1:B:279:GLY:O	1:B:290:HIS:HA	2.13	0.48
1:A:13:LEU:HA	1:A:17:TYR:CE1	2.49	0.48
1:A:167:ALA:O	1:A:169:GLN:N	2.46	0.48
1:A:43:PRO:CG	1:A:421:PHE:HB2	2.43	0.48
1:B:194:ASN:C	1:B:194:ASN:OD1	2.52	0.48
1:B:214:LEU:HD12	1:B:215:TYR:N	2.29	0.48
1:B:394:SER:HB2	1:B:407:GLU:CB	2.43	0.48
1:B:407:GLU:O	1:B:407:GLU:CG	2.62	0.48
1:B:50:VAL:CG1	1:B:414:GLY:HA3	2.43	0.48
1:B:108:PHE:HD1	1:B:118:GLY:O	1.97	0.48
1:B:158:PHE:CD2	1:B:196:LYS:HA	2.49	0.48
1:B:325:TYR:HE2	1:B:349:ASP:HB2	1.79	0.48
1:B:329:LEU:HD23	1:B:329:LEU:C	2.34	0.48
1:A:394:SER:HB2	1:A:407:GLU:CB	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:13:LEU:HA	1:B:17:TYR:CE1	2.48	0.47
1:B:204:GLN:OE1	1:B:204:GLN:HA	2.14	0.47
1:B:232:LYS:HA	1:B:270:ASN:ND2	2.30	0.47
1:A:73:ILE:O	1:A:73:ILE:CG1	2.62	0.47
1:A:173:SER:HB2	1:A:174:PRO:HD2	1.96	0.47
1:B:64:SER:HB3	1:B:165:LEU:CD2	2.44	0.47
1:B:151:ALA:HB2	1:B:233:ILE:CD1	2.44	0.47
1:A:50:VAL:CG1	1:A:414:GLY:HA3	2.42	0.47
1:A:64:SER:HB3	1:A:165:LEU:CD2	2.44	0.47
1:A:15:ARG:HD3	1:A:290:HIS:CD2	2.49	0.47
1:B:26:ASP:OD2	1:B:28:GLY:HA3	2.15	0.47
1:B:173:SER:HB2	1:B:174:PRO:HD2	1.97	0.47
1:B:33:ASN:ND2	1:B:36:LEU:CD1	2.78	0.47
1:A:26:ASP:C	1:A:28:GLY:N	2.66	0.47
1:A:151:ALA:HB2	1:A:233:ILE:CD1	2.44	0.47
1:A:329:LEU:HD23	1:A:329:LEU:C	2.35	0.47
1:B:15:ARG:HD3	1:B:290:HIS:CD2	2.49	0.47
1:A:33:ASN:ND2	1:A:36:LEU:CD1	2.78	0.47
1:A:325:TYR:HE2	1:A:349:ASP:HB2	1.80	0.47
1:B:108:PHE:CD1	1:B:118:GLY:O	2.67	0.47
1:B:214:LEU:HD12	1:B:215:TYR:H	1.80	0.47
1:B:354:ALA:O	1:B:357:ARG:HG3	2.15	0.47
1:A:232:LYS:HA	1:A:270:ASN:ND2	2.30	0.47
1:A:108:PHE:CD1	1:A:118:GLY:HA3	2.50	0.47
1:B:346:ALA:HB3	1:B:368:TRP:HB2	1.97	0.47
1:A:108:PHE:CD1	1:A:118:GLY:O	2.68	0.46
1:A:407:GLU:O	1:A:407:GLU:CG	2.62	0.46
1:B:43:PRO:CG	1:B:421:PHE:HB2	2.43	0.46
1:B:129:ASN:HB2	1:B:147:ASN:ND2	2.30	0.46
1:A:26:ASP:OD2	1:A:28:GLY:HA3	2.15	0.46
1:B:22:ALA:O	1:B:280:TYR:HB3	2.16	0.46
1:A:108:PHE:HD1	1:A:118:GLY:O	1.98	0.46
1:A:204:GLN:OE1	1:A:204:GLN:HA	2.14	0.46
1:A:364:GLN:O	1:A:365:ASP:C	2.54	0.46
1:A:376:ALA:HA	1:A:382:SER:HA	1.97	0.46
1:B:7:GLU:HA	1:B:17:TYR:CZ	2.51	0.46
1:A:22:ALA:O	1:A:280:TYR:HB3	2.15	0.46
1:B:137:ASN:HD22	1:B:137:ASN:C	2.06	0.46
1:B:364:GLN:O	1:B:365:ASP:C	2.53	0.46
1:A:26:ASP:C	1:A:28:GLY:H	2.19	0.46
1:A:136:LEU:O	1:A:137:ASN:CB	2.63	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:158:PHE:CD2	1:A:196:LYS:HA	2.49	0.46
1:A:102:TYR:O	1:A:125:THR:HB	2.15	0.46
1:A:214:LEU:HD12	1:A:215:TYR:N	2.30	0.46
1:A:217:LEU:O	1:A:218:ASP:CB	2.63	0.46
1:A:217:LEU:H	1:A:217:LEU:HD23	1.80	0.46
1:B:351:PRO:HD2	1:B:363:ASP:OD1	2.16	0.46
1:A:354:ALA:HA	1:A:357:ARG:HE	1.81	0.46
1:A:354:ALA:O	1:A:357:ARG:HG3	2.15	0.46
1:B:37:ILE:HB	1:B:133:ALA:HB3	1.98	0.46
1:B:108:PHE:CD1	1:B:118:GLY:HA3	2.51	0.46
1:B:137:ASN:ND2	1:B:137:ASN:C	2.67	0.46
1:B:217:LEU:O	1:B:218:ASP:CB	2.63	0.46
1:A:23:ILE:HG23	1:A:23:ILE:O	2.16	0.46
1:B:102:TYR:O	1:B:125:THR:HB	2.15	0.46
1:A:346:ALA:HB3	1:A:368:TRP:HB2	1.97	0.45
1:B:376:ALA:HA	1:B:382:SER:HA	1.97	0.45
1:B:377:PHE:N	1:B:381:ALA:O	2.49	0.45
1:A:20:GLU:CB	1:A:32:ARG:HD3	2.46	0.45
1:A:117:VAL:HG11	1:A:399:GLU:HG2	1.99	0.45
1:A:129:ASN:HB2	1:A:147:ASN:ND2	2.31	0.45
1:A:351:PRO:HD2	1:A:363:ASP:OD1	2.16	0.45
1:B:20:GLU:CB	1:B:32:ARG:HD3	2.46	0.45
1:B:18:SER:O	1:B:290:HIS:CD2	2.70	0.45
1:B:50:VAL:HG23	1:B:79:VAL:HB	1.98	0.45
1:A:32:ARG:HG3	1:A:32:ARG:NH1	2.31	0.45
1:A:137:ASN:ND2	1:A:137:ASN:C	2.67	0.45
1:B:217:LEU:HD23	1:B:217:LEU:H	1.80	0.45
1:A:37:ILE:HB	1:A:133:ALA:HB3	1.98	0.45
1:A:157:ARG:HA	1:A:157:ARG:HD3	1.74	0.45
1:A:377:PHE:N	1:A:381:ALA:O	2.49	0.45
1:A:50:VAL:HG23	1:A:79:VAL:HB	1.98	0.45
1:B:112:TYR:HD2	1:B:114:GLY:C	2.20	0.45
1:B:167:ALA:C	1:B:169:GLN:N	2.70	0.45
1:A:421:PHE:HE1	1:B:325:TYR:CD1	2.34	0.45
1:A:112:TYR:HD2	1:A:114:GLY:C	2.20	0.45
1:A:127:ASN:ND2	1:A:149:VAL:CG2	2.80	0.45
1:B:159:ALA:O	1:B:160:GLY:C	2.55	0.45
1:B:354:ALA:HA	1:B:357:ARG:HE	1.81	0.45
1:A:18:SER:O	1:A:290:HIS:CD2	2.70	0.45
1:B:26:ASP:C	1:B:28:GLY:H	2.20	0.45
1:B:73:ILE:O	1:B:73:ILE:CG1	2.63	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:82:MET:O	1:B:97:SER:CA	2.63	0.45
1:A:375:TYR:HD2	1:A:376:ALA:N	2.15	0.45
1:A:214:LEU:HD12	1:A:215:TYR:H	1.82	0.44
1:B:219:LYS:HD3	1:B:219:LYS:HA	1.73	0.44
1:B:273:GLU:O	1:B:296:THR:HA	2.17	0.44
1:A:273:GLU:O	1:A:296:THR:HA	2.17	0.44
1:B:161:ASP:OD1	1:B:161:ASP:C	2.56	0.44
1:B:352:VAL:HG11	1:B:357:ARG:CA	2.31	0.44
1:A:161:ASP:C	1:A:161:ASP:OD1	2.55	0.44
1:B:23:ILE:HG23	1:B:23:ILE:O	2.17	0.44
1:B:216:GLU:HB3	1:B:222:ARG:HB3	1.99	0.44
1:A:159:ALA:O	1:A:160:GLY:C	2.56	0.44
1:A:324:ALA:HB2	1:A:351:PRO:HA	1.99	0.44
1:B:117:VAL:HG11	1:B:399:GLU:HG2	1.99	0.44
1:B:134:TYR:CE2	1:B:135:ARG:O	2.70	0.44
1:B:127:ASN:ND2	1:B:149:VAL:CG2	2.80	0.44
1:A:72:ASN:C	1:A:72:ASN:HD22	2.21	0.44
1:A:219:LYS:HA	1:A:219:LYS:HD3	1.72	0.44
1:A:216:GLU:HB3	1:A:222:ARG:HB3	1.99	0.44
1:B:72:ASN:C	1:B:72:ASN:HD22	2.21	0.44
1:B:324:ALA:HB2	1:B:351:PRO:HA	1.99	0.44
1:A:73:ILE:HD11	1:A:106:THR:HB	2.00	0.44
1:A:82:MET:O	1:A:97:SER:CA	2.63	0.44
1:B:284:ASP:OD2	1:B:286:GLN:N	2.32	0.44
1:B:325:TYR:CE2	1:B:349:ASP:HB2	2.52	0.44
1:A:134:TYR:CE2	1:A:135:ARG:O	2.71	0.44
1:A:167:ALA:C	1:A:169:GLN:N	2.70	0.44
1:B:58:ILE:HB	1:B:70:ALA:HB3	1.99	0.44
1:B:81:ASN:C	1:B:81:ASN:ND2	2.69	0.44
1:B:356:ASN:HD22	1:B:356:ASN:HA	1.62	0.44
1:A:81:ASN:C	1:A:81:ASN:ND2	2.69	0.43
1:A:382:SER:O	1:A:418:ASN:N	2.48	0.43
1:B:375:TYR:HD2	1:B:376:ALA:N	2.15	0.43
1:A:13:LEU:HD22	1:A:50:VAL:HG11	2.00	0.43
1:A:121:THR:OG1	1:A:157:ARG:NH2	2.49	0.43
1:A:33:ASN:O	1:A:34:PRO:C	2.57	0.43
1:A:284:ASP:OD1	1:A:285:PRO:HD2	2.18	0.43
1:B:127:ASN:ND2	1:B:149:VAL:HG22	2.33	0.43
1:B:325:TYR:C	1:B:325:TYR:CD2	2.91	0.43
1:A:325:TYR:CE2	1:A:349:ASP:HB2	2.53	0.43
1:A:13:LEU:HD22	1:A:50:VAL:CG1	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:ILE:HB	1:A:70:ALA:HB3	2.00	0.43
1:A:127:ASN:ND2	1:A:149:VAL:HG22	2.33	0.43
1:B:117:VAL:HG23	1:B:117:VAL:O	2.18	0.43
1:B:382:SER:O	1:B:418:ASN:N	2.47	0.43
1:B:284:ASP:OD1	1:B:285:PRO:HD2	2.18	0.43
1:A:399:GLU:O	1:A:400:GLY:C	2.56	0.43
1:B:33:ASN:O	1:B:34:PRO:C	2.57	0.43
1:B:350:SER:HA	1:B:351:PRO:HD3	1.80	0.43
1:A:50:VAL:HG12	1:A:414:GLY:CA	2.47	0.43
1:A:117:VAL:HG23	1:A:117:VAL:O	2.18	0.43
1:A:325:TYR:C	1:A:325:TYR:CD2	2.91	0.43
1:A:378:ASN:O	1:A:378:ASN:CG	2.57	0.43
1:B:73:ILE:HD11	1:B:106:THR:HB	1.99	0.42
1:B:399:GLU:O	1:B:400:GLY:C	2.56	0.42
1:B:13:LEU:HD22	1:B:50:VAL:CG1	2.50	0.42
1:A:364:GLN:HE21	1:A:364:GLN:HB2	1.59	0.42
1:B:8:PHE:O	1:B:83:HIS:HE1	2.02	0.42
1:B:81:ASN:HD21	1:B:416:ASN:ND2	2.17	0.42
1:B:192:ASP:O	1:B:194:ASN:N	2.50	0.42
1:B:393:GLN:NE2	1:B:394:SER:N	2.67	0.42
1:A:6:ASN:O	1:A:8:PHE:N	2.52	0.42
1:A:179:GLN:HA	1:A:182:GLN:HB3	2.02	0.42
1:B:121:THR:OG1	1:B:157:ARG:NH2	2.50	0.42
1:B:13:LEU:HD22	1:B:50:VAL:HG11	2.01	0.42
1:B:157:ARG:HA	1:B:157:ARG:HD3	1.74	0.42
1:B:355:GLN:NE2	1:B:356:ASN:N	2.67	0.42
1:A:356:ASN:HD22	1:A:356:ASN:HA	1.62	0.42
1:A:421:PHE:CE1	1:B:325:TYR:CD1	3.07	0.42
1:B:32:ARG:HG3	1:B:32:ARG:NH1	2.31	0.42
1:B:378:ASN:O	1:B:378:ASN:CG	2.57	0.42
1:A:81:ASN:HD21	1:A:416:ASN:ND2	2.18	0.42
1:B:137:ASN:C	1:B:139:ALA:N	2.73	0.42
1:B:15:ARG:HG3	1:B:18:SER:H	1.84	0.42
1:B:179:GLN:HA	1:B:182:GLN:HB3	2.02	0.42
1:A:15:ARG:HG3	1:A:18:SER:H	1.83	0.42
1:B:7:GLU:N	1:B:7:GLU:OE2	2.36	0.42
1:A:137:ASN:C	1:A:139:ALA:N	2.72	0.41
1:B:232:LYS:HG2	1:B:270:ASN:ND2	2.32	0.41
1:A:393:GLN:NE2	1:A:394:SER:N	2.68	0.41
1:B:137:ASN:ND2	1:B:139:ALA:H	2.17	0.41
1:A:16:ALA:O	1:A:17:TYR:C	2.59	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:192:ASP:O	1:A:194:ASN:N	2.50	0.41
1:A:232:LYS:HG2	1:A:270:ASN:ND2	2.33	0.41
1:B:21:GLY:HA2	1:B:33:ASN:CB	2.50	0.41
1:B:56:VAL:C	1:B:57:ASN:HD22	2.23	0.41
1:A:34:PRO:HB2	1:A:143:GLY:C	2.41	0.41
1:B:325:TYR:CE2	1:B:349:ASP:CB	3.03	0.41
1:A:8:PHE:O	1:A:83:HIS:HE1	2.03	0.41
1:A:355:GLN:NE2	1:A:356:ASN:N	2.68	0.41
1:B:126:MET:O	1:B:149:VAL:HA	2.21	0.41
1:A:56:VAL:C	1:A:57:ASN:HD22	2.24	0.41
1:A:182:GLN:HG3	1:A:183:ALA:N	2.36	0.41
1:A:325:TYR:CE2	1:A:349:ASP:CB	3.04	0.41
1:B:73:ILE:CD1	1:B:108:PHE:CE2	2.97	0.41
1:A:21:GLY:HA2	1:A:33:ASN:CB	2.51	0.41
1:A:13:LEU:HA	1:A:13:LEU:HD12	1.94	0.41
1:B:11:SER:OG	1:B:15:ARG:NH2	2.52	0.41
1:B:16:ALA:O	1:B:17:TYR:C	2.58	0.41
1:B:34:PRO:HB2	1:B:143:GLY:C	2.41	0.41
1:B:134:TYR:CD2	1:B:135:ARG:N	2.89	0.41
1:B:151:ALA:CB	1:B:233:ILE:HD13	2.51	0.41
1:A:170:ILE:HD13	1:A:184:LEU:HB3	2.03	0.40
1:B:50:VAL:HG12	1:B:414:GLY:CA	2.47	0.40
1:A:11:SER:OG	1:A:15:ARG:NH2	2.53	0.40
1:A:126:MET:O	1:A:149:VAL:HA	2.20	0.40
1:B:8:PHE:CD1	1:B:9:SER:HB2	2.56	0.40
1:B:26:ASP:OD2	1:B:28:GLY:N	2.54	0.40
1:B:170:ILE:HD13	1:B:184:LEU:HB3	2.03	0.40
1:A:73:ILE:CD1	1:A:108:PHE:CE2	2.98	0.40
1:A:134:TYR:CD2	1:A:135:ARG:N	2.89	0.40
1:A:357:ARG:HB2	1:A:397:ILE:HG23	2.04	0.40
1:B:7:GLU:HB2	1:B:50:VAL:HG21	1.99	0.40
1:B:197:THR:O	1:B:198:ALA:HB2	2.21	0.40
1:B:348:ASP:N	1:B:348:ASP:OD2	2.53	0.40
1:A:182:GLN:CG	1:A:183:ALA:N	2.85	0.40
1:B:182:GLN:CG	1:B:183:ALA:N	2.85	0.40
1:B:209:TRP:CD1	1:B:209:TRP:N	2.88	0.40
1:A:230:GLU:OE1	1:A:232:LYS:HE3	2.22	0.40
1:B:230:GLU:OE1	1:B:232:LYS:HE3	2.21	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	357/424 (84%)	295 (83%)	41 (12%)	21 (6%)	1	11
1	B	357/424 (84%)	297 (83%)	39 (11%)	21 (6%)	1	11
All	All	714/848 (84%)	592 (83%)	80 (11%)	42 (6%)	1	11

All (42) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	7	GLU
1	A	194	ASN
1	A	354	ALA
1	A	379	LYS
1	B	194	ASN
1	B	354	ALA
1	B	379	LYS
1	A	27	ALA
1	B	27	ALA
1	A	137	ASN
1	A	187	THR
1	A	193	SER
1	A	218	ASP
1	A	348	ASP
1	B	137	ASN
1	B	187	THR
1	B	193	SER
1	B	218	ASP
1	B	348	ASP
1	A	20	GLU
1	A	118	GLY
1	A	168	GLY
1	A	198	ALA
1	B	20	GLU
1	B	118	GLY

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Mol	Chain	Res	Type
1	B	168	GLY
1	B	198	ALA
1	A	89	ASN
1	A	167	ALA
1	A	229	SER
1	B	6	ASN
1	B	89	ASN
1	B	167	ALA
1	B	229	SER
1	A	172	GLN
1	A	210	ASN
1	B	172	GLN
1	B	210	ASN
1	A	213	ILE
1	A	362	PRO
1	B	213	ILE
1	B	362	PRO

### 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	284/334 (85%)	246 (87%)	38 (13%)	4	15
1	B	284/334 (85%)	248 (87%)	36 (13%)	4	16
All	All	568/668 (85%)	494 (87%)	74 (13%)	4	16

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	7	GLU
1	A	50	VAL
1	A	72	ASN
1	A	73	ILE
1	A	80	PRO

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	81	ASN
1	A	82	MET
1	A	107	GLU
1	A	109	ASN
1	A	122	ASP
1	A	123	LEU
1	A	136	LEU
1	A	137	ASN
1	A	142	PHE
1	A	144	LEU
1	A	180	GLN
1	A	217	LEU
1	A	226	THR
1	A	267	LEU
1	A	275	TRP
1	A	297	SER
1	A	325	TYR
1	A	343	THR
1	A	345	ILE
1	A	355	GLN
1	A	356	ASN
1	A	363	ASP
1	A	364	GLN
1	A	369	LEU
1	A	373	THR
1	A	375	TYR
1	A	393	GLN
1	A	396	LYS
1	A	404	PHE
1	A	407	GLU
1	A	415	THR
1	A	416	ASN
1	B	50	VAL
1	B	72	ASN
1	B	73	ILE
1	B	80	PRO
1	B	81	ASN
1	B	82	MET
1	B	107	GLU
1	B	109	ASN
1	B	122	ASP
1	B	123	LEU

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Mol	Chain	Res	Type
1	B	136	LEU
1	B	137	ASN
1	B	142	PHE
1	B	144	LEU
1	B	180	GLN
1	B	217	LEU
1	B	226	THR
1	B	267	LEU
1	B	275	TRP
1	B	297	SER
1	B	325	TYR
1	B	343	THR
1	B	345	ILE
1	B	355	GLN
1	B	356	ASN
1	B	363	ASP
1	B	364	GLN
1	B	369	LEU
1	B	373	THR
1	B	375	TYR
1	B	393	GLN
1	B	396	LYS
1	B	404	PHE
1	B	407	GLU
1	B	415	THR
1	B	416	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (41) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	57	ASN
1	A	72	ASN
1	A	83	HIS
1	A	89	ASN
1	A	109	ASN
1	A	137	ASN
1	A	164	GLN
1	A	169	GLN
1	A	182	GLN
1	A	270	ASN
1	A	281	ASN

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Mol	Chain	Res	Type
1	A	290	HIS
1	A	355	GLN
1	A	356	ASN
1	A	364	GLN
1	A	391	HIS
1	A	393	GLN
1	A	416	ASN
1	A	418	ASN
1	B	6	ASN
1	B	57	ASN
1	B	72	ASN
1	B	83	HIS
1	B	89	ASN
1	B	101	ASN
1	B	109	ASN
1	B	137	ASN
1	B	164	GLN
1	B	169	GLN
1	B	182	GLN
1	B	270	ASN
1	B	281	ASN
1	B	290	HIS
1	B	355	GLN
1	B	356	ASN
1	B	364	GLN
1	B	391	HIS
1	B	393	GLN
1	B	416	ASN
1	B	418	ASN

### 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 monosaccharides 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	363/424 (85%)	0.03	7 (1%) 66 65	24, 49, 83, 108	0
1	B	363/424 (85%)	0.02	6 (1%) 70 68	20, 50, 82, 106	0
All	All	726/848 (85%)	0.03	13 (1%) 68 67	20, 49, 83, 108	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	123	LEU	3.6
1	A	123	LEU	3.3
1	A	298	TRP	3.2
1	A	337	ASP	2.8
1	A	232	LYS	2.8
1	B	337	ASP	2.7
1	A	362	PRO	2.6
1	B	232	LYS	2.4
1	B	359	ILE	2.2
1	A	359	ILE	2.1
1	A	233	ILE	2.1
1	B	102	TYR	2.1
1	B	233	ILE	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 monosaccharides 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.