Full wwPDB X-ray Structure Validation Report (i)

Jun 22, 2024 – 11:13 AM EDT

PDB ID	:	6JZQ
Title	:	The crystal structure of acyl-acyl carrier protein (acyl-ACP) reductase (AAR)
Authors	:	Zhang, H.M.; Li, M.; Gao, Y.
Deposited on	:	2019-05-03
Resolution	:	2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

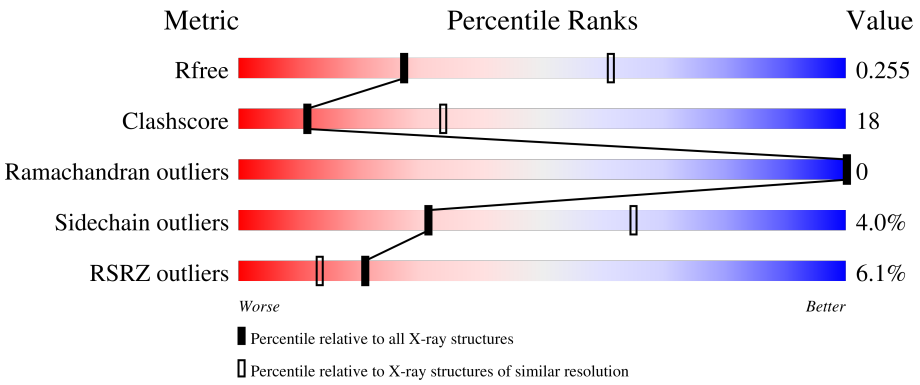
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	347	% 79% 18% • •
1	B	347	6% 76% 22% • •
1	C	347	9% 68% 29% • •
1	D	347	4% 68% 27% • •
1	E	347	11% 65% 33% • •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 12983 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 acyl-[acyl-carrier-protein] reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	340	2610	1656	448	490	16	0	0	0
1	B	345	2629	1673	444	496	16	0	0	0
1	C	344	2609	1658	452	483	16	0	0	0
1	D	339	2525	1611	426	472	16	0	0	0
1	E	345	2562	1631	432	484	15	0	1	0

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	342	GLU	-	expression tag	UNP Q54765
A	343	ASN	-	expression tag	UNP Q54765
A	344	LEU	-	expression tag	UNP Q54765
A	345	TYR	-	expression tag	UNP Q54765
A	346	PHE	-	expression tag	UNP Q54765
A	347	GLN	-	expression tag	UNP Q54765
B	342	GLU	-	expression tag	UNP Q54765
B	343	ASN	-	expression tag	UNP Q54765
B	344	LEU	-	expression tag	UNP Q54765
B	345	TYR	-	expression tag	UNP Q54765
B	346	PHE	-	expression tag	UNP Q54765
B	347	GLN	-	expression tag	UNP Q54765
C	342	GLU	-	expression tag	UNP Q54765
C	343	ASN	-	expression tag	UNP Q54765
C	344	LEU	-	expression tag	UNP Q54765
C	345	TYR	-	expression tag	UNP Q54765
C	346	PHE	-	expression tag	UNP Q54765
C	347	GLN	-	expression tag	UNP Q54765
D	342	GLU	-	expression tag	UNP Q54765

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Chain	Residue	Modelled	Actual	Comment	Reference
D	343	ASN	-	expression tag	UNP Q54765
D	344	LEU	-	expression tag	UNP Q54765
D	345	TYR	-	expression tag	UNP Q54765
D	346	PHE	-	expression tag	UNP Q54765
D	347	GLN	-	expression tag	UNP Q54765
E	342	GLU	-	expression tag	UNP Q54765
E	343	ASN	-	expression tag	UNP Q54765
E	344	LEU	-	expression tag	UNP Q54765
E	345	TYR	-	expression tag	UNP Q54765
E	346	PHE	-	expression tag	UNP Q54765
E	347	GLN	-	expression tag	UNP Q54765

- Molecule 2 is water.

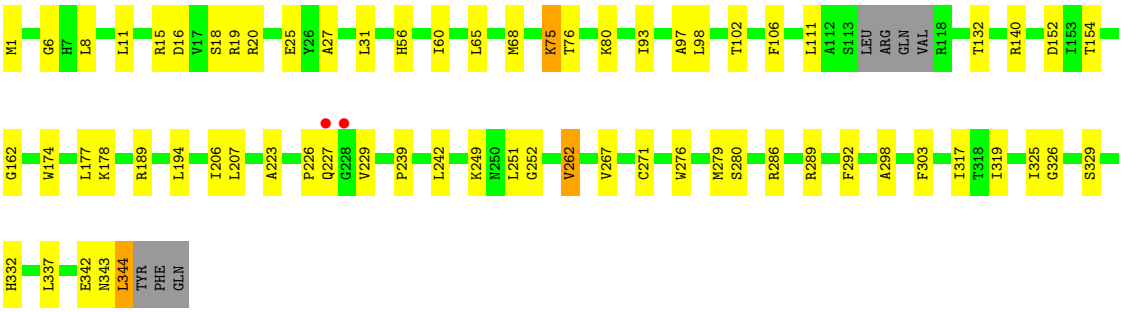
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	21	21	21	0	0
			Total	O		
2	B	12	12	12	0	0
			Total	O		
2	C	4	4	4	0	0
			Total	O		
2	D	8	8	8	0	0
			Total	O		
2	E	3	3	3	0	0

3 Residue-property plots (i)

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.

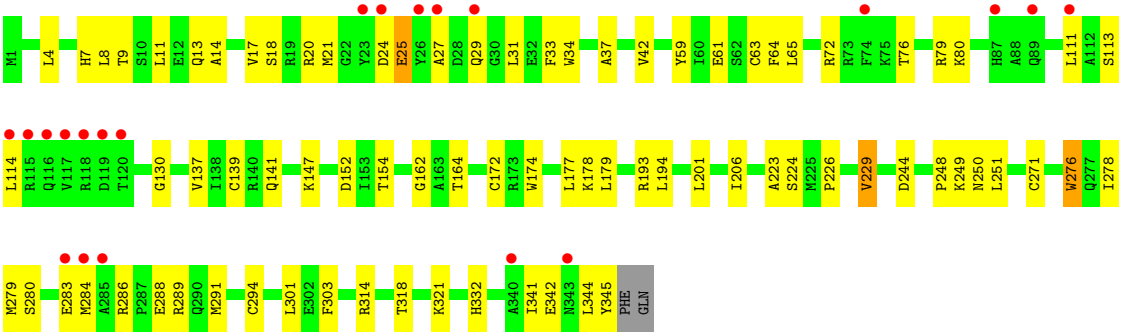
• Molecule 1: Long-chain acyl-[acyl-carrier-protein] reductase

Chain A:



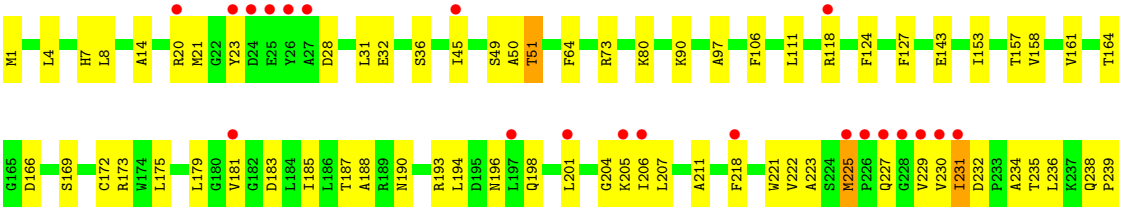
• Molecule 1: Long-chain acyl-[acyl-carrier-protein] reductase

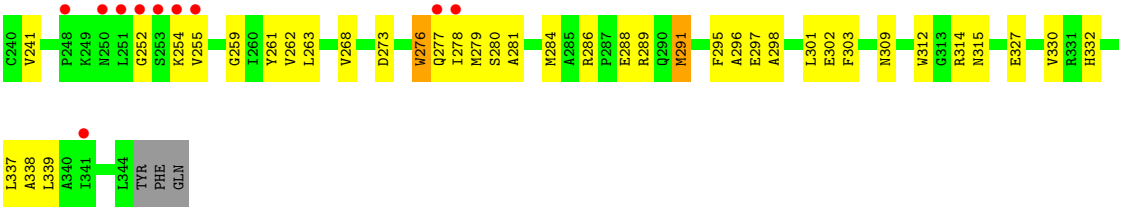
Chain B:



• Molecule 1: Long-chain acyl-[acyl-carrier-protein] reductase

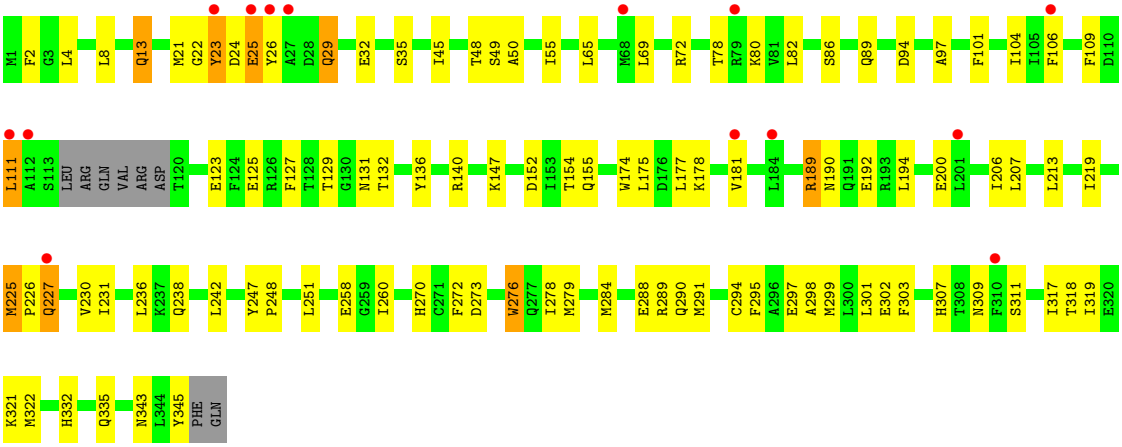
Chain C:





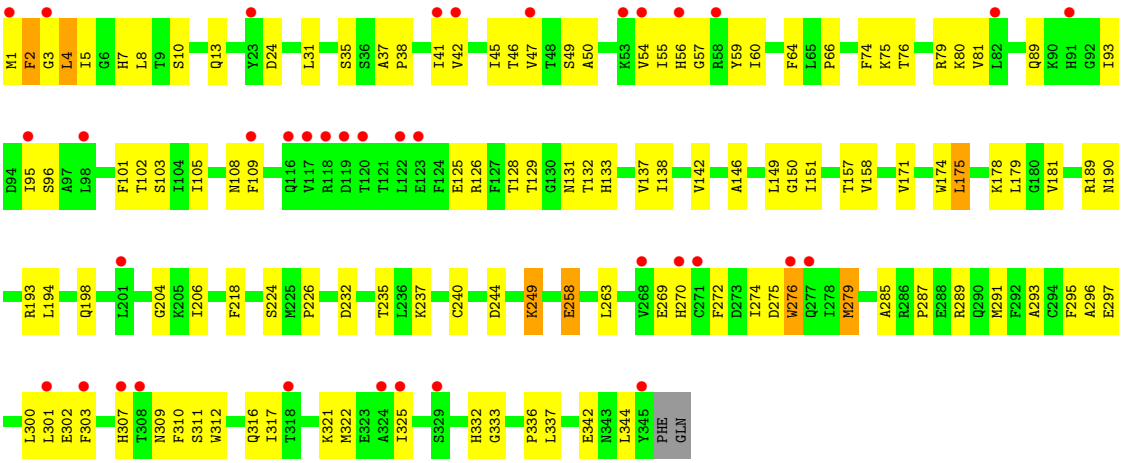
• Molecule 1: Long-chain acyl-[acyl-carrier-protein] reductase

Chain D:



• Molecule 1: Long-chain acyl-[acyl-carrier-protein] reductase

Chain E:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	92.17Å 201.44Å 124.58Å 90.00° 105.90° 90.00°	Depositor
Resolution (Å)	46.11 – 2.80 46.11 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.0 (46.11-2.80) 99.0 (46.11-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	1.87 (at 2.81Å)	Xtriage
Refinement program	PHENIX (1.15.2_3472: ???)	Depositor
R, R_{free}	0.218 , 0.254 0.222 , 0.255	Depositor DCC
R_{free} test set	1995 reflections (3.75%)	wwPDB-VP
Wilson B-factor (Å^2)	63.8	Xtriage
Anisotropy	0.600	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 60.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	12983	wwPDB-VP
Average B, all atoms (Å^2)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.71% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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	0.65	0/2661	0.76	0/3607
1	B	0.59	0/2682	0.70	0/3642
1	C	0.60	0/2660	0.73	0/3609
1	D	0.52	0/2576	0.68	0/3504
1	E	0.53	0/2615	0.69	0/3561
All	All	0.58	0/13194	0.71	0/17923

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	2610	0	2574	45	0
1	B	2629	0	2575	78	0
1	C	2609	0	2571	99	0
1	D	2525	0	2428	87	0
1	E	2562	0	2437	159	0
2	A	21	0	0	2	0
2	B	12	0	0	0	0
2	C	4	0	0	0	0
2	D	8	0	0	2	0
2	E	3	0	0	0	0
All	All	12983	0	12585	454	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 18.

All (454) 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:E:300:LEU:HD13	1:E:325:ILE:CG1	1.45	1.42
1:E:95:ILE:CG2	1:E:301:LEU:HD23	1.49	1.40
1:E:95:ILE:CG2	1:E:301:LEU:CD2	2.19	1.21
1:E:300:LEU:CD1	1:E:325:ILE:CG1	2.25	1.15
1:E:300:LEU:CD1	1:E:325:ILE:HG12	1.75	1.14
1:E:95:ILE:HG23	1:E:301:LEU:CD2	1.78	1.13
1:E:95:ILE:HG23	1:E:301:LEU:HD23	1.19	1.12
1:E:2:PHE:CE1	1:E:57:GLY:HA3	1.88	1.07
1:E:95:ILE:HG21	1:E:301:LEU:HD23	1.33	1.04
1:E:95:ILE:HG21	1:E:301:LEU:CD2	1.85	1.02
1:C:231:ILE:HD12	1:C:254:LYS:HD2	1.42	1.01
1:B:29:GLN:HB2	1:B:33:PHE:HD2	1.22	0.97
1:C:20:ARG:HH12	1:C:277:GLN:HB3	1.30	0.96
1:E:300:LEU:CD1	1:E:325:ILE:HG13	1.96	0.95
1:E:276:TRP:HH2	1:E:291:MET:SD	1.91	0.94
1:E:4:LEU:HD13	1:E:295:PHE:CD1	2.03	0.94
1:C:225:MET:SD	1:C:225:MET:N	2.39	0.94
1:E:95:ILE:HG21	1:E:301:LEU:HB3	1.48	0.94
1:E:300:LEU:HD13	1:E:325:ILE:HG12	0.95	0.93
1:E:4:LEU:HD13	1:E:295:PHE:CG	2.04	0.93
1:C:194:LEU:HB3	1:C:206:ILE:HD11	1.49	0.92
1:C:231:ILE:HD12	1:C:254:LYS:CD	2.00	0.92
1:B:29:GLN:NE2	1:B:34:TRP:CZ2	2.41	0.89
1:E:190:ASN:HD22	1:E:193:ARG:HG3	1.38	0.89
1:E:95:ILE:CD1	1:E:307:HIS:HD2	1.85	0.87
1:E:2:PHE:HD1	1:E:2:PHE:H	1.21	0.87
1:E:2:PHE:CD1	1:E:57:GLY:HA3	2.09	0.87
1:E:300:LEU:HD13	1:E:325:ILE:HG13	1.49	0.85
1:E:158:VAL:HG22	1:E:218:PHE:HB2	1.60	0.82
1:B:113:SER:C	1:B:114:LEU:HD12	2.00	0.81
1:E:310:PHE:HE2	1:E:325:ILE:HB	1.45	0.81
1:B:29:GLN:HB2	1:B:33:PHE:CD2	2.14	0.81
1:E:300:LEU:HD23	1:E:300:LEU:O	1.81	0.80
1:D:190:ASN:OD1	1:D:192:GLU:HG2	1.82	0.80
1:E:95:ILE:HG21	1:E:301:LEU:CB	2.12	0.79
1:A:252:GLY:HA2	1:A:262:VAL:HG11	1.65	0.78
1:C:232:ASP:CB	1:C:235:THR:OG1	2.33	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:29:GLN:CB	1:B:33:PHE:HD2	1.98	0.77
1:C:157:THR:HG22	1:C:183:ASP:HB2	1.67	0.77
1:C:158:VAL:HG22	1:C:218:PHE:HB2	1.68	0.76
1:E:276:TRP:CH2	1:E:291:MET:SD	2.75	0.76
1:E:55:ILE:HG21	1:E:302:GLU:HG2	1.68	0.76
1:D:89:GLN:HE22	1:D:123:GLU:H	1.33	0.76
1:B:8:LEU:HB2	1:B:65:LEU:HD22	1.68	0.75
1:C:45:ILE:HG22	1:C:276:TRP:HB3	1.71	0.73
1:C:286:ARG:HD3	1:C:289:ARG:HG3	1.68	0.73
1:E:2:PHE:CE1	1:E:57:GLY:CA	2.70	0.73
1:E:95:ILE:CG2	1:E:301:LEU:HD22	2.18	0.73
1:D:106:PHE:HA	1:D:111:LEU:HD22	1.69	0.73
1:E:190:ASN:ND2	1:E:193:ARG:HG3	2.03	0.72
1:A:286:ARG:NH2	1:A:289:ARG:HH21	1.88	0.72
1:C:221:TRP:NE1	1:C:223:ALA:O	2.23	0.72
1:D:29:GLN:HE21	1:D:29:GLN:N	1.87	0.72
1:C:205:LYS:HD3	1:C:207:LEU:HD11	1.73	0.70
1:E:178:LYS:HG3	1:E:179:LEU:HD12	1.74	0.70
1:E:95:ILE:HD11	1:E:307:HIS:HA	1.73	0.70
1:D:219:ILE:HB	1:D:242:LEU:HD23	1.73	0.70
1:D:318:THR:HG22	1:D:321:LYS:HD3	1.74	0.69
1:E:7:HIS:H	1:E:279:MET:CE	2.05	0.69
1:E:125:GLU:HG3	1:E:309:ASN:ND2	2.07	0.69
1:D:2:PHE:CZ	1:D:45:ILE:HD11	2.28	0.69
1:A:177:LEU:HD23	1:A:319:ILE:HG13	1.74	0.69
1:E:74:PHE:HB3	1:E:109:PHE:HE2	1.59	0.68
1:C:20:ARG:NH1	1:C:277:GLN:HB3	2.05	0.68
1:A:132:THR:HG23	1:A:317:ILE:HG23	1.76	0.67
1:E:291:MET:HE3	1:E:296:ALA:HB2	1.75	0.67
1:E:133:HIS:ND1	1:E:293:ALA:HB1	2.09	0.67
1:E:95:ILE:HG13	1:E:307:HIS:CD2	2.29	0.67
1:E:300:LEU:HD13	1:E:325:ILE:CD1	2.23	0.67
1:E:132:THR:HB	1:E:317:ILE:HD12	1.77	0.67
1:C:276:TRP:CZ2	1:C:278:ILE:HD13	2.29	0.66
1:D:276:TRP:HH2	1:D:291:MET:HE3	1.58	0.66
1:B:8:LEU:HD22	1:B:65:LEU:HD21	1.75	0.66
1:C:231:ILE:CD1	1:C:254:LYS:HD2	2.21	0.66
1:C:231:ILE:HD13	1:C:231:ILE:N	2.10	0.66
1:B:24:ASP:HB3	1:B:27:ALA:HB2	1.78	0.66
1:E:232:ASP:O	1:E:235:THR:HG22	1.95	0.66
1:B:114:LEU:HD12	1:B:114:LEU:N	2.10	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:224:SER:HB3	1:E:344:LEU:HB2	1.75	0.66
1:E:95:ILE:CG1	1:E:307:HIS:HA	2.26	0.65
1:B:72:ARG:HH12	1:E:258:GLU:HG3	1.61	0.65
1:B:278:ILE:HG23	1:B:280:SER:H	1.61	0.65
1:E:175:LEU:O	1:E:181:VAL:HG23	1.97	0.65
1:E:133:HIS:O	1:E:137:VAL:HG23	1.96	0.65
1:E:342:GLU:OE1	1:E:342:GLU:N	2.23	0.65
1:E:5:ILE:HG23	1:E:60:ILE:HD11	1.77	0.65
1:C:231:ILE:CD1	1:C:254:LYS:CD	2.75	0.64
1:E:297:GLU:O	1:E:301:LEU:HB2	1.96	0.64
1:E:303:PHE:HB3	1:E:332:HIS:NE2	2.11	0.64
1:D:49:SER:O	1:D:272:PHE:HE2	1.81	0.64
1:B:130:GLY:HA3	1:B:294:CYS:HB2	1.80	0.64
1:E:35:SER:O	1:E:80:LYS:NZ	2.31	0.64
1:D:194:LEU:HB3	1:D:206:ILE:HD13	1.80	0.64
1:E:270:HIS:ND1	1:E:333:GLY:O	2.29	0.63
1:B:20:ARG:HH21	1:B:42:VAL:HG11	1.63	0.63
1:B:194:LEU:HB3	1:B:206:ILE:HD13	1.80	0.63
1:D:132:THR:HG23	1:D:317:ILE:HG23	1.81	0.63
1:B:152:ASP:CG	1:C:32:GLU:HG3	2.19	0.63
1:C:166:ASP:HB3	1:C:315:ASN:H	1.62	0.63
1:A:75:LYS:O	1:A:75:LYS:HD3	1.99	0.62
1:D:21:MET:HG2	1:D:22:GLY:N	2.15	0.62
1:E:128:THR:HG22	1:E:311:SER:H	1.64	0.62
1:E:311:SER:HA	1:E:317:ILE:HD11	1.81	0.62
1:B:8:LEU:HD12	1:B:61:GLU:OE2	2.00	0.62
1:D:189:ARG:HA	1:D:189:ARG:NE	2.14	0.62
1:D:276:TRP:CH2	1:D:291:MET:HE3	2.34	0.62
1:D:78:THR:O	1:D:82:LEU:HD12	2.00	0.62
1:C:207:LEU:HD23	1:C:211:ALA:HB1	1.80	0.61
1:A:19:ARG:HB2	1:A:25:GLU:HG2	1.81	0.61
1:D:152:ASP:OD1	1:D:154:THR:HB	2.00	0.61
1:A:342:GLU:HG3	1:A:342:GLU:O	2.01	0.61
1:E:95:ILE:HG23	1:E:301:LEU:HD22	1.79	0.61
1:E:95:ILE:CD1	1:E:307:HIS:CD2	2.77	0.61
1:C:288:GLU:HG2	1:C:289:ARG:HG2	1.82	0.60
1:D:270:HIS:O	1:D:289:ARG:HG2	2.01	0.60
1:E:55:ILE:HD13	1:E:303:PHE:CE1	2.35	0.60
1:A:97:ALA:HB2	1:A:298:ALA:HB2	1.83	0.60
1:A:194:LEU:HB3	1:A:206:ILE:HD13	1.83	0.60
1:E:1:MET:C	1:E:302:GLU:OE1	2.39	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:230:VAL:O	1:D:231:ILE:HG22	2.01	0.60
1:E:7:HIS:H	1:E:279:MET:HE1	1.66	0.60
1:C:222:VAL:HG22	1:C:222:VAL:O	2.00	0.60
1:C:1:MET:N	1:C:302:GLU:OE2	2.21	0.60
1:D:175:LEU:O	1:D:181:VAL:HG23	2.02	0.59
1:D:129:THR:HG22	1:D:311:SER:HB3	1.84	0.59
1:C:231:ILE:HD13	1:C:231:ILE:H	1.66	0.59
1:A:75:LYS:HD3	1:A:75:LYS:C	2.22	0.59
1:D:35:SER:HA	1:D:80:LYS:HZ1	1.68	0.59
1:D:189:ARG:NH2	2:D:401:HOH:O	2.36	0.59
1:E:95:ILE:CD1	1:E:307:HIS:HA	2.32	0.59
1:C:45:ILE:CG2	1:C:276:TRP:HB3	2.33	0.59
1:A:8:LEU:HB2	1:A:65:LEU:HD13	1.84	0.58
1:A:162:GLY:HA3	1:A:223:ALA:HB2	1.85	0.58
1:A:280:SER:HG	1:A:292:PHE:HZ	1.50	0.58
1:E:95:ILE:HD11	1:E:307:HIS:HD2	1.67	0.58
1:B:226:PRO:HG2	1:B:229:VAL:CG2	2.34	0.58
1:D:273:ASP:OD1	1:D:289:ARG:NE	2.36	0.58
1:E:132:THR:HG22	1:E:311:SER:HB3	1.85	0.58
1:E:146:ALA:HA	1:E:151:ILE:HD11	1.86	0.58
1:E:276:TRP:CH2	1:E:291:MET:HB3	2.39	0.58
1:B:29:GLN:CD	1:B:34:TRP:CZ2	2.77	0.58
1:E:174:TRP:HB2	1:E:322:MET:HE1	1.86	0.58
1:B:152:ASP:OD1	1:B:154:THR:HB	2.04	0.57
1:C:254:LYS:O	1:C:254:LYS:HD3	2.04	0.57
1:E:300:LEU:HD23	1:E:300:LEU:C	2.24	0.57
1:A:227:GLN:NE2	1:D:343:ASN:HB3	2.19	0.57
1:C:238:GLN:O	1:C:259:GLY:HA3	2.05	0.57
1:C:252:GLY:HA2	1:C:262:VAL:HG11	1.87	0.57
1:C:8:LEU:HD22	1:C:14:ALA:HA	1.87	0.57
1:D:129:THR:HG22	1:D:311:SER:CB	2.34	0.57
1:B:31:LEU:HD13	1:E:150:GLY:HA2	1.87	0.57
1:C:190:ASN:O	1:C:194:LEU:HD12	2.05	0.57
1:C:227:GLN:O	1:C:229:VAL:HG23	2.05	0.57
1:C:303:PHE:HB3	1:C:332:HIS:CE1	2.40	0.56
1:C:314:ARG:HG2	1:C:315:ASN:ND2	2.20	0.56
1:C:273:ASP:OD1	1:C:289:ARG:HD3	2.04	0.56
1:E:1:MET:N	1:E:302:GLU:OE1	2.34	0.56
1:E:244:ASP:OD2	1:E:249:LYS:HA	2.06	0.56
1:E:47:VAL:HG13	1:E:274:ILE:HG21	1.87	0.56
1:E:54:VAL:C	1:E:55:ILE:HD12	2.26	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:95:ILE:CG1	1:E:307:HIS:HD2	2.18	0.56
1:C:97:ALA:HB2	1:C:298:ALA:HB2	1.87	0.56
1:E:7:HIS:ND1	1:E:8:LEU:O	2.38	0.56
1:C:337:LEU:HD23	1:C:338:ALA:N	2.21	0.55
1:E:125:GLU:HG3	1:E:309:ASN:HD22	1.70	0.55
1:D:55:ILE:HD13	1:D:302:GLU:HG3	1.87	0.55
1:A:60:ILE:HD11	1:A:93:ILE:HD12	1.88	0.55
1:A:271:CYS:O	1:A:289:ARG:NH1	2.29	0.55
1:B:37:ALA:O	1:B:80:LYS:NZ	2.38	0.55
1:B:137:VAL:O	1:B:141:GLN:HG3	2.07	0.55
1:C:185:ILE:HG23	1:C:207:LEU:HD13	1.87	0.55
1:E:151:ILE:HD11	1:E:218:PHE:HZ	1.71	0.55
1:D:2:PHE:HZ	1:D:45:ILE:HD11	1.71	0.55
1:C:231:ILE:CD1	1:C:231:ILE:H	2.19	0.54
1:C:45:ILE:HG21	1:C:276:TRP:HE3	1.71	0.54
1:E:95:ILE:HG21	1:E:301:LEU:CG	2.36	0.54
1:A:6:GLY:HA3	1:A:279:MET:HE2	1.88	0.54
1:B:177:LEU:HB2	1:C:118:ARG:HD3	1.88	0.54
1:A:286:ARG:NH2	1:A:289:ARG:NH2	2.55	0.54
1:D:35:SER:HA	1:D:80:LYS:NZ	2.22	0.54
1:E:76:THR:O	1:E:80:LYS:HG2	2.08	0.54
1:C:263:LEU:HD13	1:C:339:LEU:HG	1.89	0.54
1:D:32:GLU:CD	1:D:32:GLU:H	2.11	0.54
1:B:29:GLN:CB	1:B:33:PHE:CD2	2.84	0.54
1:A:242:LEU:HD23	1:A:251:LEU:HD23	1.89	0.53
1:B:7:HIS:H	1:B:279:MET:HE2	1.73	0.53
1:B:31:LEU:HD12	1:B:31:LEU:H	1.73	0.53
1:D:55:ILE:HD13	1:D:302:GLU:CG	2.39	0.53
1:D:213:LEU:HD11	1:D:231:ILE:HD11	1.90	0.53
1:B:271:CYS:O	1:B:289:ARG:CZ	2.57	0.53
1:B:8:LEU:HB2	1:B:65:LEU:CD2	2.39	0.53
1:E:95:ILE:HD12	1:E:307:HIS:HD2	1.73	0.53
1:C:31:LEU:HD23	1:C:73:ARG:CZ	2.39	0.53
1:C:169:SER:O	1:C:172:CYS:HB2	2.08	0.53
1:E:5:ILE:HA	1:E:60:ILE:HG13	1.90	0.53
1:B:9:THR:HG22	1:B:13:GLN:OE1	2.08	0.53
1:C:286:ARG:HD3	1:C:289:ARG:CG	2.38	0.53
1:E:269:GLU:HA	1:E:289:ARG:O	2.09	0.53
1:B:341:ILE:HA	1:B:344:LEU:HD12	1.92	0.52
1:C:206:ILE:O	1:C:207:LEU:HD12	2.09	0.52
1:D:97:ALA:HB2	1:D:298:ALA:HB2	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:276:TRP:CD1	1:D:278:ILE:CG1	2.92	0.52
1:B:224:SER:CB	1:E:344:LEU:HB2	2.39	0.52
1:E:237:LYS:O	1:E:240:CYS:HB2	2.10	0.52
1:B:114:LEU:N	1:B:114:LEU:CD1	2.73	0.52
1:C:20:ARG:NH1	1:C:277:GLN:CB	2.73	0.52
1:C:21:MET:HB2	1:C:23:TYR:CD2	2.45	0.52
1:E:300:LEU:HD11	1:E:325:ILE:HG13	1.84	0.52
1:D:94:ASP:HB3	1:D:307:HIS:HD2	1.74	0.52
1:A:174:TRP:CD2	1:A:178:LYS:HG3	2.45	0.52
1:C:153:ILE:HD13	1:C:179:LEU:HB3	1.92	0.51
1:C:205:LYS:HB3	1:C:207:LEU:CD1	2.41	0.51
1:E:129:THR:HG23	1:E:131:ASN:H	1.73	0.51
1:D:129:THR:HG23	1:D:131:ASN:H	1.75	0.51
1:E:95:ILE:HD11	1:E:307:HIS:CD2	2.45	0.51
1:B:20:ARG:HE	1:B:42:VAL:HG13	1.75	0.51
1:D:111:LEU:N	1:D:111:LEU:CD1	2.73	0.51
1:A:344:LEU:HD13	1:A:344:LEU:C	2.30	0.51
1:E:194:LEU:HB3	1:E:206:ILE:HD13	1.92	0.51
1:A:262:VAL:HG12	2:A:406:HOH:O	2.10	0.51
1:D:276:TRP:CD1	1:D:278:ILE:HG12	2.46	0.51
1:D:303:PHE:HB3	1:D:332:HIS:CD2	2.46	0.51
1:E:55:ILE:CG2	1:E:302:GLU:HG2	2.40	0.51
1:E:126:ARG:HH21	1:E:126:ARG:HG3	1.77	0.50
1:E:151:ILE:HD11	1:E:218:PHE:CZ	2.45	0.50
1:B:179:LEU:O	1:C:36:SER:HB3	2.11	0.50
1:B:280:SER:HB3	1:B:284:MET:CE	2.41	0.50
1:D:49:SER:OG	1:D:50:ALA:N	2.45	0.50
1:D:226:PRO:O	1:D:227:GLN:CB	2.60	0.50
1:B:174:TRP:CE2	1:B:178:LYS:HG3	2.46	0.50
1:C:231:ILE:CD1	1:C:231:ILE:N	2.74	0.50
1:D:288:GLU:OE1	1:D:288:GLU:N	2.35	0.50
1:C:193:ARG:HA	1:C:196:ASN:OD1	2.11	0.50
1:C:297:GLU:O	1:C:301:LEU:HG	2.12	0.50
1:E:2:PHE:CD1	1:E:2:PHE:N	2.72	0.50
1:B:72:ARG:NH1	1:E:258:GLU:HG3	2.26	0.50
1:B:280:SER:HB3	1:B:284:MET:HE1	1.94	0.50
1:C:20:ARG:HH12	1:C:277:GLN:CB	2.13	0.50
1:E:41:ILE:HG22	1:E:60:ILE:HG22	1.94	0.50
1:B:286:ARG:HG3	1:B:288:GLU:OE2	2.12	0.50
1:E:276:TRP:CZ2	1:E:291:MET:HB3	2.47	0.50
1:A:11:LEU:HD23	1:A:65:LEU:HD21	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:8:LEU:HD12	1:D:65:LEU:HD21	1.92	0.50
1:D:136:TYR:CE1	1:D:140:ARG:HD3	2.46	0.50
1:E:95:ILE:CG1	1:E:307:HIS:CD2	2.94	0.50
1:E:95:ILE:HG12	1:E:307:HIS:HA	1.93	0.50
1:B:29:GLN:CD	1:B:34:TRP:HZ2	2.14	0.49
1:B:271:CYS:O	1:B:289:ARG:NH2	2.45	0.49
1:C:231:ILE:CD1	1:C:254:LYS:HD3	2.42	0.49
1:D:177:LEU:HD22	1:D:319:ILE:HD12	1.93	0.49
1:E:49:SER:OG	1:E:50:ALA:N	2.45	0.49
1:E:175:LEU:HD12	1:E:181:VAL:HG21	1.93	0.49
1:E:95:ILE:HD13	1:E:301:LEU:CG	2.42	0.49
1:E:291:MET:CE	1:E:296:ALA:HB2	2.41	0.49
1:B:250:ASN:OD1	1:E:342:GLU:HG3	2.13	0.49
1:C:7:HIS:O	1:C:279:MET:HE3	2.13	0.49
1:E:2:PHE:HE1	1:E:57:GLY:N	2.10	0.49
1:A:267:VAL:HG12	1:A:337:LEU:HB2	1.94	0.49
1:B:29:GLN:OE1	1:B:34:TRP:HZ2	1.96	0.49
1:D:213:LEU:CD1	1:D:231:ILE:HD11	2.43	0.49
1:A:15:ARG:NH2	1:A:27:ALA:O	2.46	0.49
1:D:206:ILE:O	1:D:207:LEU:HD23	2.13	0.49
1:C:221:TRP:CD1	1:C:223:ALA:O	2.66	0.49
1:E:224:SER:O	1:E:226:PRO:HD3	2.13	0.48
1:D:278:ILE:HG22	1:D:279:MET:H	1.78	0.48
1:E:93:ILE:HD12	1:E:93:ILE:H	1.78	0.48
1:B:31:LEU:HD13	1:E:150:GLY:CA	2.43	0.48
1:B:139:CYS:HB3	1:B:174:TRP:CD1	2.49	0.48
1:C:143:GLU:HB2	1:C:179:LEU:HD21	1.95	0.48
1:D:278:ILE:HG22	1:D:279:MET:N	2.29	0.48
1:E:300:LEU:HD11	1:E:325:ILE:CG1	2.34	0.48
1:B:164:THR:HG21	1:B:193:ARG:HD3	1.96	0.48
1:B:226:PRO:HG2	1:B:229:VAL:HG21	1.96	0.48
1:C:172:CYS:HB3	1:C:201:LEU:HD21	1.95	0.48
1:E:59:TYR:OH	1:E:279:MET:HB3	2.14	0.48
1:E:95:ILE:HD13	1:E:301:LEU:CB	2.44	0.48
1:B:11:LEU:O	1:B:14:ALA:HB3	2.14	0.48
1:E:66:PRO:HB3	1:E:101:PHE:HE2	1.78	0.48
1:B:8:LEU:HD13	1:B:63:CYS:HA	1.96	0.48
1:C:164:THR:HG21	1:C:193:ARG:HD3	1.96	0.48
1:C:232:ASP:O	1:C:235:THR:N	2.43	0.48
1:B:164:THR:HG22	1:B:314:ARG:HH21	1.79	0.47
1:E:138:ILE:O	1:E:142:VAL:HG23	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:227:GLN:HE22	1:D:343:ASN:HB3	1.79	0.47
1:E:108:ASN:C	1:E:109:PHE:HD1	2.16	0.47
1:C:238:GLN:N	1:C:239:PRO:CD	2.78	0.47
1:E:41:ILE:O	1:E:41:ILE:HD12	2.14	0.47
1:D:294:CYS:O	1:D:297:GLU:HB3	2.14	0.47
1:E:81:VAL:HG22	1:E:102:THR:HG22	1.96	0.47
1:D:49:SER:HB2	1:D:303:PHE:CE1	2.50	0.47
1:C:7:HIS:HB2	1:C:64:PHE:O	2.14	0.47
1:C:49:SER:OG	1:C:50:ALA:N	2.46	0.47
1:A:206:ILE:O	1:A:207:LEU:HD12	2.15	0.47
1:B:276:TRP:HH2	1:B:291:MET:HB3	1.80	0.47
1:B:194:LEU:HD22	1:B:206:ILE:HG23	1.96	0.47
1:D:4:LEU:HD13	1:D:295:PHE:CD2	2.50	0.47
1:E:95:ILE:HD13	1:E:301:LEU:HB3	1.96	0.47
1:A:343:ASN:O	1:A:344:LEU:HB3	2.13	0.46
1:C:4:LEU:HD22	1:C:295:PHE:CD1	2.49	0.46
1:D:247:TYR:HB3	1:D:248:PRO:HD3	1.98	0.46
1:E:7:HIS:HB2	1:E:64:PHE:O	2.15	0.46
1:E:175:LEU:CD1	1:E:181:VAL:HG21	2.45	0.46
1:C:231:ILE:HG13	1:C:255:VAL:CB	2.45	0.46
1:B:111:LEU:HA	1:B:114:LEU:HD13	1.96	0.46
1:A:19:ARG:HD2	1:A:25:GLU:OE2	2.14	0.46
1:C:238:GLN:N	1:C:239:PRO:HD2	2.31	0.46
1:D:225:MET:HE1	1:D:251:LEU:HD13	1.98	0.46
1:E:46:THR:HA	1:E:55:ILE:O	2.14	0.46
1:B:20:ARG:HE	1:B:42:VAL:CG1	2.28	0.46
1:A:31:LEU:HD11	1:A:68:MET:HG3	1.98	0.46
1:B:17:VAL:O	1:B:21:MET:HG3	2.15	0.46
1:D:284:MET:HG3	1:D:290:GLN:HB2	1.97	0.46
1:D:29:GLN:N	1:D:29:GLN:NE2	2.59	0.46
1:E:64:PHE:CZ	1:E:105:ILE:HD11	2.51	0.46
1:D:72:ARG:HA	2:D:407:HOH:O	2.16	0.46
1:A:226:PRO:HD2	1:A:229:VAL:HG21	1.98	0.45
1:B:164:THR:CG2	1:B:193:ARG:HD3	2.47	0.45
1:B:278:ILE:HG12	1:B:279:MET:H	1.80	0.45
1:D:231:ILE:HG23	1:D:231:ILE:O	2.16	0.45
1:E:310:PHE:HZ	1:E:321:LYS:O	2.00	0.45
1:C:254:LYS:HD3	1:C:254:LYS:C	2.36	0.45
1:A:239:PRO:HG2	2:A:401:HOH:O	2.15	0.45
1:B:301:LEU:HD23	1:B:301:LEU:HA	1.71	0.45
1:C:227:GLN:HA	1:C:227:GLN:OE1	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:303:PHE:HB3	1:A:332:HIS:CE1	2.52	0.45
1:D:2:PHE:CE2	1:D:299:MET:HG2	2.52	0.45
1:B:7:HIS:H	1:B:279:MET:CE	2.28	0.45
1:A:106:PHE:HA	1:A:111:LEU:HD12	1.99	0.45
1:B:76:THR:HG22	1:B:79:ARG:NH2	2.32	0.45
1:D:32:GLU:HA	1:D:35:SER:OG	2.17	0.45
1:D:297:GLU:O	1:D:301:LEU:HG	2.16	0.45
1:B:14:ALA:O	1:B:18:SER:HB3	2.17	0.45
1:E:74:PHE:HB3	1:E:109:PHE:CE2	2.47	0.45
1:E:55:ILE:HD12	1:E:55:ILE:N	2.32	0.45
1:C:276:TRP:CE2	1:C:278:ILE:HD13	2.52	0.44
1:E:5:ILE:HD13	1:E:60:ILE:CD1	2.47	0.44
1:E:95:ILE:HD11	1:E:307:HIS:CA	2.45	0.44
1:E:171:VAL:O	1:E:175:LEU:HD23	2.16	0.44
1:C:45:ILE:HG21	1:C:276:TRP:CE3	2.52	0.44
1:A:152:ASP:OD1	1:A:154:THR:HB	2.17	0.44
1:A:344:LEU:C	1:A:344:LEU:CD1	2.85	0.44
1:C:124:PHE:HB3	1:C:312:TRP:CZ2	2.53	0.44
1:D:236:LEU:HD12	1:D:260:ILE:HG12	2.00	0.44
1:E:2:PHE:HE1	1:E:57:GLY:CA	2.28	0.44
1:E:4:LEU:CD1	1:E:295:PHE:CG	2.88	0.44
1:C:188:ALA:H	1:C:194:LEU:HD21	1.82	0.44
1:D:23:TYR:HE1	1:D:25:GLU:HG3	1.82	0.44
1:D:154:THR:HG22	1:D:155:GLN:HG3	1.99	0.44
1:E:42:VAL:HG22	1:E:59:TYR:HD1	1.83	0.44
1:B:111:LEU:C	1:B:113:SER:N	2.71	0.44
1:E:3:GLY:O	1:E:96:SER:HA	2.18	0.44
1:E:1:MET:SD	1:E:57:GLY:HA2	2.58	0.44
1:B:25:GLU:H	1:B:25:GLU:HG3	1.65	0.44
1:D:226:PRO:O	1:D:227:GLN:HG2	2.18	0.44
1:E:1:MET:HA	1:E:56:HIS:O	2.17	0.44
1:E:5:ILE:HD13	1:E:60:ILE:HG12	1.99	0.44
1:B:318:THR:OG1	1:B:321:LYS:HG3	2.18	0.43
1:E:41:ILE:CG2	1:E:60:ILE:HG22	2.48	0.43
1:C:106:PHE:HA	1:C:111:LEU:HD22	2.00	0.43
1:C:124:PHE:O	1:C:309:ASN:HB3	2.17	0.43
1:D:123:GLU:HG3	1:D:125:GLU:HG2	2.00	0.43
1:E:151:ILE:HD12	1:E:151:ILE:O	2.17	0.43
1:D:55:ILE:HG13	1:D:303:PHE:CE1	2.53	0.43
1:C:278:ILE:O	1:C:281:ALA:HB2	2.19	0.43
1:D:55:ILE:HD11	1:D:303:PHE:HA	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:300:LEU:C	1:E:300:LEU:CD2	2.86	0.43
1:B:344:LEU:O	1:B:345:TYR:HB3	2.18	0.43
1:C:198:GLN:NE2	1:C:204:GLY:O	2.41	0.43
1:E:31:LEU:HD23	1:E:31:LEU:HA	1.66	0.43
1:E:137:VAL:HG13	1:E:336:PRO:HB3	2.01	0.43
1:D:69:LEU:HD23	1:D:69:LEU:HA	1.68	0.43
1:C:127:PHE:O	1:C:309:ASN:HA	2.18	0.43
1:B:162:GLY:HA3	1:B:223:ALA:HB2	2.00	0.43
1:B:248:PRO:HG3	1:E:342:GLU:HB2	2.00	0.43
1:D:276:TRP:CD1	1:D:278:ILE:HD11	2.54	0.43
1:D:303:PHE:HB3	1:D:332:HIS:NE2	2.33	0.43
1:E:64:PHE:HE2	1:E:101:PHE:CD2	2.36	0.43
1:C:161:VAL:O	1:C:222:VAL:HG12	2.18	0.43
1:C:232:ASP:O	1:C:235:THR:HB	2.19	0.43
1:D:29:GLN:NE2	1:D:29:GLN:CA	2.82	0.43
1:B:303:PHE:HB3	1:B:332:HIS:CE1	2.54	0.42
1:C:175:LEU:O	1:C:181:VAL:HG23	2.19	0.42
1:D:174:TRP:HE3	1:D:322:MET:HE3	1.84	0.42
1:A:76:THR:O	1:A:80:LYS:HG2	2.19	0.42
1:B:276:TRP:CH2	1:B:291:MET:HB3	2.54	0.42
1:D:13:GLN:O	1:D:13:GLN:HG2	2.19	0.42
1:D:174:TRP:CZ2	1:D:178:LYS:HD3	2.55	0.42
1:E:263:LEU:HD11	1:E:337:LEU:HD23	2.00	0.42
1:C:291:MET:HE1	1:C:296:ALA:HB2	2.01	0.42
1:A:15:ARG:HA	1:A:18:SER:HB2	2.01	0.42
1:D:127:PHE:O	1:D:309:ASN:HA	2.19	0.42
1:E:64:PHE:HE2	1:E:101:PHE:HD2	1.66	0.42
1:E:95:ILE:CG1	1:E:301:LEU:HD23	2.49	0.42
1:C:166:ASP:CB	1:C:315:ASN:H	2.31	0.42
1:C:280:SER:O	1:C:284:MET:HG2	2.20	0.42
1:C:80:LYS:HD2	1:C:80:LYS:HA	1.60	0.42
1:E:3:GLY:O	1:E:4:LEU:HD23	2.20	0.42
1:E:310:PHE:CZ	1:E:321:LYS:O	2.72	0.42
1:B:172:CYS:HB3	1:B:201:LEU:HD11	2.01	0.42
1:E:89:GLN:HG3	1:E:126:ARG:HH22	1.85	0.42
1:E:198:GLN:NE2	1:E:204:GLY:O	2.53	0.42
1:E:3:GLY:HA3	1:E:93:ILE:HG21	2.01	0.42
1:E:5:ILE:HA	1:E:60:ILE:CG1	2.49	0.42
1:A:325:ILE:HA	1:A:325:ILE:HD12	1.84	0.42
1:C:327:GLU:HA	1:C:330:VAL:HG12	2.02	0.42
1:E:55:ILE:HG22	1:E:56:HIS:N	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:14:ALA:HB1	1:B:34:TRP:CZ2	2.55	0.41
1:B:342:GLU:CD	1:B:342:GLU:H	2.23	0.41
1:C:161:VAL:HA	1:C:187:THR:HB	2.02	0.41
1:D:258:GLU:O	1:D:260:ILE:HD12	2.20	0.41
1:A:140:ARG:HD2	1:A:140:ARG:HA	1.86	0.41
1:C:238:GLN:HA	1:C:238:GLN:OE1	2.21	0.41
1:C:241:VAL:HA	1:C:261:TYR:O	2.19	0.41
1:E:50:ALA:HA	1:E:272:PHE:CE2	2.54	0.41
1:B:224:SER:HB3	1:E:344:LEU:HD22	2.03	0.41
1:E:50:ALA:HA	1:E:272:PHE:HE2	1.86	0.41
1:E:126:ARG:HG3	1:E:126:ARG:NH2	2.35	0.41
1:C:49:SER:OG	1:C:51:THR:N	2.54	0.41
1:D:226:PRO:O	1:D:227:GLN:CG	2.69	0.41
1:E:75:LYS:O	1:E:79:ARG:HG2	2.20	0.41
1:A:1:MET:HA	1:A:56:HIS:O	2.20	0.41
1:B:244:ASP:OD2	1:B:251:LEU:HB2	2.21	0.41
1:C:193:ARG:O	1:C:196:ASN:HB2	2.20	0.41
1:C:206:ILE:C	1:C:207:LEU:HD12	2.40	0.41
1:A:16:ASP:OD1	1:A:16:ASP:C	2.58	0.41
1:E:95:ILE:HD13	1:E:301:LEU:HG	2.03	0.41
1:A:174:TRP:CE2	1:A:178:LYS:HG3	2.56	0.41
1:D:25:GLU:H	1:D:25:GLU:HG2	1.65	0.41
1:D:147:LYS:HB2	1:D:147:LYS:HE2	1.72	0.41
1:D:109:PHE:O	1:D:111:LEU:CD1	2.68	0.41
1:D:335:GLN:HA	1:D:335:GLN:OE1	2.21	0.41
1:C:90:LYS:HD3	1:C:118:ARG:HH22	1.86	0.40
1:C:158:VAL:HG11	1:C:175:LEU:HD23	2.02	0.40
1:C:232:ASP:C	1:C:234:ALA:N	2.70	0.40
1:D:238:GLN:OE1	1:D:238:GLN:HA	2.21	0.40
1:E:76:THR:HA	1:E:79:ARG:HD3	2.03	0.40
1:E:149:LEU:HD23	1:E:149:LEU:HA	1.91	0.40
1:E:295:PHE:HD1	1:E:295:PHE:HA	1.72	0.40
1:A:98:LEU:HB3	1:A:102:THR:CG2	2.50	0.40
1:B:8:LEU:CD1	1:B:63:CYS:HA	2.50	0.40
1:D:29:GLN:HE21	1:D:29:GLN:CA	2.34	0.40
1:D:101:PHE:HD1	1:D:104:ILE:HD12	1.85	0.40
1:D:227:GLN:O	1:D:227:GLN:HG3	2.20	0.40
1:E:37:ALA:HA	1:E:38:PRO:HD3	1.80	0.40
1:B:64:PHE:C	1:B:65:LEU:HD23	2.42	0.40
1:D:200:GLU:OE2	1:D:200:GLU:HA	2.22	0.40
1:E:285:ALA:O	1:E:287:PRO:HD3	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:326:GLY:O	1:A:329:SER:HB3	2.21	0.40
1:C:21:MET:HB2	1:C:23:TYR:CE2	2.56	0.40
1:C:153:ILE:CD1	1:C:179:LEU:HD23	2.50	0.40
1:C:229:VAL:HG12	1:C:230:VAL:N	2.37	0.40
1:E:2:PHE:HE1	1:E:56:HIS:C	2.24	0.40
1:B:4:LEU:HB2	1:B:59:TYR:HD1	1.86	0.40
1:C:268:VAL:HG23	1:C:291:MET:HE3	2.03	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	336/347 (97%)	316 (94%)	20 (6%)	0	100	100
1	B	343/347 (99%)	329 (96%)	14 (4%)	0	100	100
1	C	342/347 (99%)	322 (94%)	20 (6%)	0	100	100
1	D	335/347 (96%)	319 (95%)	16 (5%)	0	100	100
1	E	344/347 (99%)	324 (94%)	20 (6%)	0	100	100
All	All	1700/1735 (98%)	1610 (95%)	90 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	271/281 (96%)	264 (97%)	7 (3%)	46	79
1	B	271/281 (96%)	265 (98%)	6 (2%)	52	83
1	C	266/281 (95%)	258 (97%)	8 (3%)	41	75
1	D	249/281 (89%)	235 (94%)	14 (6%)	21	51
1	E	252/281 (90%)	235 (93%)	17 (7%)	16	43
All	All	1309/1405 (93%)	1257 (96%)	52 (4%)	31	65

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	ARG
1	A	75	LYS
1	A	189	ARG
1	A	249	LYS
1	A	262	VAL
1	A	276	TRP
1	A	344	LEU
1	B	25	GLU
1	B	147	LYS
1	B	229	VAL
1	B	249	LYS
1	B	276	TRP
1	B	283	GLU
1	C	28	ASP
1	C	51	THR
1	C	173	ARG
1	C	225	MET
1	C	231	ILE
1	C	236	LEU
1	C	276	TRP
1	C	291	MET
1	D	13	GLN
1	D	23	TYR
1	D	24	ASP
1	D	25	GLU
1	D	26	TYR
1	D	29	GLN
1	D	48	THR
1	D	86	SER
1	D	111	LEU
1	D	189	ARG

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Mol	Chain	Res	Type
1	D	225	MET
1	D	227	GLN
1	D	276	TRP
1	D	345	TYR
1	E	2	PHE
1	E	4	LEU
1	E	10	SER
1	E	13	GLN
1	E	24	ASP
1	E	45	ILE
1	E	103	SER
1	E	157	THR
1	E	175	LEU
1	E	189	ARG
1	E	249	LYS
1	E	258	GLU
1	E	275	ASP
1	E	276	TRP
1	E	279	MET
1	E	312	TRP
1	E	316	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	343	ASN
1	D	29	GLN
1	D	89	GLN
1	D	332	HIS
1	E	190	ASN
1	E	307	HIS
1	E	309	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, 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	340/347 (97%)	0.15	2 (0%) 89 86	38, 56, 88, 139	1 (0%)
1	B	345/347 (99%)	0.51	21 (6%) 21 13	39, 63, 105, 145	0
1	C	344/347 (99%)	0.46	31 (9%) 9 5	40, 70, 106, 145	0
1	D	339/347 (97%)	0.35	14 (4%) 37 27	24, 81, 108, 160	1 (0%)
1	E	345/347 (99%)	0.61	37 (10%) 6 3	55, 89, 109, 146	1 (0%)
All	All	1713/1735 (98%)	0.42	105 (6%) 21 13	24, 73, 106, 160	3 (0%)

All (105) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	26	TYR	6.3
1	B	114	LEU	6.1
1	D	111	LEU	5.4
1	B	29	GLN	5.3
1	C	24	ASP	5.3
1	C	201	LEU	5.0
1	E	118	ARG	4.9
1	B	26	TYR	4.3
1	D	26	TYR	4.3
1	B	117	VAL	4.2
1	C	228	GLY	4.2
1	E	56	HIS	4.2
1	C	250	ASN	4.0
1	E	3	GLY	4.0
1	C	253	SER	4.0
1	C	229	VAL	4.0
1	B	115	ARG	4.0
1	E	120	THR	4.0
1	B	74	PHE	3.9
1	E	307	HIS	3.8

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Mol	Chain	Res	Type	RSRZ
1	E	303	PHE	3.8
1	B	285	ALA	3.7
1	D	310	PHE	3.6
1	E	1	MET	3.6
1	E	117	VAL	3.6
1	B	27	ALA	3.5
1	D	27	ALA	3.5
1	C	218	PHE	3.5
1	E	301	LEU	3.4
1	E	54	VAL	3.3
1	D	25	GLU	3.3
1	E	271	CYS	3.3
1	E	325	ILE	3.2
1	A	228	GLY	3.2
1	D	184	LEU	3.2
1	C	27	ALA	3.1
1	E	277	GLN	3.1
1	C	231	ILE	3.1
1	E	47	VAL	3.0
1	C	278	ILE	3.0
1	D	201	LEU	3.0
1	B	116	GLN	2.9
1	C	248	PRO	2.8
1	C	226	PRO	2.8
1	D	23	TYR	2.8
1	B	119	ASP	2.8
1	E	318	THR	2.8
1	B	284	MET	2.7
1	B	343	ASN	2.7
1	C	206	ILE	2.7
1	D	68	MET	2.7
1	B	87	HIS	2.7
1	E	98	LEU	2.6
1	E	329	SER	2.6
1	E	109	PHE	2.5
1	E	119	ASP	2.5
1	B	120	THR	2.5
1	E	308	THR	2.5
1	E	95	ILE	2.5
1	C	252	GLY	2.4
1	C	251	LEU	2.4
1	B	340	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	118	ARG	2.4
1	C	197	LEU	2.4
1	C	25	GLU	2.4
1	E	345	TYR	2.4
1	D	106	PHE	2.4
1	E	58	ARG	2.3
1	C	45	ILE	2.3
1	C	230	VAL	2.3
1	C	227	GLN	2.3
1	C	254	LYS	2.3
1	E	201	LEU	2.3
1	D	112	ALA	2.3
1	E	41	ILE	2.3
1	D	227	GLN	2.3
1	C	23	TYR	2.2
1	E	53	LYS	2.2
1	E	276	TRP	2.2
1	D	181	VAL	2.2
1	B	283	GLU	2.2
1	E	270	HIS	2.2
1	B	111	LEU	2.2
1	E	122	LEU	2.2
1	B	24	ASP	2.2
1	B	23	TYR	2.2
1	E	123	GLU	2.2
1	E	42	VAL	2.2
1	C	20	ARG	2.2
1	E	116	GLN	2.1
1	C	181	VAL	2.1
1	C	118	ARG	2.1
1	B	89	GLN	2.1
1	D	79	ARG	2.1
1	E	82	LEU	2.1
1	A	227	GLN	2.1
1	E	324	ALA	2.1
1	C	255	VAL	2.1
1	C	277	GLN	2.1
1	C	341	ILE	2.1
1	E	268	VAL	2.1
1	C	205	LYS	2.0
1	E	23	TYR	2.0
1	C	225	MET	2.0

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Mol	Chain	Res	Type	RSRZ
1	E	91	HIS	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 (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.