



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:44 PM UTC

PDB ID : 1HZD / pdb_00001hzd
Title : CRYSTAL STRUCTURE OF HUMAN AUH PROTEIN, AN RNA-BINDING
HOMOLOGUE OF ENOYL-COA HYDRATASE
Authors : Kurimoto, K.; Fukai, S.; Nureki, O.; Muto, Y.; Yokoyama, S.; RIKEN Struc-
tural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2001-01-24
Resolution : 2.20 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

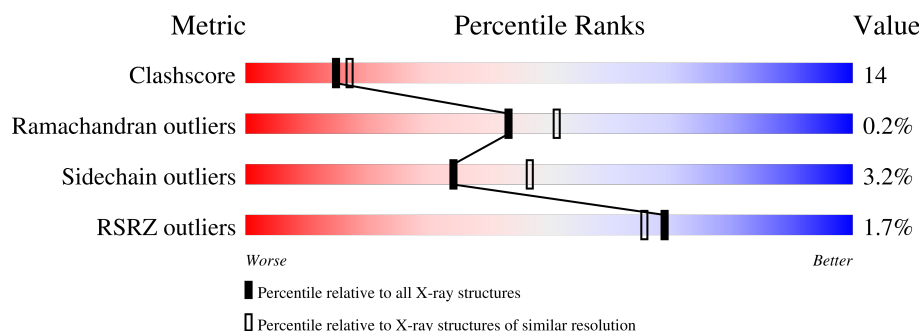
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 $\leq 5\%$. 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	272	3% 68% 28% ..
1	B	272	% 72% 24% ..
1	C	272	% 69% 26% ..
1	D	272	% 67% 27% ..
1	E	272	2% 72% 24% ..
1	F	272	2% 74% 22% ..

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 12282 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 AU-BINDING PROTEIN/ENOYL-COA HYDRATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	266	Total	C	N	O	S	0	0	0
			2000	1260	359	372	9			
1	B	265	Total	C	N	O	S	0	0	0
			1991	1255	358	369	9			
1	C	266	Total	C	N	O	S	0	0	0
			2000	1260	359	372	9			
1	D	265	Total	C	N	O	S	0	0	0
			1991	1255	358	369	9			
1	E	266	Total	C	N	O	S	0	0	0
			2000	1260	359	372	9			
1	F	265	Total	C	N	O	S	0	0	0
			1991	1255	358	369	9			

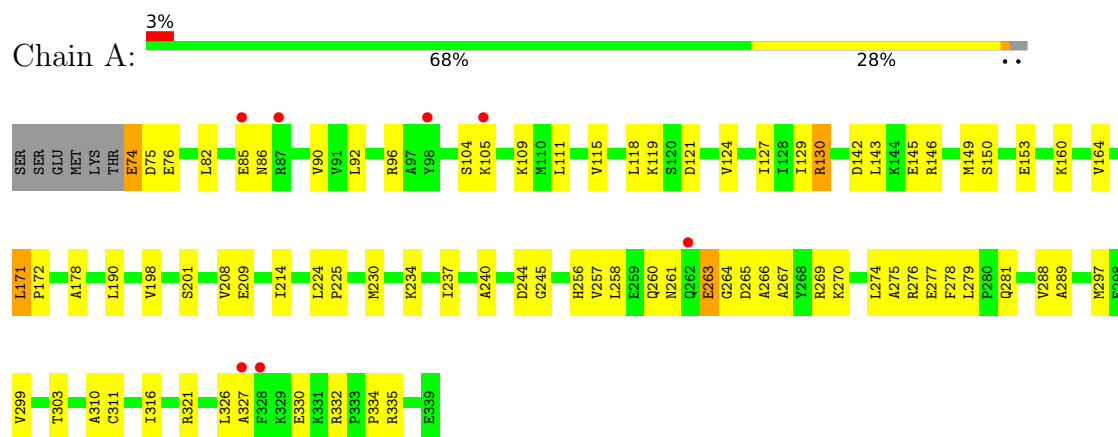
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	57	Total	O	0	0
			57	57		
2	B	53	Total	O	0	0
			53	53		
2	C	44	Total	O	0	0
			44	44		
2	D	45	Total	O	0	0
			45	45		
2	E	50	Total	O	0	0
			50	50		
2	F	60	Total	O	0	0
			60	60		

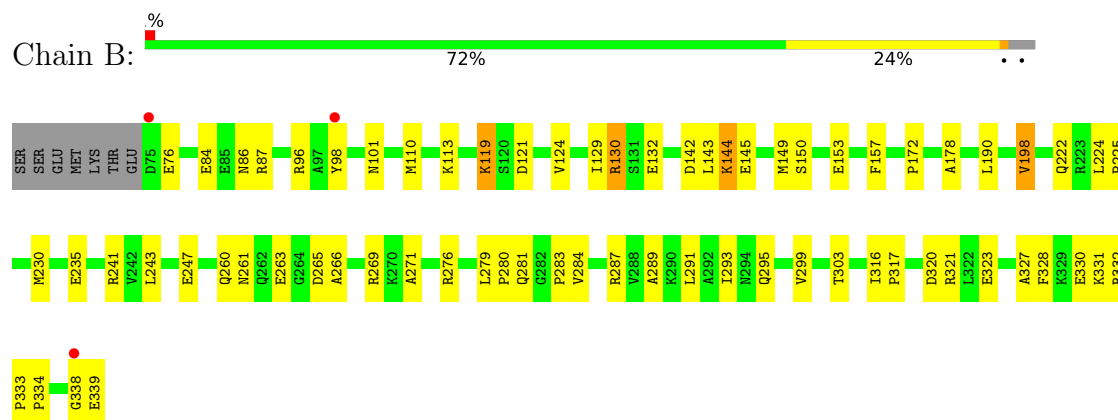
3 Residue-property plots [i](#)

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.

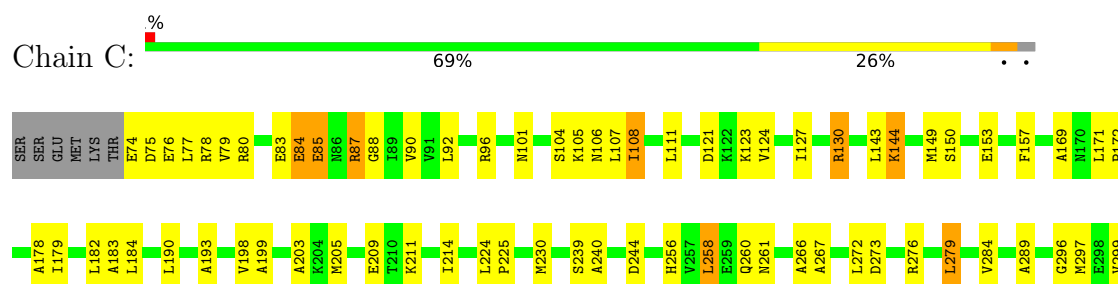
• Molecule 1: AU-BINDING PROTEIN/ENOYL-COA HYDRATASE



• Molecule 1: AU-BINDING PROTEIN/ENOYL-COA HYDRATASE

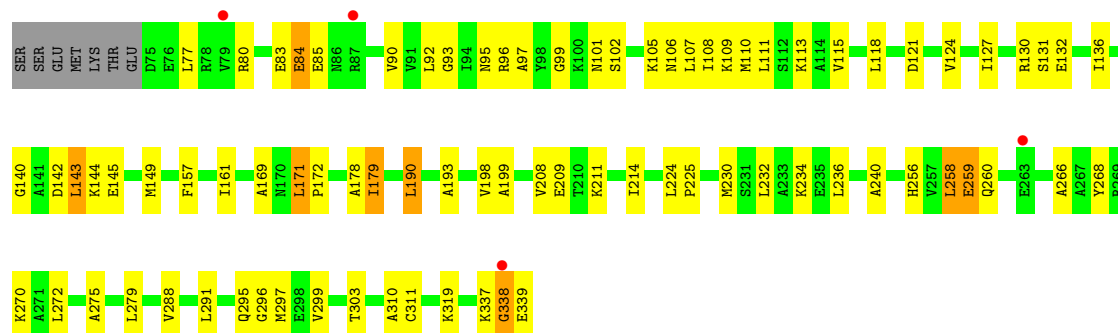


• Molecule 1: AU-BINDING PROTEIN/ENOYL-COA HYDRATASE

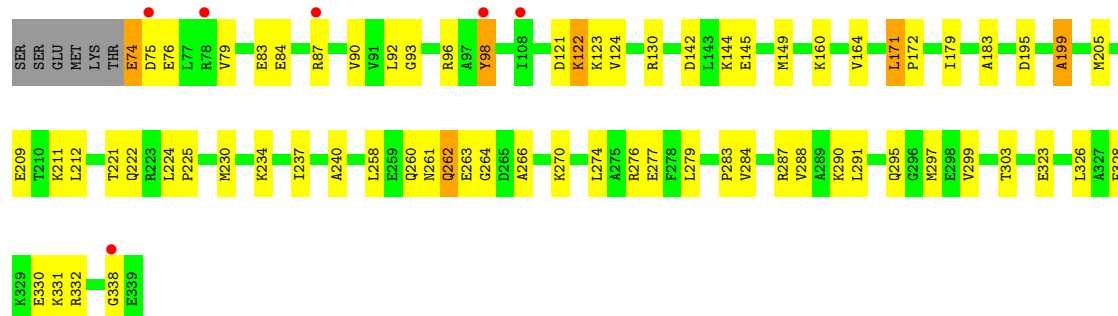




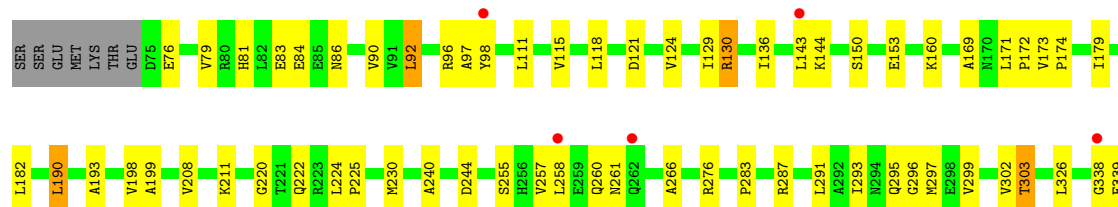
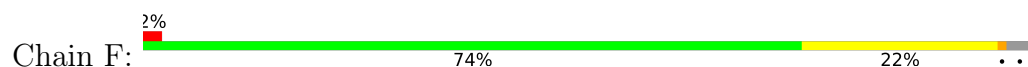
• Molecule 1: AU-BINDING PROTEIN/ENOYL-COA HYDRATASE



• Molecule 1: AU-BINDING PROTEIN/ENOYL-COA HYDRATASE



• Molecule 1: AU-BINDING PROTEIN/ENOYL-COA HYDRATASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	79.13Å 132.07Å 80.04Å 90.00° 108.14° 90.00°	Depositor
Resolution (Å)	20.00 – 2.20 20.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.5 (20.00-2.20) 96.8 (20.00-2.20)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.84 (at 2.19Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.207 , 0.253 0.216 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	22.1	Xtriage
Anisotropy	0.544	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.023 for l,-k,h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12282	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.37% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

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	0.41	0/2021	0.92	3/2717 (0.1%)
1	B	0.44	0/2012	0.90	2/2705 (0.1%)
1	C	0.40	0/2021	0.90	3/2717 (0.1%)
1	D	0.40	0/2012	0.92	6/2705 (0.2%)
1	E	0.41	0/2021	0.92	4/2717 (0.1%)
1	F	0.43	0/2012	0.91	3/2705 (0.1%)
All	All	0.42	0/12099	0.91	21/16266 (0.1%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	171	LEU	CA-C-N	8.02	127.75	119.56
1	D	171	LEU	C-N-CA	8.02	127.75	119.56
1	A	171	LEU	CA-C-N	7.62	128.00	119.32
1	A	171	LEU	C-N-CA	7.62	128.00	119.32
1	C	88	GLY	N-CA-C	-7.39	105.68	113.58
1	F	84	GLU	CB-CA-C	-5.93	109.75	116.63
1	F	303	THR	N-CA-C	-5.89	104.86	111.28
1	E	171	LEU	CA-C-N	5.68	125.36	119.05
1	E	171	LEU	C-N-CA	5.68	125.36	119.05
1	D	338	GLY	N-CA-C	-5.53	102.27	112.62
1	B	84	GLU	CB-CA-C	-5.27	110.06	117.23
1	F	220	GLY	N-CA-C	5.26	119.25	112.83
1	C	289	ALA	N-CA-C	-5.25	105.64	111.36
1	D	179	ILE	N-CA-C	5.21	115.13	107.15
1	D	259	GLU	N-CA-C	-5.19	101.24	109.96
1	E	199	ALA	N-CA-C	5.15	117.08	108.99
1	D	132	GLU	N-CA-C	-5.13	106.87	112.72
1	E	209	GLU	N-CA-C	5.10	117.51	111.33
1	B	289	ALA	N-CA-C	-5.08	105.82	111.36
1	A	289	ALA	N-CA-C	-5.06	105.84	111.36
1	C	279	LEU	N-CA-C	5.02	119.77	113.25

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2000	0	2116	62	0
1	B	1991	0	2110	60	0
1	C	2000	0	2116	65	0
1	D	1991	0	2110	66	0
1	E	2000	0	2116	62	0
1	F	1991	0	2110	50	0
2	A	57	0	0	1	0
2	B	53	0	0	0	0
2	C	44	0	0	1	0
2	D	45	0	0	0	0
2	E	50	0	0	1	0
2	F	60	0	0	1	0
All	All	12282	0	12678	340	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 14.

All (340) 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:D:258:LEU:HD21	1:D:270:LYS:HB2	1.26	1.09
1:B:86:ASN:HD21	1:B:276:ARG:HH11	1.13	0.96
1:A:146:ARG:HA	1:A:149:MET:HE3	1.50	0.92
1:B:130:ARG:HH11	1:B:130:ARG:HB2	1.36	0.90
1:F:144:LYS:HE2	1:F:144:LYS:HA	1.54	0.89
1:D:260:GLN:HE22	1:D:266:ALA:H	1.24	0.86
1:F:86:ASN:HD21	1:F:276:ARG:HH11	1.25	0.83
1:C:260:GLN:HE22	1:C:266:ALA:H	1.27	0.81
1:B:260:GLN:HE22	1:B:266:ALA:H	1.28	0.80
1:C:276:ARG:HA	1:C:279:LEU:HD13	1.60	0.80
1:D:260:GLN:NE2	1:D:266:ALA:H	1.78	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:143:LEU:HD12	1:D:143:LEU:H	1.46	0.80
1:D:258:LEU:HD21	1:D:270:LYS:CB	2.12	0.78
1:F:121:ASP:CG	1:F:124:VAL:HG23	2.10	0.75
1:A:258:LEU:HD12	1:A:267:ALA:HA	1.68	0.75
1:F:260:GLN:HE21	1:F:261:ASN:H	1.32	0.75
1:A:74:GLU:HB3	1:A:76:GLU:HG3	1.68	0.74
1:B:303:THR:HG21	1:F:299:VAL:HG11	1.70	0.73
1:A:299:VAL:HG11	1:D:303:THR:HG21	1.71	0.72
1:B:327:ALA:HA	1:B:332:ARG:NH1	2.04	0.72
1:E:276:ARG:HA	1:E:279:LEU:HD13	1.72	0.72
1:B:129:ILE:HG21	1:B:190:LEU:HD21	1.72	0.71
1:B:299:VAL:HG11	1:F:303:THR:HG21	1.74	0.70
1:B:143:LEU:HD23	1:B:143:LEU:O	1.93	0.68
1:E:121:ASP:OD1	1:E:124:VAL:HG23	1.93	0.68
1:C:299:VAL:HG11	1:E:303:THR:HG21	1.75	0.68
1:E:260:GLN:HE21	1:E:261:ASN:H	1.40	0.68
1:E:142:ASP:OD1	1:E:144:LYS:HB2	1.95	0.67
1:E:326:LEU:HD12	1:E:332:ARG:HH21	1.60	0.67
1:F:257:VAL:C	1:F:258:LEU:HD12	2.19	0.67
1:E:90:VAL:HG21	1:E:124:VAL:HG22	1.77	0.66
1:C:144:LYS:HE2	1:C:144:LYS:HA	1.78	0.66
1:D:90:VAL:CG2	1:D:127:ILE:HG12	2.25	0.65
1:B:119:LYS:HD3	1:B:172:PRO:HD3	1.78	0.65
1:F:130:ARG:HH11	1:F:130:ARG:HB2	1.62	0.65
1:D:209:GLU:HB3	1:D:214:ILE:HG13	1.79	0.65
1:A:92:LEU:HD11	1:A:118:LEU:HD11	1.78	0.65
1:A:258:LEU:CD1	1:A:267:ALA:HA	2.28	0.64
1:E:121:ASP:CG	1:E:124:VAL:HG23	2.23	0.64
1:C:104:SER:O	1:C:108:ILE:HG12	1.99	0.63
1:F:260:GLN:HE22	1:F:266:ALA:H	1.45	0.63
1:A:260:GLN:HE21	1:A:261:ASN:H	1.46	0.63
1:A:260:GLN:HE22	1:A:266:ALA:H	1.47	0.63
1:C:96:ARG:H	1:C:101:ASN:ND2	1.97	0.62
1:A:303:THR:HG21	1:D:299:VAL:HG11	1.83	0.61
1:B:76:GLU:HB2	1:B:110:MET:HE1	1.81	0.61
1:F:143:LEU:O	1:F:143:LEU:HD23	2.01	0.61
1:B:260:GLN:NE2	1:B:266:ALA:H	1.98	0.60
1:D:84:GLU:O	1:D:85:GLU:HG2	2.01	0.60
1:E:74:GLU:HG2	1:E:75:ASP:H	1.67	0.60
1:D:108:ILE:HD11	1:D:161:ILE:HG12	1.82	0.60
1:E:230:MET:HE3	1:F:297:MET:HG3	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:121:ASP:CG	1:D:124:VAL:HG13	2.27	0.59
1:E:260:GLN:HE22	1:E:266:ALA:H	1.49	0.59
1:C:90:VAL:HG22	1:C:127:ILE:HG12	1.84	0.59
1:C:182:LEU:HD11	1:C:184:LEU:HG	1.84	0.59
1:D:121:ASP:OD1	1:D:124:VAL:HG13	2.03	0.59
1:B:260:GLN:HE21	1:B:261:ASN:H	1.50	0.59
1:F:190:LEU:HD23	1:F:190:LEU:C	2.27	0.58
1:D:118:LEU:HD13	1:D:127:ILE:HD13	1.84	0.58
1:C:260:GLN:HE21	1:C:261:ASN:H	1.50	0.58
1:E:84:GLU:O	1:E:87:ARG:HG2	2.03	0.58
1:B:130:ARG:HB2	1:B:130:ARG:NH1	2.15	0.58
1:E:237:ILE:HD12	1:F:293:ILE:HD11	1.85	0.58
1:A:150:SER:OG	1:A:153:GLU:HG3	2.03	0.58
1:A:129:ILE:HG21	1:A:190:LEU:HD21	1.85	0.57
1:C:260:GLN:NE2	1:C:266:ALA:H	2.00	0.57
1:C:224:LEU:HB3	1:C:225:PRO:HD3	1.86	0.57
1:A:90:VAL:HG22	1:A:127:ILE:HG12	1.85	0.57
1:C:284:VAL:HG23	1:C:338:GLY:O	2.04	0.57
1:B:265:ASP:O	1:B:269:ARG:HG2	2.04	0.57
1:E:179:ILE:HB	1:E:199:ALA:HB2	1.84	0.57
1:B:291:LEU:HB3	1:B:295:GLN:NE2	2.19	0.57
1:D:208:VAL:HG12	1:D:240:ALA:O	2.05	0.57
1:F:76:GLU:CD	1:F:96:ARG:HH11	2.13	0.57
1:A:258:LEU:HD11	1:A:270:LYS:CB	2.34	0.57
1:E:326:LEU:HD12	1:E:332:ARG:NH2	2.18	0.57
1:C:303:THR:HG21	1:E:299:VAL:HG11	1.86	0.56
1:B:144:LYS:HE3	1:B:144:LYS:HA	1.86	0.56
1:B:145:GLU:O	1:B:149:MET:HG3	2.05	0.56
1:D:230:MET:CG	1:D:234:LYS:HE3	2.35	0.56
1:F:260:GLN:NE2	1:F:261:ASN:H	2.00	0.56
1:A:327:ALA:HB3	1:A:334:PRO:HG3	1.87	0.56
1:D:268:TYR:CE2	1:D:272:LEU:HD11	2.41	0.56
1:A:265:ASP:OD2	1:A:269:ARG:NE	2.39	0.56
1:C:79:VAL:HG22	1:C:92:LEU:CD1	2.35	0.56
1:E:122:LYS:HD3	1:E:122:LYS:N	2.20	0.56
1:C:96:ARG:HD3	1:C:107:LEU:HD22	1.88	0.56
1:A:111:LEU:O	1:A:115:VAL:HG23	2.06	0.55
1:A:258:LEU:HD11	1:A:270:LYS:HB3	1.88	0.55
1:B:86:ASN:HD21	1:B:276:ARG:NH1	1.95	0.55
1:E:76:GLU:OE2	1:E:96:ARG:HD2	2.06	0.55
1:E:145:GLU:O	1:E:149:MET:HG3	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:SER:OG	1:B:153:GLU:HG3	2.07	0.55
1:E:258:LEU:HD21	1:E:270:LYS:HB2	1.88	0.55
1:A:160:LYS:O	1:A:164:VAL:HG23	2.06	0.55
1:C:190:LEU:C	1:C:190:LEU:HD23	2.31	0.55
1:D:224:LEU:HB3	1:D:225:PRO:HD3	1.88	0.54
1:C:90:VAL:CG2	1:C:127:ILE:HG12	2.37	0.54
1:D:142:ASP:OD2	1:D:145:GLU:HG3	2.07	0.54
1:A:105:LYS:O	1:A:109:LYS:HG3	2.09	0.53
1:D:105:LYS:O	1:D:109:LYS:HG3	2.07	0.53
1:D:106:ASN:OD1	1:D:110:MET:HE2	2.08	0.53
1:D:149:MET:HE1	1:D:157:PHE:HB2	1.89	0.53
1:D:149:MET:HE1	1:D:157:PHE:CB	2.38	0.53
1:A:96:ARG:NH2	1:A:104:SER:HB3	2.23	0.53
1:B:76:GLU:HB3	1:B:96:ARG:HG2	1.90	0.53
1:C:96:ARG:H	1:C:101:ASN:HD21	1.57	0.53
1:F:79:VAL:HG22	1:F:92:LEU:CD1	2.37	0.53
1:D:275:ALA:O	1:D:279:LEU:HD13	2.09	0.53
1:D:105:LYS:HG3	1:D:106:ASN:N	2.24	0.53
1:F:121:ASP:OD2	1:F:124:VAL:HG23	2.08	0.52
1:B:230:MET:HE3	1:C:297:MET:HG3	1.91	0.52
1:F:81:HIS:HE1	1:F:121:ASP:OD1	1.93	0.52
1:F:224:LEU:HB3	1:F:225:PRO:HD3	1.91	0.52
1:E:326:LEU:CD1	1:E:332:ARG:HH21	2.22	0.52
1:C:143:LEU:HD23	1:C:143:LEU:O	2.10	0.52
1:D:258:LEU:N	1:D:258:LEU:HD12	2.25	0.52
1:A:281:GLN:HE21	1:C:211:LYS:HG2	1.75	0.52
1:D:179:ILE:HB	1:D:199:ALA:HB2	1.92	0.52
1:C:96:ARG:CD	1:C:107:LEU:HD22	2.39	0.51
1:D:108:ILE:CD1	1:D:161:ILE:HG12	2.40	0.51
1:A:327:ALA:CB	1:A:334:PRO:HG3	2.41	0.51
1:E:260:GLN:NE2	1:E:261:ASN:H	2.08	0.51
1:C:79:VAL:HG22	1:C:92:LEU:HD12	1.92	0.51
1:E:222:GLN:OE1	1:F:296:GLY:HA3	2.11	0.51
1:F:86:ASN:HD21	1:F:276:ARG:NH1	2.03	0.51
1:A:297:MET:HG3	1:C:230:MET:HE3	1.91	0.51
1:C:209:GLU:HB3	1:C:214:ILE:HG13	1.92	0.51
1:D:95:ASN:C	1:D:95:ASN:HD22	2.18	0.51
1:B:96:ARG:HB2	1:B:101:ASN:HA	1.91	0.51
1:D:297:MET:HG3	1:F:230:MET:HE3	1.92	0.51
1:F:136:ILE:HA	1:F:182:LEU:H	1.76	0.51
1:F:90:VAL:HG21	1:F:124:VAL:HG22	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:230:MET:HE3	1:E:297:MET:HG3	1.92	0.50
1:A:224:LEU:HB3	1:A:225:PRO:HD3	1.93	0.50
1:B:222:GLN:OE1	1:C:296:GLY:HA3	2.11	0.50
1:C:84:GLU:O	1:C:85:GLU:CB	2.60	0.50
1:E:122:LYS:HD3	1:E:122:LYS:H	1.75	0.50
1:E:221:THR:O	1:E:225:PRO:HG2	2.11	0.50
1:C:75:ASP:HB3	1:C:78:ARG:HE	1.77	0.50
1:D:319:LYS:HD2	1:D:339:GLU:OE2	2.12	0.49
1:E:93:GLY:HA2	1:E:130:ARG:O	2.11	0.49
1:B:121:ASP:CG	1:B:124:VAL:HG23	2.37	0.49
1:F:76:GLU:OE1	1:F:96:ARG:HD2	2.13	0.49
1:A:82:LEU:HB3	1:A:86:ASN:HB2	1.93	0.49
1:A:230:MET:HG2	1:A:234:LYS:HE3	1.93	0.49
1:D:80:ARG:HH11	1:D:80:ARG:HG2	1.77	0.49
1:A:275:ALA:O	1:A:279:LEU:HD13	2.13	0.49
1:E:195:ASP:OD1	1:E:290:LYS:HE3	2.13	0.49
1:F:169:ALA:HA	1:F:193:ALA:O	2.13	0.49
1:D:230:MET:HG2	1:D:234:LYS:HE3	1.93	0.49
1:D:338:GLY:O	1:D:339:GLU:OXT	2.31	0.49
1:E:284:VAL:HG23	1:E:338:GLY:O	2.12	0.49
1:A:208:VAL:HA	1:A:240:ALA:HB1	1.94	0.49
1:B:316:ILE:HB	1:B:317:PRO:HD3	1.94	0.49
1:D:84:GLU:O	1:D:85:GLU:CG	2.61	0.49
1:D:130:ARG:HB3	1:D:178:ALA:HB3	1.95	0.49
1:F:129:ILE:HG21	1:F:190:LEU:HD21	1.95	0.49
1:C:105:LYS:HG3	1:C:106:ASN:N	2.28	0.49
1:D:95:ASN:C	1:D:95:ASN:ND2	2.71	0.49
1:E:260:GLN:NE2	1:E:264:GLY:HA2	2.28	0.49
1:D:109:LYS:O	1:D:113:LYS:HG3	2.13	0.48
1:D:111:LEU:O	1:D:115:VAL:HG23	2.13	0.48
1:A:150:SER:HG	1:A:153:GLU:HG3	1.77	0.48
1:C:198:VAL:HG23	1:C:267:ALA:HB1	1.95	0.48
1:A:335:ARG:HG2	1:A:335:ARG:HH11	1.79	0.48
1:E:74:GLU:N	1:E:74:GLU:CD	2.72	0.48
1:F:79:VAL:HG22	1:F:92:LEU:HD12	1.96	0.48
1:E:262:GLN:HE21	1:E:262:GLN:HB3	1.51	0.48
1:C:84:GLU:O	1:C:84:GLU:HG3	2.14	0.47
1:C:87:ARG:HG3	1:C:87:ARG:O	2.14	0.47
1:B:190:LEU:C	1:B:190:LEU:HD23	2.39	0.47
1:C:211:LYS:HG3	1:C:240:ALA:CB	2.44	0.47
1:E:274:LEU:O	1:E:277:GLU:HB2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:291:LEU:HD22	1:E:295:GLN:HE22	1.78	0.47
1:C:80:ARG:HG2	1:C:80:ARG:HH11	1.79	0.47
1:F:338:GLY:O	1:F:339:GLU:OXT	2.32	0.47
1:A:85:GLU:HG2	1:A:276:ARG:HD3	1.96	0.47
1:A:118:LEU:HD13	1:A:127:ILE:HD13	1.97	0.47
1:C:121:ASP:OD2	1:C:123:LYS:HG3	2.15	0.47
1:D:169:ALA:HA	1:D:193:ALA:O	2.15	0.47
1:A:142:ASP:OD1	1:A:145:GLU:HG3	2.14	0.47
1:B:284:VAL:HG23	1:B:338:GLY:O	2.15	0.47
1:C:150:SER:HB3	1:C:153:GLU:HG3	1.97	0.47
1:C:302:VAL:HG21	1:F:302:VAL:HG11	1.97	0.47
1:D:93:GLY:HA2	1:D:130:ARG:O	2.15	0.46
1:D:143:LEU:H	1:D:143:LEU:CD1	2.23	0.46
1:E:279:LEU:HD12	1:E:279:LEU:N	2.31	0.46
1:D:84:GLU:O	1:D:84:GLU:HG3	2.14	0.46
1:C:183:ALA:O	1:C:205:MET:HA	2.16	0.46
1:F:179:ILE:HB	1:F:199:ALA:HB2	1.97	0.46
1:D:258:LEU:CD2	1:D:270:LYS:HB2	2.19	0.46
1:D:260:GLN:HE22	1:D:266:ALA:N	2.03	0.46
1:E:79:VAL:HG22	1:E:92:LEU:CD1	2.45	0.46
1:A:263:GLU:HB2	2:A:383:HOH:O	2.15	0.46
1:A:143:LEU:HD13	1:B:328:PHE:CZ	2.51	0.46
1:B:143:LEU:HD13	1:C:328:PHE:CE2	2.51	0.46
1:D:232:LEU:O	1:D:236:LEU:HG	2.15	0.46
1:A:190:LEU:HD23	1:A:190:LEU:C	2.41	0.46
1:C:273:ASP:HA	1:C:276:ARG:NH1	2.31	0.46
1:B:86:ASN:ND2	1:B:276:ARG:HH11	1.96	0.45
1:B:316:ILE:HA	1:B:321:ARG:HD3	1.98	0.45
1:D:96:ARG:HB3	1:D:101:ASN:HA	1.97	0.45
1:E:160:LYS:O	1:E:164:VAL:HG23	2.16	0.45
1:D:90:VAL:CG1	1:D:124:VAL:HG12	2.46	0.45
1:D:108:ILE:HD12	1:D:157:PHE:CZ	2.51	0.45
1:A:90:VAL:CG2	1:A:127:ILE:HG12	2.47	0.45
1:B:279:LEU:HB2	1:B:280:PRO:HD3	1.97	0.45
1:B:330:GLU:HB2	1:B:332:ARG:HH11	1.82	0.45
1:C:144:LYS:HA	1:C:144:LYS:CE	2.45	0.45
1:A:90:VAL:HG13	1:A:124:VAL:HG13	1.98	0.45
1:B:119:LYS:HB2	1:B:119:LYS:NZ	2.30	0.45
1:C:96:ARG:HB2	1:C:101:ASN:HA	1.99	0.45
1:C:190:LEU:HD23	1:C:190:LEU:O	2.17	0.45
1:A:274:LEU:O	1:A:277:GLU:HB3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:198:VAL:HG12	1:A:256:HIS:HB2	1.98	0.45
1:C:84:GLU:O	1:C:85:GLU:HB3	2.17	0.45
1:A:130:ARG:HB3	1:A:178:ALA:HB3	1.99	0.44
1:D:296:GLY:HA3	1:F:222:GLN:OE1	2.16	0.44
1:B:330:GLU:O	1:B:331:LYS:C	2.61	0.44
1:B:130:ARG:HD3	1:B:130:ARG:C	2.42	0.44
1:B:283:PRO:O	1:B:287:ARG:HG3	2.17	0.44
1:D:97:ALA:C	1:D:99:GLY:H	2.26	0.44
1:D:288:VAL:HB	1:D:311:CYS:HB3	2.00	0.44
1:F:291:LEU:HD22	1:F:295:GLN:HE22	1.81	0.44
1:E:230:MET:CG	1:E:234:LYS:HE3	2.48	0.44
1:E:284:VAL:O	1:E:288:VAL:HG22	2.18	0.44
1:C:316:ILE:HA	1:C:321:ARG:HD3	1.98	0.44
1:F:90:VAL:CG2	1:F:124:VAL:HG22	2.48	0.44
1:A:257:VAL:O	1:A:257:VAL:HG13	2.17	0.44
1:B:130:ARG:HH11	1:B:130:ARG:CB	2.20	0.44
1:B:87:ARG:HB2	1:B:87:ARG:NH1	2.33	0.44
1:B:280:PRO:HB2	1:B:281:GLN:HE22	1.83	0.44
1:C:105:LYS:HA	1:C:108:ILE:HD11	2.00	0.44
1:C:149:MET:HE1	1:C:157:PHE:CG	2.52	0.43
1:E:87:ARG:HH11	1:E:87:ARG:HG3	1.83	0.43
1:E:130:ARG:C	1:E:130:ARG:CD	2.91	0.43
1:E:130:ARG:HB2	2:E:368:HOH:O	2.18	0.43
1:B:98:TYR:HD1	1:B:98:TYR:H	1.66	0.43
1:A:330:GLU:HB2	1:A:332:ARG:HG2	2.01	0.43
1:B:98:TYR:CD1	1:B:98:TYR:N	2.87	0.43
1:E:130:ARG:C	1:E:130:ARG:HD2	2.43	0.43
1:C:284:VAL:CG2	1:C:338:GLY:O	2.66	0.43
1:E:258:LEU:HD13	1:E:266:ALA:HB1	1.99	0.43
1:B:142:ASP:OD2	1:B:144:LYS:HB2	2.18	0.43
1:F:86:ASN:ND2	1:F:276:ARG:HH11	2.03	0.43
1:F:111:LEU:O	1:F:115:VAL:HG23	2.18	0.43
1:A:74:GLU:O	1:A:75:ASP:HB3	2.19	0.43
1:E:326:LEU:HD13	1:E:326:LEU:O	2.19	0.43
1:F:173:VAL:HB	1:F:174:PRO:HD2	2.00	0.43
1:B:130:ARG:HB3	1:B:178:ALA:HB3	2.01	0.43
1:B:283:PRO:HB2	1:B:338:GLY:HA2	2.00	0.43
1:C:108:ILE:HG12	1:C:108:ILE:H	1.60	0.43
1:C:179:ILE:HB	1:C:199:ALA:HB2	2.00	0.43
1:C:273:ASP:OD1	1:C:276:ARG:NH2	2.51	0.43
1:B:144:LYS:HA	1:B:144:LYS:CE	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:338:GLY:O	1:B:339:GLU:OXT	2.37	0.43
1:D:211:LYS:HG3	1:D:240:ALA:CB	2.48	0.43
1:D:291:LEU:HD22	1:D:295:GLN:NE2	2.33	0.43
1:B:87:ARG:O	1:B:87:ARG:HG3	2.19	0.43
1:B:198:VAL:HG11	1:B:271:ALA:HB2	2.01	0.43
1:D:107:LEU:HA	1:D:110:MET:HE3	2.00	0.43
1:E:326:LEU:HD13	1:E:326:LEU:C	2.44	0.43
1:F:208:VAL:HA	1:F:240:ALA:HB1	2.01	0.43
1:A:115:VAL:O	1:A:119:LYS:HG3	2.19	0.42
1:B:241:ARG:NH2	1:B:247:GLU:OE2	2.52	0.42
1:E:179:ILE:HB	1:E:199:ALA:CB	2.47	0.42
1:D:190:LEU:HD23	1:D:190:LEU:C	2.44	0.42
1:C:79:VAL:HG22	1:C:92:LEU:HD11	2.01	0.42
1:A:130:ARG:C	1:A:130:ARG:CD	2.92	0.42
1:A:316:ILE:HA	1:A:321:ARG:HD3	2.02	0.42
1:A:146:ARG:CA	1:A:149:MET:HE3	2.35	0.42
1:A:209:GLU:HB3	1:A:214:ILE:HG13	2.02	0.42
1:F:190:LEU:HD23	1:F:190:LEU:O	2.20	0.42
1:A:288:VAL:HB	1:A:311:CYS:HB3	2.02	0.42
1:D:198:VAL:HG12	1:D:256:HIS:HB2	2.01	0.42
1:E:171:LEU:HA	1:E:172:PRO:HD3	1.86	0.42
1:F:211:LYS:HG3	1:F:240:ALA:HB3	2.02	0.42
1:A:143:LEU:HD13	1:B:328:PHE:CE2	2.54	0.42
1:B:149:MET:HE1	1:B:157:PHE:CD1	2.55	0.42
1:C:171:LEU:HA	1:C:172:PRO:HD3	1.90	0.42
1:A:260:GLN:NE2	1:A:264:GLY:HA2	2.35	0.42
1:C:203:ALA:O	1:C:244:ASP:HA	2.19	0.42
1:E:330:GLU:O	1:E:331:LYS:C	2.63	0.42
1:A:171:LEU:HA	1:A:172:PRO:HD3	1.81	0.42
1:B:178:ALA:HA	1:B:198:VAL:HG13	2.02	0.42
1:E:87:ARG:HG3	1:E:87:ARG:NH1	2.35	0.42
1:A:201:SER:HA	1:A:245:GLY:HA3	2.02	0.42
1:D:131:SER:CB	1:D:136:ILE:O	2.68	0.42
1:F:150:SER:HB2	1:F:153:GLU:HG3	2.02	0.42
1:C:178:ALA:HA	1:C:198:VAL:HG22	2.02	0.41
1:F:171:LEU:HA	1:F:172:PRO:HD3	1.93	0.41
1:A:278:PHE:HE2	1:C:239:SER:HB3	1.85	0.41
1:A:310:ALA:HB1	1:D:310:ALA:HB1	2.02	0.41
1:C:258:LEU:CD1	1:C:258:LEU:C	2.93	0.41
1:D:102:SER:HA	1:D:140:GLY:O	2.20	0.41
1:E:76:GLU:CD	1:E:96:ARG:HD2	2.45	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:237:ILE:HD12	1:B:293:ILE:HD11	2.03	0.41
1:A:326:LEU:O	1:A:330:GLU:HG3	2.19	0.41
1:A:330:GLU:OE1	1:A:332:ARG:HD3	2.19	0.41
1:C:83:GLU:O	1:C:84:GLU:O	2.38	0.41
1:C:169:ALA:HA	1:C:193:ALA:O	2.21	0.41
1:F:121:ASP:CG	1:F:124:VAL:CG2	2.88	0.41
1:E:211:LYS:NZ	1:E:240:ALA:HB3	2.36	0.41
1:C:121:ASP:CG	1:C:124:VAL:HG23	2.46	0.41
1:E:332:ARG:HG3	1:E:332:ARG:HH11	1.85	0.41
1:F:83:GLU:HB3	2:F:387:HOH:O	2.21	0.41
1:F:211:LYS:HG3	1:F:240:ALA:CB	2.51	0.41
1:A:121:ASP:CG	1:A:124:VAL:HG23	2.45	0.41
1:E:224:LEU:N	1:E:225:PRO:CD	2.84	0.41
1:C:77:LEU:CD2	1:C:111:LEU:HA	2.51	0.41
1:D:90:VAL:HG11	1:D:124:VAL:HG12	2.03	0.41
1:E:98:TYR:CD2	1:E:98:TYR:N	2.84	0.41
1:F:90:VAL:HG11	1:F:118:LEU:HD21	2.01	0.41
1:B:320:ASP:OD2	1:B:338:GLY:O	2.39	0.41
1:C:130:ARG:HB2	2:C:383:HOH:O	2.21	0.41
1:D:77:LEU:HA	1:D:93:GLY:O	2.20	0.41
1:D:171:LEU:HA	1:D:172:PRO:HD3	1.79	0.41
1:E:212:LEU:HD23	1:E:212:LEU:HA	1.93	0.41
1:E:283:PRO:O	1:E:287:ARG:HG3	2.21	0.41
1:F:97:ALA:O	1:F:98:TYR:C	2.64	0.41
1:A:121:ASP:OD1	1:A:124:VAL:HG23	2.20	0.41
1:E:291:LEU:HD22	1:E:295:GLN:NE2	2.35	0.41
1:F:173:VAL:HB	1:F:174:PRO:CD	2.51	0.41
1:C:74:GLU:HG2	1:C:76:GLU:HG3	2.01	0.40
1:D:190:LEU:HD23	1:D:190:LEU:O	2.21	0.40
1:E:122:LYS:H	1:E:122:LYS:CD	2.34	0.40
1:B:222:GLN:C	1:B:225:PRO:HD2	2.46	0.40
1:B:243:LEU:HB3	1:B:247:GLU:HB2	2.02	0.40
1:C:84:GLU:H	1:C:87:ARG:HB3	1.87	0.40
1:D:143:LEU:HD22	1:E:328:PHE:CE2	2.56	0.40
1:F:244:ASP:OD1	1:F:244:ASP:C	2.64	0.40
1:A:244:ASP:OD1	1:A:244:ASP:C	2.65	0.40
1:B:235:GLU:CD	1:C:256:HIS:HE2	2.28	0.40
1:B:333:PRO:HA	1:B:334:PRO:HD3	1.97	0.40
1:E:183:ALA:O	1:E:205:MET:HA	2.21	0.40
1:F:283:PRO:O	1:F:287:ARG:HG3	2.21	0.40
1:B:224:LEU:N	1:B:225:PRO:CD	2.84	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:83:GLU:HB2	1:E:84:GLU:H	1.76	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	264/272 (97%)	252 (96%)	12 (4%)	0	100	100
1	B	263/272 (97%)	254 (97%)	9 (3%)	0	100	100
1	C	264/272 (97%)	255 (97%)	7 (3%)	2 (1%)	16	16
1	D	263/272 (97%)	246 (94%)	16 (6%)	1 (0%)	30	34
1	E	264/272 (97%)	254 (96%)	10 (4%)	0	100	100
1	F	263/272 (97%)	250 (95%)	13 (5%)	0	100	100
All	All	1581/1632 (97%)	1511 (96%)	67 (4%)	3 (0%)	43	51

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	84	GLU
1	D	84	GLU
1	C	85	GLU

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	209/215 (97%)	206 (99%)	3 (1%)	59	75
1	B	208/215 (97%)	200 (96%)	8 (4%)	29	40
1	C	209/215 (97%)	202 (97%)	7 (3%)	33	45
1	D	208/215 (97%)	200 (96%)	8 (4%)	29	40
1	E	209/215 (97%)	202 (97%)	7 (3%)	33	45
1	F	208/215 (97%)	201 (97%)	7 (3%)	32	44
All	All	1251/1290 (97%)	1211 (97%)	40 (3%)	34	47

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	GLU
1	A	130	ARG
1	A	263	GLU
1	B	113	LYS
1	B	119	LYS
1	B	130	ARG
1	B	132	GLU
1	B	144	LYS
1	B	198	VAL
1	B	263	GLU
1	B	323	GLU
1	C	87	ARG
1	C	108	ILE
1	C	130	ARG
1	C	144	LYS
1	C	258	LEU
1	C	272	LEU
1	C	326	LEU
1	D	83	GLU
1	D	92	LEU
1	D	143	LEU
1	D	144	LYS
1	D	190	LEU
1	D	258	LEU
1	D	259	GLU
1	D	337	LYS
1	E	74	GLU
1	E	98	TYR
1	E	122	LYS
1	E	123	LYS

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Mol	Chain	Res	Type
1	E	262	GLN
1	E	263	GLU
1	E	323	GLU
1	F	92	LEU
1	F	130	ARG
1	F	160	LYS
1	F	190	LEU
1	F	198	VAL
1	F	255	SER
1	F	326	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	106	ASN
1	A	260	GLN
1	A	281	GLN
1	B	86	ASN
1	B	95	ASN
1	B	260	GLN
1	B	281	GLN
1	B	295	GLN
1	C	95	ASN
1	C	101	ASN
1	C	260	GLN
1	C	281	GLN
1	D	81	HIS
1	D	260	GLN
1	D	281	GLN
1	E	260	GLN
1	E	262	GLN
1	E	281	GLN
1	E	295	GLN
1	F	81	HIS
1	F	86	ASN
1	F	260	GLN
1	F	281	GLN
1	F	295	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	266/272 (97%)	0.03	7 (2%) 57 54	11, 23, 51, 61	0
1	B	265/272 (97%)	-0.13	3 (1%) 78 76	11, 21, 42, 51	0
1	C	266/272 (97%)	0.04	2 (0%) 82 80	13, 26, 49, 64	0
1	D	265/272 (97%)	0.18	4 (1%) 72 69	13, 29, 52, 59	0
1	E	266/272 (97%)	0.02	6 (2%) 61 58	10, 25, 49, 59	0
1	F	265/272 (97%)	-0.13	5 (1%) 66 63	12, 23, 44, 62	0
All	All	1593/1632 (97%)	0.00	27 (1%) 69 66	10, 25, 49, 64	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	328	PHE	5.1
1	B	338	GLY	4.9
1	E	98	TYR	4.9
1	C	338	GLY	4.5
1	E	338	GLY	4.5
1	F	338	GLY	4.1
1	B	98	TYR	3.6
1	F	98	TYR	3.5
1	F	258	LEU	2.9
1	F	262	GLN	2.7
1	E	75	ASP	2.7
1	A	85	GLU	2.4
1	D	87	ARG	2.4
1	A	105	LYS	2.4
1	A	87	ARG	2.4
1	A	98	TYR	2.3
1	E	78	ARG	2.3
1	E	108	ILE	2.3
1	B	75	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	263	GLU	2.1
1	D	338	GLY	2.1
1	F	143	LEU	2.1
1	A	262	GLN	2.1
1	A	327	ALA	2.0
1	D	79	VAL	2.0
1	C	335	ARG	2.0
1	E	87	ARG	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 oligosaccharides 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.