



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:59 PM GMT

PDB ID : 1DKG
Title : CRYSTAL STRUCTURE OF THE NUCLEOTIDE EXCHANGE FACTOR
GRPE BOUND TO THE ATPASE DOMAIN OF THE MOLECULAR
CHAPERONE DNAK
Authors : Harrison, C.J.; Kuriyan, J.
Deposited on : 1997-02-13
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

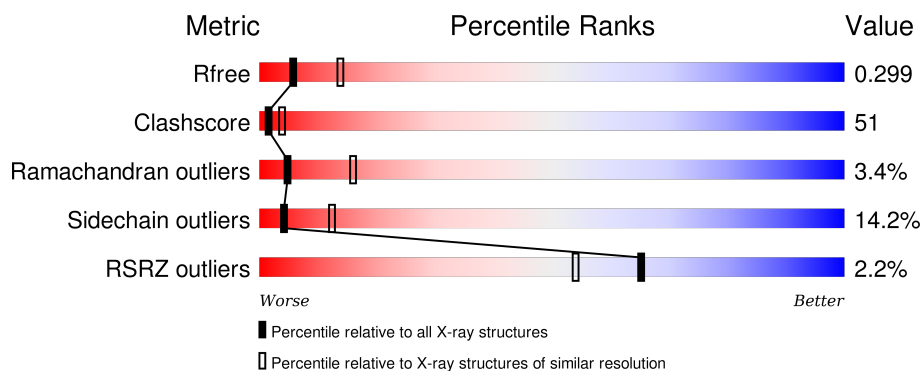
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	197	<div> <div>34%</div> <div>38%</div> <div>9%</div> <div>20%</div> </div>
1	B	197	<div> <div>34%</div> <div>37%</div> <div>6%</div> <div>23%</div> </div>
2	D	383	<div> <div>3%</div> <div>37%</div> <div>50%</div> <div>10%</div> <div>••</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5183 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 NUCLEOTIDE EXCHANGE FACTOR GRPE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	158	Total	C	N	O	S	0	0	0
			1203	750	211	235	7			
1	B	151	Total	C	N	O	S	0	0	0
			1144	716	204	217	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	122	ASP	GLY	ENGINEERED	UNP P09372
B	122	ASP	GLY	ENGINEERED	UNP P09372

- Molecule 2 is a protein called MOLECULAR CHAPERONE DNAK.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	376	Total	C	N	O	S	0	0	0
			2809	1763	483	554	9			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	319	LEU	PRO	ENGINEERED	UNP P04475

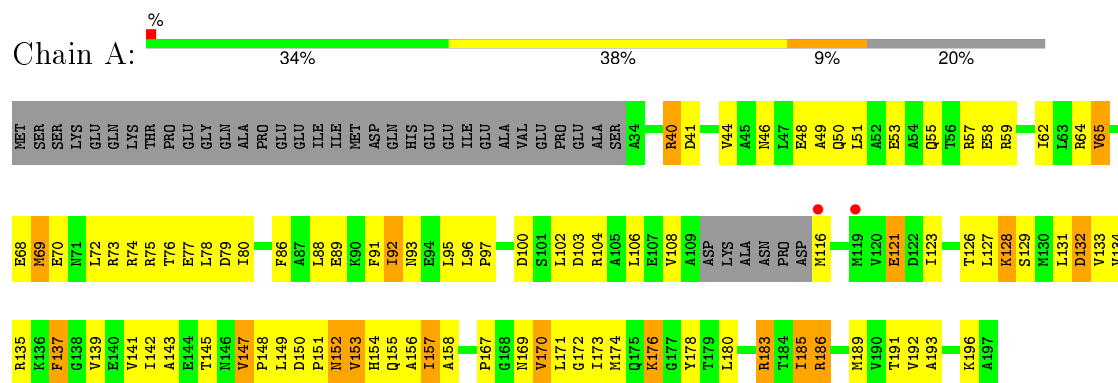
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	7	Total	O	0	0
			7	7		
3	B	4	Total	O	0	0
			4	4		
3	D	16	Total	O	0	0
			16	16		

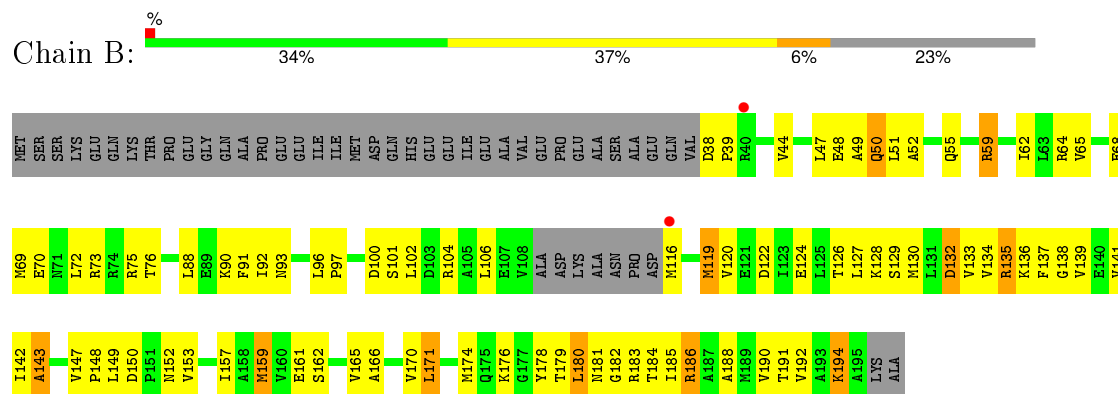
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

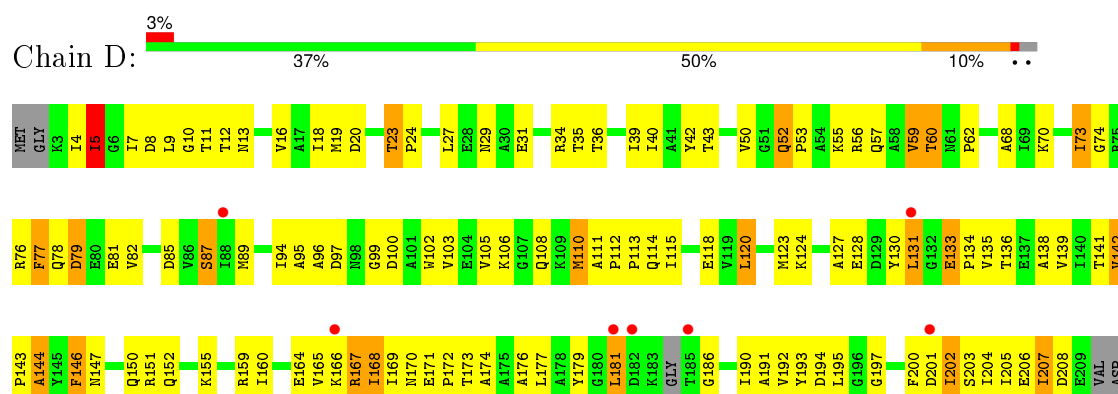
• Molecule 1: NUCLEOTIDE EXCHANGE FACTOR GRPE

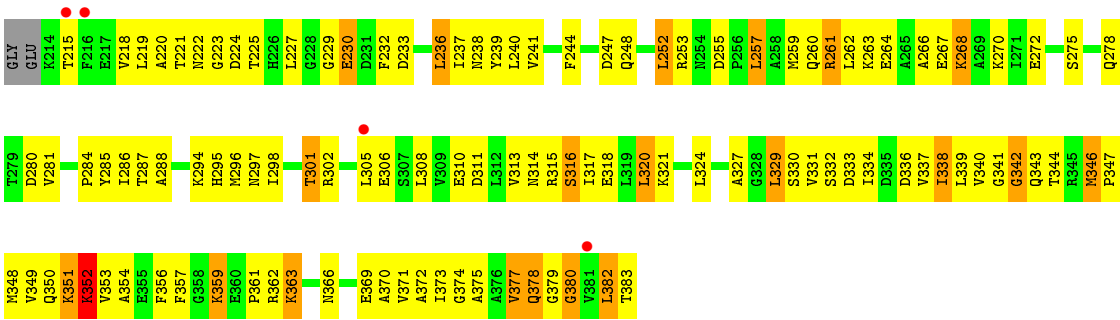


• Molecule 1: NUCLEOTIDE EXCHANGE FACTOR GRPE



• Molecule 2: MOLECULAR CHAPERONE DNAK





4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	149.43 Å 149.43 Å 49.03 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.80 29.89 – 2.80	Depositor EDS
% Data completeness (in resolution range)	87.9 (30.00-2.80) 93.9 (29.89-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.50 (at 2.80 Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.223 , 0.317 0.245 , 0.299	Depositor DCC
R_{free} test set	1967 reflections (7.70%)	DCC
Wilson B-factor (Å ²)	70.3	Xtriage
Anisotropy	0.039	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 39.8	EDS
Estimated twinning fraction	0.036 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	1 of 25559 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5183	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.80% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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	0.62	0/1212	0.81	0/1639
1	B	0.82	2/1153 (0.2%)	0.84	3/1557 (0.2%)
2	D	0.55	0/2843	0.75	0/3853
All	All	0.63	2/5208 (0.0%)	0.78	3/7049 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	132	ASP	CG-OD2	13.64	1.56	1.25
1	B	132	ASP	CG-OD1	12.77	1.54	1.25

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	132	ASP	CB-CG-OD2	-9.92	109.37	118.30
1	B	132	ASP	CB-CG-OD1	-8.79	110.39	118.30
1	B	132	ASP	OD1-CG-OD2	-5.72	112.44	123.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	132	ASP	Sidechain

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	1203	0	1213	166	0
1	B	1144	0	1158	120	0
2	D	2809	0	2807	287	0
3	A	7	0	0	0	0
3	B	4	0	0	0	0
3	D	16	0	0	2	0
All	All	5183	0	5178	525	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 51.

All (525) 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:40:ARG:HG3	1:A:40:ARG:HH11	1.03	1.12
1:A:142:ILE:HG13	1:A:153:VAL:CG2	1.77	1.12
2:D:202:ILE:HD12	2:D:203:SER:N	1.66	1.11
2:D:18:ILE:HD13	2:D:131:LEU:HD21	1.23	1.10
2:D:202:ILE:HD12	2:D:203:SER:H	1.12	1.10
1:A:142:ILE:HG13	1:A:153:VAL:HG21	1.30	1.08
1:B:141:VAL:HG12	1:B:178:TYR:CE1	1.90	1.05
1:A:186:ARG:HH11	1:A:186:ARG:HG2	1.13	1.05
1:B:171:LEU:HD23	1:B:171:LEU:O	1.59	1.03
2:D:252:LEU:HB3	2:D:259:MET:CE	1.89	1.02
1:B:157:ILE:HD11	1:B:174:MET:HG3	1.43	1.01
2:D:285:TYR:H	2:D:295:HIS:CD2	1.79	1.00
1:B:134:VAL:CG1	1:B:139:VAL:HB	1.91	1.00
1:A:150:ASP:H	1:A:154:HIS:HD2	1.08	0.99
1:A:40:ARG:HH11	1:A:40:ARG:CG	1.75	0.98
1:A:92:ILE:HD11	1:A:139:VAL:HG21	1.45	0.97
2:D:248:GLN:NE2	2:D:294:LYS:HD2	1.79	0.97
1:A:92:ILE:CD1	1:A:139:VAL:HG21	1.95	0.96
1:B:161:GLU:HA	1:B:194:LYS:O	1.65	0.96
1:B:134:VAL:HG13	1:B:139:VAL:HB	1.44	0.95
2:D:205:ILE:HD13	2:D:207:ILE:HD11	1.46	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:112:PRO:HG2	2:D:113:PRO:HD3	1.52	0.92
2:D:146:PHE:HA	2:D:150:GLN:NE2	1.85	0.92
2:D:331:VAL:HG11	2:D:356:PHE:O	1.71	0.91
1:A:116:MET:O	1:A:116:MET:HG2	1.69	0.91
2:D:52:GLN:HA	2:D:52:GLN:OE1	1.73	0.89
2:D:5:ILE:HG13	2:D:5:ILE:O	1.70	0.88
1:A:186:ARG:HG2	1:A:186:ARG:NH1	1.81	0.88
2:D:327:ALA:O	2:D:329:LEU:HG	1.73	0.88
1:A:127:LEU:HD23	1:A:127:LEU:O	1.74	0.88
2:D:77:PHE:CD2	2:D:95:ALA:HB2	2.09	0.86
1:A:153:VAL:HG22	1:A:153:VAL:O	1.71	0.86
1:A:40:ARG:NH1	1:A:40:ARG:HG3	1.81	0.86
2:D:252:LEU:HB3	2:D:259:MET:HE2	1.54	0.86
2:D:341:GLY:O	2:D:343:GLN:N	2.08	0.86
2:D:192:VAL:HG22	2:D:338:ILE:CG1	2.05	0.86
1:A:150:ASP:H	1:A:154:HIS:CD2	1.94	0.86
1:B:93:ASN:OD1	1:B:185:ILE:HD11	1.77	0.85
1:B:142:ILE:HG21	1:B:153:VAL:CG1	2.06	0.85
2:D:73:ILE:HG23	2:D:74:GLY:N	1.91	0.84
2:D:208:ASP:O	2:D:215:THR:HG22	1.78	0.84
1:B:100:ASP:OD2	1:B:186:ARG:NH2	2.10	0.83
2:D:202:ILE:HG22	2:D:222:ASN:O	1.78	0.83
1:A:129:SER:O	1:A:133:VAL:HG23	1.79	0.82
2:D:192:VAL:HG22	2:D:338:ILE:HG13	1.58	0.82
1:A:91:PHE:CD2	1:B:91:PHE:CE2	2.66	0.82
1:B:142:ILE:HG21	1:B:153:VAL:HG12	1.59	0.82
1:A:91:PHE:CD2	1:B:91:PHE:HE2	1.96	0.82
2:D:205:ILE:CD1	2:D:207:ILE:HD11	2.07	0.82
1:B:180:LEU:HD12	1:B:180:LEU:O	1.79	0.82
1:A:64:ARG:O	1:A:68:GLU:HB2	1.80	0.82
2:D:18:ILE:HD13	2:D:131:LEU:CD2	2.09	0.81
1:A:145:THR:HA	1:A:170:VAL:CG2	2.10	0.81
1:A:186:ARG:HH11	1:A:186:ARG:CG	1.93	0.81
2:D:11:THR:OG1	2:D:70:LYS:HD2	1.81	0.80
1:B:92:ILE:HD13	1:B:139:VAL:CG2	2.11	0.80
2:D:281:VAL:HG11	2:D:296:MET:HE2	1.62	0.80
1:B:88:LEU:O	1:B:92:ILE:HG13	1.82	0.80
2:D:310:GLU:HA	2:D:348:MET:HE1	1.64	0.80
2:D:146:PHE:HA	2:D:150:GLN:HE22	1.44	0.79
1:A:131:LEU:O	1:A:135:ARG:HG3	1.82	0.79
2:D:124:LYS:O	2:D:128:GLU:HG3	1.82	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:119:MET:O	1:B:122:ASP:HB2	1.84	0.78
2:D:268:LYS:O	2:D:272:GLU:HG3	1.83	0.78
2:D:236:LEU:HD22	2:D:305:LEU:HD13	1.64	0.77
1:B:141:VAL:HG12	1:B:178:TYR:HE1	1.49	0.77
2:D:142:VAL:CG1	2:D:146:PHE:CD2	2.68	0.77
1:A:147:VAL:C	1:A:169:ASN:HD22	1.88	0.77
1:B:150:ASP:OD1	1:B:152:ASN:HB3	1.83	0.77
2:D:252:LEU:HB3	2:D:259:MET:HE3	1.66	0.76
1:B:141:VAL:HG12	1:B:178:TYR:CD1	2.19	0.76
1:A:89:GLU:O	1:A:93:ASN:HB2	1.86	0.76
2:D:205:ILE:HG22	2:D:218:VAL:HA	1.68	0.76
1:B:64:ARG:O	1:B:68:GLU:HB2	1.86	0.76
1:A:142:ILE:CG1	1:A:153:VAL:HG21	2.14	0.76
1:A:91:PHE:HD1	1:B:137:PHE:CZ	2.03	0.75
2:D:313:VAL:HG21	2:D:348:MET:HE2	1.68	0.75
1:A:91:PHE:CE2	1:B:91:PHE:HE2	2.03	0.75
2:D:5:ILE:HD11	2:D:7:ILE:HD11	1.67	0.75
1:A:147:VAL:C	1:A:169:ASN:ND2	2.40	0.75
1:A:178:TYR:HB2	1:A:185:ILE:HD11	1.66	0.75
1:B:106:LEU:HD21	1:B:124:GLU:HG2	1.67	0.75
2:D:204:ILE:HD12	2:D:220:ALA:HB3	1.68	0.74
2:D:318:GLU:OE1	2:D:321:LYS:HD2	1.86	0.74
2:D:105:VAL:CG2	2:D:110:MET:HE2	2.18	0.74
2:D:112:PRO:CG	2:D:113:PRO:HD3	2.16	0.74
2:D:329:LEU:HB3	2:D:333:ASP:OD2	1.87	0.74
2:D:59:VAL:O	2:D:62:PRO:HD3	1.87	0.74
1:A:91:PHE:HD2	1:B:91:PHE:CE2	2.04	0.74
1:A:69:MET:HB2	1:B:69:MET:CG	2.18	0.74
2:D:302:ARG:O	2:D:306:GLU:HG3	1.87	0.74
2:D:18:ILE:HG13	2:D:18:ILE:O	1.87	0.74
2:D:179:TYR:CD2	2:D:338:ILE:HG21	2.23	0.73
1:A:88:LEU:CD1	1:A:137:PHE:HD2	2.00	0.73
1:A:91:PHE:CE2	1:B:91:PHE:CE2	2.75	0.73
1:B:133:VAL:O	1:B:136:LYS:HB2	1.89	0.73
2:D:378:GLN:HE21	2:D:378:GLN:HA	1.54	0.73
1:B:138:GLY:O	1:B:180:LEU:HA	1.89	0.73
1:B:106:LEU:HD11	1:B:120:VAL:HG13	1.70	0.73
2:D:285:TYR:N	2:D:295:HIS:CD2	2.55	0.72
1:A:92:ILE:HD11	1:A:134:VAL:HB	1.72	0.72
2:D:18:ILE:CD1	2:D:131:LEU:HD21	2.14	0.72
2:D:10:GLY:HA3	2:D:13:ASN:O	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:178:TYR:O	1:B:185:ILE:HG22	1.90	0.72
2:D:27:LEU:HD21	2:D:130:TYR:CD2	2.26	0.71
1:A:171:LEU:HB3	1:A:191:THR:HG23	1.72	0.71
2:D:39:ILE:C	2:D:40:ILE:HG13	2.10	0.71
2:D:281:VAL:HG11	2:D:296:MET:CE	2.21	0.70
2:D:248:GLN:HE22	2:D:294:LYS:HD2	1.57	0.70
1:A:86:PHE:HE1	1:A:180:LEU:O	1.75	0.70
1:A:89:GLU:OE1	1:A:183:ARG:NH1	2.25	0.70
1:A:102:LEU:HD13	1:B:126:THR:HG21	1.73	0.70
2:D:105:VAL:CG2	2:D:110:MET:CE	2.70	0.69
1:B:104:ARG:HB3	1:B:104:ARG:HH11	1.58	0.69
2:D:181:LEU:HD22	2:D:181:LEU:N	2.08	0.69
2:D:142:VAL:HG11	2:D:146:PHE:HD2	1.58	0.69
2:D:171:GLU:O	2:D:174:ALA:HB3	1.91	0.69
2:D:73:ILE:HG23	2:D:74:GLY:H	1.58	0.69
2:D:313:VAL:CG2	2:D:348:MET:HE2	2.23	0.69
2:D:152:GLN:O	2:D:155:LYS:HB2	1.93	0.69
2:D:39:ILE:HD12	2:D:55:LYS:HB2	1.74	0.69
1:A:191:THR:O	1:A:191:THR:HG23	1.93	0.69
1:A:150:ASP:N	1:A:154:HIS:HD2	1.87	0.69
2:D:105:VAL:HG21	2:D:110:MET:CE	2.23	0.69
1:A:102:LEU:HD21	1:A:126:THR:HG22	1.74	0.68
2:D:351:LYS:O	2:D:354:ALA:N	2.28	0.67
2:D:78:GLN:O	2:D:79:ASP:HB2	1.95	0.67
1:B:90:LYS:O	1:B:93:ASN:HB2	1.94	0.67
2:D:146:PHE:O	2:D:151:ARG:HD2	1.95	0.67
1:B:101:SER:HA	1:B:104:ARG:HH12	1.58	0.67
2:D:73:ILE:CG2	2:D:74:GLY:N	2.57	0.67
2:D:40:ILE:HG23	2:D:50:VAL:HG22	1.77	0.67
1:B:72:LEU:O	1:B:76:THR:HG23	1.93	0.67
1:B:104:ARG:HB3	1:B:104:ARG:NH1	2.09	0.67
2:D:139:VAL:HG13	2:D:167:ARG:O	1.95	0.67
1:B:101:SER:HA	1:B:104:ARG:NH1	2.10	0.67
1:A:76:THR:O	1:A:80:ILE:HG13	1.96	0.66
2:D:317:ILE:HA	2:D:320:LEU:HB2	1.77	0.66
2:D:351:LYS:O	2:D:353:VAL:N	2.29	0.66
2:D:356:PHE:O	2:D:356:PHE:CD1	2.47	0.66
2:D:174:ALA:O	2:D:374:GLY:HA3	1.95	0.66
2:D:142:VAL:HG12	2:D:146:PHE:CE2	2.29	0.66
1:A:91:PHE:HD1	1:B:137:PHE:CE2	2.14	0.66
1:A:123:ILE:HG22	1:A:123:ILE:O	1.94	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:102:LEU:HD12	1:B:127:LEU:HD12	1.76	0.66
1:B:142:ILE:CG2	1:B:153:VAL:CG1	2.74	0.66
1:A:41:ASP:O	1:A:44:VAL:HG12	1.96	0.65
2:D:195:LEU:O	2:D:341:GLY:HA3	1.96	0.65
1:A:51:LEU:O	1:A:51:LEU:HG	1.96	0.65
1:A:135:ARG:HA	1:A:139:VAL:O	1.97	0.65
1:A:172:GLY:N	1:A:191:THR:HG22	2.11	0.65
1:A:128:LYS:O	1:A:132:ASP:HB2	1.96	0.64
2:D:159:ARG:HH12	2:D:164:GLU:HG2	1.60	0.64
1:A:147:VAL:CA	1:A:169:ASN:HD22	2.11	0.64
1:A:155:GLN:HE21	1:A:157:ILE:HD13	1.61	0.64
1:B:179:THR:HG22	1:B:184:THR:HA	1.79	0.64
1:A:174:MET:CE	2:D:261:ARG:HD2	2.28	0.64
1:A:174:MET:HE2	2:D:261:ARG:HD2	1.79	0.64
2:D:285:TYR:N	2:D:295:HIS:HD2	1.95	0.64
1:A:150:ASP:O	1:A:154:HIS:HB2	1.97	0.64
1:A:116:MET:O	1:A:116:MET:CG	2.42	0.64
1:B:106:LEU:HD21	1:B:124:GLU:CG	2.27	0.64
1:B:92:ILE:HD13	1:B:139:VAL:HG21	1.79	0.63
2:D:207:ILE:CD1	2:D:207:ILE:N	2.60	0.63
1:A:191:THR:O	1:A:191:THR:CG2	2.46	0.63
2:D:281:VAL:CG1	2:D:296:MET:HE2	2.29	0.63
2:D:285:TYR:H	2:D:295:HIS:HD2	1.42	0.63
1:B:149:LEU:HD13	1:B:192:VAL:CG2	2.28	0.63
1:B:141:VAL:CG1	1:B:178:TYR:CE1	2.77	0.63
1:A:92:ILE:HD13	1:A:139:VAL:HG21	1.78	0.63
2:D:31:GLU:OE1	2:D:31:GLU:N	2.31	0.63
2:D:263:LYS:O	2:D:267:GLU:HG3	1.99	0.63
2:D:379:GLY:O	2:D:380:GLY:O	2.17	0.62
1:A:92:ILE:CD1	1:A:139:VAL:CG2	2.76	0.62
2:D:40:ILE:CD1	2:D:115:ILE:HG22	2.29	0.62
1:B:102:LEU:HD12	1:B:127:LEU:CD1	2.28	0.62
2:D:105:VAL:HG23	2:D:110:MET:HE2	1.81	0.62
1:B:143:ALA:HB2	1:B:176:LYS:HG2	1.81	0.62
2:D:318:GLU:OE2	2:D:321:LYS:HE3	2.00	0.62
1:A:143:ALA:HB1	1:A:173:ILE:CD1	2.29	0.62
1:A:96:LEU:N	1:A:97:PRO:HD2	2.14	0.61
2:D:73:ILE:CG2	2:D:74:GLY:H	2.12	0.61
1:B:102:LEU:HB3	1:B:127:LEU:HD13	1.80	0.61
2:D:207:ILE:N	2:D:207:ILE:HD13	2.15	0.61
2:D:152:GLN:O	2:D:155:LYS:N	2.31	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:147:ASN:H	2:D:150:GLN:NE2	1.98	0.61
1:A:149:LEU:HD13	1:A:192:VAL:CG2	2.31	0.61
2:D:200:PHE:O	2:D:224:ASP:N	2.31	0.61
2:D:176:ALA:O	2:D:181:LEU:HD23	2.00	0.60
2:D:346:MET:O	2:D:349:VAL:HG22	2.01	0.60
2:D:133:GLU:OE1	2:D:134:PRO:O	2.19	0.60
1:A:147:VAL:N	1:A:169:ASN:HD22	1.99	0.60
1:A:172:GLY:H	1:A:191:THR:HG22	1.65	0.60
2:D:142:VAL:HG12	2:D:146:PHE:CD2	2.35	0.60
2:D:241:VAL:HG11	2:D:253:ARG:HG3	1.84	0.60
2:D:240:LEU:HD11	2:D:281:VAL:HG21	1.84	0.60
1:B:92:ILE:HD13	1:B:139:VAL:HG23	1.82	0.59
2:D:142:VAL:CG1	2:D:146:PHE:HD2	2.13	0.59
2:D:240:LEU:HD13	2:D:296:MET:HE1	1.84	0.59
2:D:331:VAL:CG1	2:D:356:PHE:O	2.46	0.59
1:A:171:LEU:O	1:A:171:LEU:HD12	2.02	0.59
1:A:91:PHE:HD1	1:B:137:PHE:CE1	2.21	0.59
2:D:105:VAL:HG21	2:D:110:MET:HE1	1.83	0.59
1:B:149:LEU:HD13	1:B:192:VAL:HG21	1.83	0.59
1:A:156:ALA:CB	2:D:60:THR:CG2	2.80	0.58
1:A:137:PHE:CZ	1:B:91:PHE:HA	2.37	0.58
1:A:151:PRO:O	2:D:56:ARG:NH1	2.35	0.58
2:D:252:LEU:CB	2:D:259:MET:HE3	2.33	0.58
1:A:155:GLN:NE2	2:D:264:GLU:OE2	2.36	0.58
2:D:143:PRO:HD2	2:D:146:PHE:CG	2.38	0.58
2:D:5:ILE:CG1	2:D:5:ILE:O	2.48	0.58
1:B:188:ALA:O	1:B:190:VAL:HG13	2.03	0.58
2:D:344:THR:HA	2:D:349:VAL:HG21	1.85	0.58
2:D:20:ASP:O	2:D:23:THR:HG23	2.04	0.58
2:D:295:HIS:HB3	3:D:935:HOH:O	2.03	0.57
2:D:201:ASP:HA	2:D:223:GLY:HA3	1.85	0.57
2:D:4:ILE:O	2:D:5:ILE:C	2.43	0.57
2:D:301:THR:O	2:D:302:ARG:C	2.41	0.57
2:D:77:PHE:CG	2:D:95:ALA:HB2	2.39	0.57
1:A:121:GLU:O	1:A:121:GLU:HG2	2.01	0.57
2:D:105:VAL:O	2:D:108:GLN:HG2	2.05	0.57
2:D:202:ILE:HG12	2:D:320:LEU:HD13	1.87	0.57
2:D:190:ILE:CG2	2:D:191:ALA:N	2.68	0.57
1:A:131:LEU:HD22	1:A:141:VAL:HG21	1.87	0.57
2:D:281:VAL:CG1	2:D:296:MET:CE	2.84	0.56
1:A:91:PHE:CD1	1:B:137:PHE:CE2	2.92	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:159:MET:HE3	1:B:194:LYS:HB3	1.86	0.56
2:D:147:ASN:OD1	2:D:150:GLN:HG3	2.06	0.56
2:D:278:GLN:HB3	2:D:301:THR:HG22	1.87	0.56
1:A:139:VAL:HG22	1:A:180:LEU:HD13	1.86	0.56
2:D:146:PHE:CA	2:D:150:GLN:NE2	2.64	0.56
1:A:72:LEU:O	1:A:76:THR:HG23	2.06	0.56
1:A:143:ALA:HB1	1:A:173:ILE:HD11	1.88	0.56
2:D:36:THR:HB	2:D:123:MET:CE	2.34	0.56
1:B:142:ILE:HD11	1:B:179:THR:CG2	2.35	0.56
1:B:59:ARG:HH11	1:B:59:ARG:HG2	1.71	0.56
1:A:137:PHE:CE2	1:B:91:PHE:HB2	2.41	0.56
1:A:137:PHE:N	1:A:137:PHE:CD1	2.74	0.56
2:D:305:LEU:O	2:D:308:LEU:HB2	2.05	0.56
2:D:311:ASP:O	2:D:315:ARG:HB2	2.06	0.56
2:D:193:TYR:HB3	2:D:339:LEU:HD23	1.87	0.55
1:B:165:VAL:O	1:B:166:ALA:C	2.44	0.55
1:A:92:ILE:HD13	1:A:139:VAL:CG2	2.35	0.55
2:D:74:GLY:HA3	2:D:150:GLN:HA	1.88	0.55
2:D:324:LEU:HD11	2:D:357:PHE:HE1	1.71	0.55
2:D:40:ILE:HD12	2:D:115:ILE:HG22	1.87	0.55
2:D:193:TYR:CE1	2:D:200:PHE:CE1	2.94	0.55
1:A:73:ARG:O	1:A:77:GLU:HG3	2.07	0.55
1:A:142:ILE:HG13	1:A:153:VAL:HG22	1.80	0.55
1:A:178:TYR:O	1:A:185:ILE:HG12	2.06	0.55
2:D:146:PHE:CD1	2:D:150:GLN:NE2	2.75	0.55
2:D:27:LEU:CD2	2:D:130:TYR:CD2	2.90	0.55
1:A:100:ASP:O	1:A:104:ARG:HB2	2.07	0.55
1:A:96:LEU:N	1:A:97:PRO:CD	2.70	0.55
1:A:69:MET:HB2	1:B:69:MET:HG2	1.88	0.54
1:A:76:THR:HB	1:B:76:THR:HG21	1.90	0.54
2:D:220:ALA:HB2	2:D:327:ALA:HB2	1.89	0.54
1:A:127:LEU:CD2	1:A:127:LEU:O	2.51	0.54
2:D:143:PRO:HD2	2:D:146:PHE:CD2	2.43	0.54
2:D:233:ASP:OD2	2:D:267:GLU:HG2	2.07	0.54
1:B:102:LEU:CD1	1:B:127:LEU:CD1	2.86	0.54
1:A:69:MET:HG3	1:A:69:MET:O	2.05	0.54
2:D:147:ASN:N	2:D:150:GLN:NE2	2.55	0.54
1:A:148:PRO:N	1:A:169:ASN:ND2	2.55	0.54
1:B:44:VAL:O	1:B:48:GLU:HG3	2.08	0.54
2:D:337:VAL:HG13	2:D:361:PRO:HA	1.90	0.54
2:D:229:GLY:O	2:D:232:PHE:N	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:LEU:HD21	1:A:126:THR:CG2	2.38	0.54
1:A:88:LEU:CD1	1:A:137:PHE:CD2	2.87	0.53
2:D:173:THR:HG22	2:D:177:LEU:CD1	2.39	0.53
1:B:141:VAL:CG1	1:B:178:TYR:HE1	2.18	0.53
1:B:159:MET:HE3	1:B:194:LYS:CB	2.39	0.53
1:B:59:ARG:NH1	1:B:59:ARG:HG2	2.23	0.53
2:D:192:VAL:O	2:D:202:ILE:HD13	2.09	0.53
1:B:102:LEU:CB	1:B:127:LEU:HD13	2.39	0.53
2:D:5:ILE:HA	2:D:18:ILE:HG22	1.90	0.53
2:D:5:ILE:HD12	2:D:16:VAL:HG21	1.89	0.53
2:D:190:ILE:HG22	2:D:191:ALA:N	2.22	0.53
2:D:205:ILE:CD1	2:D:207:ILE:CD1	2.84	0.53
2:D:207:ILE:HA	2:D:215:THR:O	2.09	0.53
1:A:158:ALA:CB	2:D:257:LEU:HD23	2.39	0.53
1:A:91:PHE:HE2	1:B:91:PHE:CE2	2.25	0.53
2:D:144:ALA:HB2	2:D:172:PRO:HG2	1.91	0.53
2:D:36:THR:HG22	2:D:36:THR:O	2.09	0.52
1:A:126:THR:OG1	1:B:101:SER:HB3	2.10	0.52
1:A:157:ILE:HD11	2:D:264:GLU:OE2	2.10	0.52
1:A:157:ILE:HG13	2:D:260:GLN:HG3	1.92	0.52
2:D:366:ASN:HB3	2:D:369:GLU:HB2	1.91	0.52
1:A:178:TYR:CB	1:A:185:ILE:HD11	2.37	0.52
1:B:142:ILE:HG12	1:B:153:VAL:HG13	1.92	0.52
2:D:224:ASP:OD1	2:D:225:THR:N	2.43	0.52
2:D:259:MET:HE1	2:D:262:LEU:HD23	1.91	0.52
2:D:329:LEU:HB3	2:D:333:ASP:CB	2.40	0.52
1:A:91:PHE:CD1	1:B:137:PHE:CD2	2.97	0.52
1:A:62:ILE:HD13	1:B:62:ILE:HD13	1.90	0.52
2:D:141:THR:CG2	2:D:375:ALA:HB2	2.40	0.52
2:D:208:ASP:CB	2:D:215:THR:CG2	2.88	0.52
1:A:142:ILE:HG21	1:A:153:VAL:HG22	1.91	0.52
2:D:349:VAL:HG23	2:D:350:GLN:N	2.24	0.52
2:D:105:VAL:HG23	2:D:110:MET:CE	2.38	0.51
1:A:72:LEU:O	1:A:72:LEU:HG	2.08	0.51
2:D:42:TYR:CD1	2:D:105:VAL:HG11	2.46	0.51
2:D:191:ALA:O	2:D:337:VAL:HG23	2.10	0.51
1:B:135:ARG:HH11	1:B:135:ARG:HG3	1.75	0.51
1:A:89:GLU:HB2	1:A:180:LEU:HD23	1.92	0.51
1:A:143:ALA:CB	1:A:173:ILE:CD1	2.88	0.51
1:A:70:GLU:OE1	1:A:70:GLU:HA	2.10	0.51
2:D:237:ILE:O	2:D:240:LEU:HB2	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:PHE:HD1	1:A:137:PHE:N	2.08	0.51
1:A:157:ILE:CD1	2:D:260:GLN:HG3	2.40	0.51
2:D:166:LYS:O	2:D:167:ARG:CB	2.59	0.51
2:D:171:GLU:N	2:D:172:PRO:CD	2.73	0.51
2:D:286:ILE:O	2:D:287:THR:HG23	2.12	0.50
2:D:190:ILE:HA	2:D:336:ASP:O	2.11	0.50
2:D:142:VAL:HG11	2:D:146:PHE:CD2	2.37	0.50
1:A:145:THR:HA	1:A:170:VAL:HG23	1.94	0.50
2:D:317:ILE:HD13	2:D:320:LEU:HD23	1.94	0.50
2:D:192:VAL:O	2:D:202:ILE:CD1	2.59	0.50
2:D:374:GLY:O	2:D:377:VAL:HG12	2.12	0.50
1:B:49:ALA:O	1:B:51:LEU:N	2.45	0.50
2:D:346:MET:HE1	2:D:349:VAL:HG13	1.94	0.50
2:D:193:TYR:CE1	2:D:200:PHE:CZ	3.00	0.50
1:A:40:ARG:NH1	1:A:40:ARG:CG	2.46	0.49
1:B:49:ALA:O	1:B:50:GLN:C	2.50	0.49
1:B:171:LEU:HD23	1:B:171:LEU:C	2.31	0.49
2:D:240:LEU:HD13	2:D:296:MET:CE	2.42	0.49
1:B:157:ILE:CD1	1:B:174:MET:HG3	2.30	0.48
2:D:252:LEU:HD12	2:D:252:LEU:HA	1.59	0.48
2:D:270:LYS:NZ	2:D:343:GLN:HE22	2.10	0.48
2:D:36:THR:HB	2:D:123:MET:HE1	1.94	0.48
2:D:278:GLN:HB3	2:D:301:THR:CG2	2.43	0.48
2:D:76:ARG:HA	2:D:100:ASP:HA	1.95	0.48
2:D:173:THR:HG22	2:D:177:LEU:HD11	1.95	0.48
2:D:112:PRO:N	2:D:113:PRO:CD	2.76	0.48
1:A:172:GLY:O	1:A:191:THR:HG22	2.13	0.48
1:A:57:ARG:HH22	1:B:55:GLN:HE22	1.62	0.48
2:D:94:ILE:HD12	2:D:94:ILE:C	2.34	0.48
1:A:157:ILE:HA	1:A:157:ILE:HD12	1.65	0.48
1:A:157:ILE:O	1:A:158:ALA:HB2	2.14	0.48
1:A:91:PHE:C	1:A:93:ASN:H	2.17	0.48
1:A:155:GLN:HE21	1:A:157:ILE:CD1	2.27	0.48
2:D:317:ILE:O	2:D:320:LEU:N	2.47	0.48
1:A:96:LEU:HD11	1:A:185:ILE:HG13	1.95	0.48
1:B:127:LEU:HD12	1:B:127:LEU:HA	1.67	0.48
1:B:135:ARG:CG	1:B:135:ARG:HH11	2.27	0.48
1:B:135:ARG:NH1	1:B:135:ARG:HG3	2.28	0.48
2:D:287:THR:HB	2:D:288:ALA:H	1.48	0.47
2:D:378:GLN:C	2:D:380:GLY:N	2.64	0.47
1:A:103:ASP:OD2	1:A:176:LYS:NZ	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:LEU:HD13	1:A:192:VAL:HG22	1.96	0.47
2:D:321:LYS:HG2	2:D:356:PHE:CE2	2.49	0.47
2:D:181:LEU:HG	2:D:190:ILE:HD11	1.96	0.47
1:A:91:PHE:CE2	1:B:91:PHE:CD2	3.02	0.47
2:D:205:ILE:O	2:D:205:ILE:HD12	2.15	0.47
2:D:168:ILE:O	2:D:168:ILE:HG23	2.14	0.47
2:D:255:ASP:C	2:D:255:ASP:OD1	2.53	0.47
1:A:153:VAL:CG2	1:A:153:VAL:O	2.45	0.47
2:D:240:LEU:HD22	2:D:298:ILE:HG13	1.96	0.47
2:D:171:GLU:HB2	2:D:172:PRO:HD3	1.95	0.47
1:B:70:GLU:HG3	1:B:70:GLU:O	2.14	0.47
2:D:143:PRO:HG2	2:D:146:PHE:CE1	2.49	0.47
1:A:157:ILE:HG12	1:A:189:MET:CE	2.45	0.47
1:A:78:LEU:O	1:A:79:ASP:C	2.52	0.47
1:A:100:ASP:O	1:A:103:ASP:OD1	2.32	0.47
2:D:135:VAL:CG1	2:D:136:THR:N	2.78	0.47
2:D:5:ILE:HD12	2:D:16:VAL:CG2	2.45	0.46
2:D:194:ASP:HB3	2:D:201:ASP:O	2.16	0.46
1:B:49:ALA:O	1:B:52:ALA:N	2.49	0.46
2:D:68:ALA:HB3	2:D:89:MET:CE	2.45	0.46
1:A:74:ARG:O	1:A:78:LEU:HG	2.15	0.46
1:B:65:VAL:O	1:B:65:VAL:HG12	2.15	0.46
1:A:40:ARG:CB	1:A:40:ARG:HH11	2.26	0.46
2:D:341:GLY:O	2:D:342:GLY:C	2.51	0.46
2:D:76:ARG:O	2:D:78:GLN:N	2.47	0.46
1:A:127:LEU:HD23	1:A:127:LEU:C	2.34	0.46
1:A:158:ALA:HB2	2:D:257:LEU:HD23	1.98	0.46
1:A:92:ILE:HG21	1:A:180:LEU:HD22	1.97	0.46
2:D:244:PHE:CZ	2:D:286:ILE:HG21	2.51	0.46
1:A:91:PHE:CD1	1:B:137:PHE:CE1	3.03	0.46
1:A:123:ILE:O	1:A:123:ILE:CG2	2.63	0.46
2:D:181:LEU:CD2	2:D:181:LEU:N	2.78	0.46
2:D:166:LYS:HB3	2:D:383:THR:HG21	1.98	0.46
2:D:18:ILE:CD1	2:D:131:LEU:CD2	2.86	0.46
2:D:27:LEU:HD13	2:D:127:ALA:HA	1.98	0.46
2:D:346:MET:HE3	2:D:346:MET:HB3	1.63	0.45
1:A:95:LEU:HD23	1:B:133:VAL:HG11	1.97	0.45
2:D:97:ASP:C	2:D:99:GLY:N	2.69	0.45
1:B:178:TYR:C	1:B:185:ILE:HG22	2.36	0.45
2:D:247:ASP:O	2:D:248:GLN:HG3	2.15	0.45
2:D:170:ASN:HB3	2:D:172:PRO:HD2	1.96	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:197:GLY:O	2:D:230:GLU:HB3	2.15	0.45
2:D:9:LEU:HD13	2:D:120:LEU:HD13	1.98	0.45
2:D:351:LYS:C	2:D:353:VAL:N	2.68	0.45
1:B:180:LEU:CD1	1:B:180:LEU:O	2.59	0.45
2:D:24:PRO:HB3	2:D:373:ILE:HG23	1.97	0.45
1:A:116:MET:HB2	1:B:116:MET:HE3	1.97	0.45
2:D:337:VAL:HG13	2:D:337:VAL:O	2.17	0.45
2:D:352:LYS:NZ	3:D:922:HOH:O	2.49	0.45
2:D:7:ILE:HG12	2:D:16:VAL:HG23	1.97	0.45
1:A:75:ARG:O	1:A:76:THR:C	2.55	0.45
1:A:149:LEU:HD13	1:A:192:VAL:HG21	1.98	0.45
2:D:207:ILE:HG22	2:D:207:ILE:O	2.17	0.45
2:D:96:ALA:HB2	2:D:102:TRP:CD1	2.51	0.45
2:D:4:ILE:HG22	2:D:138:ALA:HA	1.99	0.45
1:B:174:MET:SD	1:B:191:THR:HG23	2.57	0.45
1:A:91:PHE:CD2	1:B:91:PHE:CD2	3.05	0.45
2:D:204:ILE:HD13	2:D:324:LEU:HD23	1.98	0.45
1:A:127:LEU:CD2	1:A:127:LEU:C	2.85	0.45
1:A:91:PHE:CD1	1:B:137:PHE:CG	3.05	0.45
2:D:232:PHE:CD2	2:D:346:MET:HE2	2.51	0.45
1:A:150:ASP:OD1	1:A:152:ASN:HB2	2.17	0.44
1:A:50:GLN:O	1:A:53:GLU:HB2	2.18	0.44
2:D:147:ASN:N	2:D:150:GLN:HE21	2.15	0.44
1:A:48:GLU:HG3	1:B:47:LEU:HD11	2.00	0.44
1:B:88:LEU:HD11	1:B:137:PHE:HD2	1.82	0.44
2:D:146:PHE:HB3	2:D:151:ARG:HG2	1.98	0.44
2:D:378:GLN:C	2:D:380:GLY:H	2.19	0.44
1:A:157:ILE:HD12	2:D:260:GLN:CG	2.47	0.44
1:B:49:ALA:C	1:B:51:LEU:N	2.70	0.44
1:B:51:LEU:O	1:B:55:GLN:HB2	2.17	0.44
2:D:77:PHE:CE1	2:D:94:ILE:C	2.91	0.44
1:A:41:ASP:O	1:A:44:VAL:CG1	2.66	0.44
2:D:379:GLY:HA2	2:D:383:THR:HB	1.99	0.44
1:B:170:VAL:HG13	1:B:190:VAL:HB	1.98	0.44
2:D:124:LYS:NZ	2:D:128:GLU:OE2	2.48	0.44
2:D:115:ILE:O	2:D:118:GLU:HB2	2.17	0.44
1:B:165:VAL:O	1:B:165:VAL:HG23	2.17	0.44
2:D:363:LYS:HB3	2:D:363:LYS:HE2	1.89	0.44
1:B:96:LEU:N	1:B:97:PRO:HD2	2.33	0.44
2:D:12:THR:HG22	2:D:13:ASN:ND2	2.32	0.44
2:D:232:PHE:CD2	2:D:343:GLN:OE1	2.71	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:380:GLY:O	2:D:382:LEU:N	2.51	0.44
2:D:19:MET:SD	2:D:377:VAL:HG23	2.58	0.44
2:D:96:ALA:HB3	2:D:100:ASP:H	1.83	0.44
2:D:112:PRO:CB	2:D:113:PRO:HD3	2.47	0.43
1:B:171:LEU:CD2	1:B:171:LEU:O	2.48	0.43
2:D:78:GLN:O	2:D:79:ASP:CB	2.66	0.43
2:D:85:ASP:C	2:D:87:SER:N	2.69	0.43
1:A:141:VAL:HG22	1:A:178:TYR:CE2	2.54	0.43
1:A:73:ARG:HG2	1:B:72:LEU:HD21	2.00	0.43
1:B:102:LEU:CD1	1:B:127:LEU:HD13	2.47	0.43
2:D:353:VAL:HG12	2:D:361:PRO:HG3	2.00	0.43
1:A:50:GLN:O	1:A:53:GLU:N	2.51	0.43
1:A:91:PHE:CD1	1:B:137:PHE:CD1	3.07	0.43
1:B:180:LEU:O	1:B:181:ASN:HB2	2.18	0.43
2:D:380:GLY:C	2:D:382:LEU:N	2.71	0.43
1:A:157:ILE:CD1	2:D:260:GLN:CG	2.97	0.43
1:B:38:ASP:N	1:B:39:PRO:CD	2.81	0.43
1:A:91:PHE:CE1	1:B:137:PHE:CG	3.06	0.43
2:D:329:LEU:HB3	2:D:333:ASP:HB2	1.99	0.43
2:D:378:GLN:HE21	2:D:378:GLN:CA	2.21	0.43
1:A:157:ILE:HG12	1:A:189:MET:HE2	2.01	0.43
2:D:179:TYR:CD2	2:D:338:ILE:CG2	2.99	0.43
2:D:316:SER:OG	2:D:317:ILE:N	2.52	0.43
2:D:331:VAL:HG12	2:D:357:PHE:CD1	2.54	0.43
1:A:46:ASN:O	1:A:49:ALA:N	2.52	0.43
2:D:351:LYS:O	2:D:352:LYS:C	2.56	0.42
1:B:104:ARG:HH11	1:B:104:ARG:CB	2.29	0.42
1:A:91:PHE:CE1	1:B:137:PHE:CD1	3.07	0.42
2:D:343:GLN:HG3	2:D:343:GLN:O	2.19	0.42
1:A:167:PRO:HA	1:A:193:ALA:O	2.20	0.42
2:D:314:ASN:OD1	2:D:352:LYS:NZ	2.49	0.42
2:D:347:PRO:O	2:D:350:GLN:HB2	2.18	0.42
2:D:236:LEU:HD12	2:D:236:LEU:O	2.20	0.42
2:D:171:GLU:N	2:D:172:PRO:HD2	2.34	0.42
2:D:173:THR:O	2:D:177:LEU:HD12	2.18	0.42
1:A:106:LEU:HA	1:A:106:LEU:HD23	1.77	0.42
2:D:238:ASN:O	2:D:239:TYR:C	2.57	0.42
2:D:334:ILE:O	2:D:359:LYS:NZ	2.51	0.42
1:A:72:LEU:HD13	1:B:73:ARG:CZ	2.50	0.42
2:D:193:TYR:CD1	2:D:200:PHE:CE1	3.07	0.42
2:D:8:ASP:HB3	2:D:372:ALA:HB2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:252:LEU:O	2:D:259:MET:HG2	2.20	0.42
2:D:331:VAL:HG12	2:D:357:PHE:HD1	1.84	0.42
2:D:356:PHE:CG	2:D:356:PHE:O	2.72	0.42
1:A:171:LEU:HB3	1:A:191:THR:O	2.19	0.42
2:D:334:ILE:O	2:D:359:LYS:HE2	2.20	0.42
1:B:147:VAL:O	1:B:148:PRO:C	2.57	0.42
2:D:179:TYR:HB2	2:D:181:LEU:HD21	2.02	0.42
1:A:86:PHE:CE1	1:A:180:LEU:O	2.65	0.42
2:D:142:VAL:CG2	2:D:168:ILE:HD11	2.49	0.42
2:D:310:GLU:HA	2:D:348:MET:CE	2.39	0.42
2:D:174:ALA:CB	2:D:371:VAL:HG23	2.50	0.42
2:D:181:LEU:HG	2:D:190:ILE:CD1	2.50	0.41
1:A:58:GLU:HG2	1:A:59:ARG:N	2.35	0.41
1:A:143:ALA:HA	1:A:173:ILE:HD13	2.02	0.41
2:D:52:GLN:N	2:D:53:PRO:HD2	2.35	0.41
1:A:172:GLY:N	1:A:191:THR:CG2	2.80	0.41
1:B:182:GLY:O	1:B:183:ARG:C	2.56	0.41
1:A:171:LEU:HB2	1:A:193:ALA:HB2	2.02	0.41
2:D:111:ALA:O	2:D:114:GLN:HB2	2.20	0.41
2:D:330:SER:C	2:D:332:SER:N	2.73	0.41
1:A:65:VAL:O	1:B:65:VAL:HG11	2.20	0.41
2:D:192:VAL:HG22	2:D:338:ILE:CD1	2.50	0.41
1:A:95:LEU:C	1:A:97:PRO:HD2	2.40	0.41
2:D:85:ASP:O	2:D:87:SER:N	2.54	0.41
2:D:202:ILE:N	2:D:222:ASN:O	2.35	0.41
1:A:141:VAL:CG1	1:A:176:LYS:HG2	2.50	0.41
1:A:88:LEU:O	1:A:92:ILE:HG22	2.21	0.41
1:B:134:VAL:HG12	1:B:139:VAL:HB	1.92	0.41
1:A:137:PHE:HE2	1:B:91:PHE:HB2	1.85	0.41
2:D:160:ILE:O	2:D:160:ILE:CG2	2.67	0.41
2:D:294:LYS:HD3	2:D:294:LYS:HA	1.66	0.41
1:A:40:ARG:NH1	1:A:40:ARG:HB2	2.36	0.41
2:D:244:PHE:O	2:D:248:GLN:N	2.54	0.41
2:D:115:ILE:N	2:D:115:ILE:HD13	2.35	0.41
1:A:156:ALA:CB	2:D:60:THR:HG21	2.50	0.41
2:D:141:THR:HA	2:D:169:ILE:O	2.20	0.41
1:B:92:ILE:O	1:B:92:ILE:HG22	2.20	0.41
2:D:194:ASP:OD1	2:D:340:VAL:CG1	2.69	0.41
1:A:89:GLU:CD	1:A:183:ARG:HH11	2.24	0.40
1:A:62:ILE:HD13	1:B:62:ILE:CD1	2.49	0.40
2:D:68:ALA:HB3	2:D:89:MET:HE3	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:237:ILE:HD12	2:D:266:ALA:CB	2.52	0.40
1:B:65:VAL:CG1	1:B:65:VAL:O	2.69	0.40
2:D:240:LEU:CD1	2:D:262:LEU:HD11	2.51	0.40
2:D:205:ILE:HD13	2:D:207:ILE:CD1	2.32	0.40
2:D:329:LEU:HB3	2:D:333:ASP:CG	2.41	0.40
2:D:40:ILE:HD13	2:D:115:ILE:HG22	2.00	0.40
2:D:131:LEU:HG	2:D:131:LEU:H	1.63	0.40
2:D:284:PRO:O	2:D:285:TYR:C	2.60	0.40
1:B:92:ILE:CG2	1:B:139:VAL:HG21	2.52	0.40
2:D:52:GLN:HB3	2:D:53:PRO:CD	2.52	0.40
1:B:182:GLY:O	1:B:184:THR:N	2.55	0.40
2:D:330:SER:O	2:D:332:SER:N	2.54	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	154/197 (78%)	134 (87%)	17 (11%)	3 (2%)	10	32
1	B	147/197 (75%)	127 (86%)	18 (12%)	2 (1%)	14	42
2	D	370/383 (97%)	307 (83%)	45 (12%)	18 (5%)	3	8
All	All	671/777 (86%)	568 (85%)	80 (12%)	23 (3%)	5	16

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	196	LYS
2	D	167	ARG
2	D	342	GLY
2	D	352	LYS
1	B	143	ALA

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Mol	Chain	Res	Type
2	D	5	ILE
2	D	73	ILE
2	D	77	PHE
2	D	79	ASP
2	D	81	GLU
2	D	144	ALA
2	D	380	GLY
1	B	50	GLN
2	D	29	ASN
2	D	165	VAL
2	D	370	ALA
1	A	92	ILE
2	D	351	LYS
2	D	382	LEU
1	A	153	VAL
2	D	59	VAL
2	D	82	VAL
2	D	186	GLY

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	126/166 (76%)	109 (86%)	17 (14%)	5	14
1	B	118/166 (71%)	105 (89%)	13 (11%)	8	23
2	D	292/312 (94%)	246 (84%)	46 (16%)	3	9
All	All	536/644 (83%)	460 (86%)	76 (14%)	4	12

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	ARG
1	A	55	GLN
1	A	65	VAL
1	A	69	MET

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Mol	Chain	Res	Type
1	A	108	VAL
1	A	121	GLU
1	A	128	LYS
1	A	132	ASP
1	A	137	PHE
1	A	147	VAL
1	A	152	ASN
1	A	157	ILE
1	A	170	VAL
1	A	176	LYS
1	A	183	ARG
1	A	185	ILE
1	A	186	ARG
1	B	59	ARG
1	B	75	ARG
1	B	119	MET
1	B	128	LYS
1	B	129	SER
1	B	130	MET
1	B	135	ARG
1	B	159	MET
1	B	162	SER
1	B	171	LEU
1	B	180	LEU
1	B	186	ARG
1	B	194	LYS
2	D	5	ILE
2	D	23	THR
2	D	34	ARG
2	D	35	THR
2	D	43	THR
2	D	52	GLN
2	D	57	GLN
2	D	60	THR
2	D	87	SER
2	D	103	VAL
2	D	106	LYS
2	D	110	MET
2	D	120	LEU
2	D	131	LEU
2	D	133	GLU
2	D	142	VAL

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Mol	Chain	Res	Type
2	D	146	PHE
2	D	168	ILE
2	D	181	LEU
2	D	202	ILE
2	D	206	GLU
2	D	207	ILE
2	D	219	LEU
2	D	221	THR
2	D	227	LEU
2	D	230	GLU
2	D	236	LEU
2	D	252	LEU
2	D	257	LEU
2	D	261	ARG
2	D	268	LYS
2	D	275	SER
2	D	280	ASP
2	D	297	ASN
2	D	301	THR
2	D	316	SER
2	D	320	LEU
2	D	329	LEU
2	D	338	ILE
2	D	346	MET
2	D	352	LYS
2	D	359	LYS
2	D	362	ARG
2	D	363	LYS
2	D	377	VAL
2	D	378	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	GLN
1	A	154	HIS
1	A	155	GLN
1	A	169	ASN
1	B	84	HIS
1	B	146	ASN
2	D	13	ASN
2	D	98	ASN

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Mol	Chain	Res	Type
2	D	150	GLN
2	D	248	GLN
2	D	295	HIS
2	D	297	ASN
2	D	343	GLN
2	D	378	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 carbohydrates 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	158/197 (80%)	-0.13	2 (1%) 79 71	27, 44, 64, 71	0
1	B	151/197 (76%)	-0.32	2 (1%) 79 71	30, 51, 72, 87	0
2	D	376/383 (98%)	-0.09	11 (2%) 55 43	27, 52, 79, 98	0
All	All	685/777 (88%)	-0.15	15 (2%) 65 54	27, 51, 73, 98	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	185	THR	4.7
1	B	116	MET	3.5
2	D	381	VAL	3.1
2	D	88	ILE	3.0
1	A	119	MET	2.7
2	D	216	PHE	2.7
2	D	305	LEU	2.7
2	D	215	THR	2.6
2	D	181	LEU	2.6
2	D	182	ASP	2.6
2	D	166	LYS	2.5
1	B	40	ARG	2.2
2	D	131	LEU	2.1
1	A	116	MET	2.1
2	D	201	ASP	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.