



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 05:00 PM UTC

PDB ID : 3K53 / pdb\_00003k53  
Title : Crystal Structure of NFeoB from *P. furiosus*  
Authors : Eng, E.T.; Dong, G.; Unger, V.M.  
Deposited on : 2009-10-06  
Resolution : 2.70 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 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

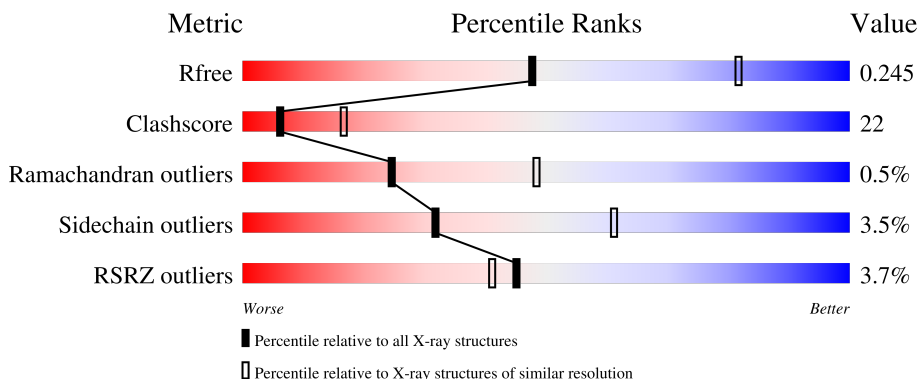
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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))
$R_{free}$	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	271	 4% 55% 40% ..
1	B	271	 3% 55% 41% ..
1	C	271	 4% 58% 37% ..
1	D	271	 4% 59% 37% .

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8696 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 Ferrous iron transport protein b.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	268	Total 2163	C 1401	N 362	O 391	S 9	0	0	0
1	B	265	Total 2139	C 1386	N 356	O 388	S 9	0	0	0
1	C	267	Total 2159	C 1397	N 360	O 393	S 9	0	0	0
1	D	270	Total 2180	C 1412	N 366	O 393	S 9	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q8U2H8
A	2	VAL	-	expression tag	UNP Q8U2H8
B	1	MET	-	expression tag	UNP Q8U2H8
B	2	VAL	-	expression tag	UNP Q8U2H8
C	1	MET	-	expression tag	UNP Q8U2H8
C	2	VAL	-	expression tag	UNP Q8U2H8
D	1	MET	-	expression tag	UNP Q8U2H8
D	2	VAL	-	expression tag	UNP Q8U2H8

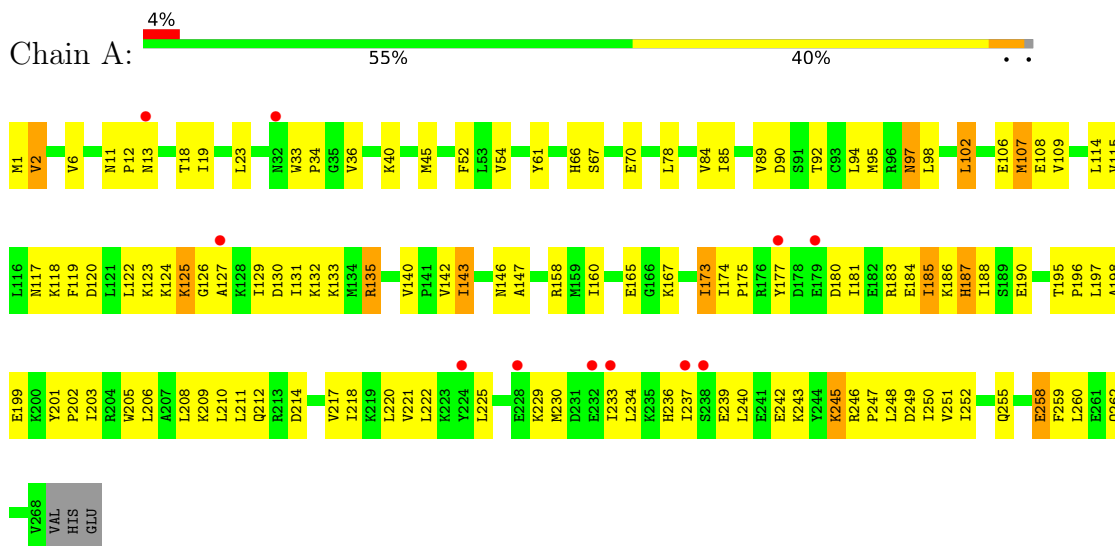
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total 3	O 3	0	0
2	B	11	Total 11	O 11	0	0
2	C	20	Total 20	O 20	0	0
2	D	21	Total 21	O 21	0	0

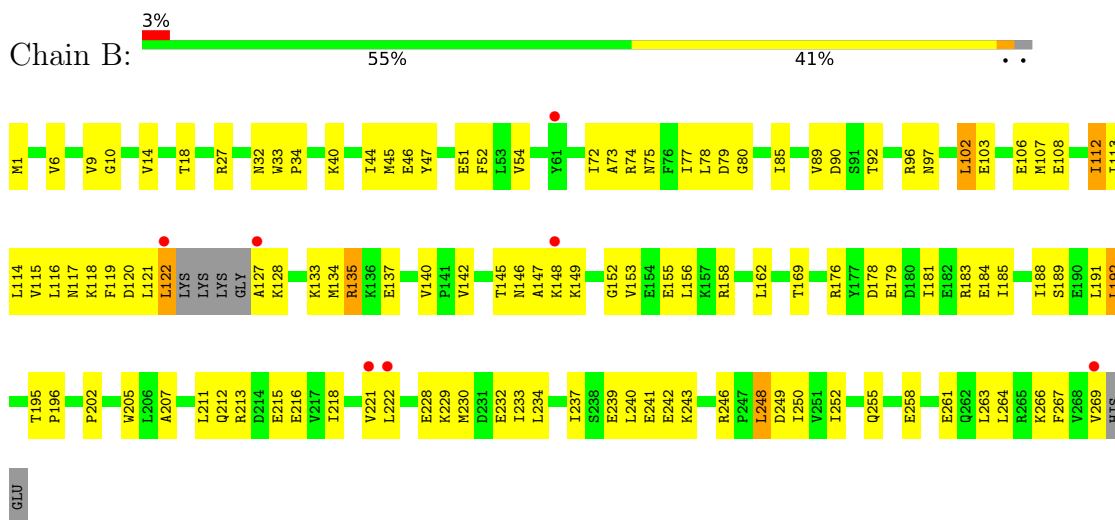
### 3 Residue-property plots [i](#)

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: Ferrous iron transport protein b

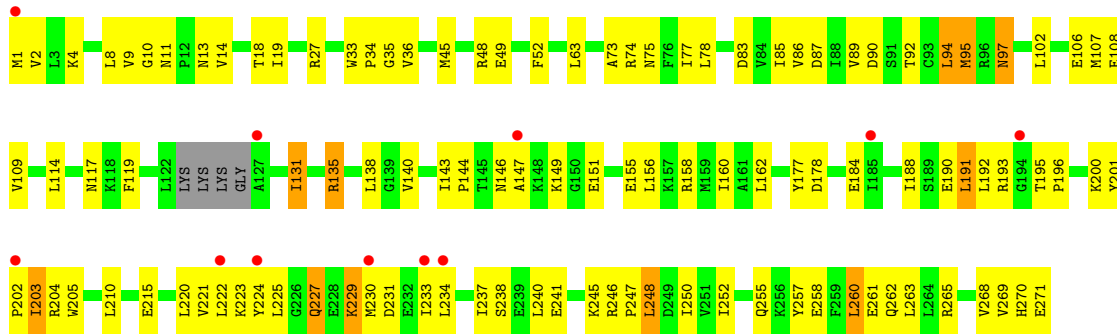


- Molecule 1: Ferrous iron transport protein b

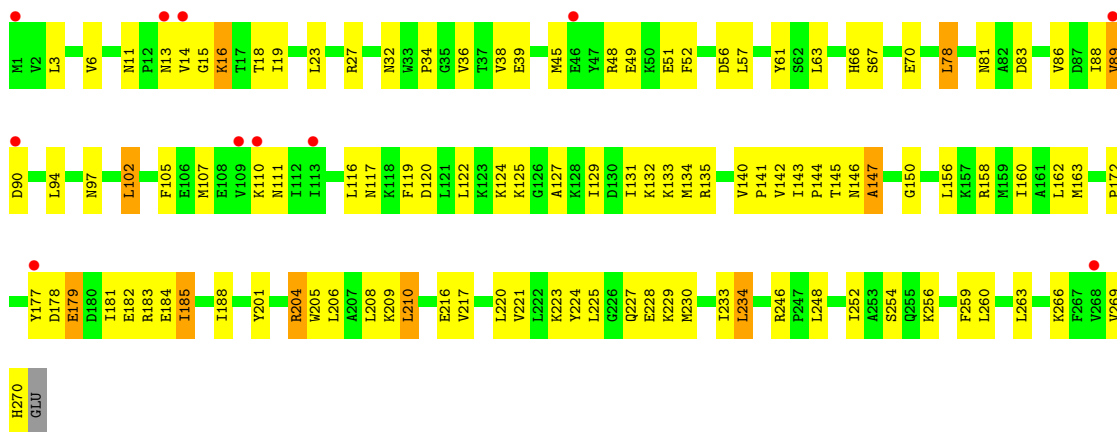


- Molecule 1: Ferrous iron transport protein b





● Molecule 1: Ferrous iron transport protein b



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	55.26Å 106.11Å 253.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.70 50.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	94.1 (50.00-2.70) 94.1 (50.00-2.70)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.84 (at 2.51Å)	Xtrriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.230 , 0.286 0.242 , 0.245	Depositor DCC
$R_{free}$ test set	3890 reflections (3.97%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	55.6	Xtrriage
Anisotropy	0.685	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 42.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	8696	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	72.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% of the height of the origin peak. No significant pseudotranslation is detected.*

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

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

## 5 Model quality i

### 5.1 Standard geometry i

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z  > 5$	RMSZ	# $ Z  > 5$
1	A	0.48	0/2195	1.00	9/2955 (0.3%)
1	B	0.44	0/2170	0.98	7/2924 (0.2%)
1	C	0.47	0/2191	0.97	11/2951 (0.4%)
1	D	0.48	0/2213	1.01	14/2980 (0.5%)
All	All	0.47	0/8769	0.99	41/11810 (0.3%)

There are no bond length outliers.

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	14	VAL	N-CA-C	-9.33	104.05	113.10
1	B	248	LEU	N-CA-C	8.94	121.99	111.71
1	D	143	ILE	N-CA-C	8.67	115.45	107.56
1	D	14	VAL	N-CA-C	-7.82	102.44	113.07
1	D	89	VAL	N-CA-C	7.25	118.34	107.75
1	D	110	LYS	N-CA-C	7.00	118.99	111.36
1	A	94	LEU	N-CA-C	6.83	119.30	111.11
1	C	131	ILE	N-CA-C	6.82	117.32	110.23
1	C	19	ILE	N-CA-C	-6.66	104.00	110.72
1	D	16	LYS	N-CA-C	-6.63	104.13	111.82
1	C	94	LEU	N-CA-C	6.57	119.04	111.02
1	B	240	LEU	N-CA-C	-6.17	104.56	111.28
1	C	33	TRP	CA-C-N	6.13	126.14	119.89
1	C	33	TRP	C-N-CA	6.13	126.14	119.89
1	B	216	GLU	N-CA-C	6.09	117.71	111.14
1	D	179	GLU	N-CA-C	6.00	119.43	111.75
1	C	97	ASN	N-CA-C	5.84	117.72	111.36
1	C	14	VAL	N-CA-C	-5.78	107.49	113.10
1	D	227	GLN	N-CA-C	-5.74	105.38	112.38
1	C	227	GLN	N-CA-C	-5.62	105.92	112.89
1	A	258	GLU	N-CA-C	-5.58	105.19	111.28
1	B	107	MET	N-CA-C	-5.58	106.55	113.19
1	A	185	ILE	N-CA-C	-5.47	105.20	110.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	191	LEU	N-CA-C	-5.44	106.95	113.21
1	A	97	ASN	N-CA-C	5.44	117.91	111.33
1	A	143	ILE	N-CA-C	5.44	113.93	107.73
1	A	66	HIS	N-CA-C	5.32	122.14	110.80
1	D	143	ILE	CA-C-N	5.28	125.44	119.90
1	D	143	ILE	C-N-CA	5.28	125.44	119.90
1	D	256	LYS	N-CA-C	-5.27	105.58	112.23
1	D	145	THR	N-CA-C	5.25	117.05	109.07
1	B	72	ILE	N-CA-C	-5.22	105.45	110.72
1	D	147	ALA	N-CA-C	5.20	116.76	111.14
1	D	66	HIS	N-CA-C	5.16	119.43	113.18
1	A	107	MET	N-CA-C	-5.14	107.03	113.15
1	C	248	LEU	N-CA-C	5.14	117.55	111.33
1	A	2	VAL	N-CA-C	-5.13	104.45	110.05
1	A	19	ILE	N-CA-C	-5.08	105.59	110.72
1	C	13	ASN	N-CA-C	5.02	118.32	111.39
1	D	185	ILE	N-CA-C	-5.02	105.65	110.72
1	B	192	LEU	N-CA-C	-5.02	107.83	114.31

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	2163	0	2302	128	0
1	B	2139	0	2268	100	0
1	C	2159	0	2281	102	0
1	D	2180	0	2318	101	0
2	A	3	0	0	1	0
2	B	11	0	0	0	0
2	C	20	0	0	0	0
2	D	21	0	0	2	0
All	All	8696	0	9169	399	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 22.

All (399) 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:23:LEU:HD22	1:A:45:MET:HE2	1.39	1.03
1:C:195:THR:HB	1:C:225:LEU:HD11	1.41	1.03
1:A:245:LYS:H	1:A:245:LYS:HD3	1.21	1.01
1:A:78:LEU:HG	1:A:107:MET:HE3	1.46	0.98
1:A:245:LYS:HE2	1:A:246:ARG:HG3	1.51	0.90
1:C:195:THR:HB	1:C:196:PRO:HD2	1.60	0.84
1:B:213:ARG:HH12	1:B:234:LEU:HD22	1.44	0.83
1:A:245:LYS:HD3	1:A:245:LYS:N	1.94	0.82
1:A:245:LYS:HG3	1:B:179:GLU:HG3	1.61	0.82
1:D:177:TYR:CE1	1:D:252:ILE:HG23	2.14	0.82
1:A:201:TYR:CE2	1:A:220:LEU:HD21	2.17	0.79
1:B:228:GLU:HG3	1:B:229:LYS:H	1.48	0.79
1:B:120:ASP:C	1:B:121:LEU:HD12	2.06	0.78
1:D:45:MET:HE1	1:D:160:ILE:HG21	1.65	0.78
1:B:122:LEU:HB3	1:B:127:ALA:N	1.99	0.78
1:C:229:LYS:HE3	1:C:229:LYS:HA	1.64	0.78
1:D:225:LEU:HD13	1:D:229:LYS:HG2	1.64	0.78
1:C:11:ASN:HD21	1:C:97:ASN:HA	1.49	0.77
1:A:221:VAL:HG11	1:A:230:MET:SD	2.25	0.77
1:B:212:GLN:HE21	1:B:249:ASP:HA	1.50	0.77
1:C:90:ASP:H	1:C:97:ASN:ND2	1.83	0.76
1:B:213:ARG:NH1	1:B:234:LEU:HD22	2.00	0.75
1:D:122:LEU:HG	1:D:127:ALA:HB3	1.68	0.75
1:C:223:LYS:HG3	1:C:224:TYR:H	1.52	0.75
1:D:183:ARG:HG2	1:D:183:ARG:HH11	1.52	0.74
1:C:146:ASN:HD21	1:C:149:LYS:HD3	1.52	0.73
1:D:89:VAL:HG13	1:D:97:ASN:HB3	1.71	0.73
1:B:146:ASN:ND2	1:B:149:LYS:HD3	2.04	0.72
1:B:184:GLU:O	1:B:188:ILE:HG12	1.89	0.72
1:A:177:TYR:CE1	1:A:252:ILE:HG23	2.24	0.72
1:A:165:GLU:OE1	1:A:167:LYS:HE3	1.89	0.72
1:A:78:LEU:HG	1:A:107:MET:CE	2.20	0.71
1:C:191:LEU:HD23	1:C:233:ILE:HD13	1.70	0.71
1:C:222:LEU:HD22	1:C:227:GLN:HG2	1.71	0.71
1:D:221:VAL:HB	1:D:230:MET:HE1	1.73	0.70
1:A:262:GLN:HE22	1:B:246:ARG:HE	1.38	0.70
1:C:195:THR:HG21	1:C:225:LEU:HD21	1.74	0.70
1:B:34:PRO:HG2	1:D:34:PRO:HG2	1.72	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:18:THR:CG2	1:D:147:ALA:HB1	2.21	0.70
1:B:211:LEU:HB3	1:B:248:LEU:HD23	1.73	0.70
1:C:230:MET:O	1:C:234:LEU:HG	1.92	0.69
1:D:124:LYS:HZ1	1:D:125:LYS:HE2	1.55	0.69
1:A:6:VAL:HG22	1:A:84:VAL:HB	1.74	0.69
1:A:117:ASN:HD21	1:A:147:ALA:H	1.39	0.69
1:C:231:ASP:HA	1:C:234:LEU:HD12	1.74	0.69
1:D:131:ILE:HD13	1:D:144:PRO:CG	2.22	0.69
1:C:230:MET:HE3	1:C:234:LEU:HD21	1.73	0.69
1:B:44:ILE:HG12	1:D:3:LEU:HD22	1.73	0.69
1:A:95:MET:CE	1:A:260:LEU:HB3	2.22	0.69
1:D:177:TYR:CE2	1:D:208:LEU:HD21	2.28	0.68
1:B:46:GLU:OE1	1:B:51:GLU:HG2	1.93	0.68
1:C:196:PRO:HD2	1:C:225:LEU:HD11	1.74	0.67
1:A:209:LYS:HE2	1:A:209:LYS:HA	1.76	0.67
1:D:89:VAL:HG13	1:D:97:ASN:CB	2.24	0.67
1:D:16:LYS:HG3	1:D:88:ILE:HD12	1.75	0.67
1:C:77:ILE:HD12	1:C:85:ILE:HD13	1.76	0.66
1:C:85:ILE:HD12	1:C:109:VAL:HG11	1.75	0.66
1:A:180:ASP:OD2	1:A:243:LYS:HD3	1.94	0.66
1:D:201:TYR:CE2	1:D:220:LEU:HD21	2.30	0.66
1:A:248:LEU:HA	1:A:251:VAL:HG12	1.76	0.66
1:C:89:VAL:HG13	1:C:97:ASN:HB3	1.78	0.66
1:D:27:ARG:HG2	1:D:27:ARG:HH11	1.61	0.66
1:B:228:GLU:HG3	1:B:229:LYS:N	2.12	0.65
1:A:258:GLU:OE2	1:B:246:ARG:HD2	1.96	0.65
1:D:188:ILE:HD12	1:D:233:ILE:HG23	1.77	0.65
1:C:223:LYS:HG3	1:C:224:TYR:N	2.11	0.64
1:B:181:ILE:O	1:B:185:ILE:HG12	1.97	0.64
1:C:87:ASP:HB3	1:C:114:LEU:HD12	1.79	0.64
1:C:45:MET:HE2	1:C:52:PHE:CD2	2.32	0.64
1:D:63:LEU:HD13	1:D:107:MET:HE2	1.80	0.64
1:A:245:LYS:H	1:A:245:LYS:CD	2.04	0.64
1:B:120:ASP:O	1:B:121:LEU:HD12	1.97	0.64
1:C:117:ASN:HD21	1:C:147:ALA:H	1.44	0.64
1:C:135:ARG:HD3	1:C:135:ARG:C	2.23	0.64
1:C:262:GLN:HE22	1:D:246:ARG:NH1	1.94	0.64
1:A:158:ARG:HH11	1:A:158:ARG:HG2	1.63	0.64
1:D:188:ILE:HG21	1:D:210:LEU:HD23	1.80	0.64
1:A:95:MET:HE1	1:A:260:LEU:HB3	1.80	0.63
1:C:146:ASN:ND2	1:C:149:LYS:HD3	2.14	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:188:ILE:O	1:B:192:LEU:HB2	1.98	0.62
1:C:131:ILE:HD13	1:C:144:PRO:HG3	1.81	0.62
1:A:201:TYR:HB2	1:A:206:LEU:HD11	1.81	0.62
1:B:137:GLU:OE2	1:B:266:LYS:HD3	1.99	0.62
1:A:67:SER:OG	1:A:70:GLU:HG3	1.99	0.62
1:D:131:ILE:HD13	1:D:144:PRO:HG2	1.81	0.62
1:C:234:LEU:C	1:C:237:ILE:HG22	2.24	0.62
1:D:184:GLU:O	1:D:188:ILE:HG12	2.00	0.62
1:C:230:MET:HE3	1:C:234:LEU:CD2	2.30	0.62
1:C:200:LYS:HG2	1:C:201:TYR:CE2	2.35	0.61
1:A:119:PHE:CZ	1:A:131:ILE:HD11	2.35	0.61
1:A:135:ARG:HG2	1:A:140:VAL:O	2.01	0.61
1:B:135:ARG:HD3	1:B:135:ARG:C	2.25	0.61
1:A:247:PRO:HD2	1:A:250:ILE:HD12	1.83	0.60
1:B:192:LEU:HD11	1:B:221:VAL:HG13	1.83	0.60
1:A:85:ILE:HD12	1:A:109:VAL:HG11	1.83	0.60
1:A:262:GLN:NE2	1:B:246:ARG:HE	1.99	0.60
1:D:32:ASN:ND2	1:D:36:VAL:O	2.29	0.60
1:B:47:TYR:OH	1:B:158:ARG:HG3	2.02	0.60
1:B:239:GLU:O	1:B:242:GLU:HB2	2.01	0.60
1:B:222:LEU:HD13	1:B:230:MET:HG2	1.82	0.60
1:A:222:LEU:HD21	1:A:230:MET:HG2	1.84	0.60
1:D:18:THR:HG22	1:D:147:ALA:HB1	1.83	0.60
1:A:108:GLU:CG	1:A:202:PRO:HB3	2.32	0.60
1:D:116:LEU:HD11	1:D:134:MET:HE1	1.82	0.60
1:B:32:ASN:H	1:D:81:ASN:HD21	1.49	0.60
1:C:257:TYR:O	1:C:261:GLU:HG3	2.02	0.59
1:C:177:TYR:CE1	1:C:252:ILE:HG23	2.38	0.59
1:D:117:ASN:HD21	1:D:147:ALA:H	1.51	0.59
1:D:188:ILE:CG2	1:D:210:LEU:HD23	2.32	0.59
1:A:225:LEU:HD23	1:A:229:LYS:HG2	1.84	0.59
1:B:215:GLU:O	1:B:218:ILE:HG22	2.03	0.58
1:C:74:ARG:HG3	1:C:75:ASN:N	2.17	0.58
1:D:158:ARG:O	1:D:162:LEU:HD13	2.03	0.58
1:C:270:HIS:O	1:C:271:GLU:HB2	2.04	0.58
1:A:221:VAL:CG1	1:A:230:MET:SD	2.91	0.57
1:C:190:GLU:C	1:C:192:LEU:H	2.11	0.57
1:A:184:GLU:HG3	1:A:240:LEU:HB2	1.87	0.57
1:C:193:ARG:HE	1:C:203:ILE:CD1	2.17	0.57
1:C:223:LYS:HG3	1:C:224:TYR:HD1	1.69	0.57
1:B:89:VAL:HG13	1:B:97:ASN:HB3	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:269:VAL:O	1:D:270:HIS:HB2	2.04	0.57
1:B:89:VAL:HG13	1:B:97:ASN:CB	2.35	0.57
1:D:225:LEU:CD1	1:D:229:LYS:HG2	2.34	0.57
1:D:13:ASN:HA	2:D:278:HOH:O	2.04	0.56
1:C:270:HIS:HD2	1:D:61:TYR:CE2	2.24	0.56
1:A:34:PRO:O	1:A:36:VAL:HG22	2.04	0.56
1:B:134:MET:HA	1:B:267:PHE:CE1	2.40	0.56
1:A:45:MET:HE1	1:A:160:ILE:CD1	2.36	0.56
1:B:73:ALA:O	1:B:77:ILE:HG12	2.06	0.56
1:B:90:ASP:OD1	1:B:118:LYS:HD2	2.06	0.56
1:A:209:LYS:CB	1:A:217:VAL:HG21	2.35	0.56
1:A:85:ILE:HD12	1:A:109:VAL:CG1	2.36	0.56
1:A:92:THR:O	1:B:96:ARG:NH2	2.39	0.55
1:D:16:LYS:HE3	1:D:57:LEU:O	2.07	0.55
1:D:177:TYR:CZ	1:D:252:ILE:HG23	2.40	0.55
1:C:4:LYS:HB2	1:C:52:PHE:HD1	1.71	0.55
1:A:245:LYS:CG	1:B:179:GLU:HG3	2.33	0.55
1:C:4:LYS:HB2	1:C:52:PHE:CD1	2.42	0.55
1:C:156:LEU:O	1:C:160:ILE:HG13	2.06	0.55
1:A:209:LYS:HB3	1:A:217:VAL:HG21	1.88	0.55
1:C:90:ASP:OD1	1:C:92:THR:OG1	2.23	0.55
1:C:234:LEU:O	1:C:237:ILE:HG22	2.07	0.55
1:B:176:ARG:HH11	1:B:176:ARG:HG2	1.71	0.55
1:A:230:MET:O	1:A:234:LEU:HG	2.07	0.55
1:C:77:ILE:HD12	1:C:85:ILE:CD1	2.37	0.55
1:C:237:ILE:O	1:C:241:GLU:HG3	2.07	0.55
1:D:216:GLU:HG3	2:D:272:HOH:O	2.06	0.55
1:A:205:TRP:CZ2	1:A:209:LYS:HD2	2.42	0.54
1:D:67:SER:OG	1:D:70:GLU:HG3	2.06	0.54
1:C:191:LEU:CD2	1:C:233:ILE:HD13	2.35	0.54
1:D:124:LYS:HZ1	1:D:125:LYS:CE	2.20	0.54
1:B:147:ALA:C	1:B:149:LYS:H	2.15	0.54
1:C:146:ASN:HB3	1:C:151:GLU:HB2	1.89	0.54
1:A:142:VAL:C	1:A:143:ILE:HD13	2.33	0.54
1:A:245:LYS:HG2	1:A:246:ARG:N	2.23	0.54
1:B:202:PRO:HG2	1:B:205:TRP:HB3	1.90	0.54
1:C:78:LEU:HD11	1:C:205:TRP:CH2	2.43	0.54
1:A:173:ILE:HD13	1:A:174:ILE:N	2.22	0.54
1:C:1:MET:HG2	1:C:2:VAL:O	2.06	0.54
1:A:175:PRO:HA	1:A:259:PHE:CE1	2.43	0.53
1:B:183:ARG:HG3	1:B:183:ARG:HH11	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:LYS:HE2	1:A:127:ALA:HB2	1.90	0.53
1:C:262:GLN:NE2	1:D:246:ARG:NH1	2.56	0.53
1:C:270:HIS:O	1:C:271:GLU:CB	2.57	0.53
1:B:51:GLU:OE1	1:D:27:ARG:NH2	2.42	0.53
1:D:122:LEU:HD21	1:D:129:ILE:HD12	1.91	0.53
1:A:195:THR:HG21	1:A:225:LEU:HD21	1.91	0.52
1:D:131:ILE:HD13	1:D:144:PRO:HG3	1.90	0.52
1:A:187:HIS:O	1:A:190:GLU:N	2.42	0.52
1:B:249:ASP:OD1	1:B:250:ILE:N	2.43	0.52
1:C:245:LYS:CD	1:D:179:GLU:HG2	2.40	0.52
1:A:117:ASN:ND2	1:A:118:LYS:H	2.07	0.52
1:C:191:LEU:HD11	1:C:229:LYS:HE2	1.91	0.52
1:B:195:THR:HG22	1:B:196:PRO:HD2	1.91	0.52
1:D:102:LEU:HD21	1:D:259:PHE:HD2	1.74	0.52
1:A:195:THR:HB	1:A:196:PRO:HD2	1.90	0.52
1:B:237:ILE:O	1:B:241:GLU:HG3	2.09	0.52
1:A:158:ARG:HG2	1:A:158:ARG:NH1	2.24	0.51
1:D:122:LEU:CD2	1:D:129:ILE:HD12	2.40	0.51
1:B:116:LEU:O	1:B:119:PHE:HB2	2.10	0.51
1:B:45:MET:HE2	1:B:52:PHE:CD2	2.45	0.51
1:A:18:THR:CG2	1:A:147:ALA:HB1	2.40	0.51
1:A:67:SER:HG	1:A:70:GLU:HG3	1.75	0.51
1:A:130:ASP:OD2	1:A:133:LYS:HD2	2.10	0.51
1:A:173:ILE:HD12	1:A:259:PHE:HZ	1.76	0.51
1:A:197:LEU:C	1:A:199:GLU:H	2.17	0.51
1:C:48:ARG:O	1:C:49:GLU:HB2	2.08	0.51
1:B:6:VAL:HB	1:B:54:VAL:HG22	1.92	0.51
1:C:143:ILE:HD12	1:C:155:GLU:HG3	1.92	0.51
1:D:45:MET:HE2	1:D:52:PHE:CD2	2.46	0.51
1:C:224:TYR:O	1:C:225:LEU:HD12	2.11	0.51
1:A:214:ASP:OD1	1:A:217:VAL:HG23	2.10	0.51
1:D:183:ARG:HG2	1:D:183:ARG:NH1	2.23	0.51
1:B:118:LYS:O	1:B:121:LEU:HD13	2.10	0.51
1:B:261:GLU:O	1:B:261:GLU:HG2	2.11	0.51
1:C:247:PRO:HG2	1:C:250:ILE:HG12	1.92	0.51
1:D:45:MET:HE1	1:D:160:ILE:CG2	2.38	0.51
1:B:77:ILE:HD12	1:B:85:ILE:HD12	1.93	0.50
1:D:209:LYS:CB	1:D:217:VAL:HG21	2.42	0.50
1:D:11:ASN:ND2	1:D:97:ASN:CG	2.69	0.50
1:A:6:VAL:O	1:A:54:VAL:HA	2.11	0.50
1:A:123:LYS:HZ3	1:A:129:ILE:H	1.60	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:223:LYS:C	1:D:224:TYR:HD1	2.19	0.50
1:A:236:HIS:O	1:A:239:GLU:HB3	2.11	0.50
1:D:210:LEU:HD13	1:D:217:VAL:HG12	1.94	0.50
1:A:13:ASN:ND2	1:B:122:LEU:HD12	2.27	0.50
1:A:45:MET:HE1	1:A:160:ILE:HD13	1.93	0.50
1:A:108:GLU:HG2	1:A:202:PRO:HB3	1.92	0.50
1:A:218:ILE:HG23	1:A:230:MET:SD	2.52	0.50
1:C:201:TYR:CE2	1:C:220:LEU:HD21	2.47	0.50
1:A:119:PHE:O	1:A:122:LEU:HB3	2.11	0.49
1:A:198:ALA:HB1	1:A:203:ILE:HG12	1.94	0.49
1:C:73:ALA:O	1:C:77:ILE:HG12	2.12	0.49
1:A:114:LEU:HD23	1:A:115:VAL:N	2.28	0.49
1:A:124:LYS:C	1:A:126:GLY:N	2.68	0.49
1:D:122:LEU:HD21	1:D:129:ILE:CD1	2.43	0.49
1:D:260:LEU:HD23	1:D:263:LEU:HD12	1.95	0.49
1:C:195:THR:CB	1:C:225:LEU:HD11	2.26	0.49
1:A:230:MET:O	1:A:230:MET:HE3	2.13	0.49
1:B:230:MET:O	1:B:234:LEU:HG	2.12	0.49
1:D:210:LEU:HD13	1:D:217:VAL:CG1	2.43	0.49
1:A:13:ASN:CG	1:B:122:LEU:HD12	2.37	0.49
1:B:9:VAL:HG22	1:B:10:GLY:N	2.28	0.49
1:A:258:GLU:HG2	1:B:246:ARG:CD	2.43	0.49
1:D:117:ASN:HD21	1:D:147:ALA:N	2.10	0.49
1:A:184:GLU:OE2	1:A:184:GLU:HA	2.12	0.48
1:B:79:ASP:HB3	1:D:32:ASN:OD1	2.12	0.48
1:D:48:ARG:O	1:D:49:GLU:C	2.54	0.48
1:C:89:VAL:HG13	1:C:97:ASN:CB	2.41	0.48
1:A:124:LYS:C	1:A:126:GLY:H	2.19	0.48
1:D:140:VAL:HB	1:D:141:PRO:HD2	1.94	0.48
1:A:132:LYS:HB2	2:A:273:HOH:O	2.12	0.48
1:C:8:LEU:HD23	1:C:86:VAL:HB	1.95	0.48
1:C:119:PHE:CE2	1:C:131:ILE:HD11	2.48	0.48
1:D:124:LYS:HZ2	1:D:125:LYS:HG2	1.77	0.48
1:B:74:ARG:HG3	1:B:75:ASN:N	2.29	0.48
1:A:13:ASN:HD21	1:B:122:LEU:CD1	2.26	0.48
1:A:218:ILE:O	1:A:221:VAL:HG12	2.12	0.48
1:B:106:GLU:C	1:B:108:GLU:H	2.21	0.48
1:B:122:LEU:HD13	1:B:127:ALA:N	2.28	0.48
1:C:246:ARG:HH11	1:D:178:ASP:HB3	1.78	0.48
1:A:181:ILE:HD12	1:A:255:GLN:HE22	1.78	0.48
1:B:112:ILE:HG13	1:B:113:ILE:N	2.28	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:147:ALA:C	1:B:149:LYS:N	2.72	0.47
1:B:33:TRP:CE2	1:B:40:LYS:HG3	2.49	0.47
1:D:119:PHE:CZ	1:D:131:ILE:HD11	2.50	0.47
1:B:18:THR:HG22	1:B:147:ALA:HB1	1.97	0.47
1:B:128:LYS:O	1:B:269:VAL:HG23	2.14	0.47
1:C:202:PRO:HG2	1:C:205:TRP:HB3	1.96	0.47
1:C:222:LEU:CD2	1:C:227:GLN:HG2	2.41	0.47
1:A:102:LEU:HD12	1:A:102:LEU:HA	1.79	0.47
1:A:258:GLU:HG2	1:B:246:ARG:HD2	1.97	0.47
1:B:102:LEU:O	1:B:106:GLU:HG3	2.15	0.47
1:C:203:ILE:O	1:C:204:ARG:C	2.57	0.47
1:C:138:LEU:HD21	1:C:263:LEU:HD21	1.97	0.47
1:A:123:LYS:NZ	1:A:129:ILE:H	2.13	0.46
1:C:94:LEU:O	1:C:95:MET:C	2.58	0.46
1:D:105:PHE:CG	1:D:172:PRO:HB3	2.50	0.46
1:D:124:LYS:HD2	1:D:124:LYS:C	2.40	0.46
1:A:240:LEU:HD22	1:A:248:LEU:HD21	1.96	0.46
1:A:1:MET:HG2	1:A:2:VAL:O	2.15	0.46
1:A:225:LEU:HB3	1:A:229:LYS:HG2	1.96	0.46
1:C:210:LEU:HD13	1:C:233:ILE:HG21	1.98	0.46
1:A:181:ILE:O	1:A:185:ILE:HG13	2.15	0.46
1:A:210:LEU:HD13	1:A:233:ILE:HG21	1.97	0.46
1:A:120:ASP:CG	1:A:146:ASN:ND2	2.73	0.46
1:C:229:LYS:HA	1:C:229:LYS:CE	2.41	0.46
1:D:78:LEU:HD11	1:D:205:TRP:CZ3	2.51	0.46
1:D:209:LYS:HB3	1:D:217:VAL:HG21	1.98	0.46
1:A:108:GLU:HG3	1:A:202:PRO:HB3	1.97	0.46
1:A:187:HIS:O	1:A:188:ILE:C	2.59	0.46
1:D:120:ASP:HB3	1:D:146:ASN:HD22	1.80	0.46
1:C:131:ILE:CD1	1:C:144:PRO:HG3	2.44	0.46
1:C:158:ARG:HH12	1:C:162:LEU:HD22	1.80	0.46
1:D:86:VAL:HG21	1:D:160:ILE:HD11	1.98	0.46
1:D:181:ILE:O	1:D:185:ILE:HG13	2.16	0.46
1:A:11:ASN:HD21	1:A:97:ASN:HA	1.81	0.45
1:A:120:ASP:HB3	1:A:146:ASN:HD22	1.80	0.45
1:B:133:LYS:HD3	1:B:266:LYS:O	2.16	0.45
1:C:270:HIS:ND1	1:C:270:HIS:N	2.64	0.45
1:D:160:ILE:O	1:D:163:MET:HB3	2.16	0.45
1:B:34:PRO:CG	1:D:34:PRO:HG2	2.44	0.45
1:A:175:PRO:HA	1:A:259:PHE:CZ	2.52	0.45
1:B:189:SER:C	1:B:191:LEU:H	2.23	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:ARG:HD3	1:A:135:ARG:C	2.41	0.45
1:A:195:THR:HB	1:A:225:LEU:HD11	1.98	0.45
1:B:146:ASN:HD21	1:B:149:LYS:HD3	1.79	0.45
1:C:210:LEU:HD21	1:C:221:VAL:HG21	1.98	0.45
1:A:106:GLU:C	1:A:108:GLU:H	2.23	0.45
1:A:184:GLU:HB2	1:A:240:LEU:HD12	1.99	0.45
1:A:248:LEU:O	1:A:251:VAL:HG12	2.17	0.45
1:C:234:LEU:HA	1:C:237:ILE:HG22	1.99	0.45
1:B:153:VAL:O	1:B:156:LEU:HB3	2.17	0.45
1:B:207:ALA:O	1:B:211:LEU:HD13	2.17	0.45
1:A:13:ASN:OD1	1:B:122:LEU:HD12	2.17	0.45
1:A:211:LEU:HD13	1:A:240:LEU:CD1	2.46	0.45
1:C:95:MET:HE2	1:C:260:LEU:HB3	1.98	0.44
1:D:19:ILE:O	1:D:23:LEU:HG	2.16	0.44
1:D:34:PRO:O	1:D:36:VAL:HG22	2.16	0.44
1:A:114:LEU:HD23	1:A:114:LEU:C	2.42	0.44
1:D:209:LYS:HA	1:D:209:LYS:HD3	1.74	0.44
1:A:233:ILE:HG22	1:A:237:ILE:HD11	1.98	0.44
1:B:229:LYS:O	1:B:233:ILE:HG13	2.17	0.44
1:C:9:VAL:HG22	1:C:10:GLY:N	2.32	0.44
1:D:27:ARG:HG2	1:D:27:ARG:NH1	2.31	0.44
1:D:90:ASP:OD1	1:D:90:ASP:C	2.60	0.44
1:D:105:PHE:CD1	1:D:172:PRO:HB3	2.52	0.44
1:D:221:VAL:HG12	1:D:230:MET:HE2	1.97	0.44
1:A:45:MET:HB3	1:A:52:PHE:HB2	2.00	0.44
1:A:122:LEU:HD23	1:A:123:LYS:HE2	1.99	0.44
1:C:268:VAL:HG12	1:C:269:VAL:N	2.31	0.44
1:B:122:LEU:O	1:B:122:LEU:HD23	2.18	0.44
1:D:13:ASN:C	1:D:15:GLY:N	2.74	0.44
1:D:16:LYS:HE2	1:D:56:ASP:OD1	2.17	0.44
1:A:245:LYS:HG2	1:A:246:ARG:H	1.82	0.44
1:D:221:VAL:HB	1:D:230:MET:CE	2.43	0.44
1:A:262:GLN:OE1	1:B:246:ARG:HG2	2.18	0.44
1:B:212:GLN:CG	1:B:252:ILE:HD12	2.48	0.44
1:D:182:GLU:OE2	1:D:204:ARG:NH2	2.48	0.43
1:D:230:MET:O	1:D:234:LEU:HB2	2.17	0.43
1:A:33:TRP:CD1	1:A:40:LYS:HB2	2.53	0.43
1:B:122:LEU:HD22	1:B:127:ALA:HA	2.00	0.43
1:C:258:GLU:HG3	1:D:254:SER:OG	2.18	0.43
1:B:77:ILE:HD12	1:B:85:ILE:CD1	2.48	0.43
1:B:114:LEU:O	1:B:142:VAL:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:258:GLU:HG3	1:D:254:SER:CB	2.49	0.43
1:D:133:LYS:HG2	1:D:266:LYS:O	2.19	0.43
1:A:197:LEU:C	1:A:199:GLU:N	2.77	0.43
1:D:6:VAL:HG11	1:D:160:ILE:HD12	2.00	0.43
1:A:102:LEU:O	1:A:106:GLU:HG3	2.19	0.43
1:C:89:VAL:HG21	1:C:114:LEU:HD11	2.01	0.43
1:C:4:LYS:HA	1:C:83:ASP:OD2	2.18	0.42
1:A:12:PRO:HD2	1:A:61:TYR:CZ	2.54	0.42
1:B:135:ARG:HG2	1:B:140:VAL:O	2.19	0.42
1:C:245:LYS:HD2	1:D:179:GLU:HG2	2.01	0.42
1:C:248:LEU:HA	1:C:248:LEU:HD23	1.80	0.42
1:C:245:LYS:HD3	1:D:179:GLU:HG2	2.00	0.42
1:A:188:ILE:HG21	1:A:210:LEU:HD12	2.01	0.42
1:A:188:ILE:HD11	1:A:237:ILE:CD1	2.49	0.42
1:A:212:GLN:HG2	1:A:249:ASP:HA	2.01	0.42
1:A:239:GLU:O	1:A:242:GLU:HB3	2.19	0.42
1:B:90:ASP:H	1:B:97:ASN:ND2	2.17	0.42
1:B:92:THR:HG21	1:B:121:LEU:HD22	2.01	0.42
1:D:83:ASP:O	1:D:111:ASN:ND2	2.35	0.42
1:D:156:LEU:O	1:D:160:ILE:HG12	2.19	0.42
1:C:178:ASP:OD2	1:C:255:GLN:NE2	2.51	0.42
1:C:234:LEU:HA	1:C:237:ILE:CG2	2.50	0.42
1:B:178:ASP:OD2	1:B:255:GLN:NE2	2.53	0.42
1:C:190:GLU:C	1:C:192:LEU:N	2.77	0.42
1:D:210:LEU:HD12	1:D:210:LEU:HA	1.90	0.42
1:A:89:VAL:HG13	1:A:97:ASN:HB3	2.01	0.42
1:A:262:GLN:HG3	1:B:250:ILE:HD13	2.02	0.42
1:D:13:ASN:C	1:D:15:GLY:H	2.28	0.42
1:A:212:GLN:HE21	1:A:249:ASP:HA	1.85	0.42
1:B:117:ASN:ND2	1:B:118:LYS:H	2.18	0.42
1:C:45:MET:HE2	1:C:52:PHE:HD2	1.83	0.42
1:C:63:LEU:HD13	1:C:107:MET:HE2	2.02	0.42
1:C:234:LEU:CA	1:C:237:ILE:HG22	2.49	0.42
1:D:248:LEU:HD23	1:D:248:LEU:HA	1.76	0.42
1:B:115:VAL:HG13	1:B:145:THR:HG23	2.02	0.41
1:C:196:PRO:HD2	1:C:225:LEU:CD1	2.45	0.41
1:A:188:ILE:HG23	1:A:233:ILE:HD12	2.02	0.41
1:D:32:ASN:HA	1:D:38:VAL:O	2.20	0.41
1:A:33:TRP:HA	1:A:34:PRO:HD3	1.95	0.41
1:A:98:LEU:HA	1:A:98:LEU:HD23	1.84	0.41
1:C:231:ASP:HA	1:C:234:LEU:CD1	2.47	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:237:ILE:HG23	1:C:238:SER:N	2.35	0.41
1:C:106:GLU:C	1:C:108:GLU:H	2.27	0.41
1:C:158:ARG:NH1	1:C:162:LEU:HD22	2.34	0.41
1:C:246:ARG:HD2	1:D:178:ASP:HA	2.03	0.41
1:A:106:GLU:OE2	1:A:175:PRO:HG2	2.21	0.41
1:A:125:LYS:HD2	1:A:125:LYS:O	2.21	0.41
1:B:146:ASN:CG	1:B:149:LYS:HD3	2.44	0.41
1:B:78:LEU:C	1:B:80:GLY:H	2.29	0.41
1:B:106:GLU:C	1:B:108:GLU:N	2.78	0.41
1:A:13:ASN:ND2	1:B:122:LEU:CD1	2.84	0.41
1:B:152:GLY:O	1:B:155:GLU:N	2.54	0.41
1:B:179:GLU:O	1:B:183:ARG:N	2.48	0.41
1:B:258:GLU:OE1	1:B:258:GLU:HA	2.21	0.41
1:C:34:PRO:O	1:C:36:VAL:N	2.54	0.41
1:C:95:MET:CE	1:C:260:LEU:HB3	2.51	0.41
1:C:135:ARG:HG2	1:C:140:VAL:O	2.21	0.41
1:B:195:THR:CG2	1:B:196:PRO:HD2	2.51	0.41
1:A:177:TYR:CE2	1:A:208:LEU:HD11	2.56	0.40
1:B:147:ALA:O	1:B:149:LYS:N	2.54	0.40
1:D:94:LEU:HD11	1:D:129:ILE:HD13	2.01	0.40
1:B:229:LYS:HD3	1:B:232:GLU:OE2	2.21	0.40
1:D:120:ASP:HB3	1:D:146:ASN:ND2	2.36	0.40
1:B:27:ARG:NH2	1:D:51:GLU:OE1	2.41	0.40
1:B:133:LYS:O	1:B:134:MET:C	2.64	0.40
1:C:184:GLU:HG3	1:C:240:LEU:HB2	2.04	0.40
1:C:188:ILE:HG23	1:C:233:ILE:HD12	2.04	0.40
1:D:116:LEU:HD13	1:D:142:VAL:CG1	2.52	0.40
1:A:90:ASP:OD1	1:A:92:THR:OG1	2.38	0.40
1:A:183:ARG:O	1:A:186:LYS:HB3	2.21	0.40
1:B:242:GLU:O	1:B:243:LYS:C	2.62	0.40
1:A:122:LEU:C	1:A:124:LYS:H	2.30	0.40
1:A:249:ASP:OD1	1:A:250:ILE:N	2.55	0.40
1:C:268:VAL:CG1	1:C:269:VAL:N	2.84	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	266/271 (98%)	236 (89%)	29 (11%)	1 (0%)	30	54
1	B	261/271 (96%)	235 (90%)	25 (10%)	1 (0%)	30	54
1	C	263/271 (97%)	241 (92%)	20 (8%)	2 (1%)	16	37
1	D	235/271 (87%)	221 (94%)	13 (6%)	1 (0%)	30	54
All	All	1025/1084 (95%)	933 (91%)	87 (8%)	5 (0%)	24	48

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	35	GLY
1	D	150	GLY
1	A	187	HIS
1	C	27	ARG
1	B	148	LYS

### 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	241/244 (99%)	236 (98%)	5 (2%)	47	75
1	B	239/244 (98%)	229 (96%)	10 (4%)	26	55
1	C	241/244 (99%)	232 (96%)	9 (4%)	30	59
1	D	243/244 (100%)	233 (96%)	10 (4%)	27	56
All	All	964/976 (99%)	930 (96%)	34 (4%)	32	61

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	102	LEU
1	A	125	LYS
1	A	135	ARG
1	A	173	ILE
1	A	245	LYS
1	B	1	MET
1	B	102	LEU
1	B	103	GLU
1	B	112	ILE
1	B	122	LEU
1	B	135	ARG
1	B	162	LEU
1	B	169	THR
1	B	263	LEU
1	B	264	LEU
1	C	18	THR
1	C	95	MET
1	C	102	LEU
1	C	135	ARG
1	C	203	ILE
1	C	215	GLU
1	C	229	LYS
1	C	260	LEU
1	C	265	ARG
1	D	39	GLU
1	D	78	LEU
1	D	102	LEU
1	D	132	LYS
1	D	135	ARG
1	D	204	ARG
1	D	206	LEU
1	D	210	LEU
1	D	228	GLU
1	D	234	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	13	ASN
1	A	75	ASN

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Mol	Chain	Res	Type
1	A	97	ASN
1	A	117	ASN
1	A	146	ASN
1	A	187	HIS
1	A	212	GLN
1	A	262	GLN
1	B	11	ASN
1	B	28	GLN
1	B	75	ASN
1	B	97	ASN
1	B	117	ASN
1	B	212	GLN
1	B	236	HIS
1	C	11	ASN
1	C	28	GLN
1	C	75	ASN
1	C	81	ASN
1	C	97	ASN
1	C	117	ASN
1	C	236	HIS
1	C	262	GLN
1	C	270	HIS
1	D	11	ASN
1	D	28	GLN
1	D	81	ASN
1	D	97	ASN
1	D	117	ASN
1	D	146	ASN
1	D	212	GLN
1	D	236	HIS
1	D	262	GLN
1	D	270	HIS

### 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, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	268/271 (98%)	0.28	11 (4%) 41 37	43, 63, 103, 111	0
1	B	265/271 (97%)	0.36	7 (2%) 57 54	49, 74, 103, 109	0
1	C	267/271 (98%)	0.33	11 (4%) 41 37	41, 69, 109, 121	0
1	D	270/271 (99%)	0.25	11 (4%) 41 37	26, 67, 100, 119	0
All	All	1070/1084 (98%)	0.30	40 (3%) 45 41	26, 69, 104, 121	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	109	VAL	3.9
1	C	224	TYR	3.4
1	A	177	TYR	3.3
1	D	14	VAL	3.3
1	C	127	ALA	3.1
1	D	177	TYR	3.1
1	B	122	LEU	3.1
1	B	222	LEU	2.9
1	A	232	GLU	2.9
1	D	268	VAL	2.9
1	D	1	MET	2.7
1	B	269	VAL	2.6
1	C	1	MET	2.6
1	A	237	ILE	2.6
1	C	233	ILE	2.6
1	B	221	VAL	2.5
1	A	32	ASN	2.5
1	A	224	TYR	2.5
1	D	46	GLU	2.4
1	C	230	MET	2.4
1	A	127	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	13	ASN	2.4
1	B	61	TYR	2.3
1	A	233	ILE	2.3
1	D	90	ASP	2.3
1	D	89	VAL	2.3
1	A	179	GLU	2.3
1	C	202	PRO	2.2
1	D	110	LYS	2.2
1	D	113	ILE	2.2
1	C	147	ALA	2.2
1	C	194	GLY	2.2
1	A	238	SER	2.2
1	A	13	ASN	2.1
1	B	127	ALA	2.1
1	B	148	LYS	2.1
1	C	222	LEU	2.1
1	C	185	ILE	2.1
1	C	234	LEU	2.1
1	A	228	GLU	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.