



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2018 – 02:10 am GMT

PDB ID : 3F4L  
Title : Crystal structure of a probable oxidoreductase yhhX in Triclinic form. Northeast Structural Genomics target ER647  
Authors : Seetharaman, J.; Abashidze, M.; Wang, H.; Janjua, H.; Foote, E.L.; Xiao, R.; Nair, R.; Everett, J.K.; Acton, T.B.; Rost, B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)  
Deposited on : 2008-10-31  
Resolution : 2.00 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.

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

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : trunk30967  
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)  
Refmac : 5.8.0158  
CCP4 : 7.0 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk30967

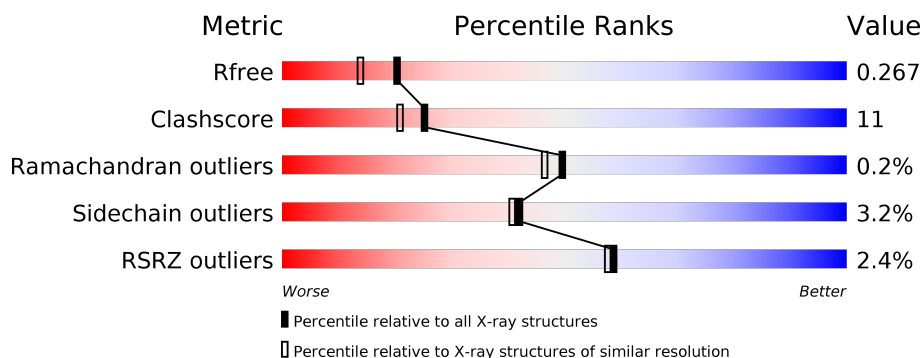
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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))
$R_{free}$	111664	7193 (2.00-2.00)
Clashscore	122126	8267 (2.00-2.00)
Ramachandran outliers	120053	8166 (2.00-2.00)
Sidechain outliers	120020	8165 (2.00-2.00)
RSRZ outliers	108989	7011 (2.00-2.00)

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  $\geq 3$ , 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	345	<div> <div>82%</div> <div>16%</div> <div>.</div> </div>
1	B	345	<div> <div>%</div> <div>76%</div> <div>21%</div> <div>.</div> </div>
1	C	345	<div> <div>3%</div> <div>77%</div> <div>21%</div> <div>.</div> </div>
1	D	345	<div> <div>%</div> <div>79%</div> <div>19%</div> <div>.</div> </div>
1	E	345	<div> <div>4%</div> <div>77%</div> <div>19%</div> <div>.</div> </div>
1	F	345	<div> <div>5%</div> <div>78%</div> <div>19%</div> <div>.</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 16970 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 Putative oxidoreductase yhhX.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	344	Total	C	N	O	S	Se	0	0	0
			2731	1747	462	515	3	4			
1	B	344	Total	C	N	O	S	Se	0	0	0
			2731	1747	462	515	3	4			
1	C	344	Total	C	N	O	S	Se	0	0	0
			2731	1747	462	515	3	4			
1	D	344	Total	C	N	O	S	Se	0	0	0
			2731	1747	462	515	3	4			
1	E	344	Total	C	N	O	S	Se	0	0	0
			2731	1747	462	515	3	4			
1	F	344	Total	C	N	O	S	Se	0	0	0
			2731	1747	462	515	3	4			

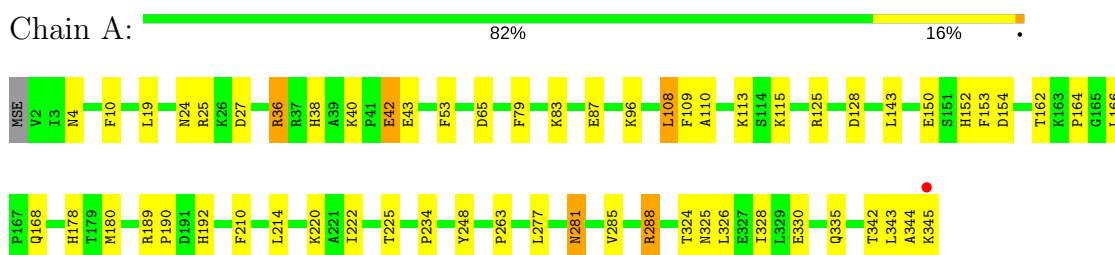
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	123	Total	O	0	0
			123	123		
2	B	96	Total	O	0	0
			96	96		
2	C	106	Total	O	0	0
			106	106		
2	D	87	Total	O	0	0
			87	87		
2	E	88	Total	O	0	0
			88	88		
2	F	84	Total	O	0	0
			84	84		

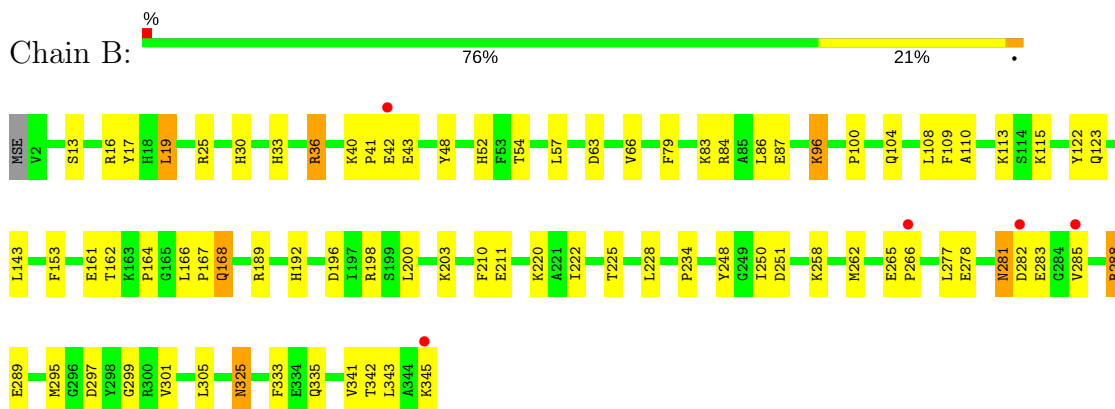
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

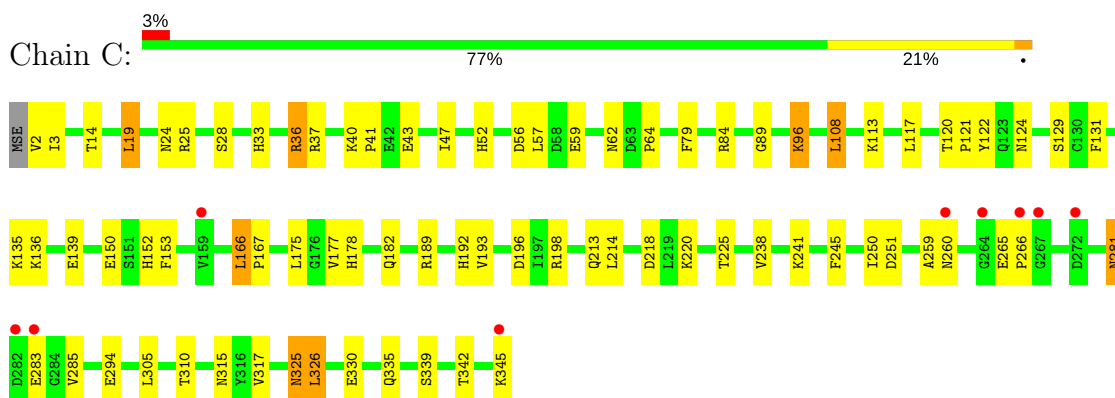
#### • Molecule 1: Putative oxidoreductase yhhX



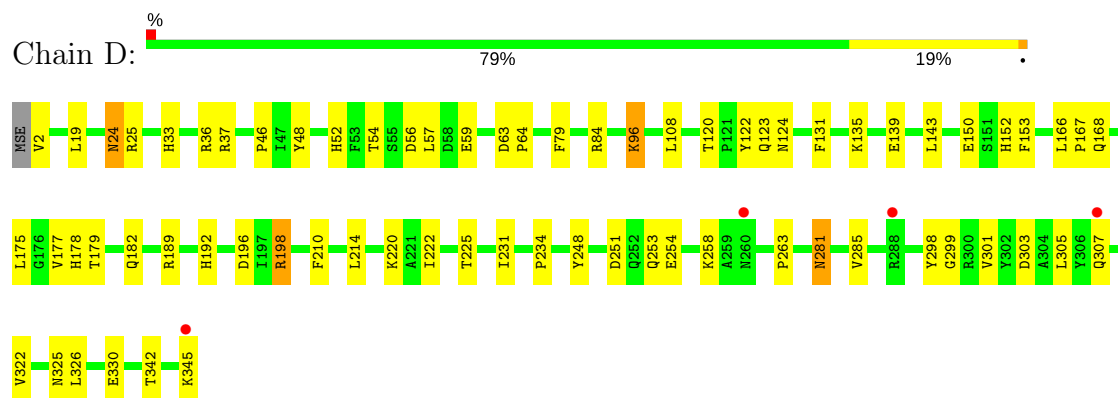
#### • Molecule 1: Putative oxidoreductase yhhX



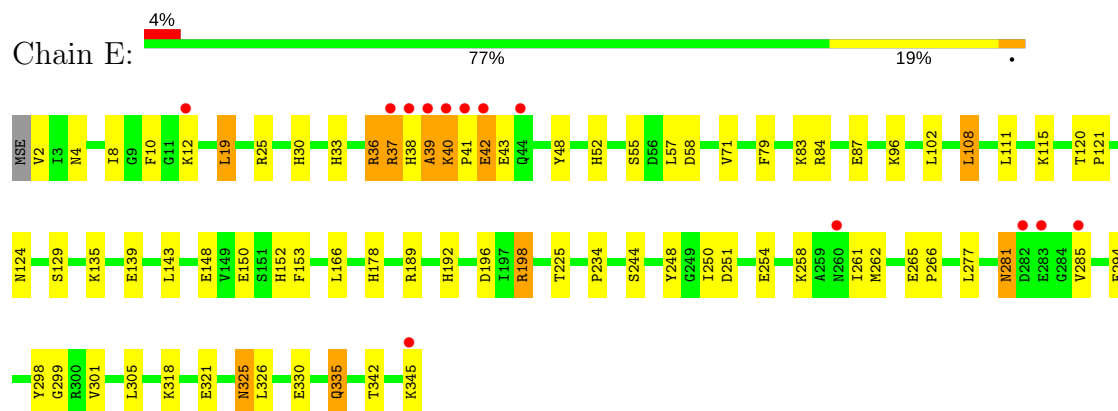
#### • Molecule 1: Putative oxidoreductase yhhX



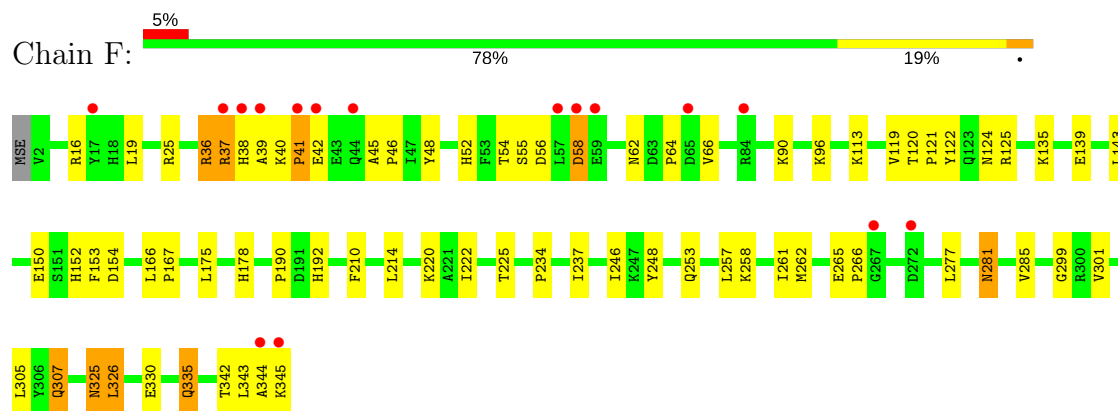
#### • Molecule 1: Putative oxidoreductase yhhX



• Molecule 1: Putative oxidoreductase yhhX



• Molecule 1: Putative oxidoreductase yhhX



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	66.19Å 100.78Å 102.17Å 114.40° 103.31° 101.62°	Depositor
Resolution (Å)	42.60 – 2.00 42.59 – 2.00	Depositor EDS
% Data completeness (in resolution range)	86.4 (42.60-2.00) 91.8 (42.59-2.00)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.04	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.50 (at 2.00Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.231 , 0.264 0.234 , 0.267	Depositor DCC
$R_{free}$ test set	11693 reflections (4.26%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	22.9	Xtriage
Anisotropy	0.210	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 39.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.018 for h,l,-h-k-l 0.018 for h,-h-k-l,k 0.002 for -h,-l,-k 0.001 for -h,-k,h+k+l 0.004 for -h,h+k+l,-l	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	16970	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.01% 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	0.36	0/2795	0.62	0/3780
1	B	0.34	0/2795	0.61	0/3780
1	C	0.34	0/2795	0.60	0/3780
1	D	0.34	0/2795	0.61	0/3780
1	E	0.34	0/2795	0.59	0/3780
1	F	0.34	0/2795	0.60	0/3780
All	All	0.35	0/16770	0.61	0/22680

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2731	0	2684	51	0
1	B	2731	0	2684	77	0
1	C	2731	0	2684	65	0
1	D	2731	0	2684	56	0
1	E	2731	0	2684	67	0
1	F	2731	0	2684	62	0
2	A	123	0	0	2	0
2	B	96	0	0	7	0
2	C	106	0	0	4	0
2	D	87	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	88	0	0	2	0
2	F	84	0	0	3	0
All	All	16970	0	16104	373	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 11.

All (373) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:12:LYS:H	1:E:12:LYS:HD2	1.14	1.08
1:B:40:LYS:HZ2	1:B:42:GLU:H	1.15	0.95
1:B:143:LEU:HD11	1:B:277:LEU:HD11	1.56	0.88
1:F:192:HIS:ND1	1:F:342:THR:HG22	1.89	0.87
1:A:96:LYS:H	1:A:96:LYS:HD3	1.41	0.86
1:D:301:VAL:O	1:D:305:LEU:HD23	1.77	0.84
1:B:189:ARG:HD3	1:B:345:LYS:OXT	1.81	0.81
1:C:192:HIS:ND1	1:C:342:THR:HG22	1.96	0.80
1:B:189:ARG:HD3	1:B:345:LYS:C	2.02	0.80
1:E:12:LYS:H	1:E:12:LYS:CD	1.95	0.80
1:D:96:LYS:HD3	1:D:96:LYS:H	1.47	0.78
1:E:40:LYS:HG3	1:E:41:PRO:HD2	1.64	0.78
1:B:301:VAL:O	1:B:305:LEU:HD23	1.83	0.78
1:B:192:HIS:ND1	1:B:342:THR:HG22	1.98	0.78
1:D:322:VAL:O	1:D:326:LEU:HD13	1.84	0.77
1:D:281:ASN:HD21	1:D:285:VAL:H	1.30	0.76
1:A:192:HIS:ND1	1:A:342:THR:HG22	2.00	0.76
1:F:37:ARG:HG2	1:F:38:HIS:H	1.51	0.76
1:E:12:LYS:N	1:E:12:LYS:HD2	1.98	0.75
1:A:110:ALA:HA	1:A:113:LYS:HE3	1.67	0.74
1:A:143:LEU:HD11	1:A:277:LEU:HD11	1.70	0.74
1:E:192:HIS:ND1	1:E:342:THR:HG22	2.03	0.74
1:D:281:ASN:ND2	1:D:285:VAL:H	1.85	0.73
1:E:150:GLU:OE2	1:E:152:HIS:HE1	1.71	0.73
1:D:189:ARG:HD3	1:D:345:LYS:C	2.09	0.73
1:A:153:PHE:O	1:A:225:THR:HB	1.88	0.72
1:B:83:LYS:O	1:B:87:GLU:HG3	1.90	0.72
1:A:189:ARG:CZ	1:A:345:LYS:HB3	2.20	0.72
1:A:36:ARG:HD3	1:A:43:GLU:OE1	1.90	0.72
1:B:96:LYS:H	1:B:96:LYS:HE2	1.56	0.70
1:B:96:LYS:HD3	2:B:526:HOH:O	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:281:ASN:C	1:C:281:ASN:HD22	1.95	0.70
1:A:189:ARG:NH1	1:A:345:LYS:HB3	2.07	0.69
1:B:40:LYS:HD2	1:B:41:PRO:HD2	1.72	0.69
1:B:96:LYS:H	1:B:96:LYS:CE	2.06	0.69
1:E:96:LYS:HE3	1:E:178:HIS:NE2	2.08	0.69
1:B:36:ARG:HD3	1:B:43:GLU:OE1	1.93	0.68
1:E:33:HIS:ND1	1:E:52:HIS:HD2	1.92	0.68
1:F:56:ASP:OD1	1:F:58:ASP:HB2	1.94	0.68
1:E:87:GLU:HA	1:E:115:LYS:HE2	1.74	0.68
1:E:40:LYS:CG	1:E:41:PRO:HD2	2.24	0.68
1:F:96:LYS:HE3	1:F:178:HIS:NE2	2.08	0.68
1:D:135:LYS:O	1:D:139:GLU:HG3	1.94	0.67
1:C:96:LYS:H	1:C:96:LYS:HE3	1.57	0.67
1:D:143:LEU:N	1:D:143:LEU:HD12	2.09	0.67
1:F:345:LYS:HB3	2:F:396:HOH:O	1.94	0.66
1:F:40:LYS:HD2	1:F:41:PRO:HD2	1.77	0.65
1:E:281:ASN:C	1:E:281:ASN:HD22	2.00	0.65
1:C:281:ASN:ND2	1:C:285:VAL:H	1.94	0.65
1:A:220:LYS:HE3	1:A:222:ILE:HD11	1.77	0.65
1:F:66:VAL:O	1:F:90:LYS:HD2	1.97	0.64
1:A:210:PHE:CZ	1:A:225:THR:HG23	2.33	0.64
1:B:281:ASN:ND2	1:B:285:VAL:H	1.95	0.64
1:E:254:GLU:HG2	1:E:258:LYS:HE2	1.79	0.64
1:F:150:GLU:OE2	1:F:152:HIS:HE1	1.81	0.64
1:D:303:ASP:O	1:D:307:GLN:HG3	1.98	0.63
1:B:153:PHE:O	1:B:225:THR:HB	1.98	0.63
1:F:281:ASN:HD22	1:F:281:ASN:C	2.02	0.63
1:A:189:ARG:HG3	1:A:324:THR:OG1	1.98	0.63
1:E:57:LEU:HD23	1:E:84:ARG:CZ	2.29	0.62
1:D:96:LYS:CD	1:D:96:LYS:H	2.12	0.62
1:D:33:HIS:ND1	1:D:52:HIS:HD2	1.97	0.62
1:F:40:LYS:C	1:F:42:GLU:H	2.02	0.62
1:F:143:LEU:HD12	1:F:143:LEU:N	2.15	0.62
1:F:281:ASN:HD21	1:F:285:VAL:H	1.48	0.62
1:A:150:GLU:OE2	1:A:152:HIS:HE1	1.83	0.62
1:E:83:LYS:O	1:E:87:GLU:HG3	2.00	0.61
1:F:37:ARG:HG2	1:F:38:HIS:N	2.14	0.61
1:D:196:ASP:OD1	1:D:198:ARG:HD3	2.01	0.61
1:C:96:LYS:H	1:C:96:LYS:CE	2.14	0.61
1:D:153:PHE:HB3	1:D:225:THR:HG22	1.81	0.61
1:F:143:LEU:HD11	1:F:277:LEU:HD11	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:150:GLU:OE2	1:D:152:HIS:HE1	1.83	0.61
1:B:281:ASN:HD21	1:B:285:VAL:H	1.49	0.61
1:D:220:LYS:HE3	1:D:222:ILE:HD11	1.82	0.60
1:D:96:LYS:N	1:D:96:LYS:HD3	2.15	0.60
1:A:96:LYS:HE3	1:A:178:HIS:CE1	2.37	0.60
1:A:83:LYS:O	1:A:87:GLU:HG3	2.01	0.60
1:F:281:ASN:ND2	1:F:285:VAL:H	1.99	0.60
1:A:110:ALA:HA	1:A:113:LYS:CE	2.30	0.60
1:D:189:ARG:HD3	1:D:345:LYS:OXT	2.02	0.60
1:E:301:VAL:O	1:E:305:LEU:HD23	2.02	0.60
1:E:189:ARG:HD3	1:E:345:LYS:OXT	2.03	0.59
1:A:96:LYS:N	1:A:96:LYS:HD3	2.15	0.59
1:B:40:LYS:NZ	1:B:42:GLU:HB2	2.19	0.58
1:C:326:LEU:O	1:C:330:GLU:HG3	2.04	0.58
1:B:40:LYS:NZ	1:B:42:GLU:H	1.95	0.58
1:A:87:GLU:HA	1:A:115:LYS:HE2	1.84	0.58
1:C:40:LYS:HD2	1:C:41:PRO:HD2	1.84	0.58
1:C:56:ASP:OD2	1:C:59:GLU:HG3	2.04	0.58
1:E:4:ASN:HD22	1:E:30:HIS:HB3	1.68	0.58
1:F:325:ASN:HD22	1:F:325:ASN:C	2.07	0.58
1:D:281:ASN:C	1:D:281:ASN:HD22	2.06	0.58
1:F:37:ARG:CG	1:F:38:HIS:H	2.15	0.57
1:C:96:LYS:HD3	1:C:96:LYS:N	2.18	0.57
1:C:135:LYS:O	1:C:139:GLU:HG3	2.04	0.57
1:D:19:LEU:HD23	1:D:48:TYR:CE1	2.39	0.57
1:D:96:LYS:HE3	1:D:178:HIS:NE2	2.20	0.57
1:B:220:LYS:HE3	1:B:222:ILE:HD11	1.87	0.57
1:C:96:LYS:HG3	2:C:439:HOH:O	2.05	0.56
1:D:56:ASP:OD2	1:D:59:GLU:HG3	2.05	0.56
1:A:281:ASN:HD22	1:A:281:ASN:C	2.08	0.56
1:B:40:LYS:HZ2	1:B:42:GLU:N	1.95	0.56
1:C:281:ASN:ND2	1:C:283:GLU:H	2.02	0.56
1:E:39:ALA:O	1:E:40:LYS:HB2	2.05	0.56
1:A:234:PRO:HG3	1:A:248:TYR:CZ	2.40	0.56
1:B:281:ASN:C	1:B:281:ASN:HD22	2.09	0.56
1:E:19:LEU:HD23	1:E:48:TYR:CE1	2.41	0.56
1:C:281:ASN:HD21	1:C:285:VAL:H	1.53	0.56
1:F:25:ARG:HH21	1:F:299:GLY:HA3	1.72	0.55
1:C:96:LYS:HG2	1:C:178:HIS:NE2	2.21	0.55
1:D:231:ILE:HG13	1:E:244:SER:HB2	1.87	0.55
1:E:96:LYS:HD3	1:E:96:LYS:N	2.22	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:150:GLU:OE2	1:C:152:HIS:HE1	1.90	0.55
1:F:125:ARG:HB2	2:F:359:HOH:O	2.07	0.55
1:E:143:LEU:HD21	1:E:277:LEU:CD1	2.37	0.55
1:A:281:ASN:HD21	1:A:285:VAL:H	1.56	0.54
1:A:326:LEU:O	1:A:330:GLU:HG3	2.08	0.54
1:A:143:LEU:HD12	1:A:143:LEU:N	2.23	0.54
1:D:131:PHE:CE1	1:D:182:GLN:HB3	2.43	0.54
1:E:42:GLU:HB3	1:E:48:TYR:CE2	2.43	0.54
1:F:166:LEU:HD22	1:F:166:LEU:N	2.23	0.54
1:C:25:ARG:CZ	2:C:366:HOH:O	2.56	0.53
1:C:96:LYS:HD2	1:C:122:TYR:O	2.08	0.53
1:F:125:ARG:HH11	1:F:125:ARG:HG3	1.72	0.53
1:E:57:LEU:HD23	1:E:84:ARG:NH2	2.22	0.53
1:D:326:LEU:O	1:D:330:GLU:HG3	2.08	0.53
1:E:335:GLN:HG2	2:E:364:HOH:O	2.09	0.53
1:E:318:LYS:HB2	1:E:321:GLU:HG3	1.89	0.53
1:E:42:GLU:O	1:E:48:TYR:HD2	1.91	0.53
1:F:153:PHE:HB3	1:F:225:THR:HG22	1.91	0.53
1:B:109:PHE:O	1:B:113:LYS:HE2	2.08	0.53
1:B:86:LEU:O	1:B:115:LYS:HE2	2.09	0.53
1:C:96:LYS:H	1:C:96:LYS:CD	2.22	0.53
1:E:143:LEU:HD21	1:E:277:LEU:HD11	1.91	0.53
1:B:234:PRO:HG3	1:B:248:TYR:CZ	2.43	0.53
1:C:33:HIS:ND1	1:C:52:HIS:HD2	2.07	0.53
1:B:278:GLU:HB2	1:B:288:ARG:NH1	2.24	0.52
1:A:96:LYS:H	1:A:96:LYS:CD	2.16	0.52
1:B:30:HIS:HD2	2:B:463:HOH:O	1.91	0.52
1:C:136:LYS:HE3	1:C:294:GLU:OE2	2.09	0.52
1:D:25:ARG:HH21	1:D:299:GLY:HA3	1.74	0.52
1:D:123:GLN:NE2	2:D:392:HOH:O	2.42	0.52
1:D:168:GLN:H	1:D:168:GLN:CD	2.12	0.52
1:A:288:ARG:HH11	1:A:288:ARG:HG3	1.74	0.52
1:B:110:ALA:HA	1:B:113:LYS:CE	2.40	0.52
1:B:153:PHE:HB3	1:B:225:THR:HG22	1.91	0.52
1:D:96:LYS:HE3	1:D:178:HIS:CE1	2.44	0.52
1:B:25:ARG:HH21	1:B:299:GLY:HA3	1.76	0.51
1:F:190:PRO:HG2	1:F:343:LEU:HD22	1.91	0.51
1:E:37:ARG:NH2	1:E:55:SER:O	2.43	0.51
1:C:14:THR:HG22	1:C:19:LEU:HD22	1.91	0.51
1:F:37:ARG:H	1:F:37:ARG:HD2	1.76	0.51
1:A:189:ARG:HG2	1:A:345:LYS:OXT	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:343:LEU:HB3	1:B:345:LYS:OXT	2.10	0.51
1:C:113:LYS:NZ	1:C:113:LYS:HB2	2.26	0.51
1:C:96:LYS:N	1:C:96:LYS:CD	2.72	0.51
1:C:166:LEU:HB3	1:C:167:PRO:HD2	1.92	0.51
1:C:189:ARG:HD3	1:C:345:LYS:OXT	2.11	0.50
1:B:203:LYS:HZ2	1:C:218:ASP:CG	2.14	0.50
1:C:325:ASN:HD22	1:C:325:ASN:C	2.14	0.50
1:F:52:HIS:CE1	1:F:54:THR:HB	2.46	0.50
1:C:335:GLN:HG2	2:C:399:HOH:O	2.11	0.50
1:D:166:LEU:HB3	1:D:167:PRO:HD2	1.93	0.50
1:A:281:ASN:ND2	1:A:285:VAL:H	2.09	0.50
1:B:40:LYS:HZ1	1:B:42:GLU:HB2	1.77	0.50
1:C:57:LEU:HD23	1:C:84:ARG:CZ	2.41	0.50
1:D:37:ARG:HG3	2:D:389:HOH:O	2.12	0.50
1:E:40:LYS:O	1:E:43:GLU:HG3	2.11	0.50
1:D:175:LEU:N	1:D:175:LEU:HD22	2.26	0.50
1:E:261:ILE:C	1:E:262:MSE:HE2	2.32	0.49
1:E:40:LYS:CD	1:E:41:PRO:HD2	2.41	0.49
1:F:234:PRO:HG3	1:F:248:TYR:CZ	2.46	0.49
1:C:265:GLU:HB3	1:C:266:PRO:HD2	1.94	0.49
1:C:136:LYS:HE3	1:C:294:GLU:CD	2.32	0.49
1:D:166:LEU:HB3	1:D:167:PRO:CD	2.42	0.49
1:E:37:ARG:HE	1:E:37:ARG:HA	1.76	0.49
1:F:96:LYS:HE3	1:F:178:HIS:CE1	2.47	0.49
1:F:40:LYS:C	1:F:42:GLU:N	2.66	0.49
1:C:259:ALA:O	1:C:260:ASN:HB2	2.13	0.49
1:C:281:ASN:C	1:C:281:ASN:ND2	2.64	0.49
1:D:143:LEU:CD1	1:D:143:LEU:N	2.74	0.49
1:F:120:THR:OG1	1:F:121:PRO:HD2	2.13	0.49
1:F:122:TYR:CZ	1:F:124:ASN:HB3	2.47	0.49
1:B:123:GLN:NE2	2:B:361:HOH:O	2.46	0.49
1:C:47:ILE:HG23	2:C:365:HOH:O	2.11	0.49
1:E:111:LEU:HD11	1:E:115:LYS:NZ	2.28	0.49
1:A:4:ASN:HB2	1:A:65:ASP:O	2.13	0.49
1:B:16:ARG:HD2	1:B:17:TYR:CZ	2.48	0.49
1:B:96:LYS:HG3	2:B:361:HOH:O	2.11	0.49
1:F:301:VAL:O	1:F:305:LEU:HD23	2.12	0.49
1:B:110:ALA:HA	1:B:113:LYS:HE3	1.94	0.49
1:B:295:MSE:HE2	2:B:365:HOH:O	2.12	0.49
1:F:307:GLN:HE21	1:F:307:GLN:CA	2.25	0.49
1:B:25:ARG:NE	2:B:351:HOH:O	2.44	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:129:SER:HB2	1:E:294:GLU:O	2.13	0.49
1:B:210:PHE:CZ	1:B:225:THR:HG23	2.48	0.49
1:B:16:ARG:HH12	1:B:258:LYS:NZ	2.11	0.48
1:B:25:ARG:CZ	2:B:351:HOH:O	2.61	0.48
1:B:262:MSE:HB2	1:B:265:GLU:OE2	2.13	0.48
1:D:179:THR:HG21	2:D:388:HOH:O	2.13	0.48
1:A:335:GLN:HG2	2:A:399:HOH:O	2.13	0.48
1:E:96:LYS:H	1:E:96:LYS:CE	2.25	0.48
1:B:96:LYS:CD	1:B:96:LYS:H	2.26	0.48
1:A:143:LEU:N	1:A:143:LEU:CD1	2.77	0.48
1:F:62:ASN:O	1:F:64:PRO:HD3	2.14	0.48
1:B:19:LEU:HD23	1:B:48:TYR:CE1	2.48	0.48
1:C:40:LYS:HD2	1:C:41:PRO:CD	2.44	0.48
1:F:220:LYS:HE3	1:F:222:ILE:HD11	1.96	0.48
1:F:253:GLN:O	1:F:257:LEU:HG	2.13	0.48
1:C:89:GLY:HA2	1:C:117:LEU:HD11	1.95	0.48
1:E:40:LYS:HA	1:E:40:LYS:HD2	1.56	0.48
1:C:129:SER:HB2	1:C:294:GLU:O	2.14	0.48
1:E:166:LEU:HD22	1:E:166:LEU:N	2.28	0.48
1:E:254:GLU:CG	1:E:258:LYS:HE2	2.42	0.48
1:B:143:LEU:CD1	1:B:143:LEU:N	2.77	0.47
1:B:228:LEU:HD13	1:C:241:LYS:HG2	1.96	0.47
1:C:79:PHE:CE1	1:C:108:LEU:HD13	2.49	0.47
1:B:282:ASP:HB2	1:B:283:GLU:OE2	2.13	0.47
1:C:120:THR:OG1	1:C:121:PRO:HD2	2.15	0.47
1:D:153:PHE:O	1:D:225:THR:HB	2.14	0.47
1:E:4:ASN:ND2	1:E:30:HIS:HB3	2.29	0.47
1:C:131:PHE:CE1	1:C:182:GLN:HB3	2.49	0.47
1:F:19:LEU:HD23	1:F:48:TYR:CE1	2.49	0.47
1:E:265:GLU:HB3	1:E:266:PRO:HD2	1.97	0.47
1:F:143:LEU:CD1	1:F:143:LEU:N	2.76	0.47
1:A:180:MSE:SE	1:A:328:ILE:HG21	2.64	0.47
1:C:3:ILE:HD11	1:C:310:THR:OG1	2.13	0.47
1:E:234:PRO:HG3	1:E:248:TYR:CZ	2.50	0.47
1:B:166:LEU:HD22	1:B:166:LEU:N	2.29	0.47
1:B:250:ILE:HG12	1:B:251:ASP:N	2.30	0.47
1:F:40:LYS:O	1:F:42:GLU:N	2.48	0.47
1:B:295:MSE:HE3	1:B:297:ASP:HB2	1.97	0.47
1:F:153:PHE:O	1:F:225:THR:HB	2.15	0.47
1:E:196:ASP:OD1	1:E:198:ARG:HD3	2.15	0.47
1:C:238:VAL:HB	1:C:245:PHE:HB3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:178:HIS:HD2	2:E:352:HOH:O	1.98	0.47
1:F:96:LYS:H	1:F:96:LYS:HE2	1.79	0.47
1:B:200:LEU:CD2	1:C:218:ASP:HB3	2.45	0.46
1:A:214:LEU:N	1:A:214:LEU:HD12	2.31	0.46
1:F:326:LEU:O	1:F:330:GLU:HG3	2.15	0.46
1:E:40:LYS:HE3	1:E:41:PRO:HD2	1.98	0.46
1:B:162:THR:HG22	1:B:164:PRO:HD3	1.97	0.46
1:D:96:LYS:HD2	1:D:122:TYR:O	2.16	0.46
1:D:251:ASP:OD2	1:D:253:GLN:HG3	2.16	0.46
1:E:281:ASN:C	1:E:281:ASN:ND2	2.69	0.46
1:F:96:LYS:CE	1:F:96:LYS:H	2.29	0.46
1:D:166:LEU:HD22	1:D:166:LEU:N	2.31	0.45
1:F:335:GLN:HG2	2:F:356:HOH:O	2.16	0.45
1:B:96:LYS:NZ	1:B:122:TYR:HB3	2.31	0.45
1:F:175:LEU:HD22	1:F:175:LEU:N	2.32	0.45
1:A:344:ALA:O	1:A:345:LYS:HG3	2.17	0.45
1:E:96:LYS:CD	1:E:96:LYS:H	2.29	0.45
1:F:154:ASP:OD1	1:F:225:THR:HA	2.17	0.45
1:A:96:LYS:HE3	1:A:178:HIS:NE2	2.32	0.45
1:B:79:PHE:CE1	1:B:108:LEU:HD13	2.51	0.45
1:B:281:ASN:C	1:B:281:ASN:ND2	2.69	0.45
1:A:154:ASP:OD1	1:A:225:THR:HA	2.17	0.45
1:B:13:SER:HA	1:B:17:TYR:HD2	1.81	0.45
1:C:175:LEU:HD22	1:C:175:LEU:N	2.32	0.45
1:D:254:GLU:CG	1:D:258:LYS:HE2	2.46	0.45
1:E:25:ARG:HH21	1:E:299:GLY:HA3	1.82	0.45
1:A:125:ARG:NE	1:A:128:ASP:OD2	2.46	0.45
1:B:96:LYS:CD	1:B:96:LYS:N	2.80	0.45
1:E:96:LYS:CD	1:E:96:LYS:N	2.80	0.45
1:D:254:GLU:HG2	1:D:258:LYS:HE2	1.99	0.45
1:D:57:LEU:HD23	1:D:84:ARG:NH1	2.32	0.45
1:A:162:THR:HG22	1:A:164:PRO:HD3	1.98	0.44
1:B:143:LEU:N	1:B:143:LEU:HD12	2.31	0.44
1:E:102:LEU:HA	1:E:326:LEU:HD23	1.98	0.44
1:A:113:LYS:HG3	2:A:407:HOH:O	2.15	0.44
1:B:109:PHE:O	1:B:113:LYS:HG3	2.18	0.44
1:B:40:LYS:HD2	1:B:41:PRO:CD	2.46	0.44
1:F:119:VAL:O	1:F:119:VAL:HG13	2.16	0.44
1:F:56:ASP:CG	1:F:58:ASP:HB2	2.38	0.44
1:B:335:GLN:NE2	1:B:341:VAL:HG12	2.33	0.44
1:D:122:TYR:CZ	1:D:124:ASN:HB3	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:189:ARG:CD	1:C:345:LYS:OXT	2.66	0.44
1:E:135:LYS:O	1:E:139:GLU:HG3	2.17	0.44
1:A:166:LEU:N	1:A:166:LEU:HD22	2.33	0.44
1:E:36:ARG:HG2	1:E:36:ARG:HH11	1.83	0.44
1:C:250:ILE:HG12	1:C:251:ASP:N	2.33	0.44
1:D:2:VAL:HG13	1:D:2:VAL:O	2.17	0.44
1:B:96:LYS:HD3	1:B:96:LYS:N	2.33	0.44
1:E:124:ASN:HB2	1:E:298:TYR:CD1	2.53	0.44
1:E:36:ARG:NE	1:E:43:GLU:OE2	2.42	0.44
1:B:168:GLN:H	1:B:168:GLN:CD	2.21	0.44
1:B:33:HIS:ND1	1:B:52:HIS:HD2	2.15	0.44
1:C:214:LEU:HD12	1:C:214:LEU:N	2.33	0.44
1:D:124:ASN:HB2	1:D:298:TYR:CD1	2.53	0.44
1:D:46:PRO:HD2	2:D:468:HOH:O	2.18	0.44
1:E:8:ILE:HB	1:E:71:VAL:HA	1.99	0.44
1:F:45:ALA:HA	1:F:46:PRO:HD3	1.86	0.43
1:A:24:ASN:ND2	1:A:263:PRO:O	2.52	0.43
1:C:122:TYR:CZ	1:C:124:ASN:HB3	2.52	0.43
1:C:40:LYS:CD	1:C:41:PRO:HD2	2.49	0.43
1:A:190:PRO:HG2	1:A:343:LEU:HD22	1.99	0.43
1:E:281:ASN:HD21	1:E:285:VAL:H	1.66	0.43
1:F:36:ARG:C	1:F:36:ARG:CD	2.87	0.43
1:B:211:GLU:OE1	1:C:220:LYS:NZ	2.49	0.43
1:B:57:LEU:HD23	1:B:84:ARG:CZ	2.49	0.43
1:C:177:VAL:HG13	1:C:178:HIS:N	2.33	0.43
1:C:2:VAL:HG23	1:C:28:SER:C	2.39	0.43
1:E:153:PHE:HB3	1:E:225:THR:HG22	1.99	0.43
1:A:109:PHE:C	1:A:113:LYS:HE2	2.39	0.43
1:F:210:PHE:CZ	1:F:225:THR:HG23	2.53	0.43
1:F:96:LYS:H	1:F:96:LYS:CD	2.32	0.43
1:B:325:ASN:HD22	1:B:325:ASN:C	2.22	0.42
1:C:166:LEU:HB3	1:C:167:PRO:CD	2.49	0.42
1:E:281:ASN:ND2	1:E:285:VAL:H	2.17	0.42
1:E:10:PHE:HE2	1:E:42:GLU:HB2	1.83	0.42
1:C:193:VAL:HA	1:C:213:GLN:O	2.19	0.42
1:A:43:GLU:HG2	1:A:53:PHE:CE2	2.54	0.42
1:D:24:ASN:ND2	1:D:263:PRO:O	2.53	0.42
1:F:214:LEU:N	1:F:214:LEU:HD12	2.33	0.42
1:D:166:LEU:HB3	1:D:168:GLN:HE22	1.85	0.42
1:A:25:ARG:HG2	1:A:25:ARG:HH11	1.85	0.42
1:D:210:PHE:CE1	1:D:225:THR:OG1	2.64	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:261:ILE:C	1:F:262:MSE:HE2	2.40	0.42
1:B:265:GLU:HB3	1:B:266:PRO:HD2	2.01	0.42
1:D:79:PHE:CE1	1:D:108:LEU:HD13	2.54	0.42
1:E:41:PRO:HG2	1:E:42:GLU:OE1	2.20	0.42
1:A:10:PHE:O	1:A:40:LYS:HD2	2.20	0.42
1:C:25:ARG:HH11	1:C:25:ARG:HG2	1.85	0.42
1:C:335:GLN:HG3	1:C:339:SER:HB3	2.02	0.42
1:D:214:LEU:HD12	1:D:214:LEU:N	2.35	0.42
1:A:189:ARG:HD3	1:A:345:LYS:C	2.40	0.41
1:B:100:PRO:HD2	1:B:104:GLN:OE1	2.20	0.41
1:B:40:LYS:CD	1:B:41:PRO:HD2	2.45	0.41
1:F:96:LYS:HD3	1:F:96:LYS:N	2.34	0.41
1:B:167:PRO:HB3	1:B:333:PHE:CE1	2.54	0.41
1:B:166:LEU:HB3	1:B:167:PRO:HD2	2.01	0.41
1:B:196:ASP:OD2	1:B:198:ARG:NH1	2.53	0.41
1:C:189:ARG:CZ	1:C:345:LYS:HA	2.49	0.41
1:D:166:LEU:CB	1:D:168:GLN:HE22	2.32	0.41
1:E:326:LEU:O	1:E:330:GLU:HG3	2.19	0.41
1:C:196:ASP:OD1	1:C:198:ARG:HD3	2.20	0.41
1:D:234:PRO:HG3	1:D:248:TYR:CZ	2.55	0.41
1:E:120:THR:OG1	1:E:121:PRO:HD2	2.21	0.41
1:F:113:LYS:HE3	1:F:113:LYS:HB2	1.78	0.41
1:A:36:ARG:HD2	1:A:38:HIS:O	2.21	0.41
1:B:161:GLU:OE2	1:B:161:GLU:HA	2.21	0.41
1:F:265:GLU:HB3	1:F:266:PRO:HD2	2.02	0.41
1:A:189:ARG:NE	1:A:345:LYS:O	2.53	0.41
1:D:166:LEU:HB2	1:D:168:GLN:NE2	2.36	0.41
1:D:177:VAL:HG13	1:D:178:HIS:N	2.36	0.41
1:B:109:PHE:C	1:B:113:LYS:HE2	2.41	0.41
1:F:39:ALA:O	1:F:40:LYS:HB2	2.21	0.41
1:E:250:ILE:HG12	1:E:251:ASP:N	2.36	0.41
1:F:192:HIS:CE1	1:F:342:THR:HG22	2.54	0.41
1:A:10:PHE:CE2	1:A:42:GLU:HG3	2.54	0.41
1:C:153:PHE:O	1:C:225:THR:HB	2.21	0.41
1:C:315:ASN:ND2	1:C:317:VAL:H	2.19	0.41
1:C:62:ASN:O	1:C:64:PRO:HD3	2.20	0.41
1:E:79:PHE:CE1	1:E:108:LEU:HD13	2.56	0.41
1:D:192:HIS:ND1	1:D:342:THR:HG22	2.36	0.40
1:D:63:ASP:HA	1:D:64:PRO:HD3	1.87	0.40
1:E:325:ASN:HD22	1:E:325:ASN:C	2.25	0.40
1:E:38:HIS:O	1:E:39:ALA:HB2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:36:ARG:HD2	1:C:43:GLU:OE1	2.21	0.40
1:F:166:LEU:HB3	1:F:167:PRO:HD2	2.03	0.40
1:A:79:PHE:CE1	1:A:108:LEU:HD13	2.56	0.40
1:B:16:ARG:NH1	1:B:258:LYS:NZ	2.69	0.40
1:C:305:LEU:HA	1:C:305:LEU:HD13	1.96	0.40
1:C:335:GLN:HG3	1:C:339:SER:CB	2.52	0.40
1:E:2:VAL:O	1:E:2:VAL:HG13	2.21	0.40
1:F:135:LYS:O	1:F:139:GLU:HG3	2.22	0.40
1:F:237:ILE:HG23	1:F:246:ILE:CD1	2.52	0.40
1:A:109:PHE:O	1:A:113:LYS:HE2	2.22	0.40
1:B:96:LYS:HZ3	1:B:122:TYR:HB3	1.86	0.40
1:F:16:ARG:NH1	1:F:258:LYS:NZ	2.69	0.40
1:A:288:ARG:NH1	1:A:288:ARG:HG3	2.37	0.40
1:B:63:ASP:HB3	1:B:66:VAL:HG23	2.03	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	342/345 (99%)	334 (98%)	8 (2%)	0	100	100
1	B	342/345 (99%)	338 (99%)	4 (1%)	0	100	100
1	C	342/345 (99%)	336 (98%)	6 (2%)	0	100	100
1	D	342/345 (99%)	334 (98%)	8 (2%)	0	100	100
1	E	342/345 (99%)	328 (96%)	12 (4%)	2 (1%)	27	21
1	F	342/345 (99%)	323 (94%)	17 (5%)	2 (1%)	27	21
All	All	2052/2070 (99%)	1993 (97%)	55 (3%)	4 (0%)	49	46

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	39	ALA
1	F	344	ALA
1	F	41	PRO
1	E	40	LYS

### 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	294/290 (101%)	285 (97%)	9 (3%)	43	42
1	B	294/290 (101%)	285 (97%)	9 (3%)	43	42
1	C	294/290 (101%)	284 (97%)	10 (3%)	40	38
1	D	294/290 (101%)	286 (97%)	8 (3%)	48	49
1	E	294/290 (101%)	283 (96%)	11 (4%)	37	35
1	F	294/290 (101%)	285 (97%)	9 (3%)	43	42
All	All	1764/1740 (101%)	1708 (97%)	56 (3%)	42	41

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LEU
1	A	27	ASP
1	A	36	ARG
1	A	42	GLU
1	A	108	LEU
1	A	168	GLN
1	A	281	ASN
1	A	288	ARG
1	A	325	ASN
1	B	19	LEU
1	B	36	ARG
1	B	54	THR
1	B	96	LYS
1	B	168	GLN
1	B	281	ASN

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Mol	Chain	Res	Type
1	B	288	ARG
1	B	289	GLU
1	B	325	ASN
1	C	19	LEU
1	C	24	ASN
1	C	36	ARG
1	C	37	ARG
1	C	96	LYS
1	C	108	LEU
1	C	166	LEU
1	C	281	ASN
1	C	325	ASN
1	C	326	LEU
1	D	24	ASN
1	D	36	ARG
1	D	54	THR
1	D	96	LYS
1	D	120	THR
1	D	198	ARG
1	D	281	ASN
1	D	325	ASN
1	E	19	LEU
1	E	36	ARG
1	E	37	ARG
1	E	42	GLU
1	E	58	ASP
1	E	108	LEU
1	E	148	GLU
1	E	198	ARG
1	E	281	ASN
1	E	325	ASN
1	E	335	GLN
1	F	36	ARG
1	F	37	ARG
1	F	55	SER
1	F	58	ASP
1	F	281	ASN
1	F	307	GLN
1	F	325	ASN
1	F	326	LEU
1	F	335	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (49) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	ASN
1	A	52	HIS
1	A	152	HIS
1	A	213	GLN
1	A	281	ASN
1	A	307	GLN
1	A	325	ASN
1	B	4	ASN
1	B	24	ASN
1	B	38	HIS
1	B	44	GLN
1	B	52	HIS
1	B	123	GLN
1	B	213	GLN
1	B	281	ASN
1	B	325	ASN
1	B	335	GLN
1	C	4	ASN
1	C	30	HIS
1	C	52	HIS
1	C	152	HIS
1	C	213	GLN
1	C	281	ASN
1	C	307	GLN
1	C	325	ASN
1	D	4	ASN
1	D	24	ASN
1	D	30	HIS
1	D	44	GLN
1	D	52	HIS
1	D	152	HIS
1	D	168	GLN
1	D	213	GLN
1	D	281	ASN
1	D	307	GLN
1	D	325	ASN
1	E	4	ASN
1	E	44	GLN
1	E	52	HIS
1	E	152	HIS
1	E	213	GLN
1	E	281	ASN

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Mol	Chain	Res	Type
1	E	307	GLN
1	E	325	ASN
1	F	152	HIS
1	F	213	GLN
1	F	281	ASN
1	F	307	GLN
1	F	325	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates [i](#)

There are no 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

### 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	340/345 (98%)	-0.15	1 (0%) 93 93	13, 22, 36, 52	0
1	B	340/345 (98%)	0.01	5 (1%) 73 72	12, 25, 43, 52	0
1	C	340/345 (98%)	0.13	9 (2%) 56 55	14, 27, 46, 55	0
1	D	340/345 (98%)	0.04	4 (1%) 79 78	14, 27, 41, 58	0
1	E	340/345 (98%)	0.16	13 (3%) 40 40	15, 29, 48, 67	0
1	F	340/345 (98%)	0.19	16 (4%) 31 31	15, 29, 48, 65	0
All	All	2040/2070 (98%)	0.06	48 (2%) 59 58	12, 26, 44, 67	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	41	PRO	10.6
1	F	345	LYS	6.8
1	F	41	PRO	6.5
1	F	42	GLU	5.0
1	E	38	HIS	4.6
1	F	44	GLN	4.6
1	F	38	HIS	4.2
1	D	345	LYS	4.1
1	F	37	ARG	4.0
1	C	267	GLY	3.9
1	B	345	LYS	3.8
1	E	345	LYS	3.6
1	E	42	GLU	3.6
1	C	266	PRO	3.5
1	C	345	LYS	3.4
1	E	39	ALA	3.3
1	E	44	GLN	3.2
1	A	345	LYS	3.2
1	E	37	ARG	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	282	ASP	3.0
1	C	283	GLU	3.0
1	F	344	ALA	3.0
1	F	57	LEU	2.9
1	F	65	ASP	2.9
1	F	17	TYR	2.9
1	E	40	LYS	2.9
1	E	283	GLU	2.8
1	C	264	GLY	2.8
1	F	267	GLY	2.7
1	B	266	PRO	2.6
1	D	288	ARG	2.5
1	E	12	LYS	2.5
1	F	58	ASP	2.4
1	F	59	GLU	2.3
1	F	272	ASP	2.3
1	B	282	ASP	2.2
1	C	260	ASN	2.2
1	E	260	ASN	2.2
1	D	260	ASN	2.2
1	F	39	ALA	2.2
1	B	42	GLU	2.1
1	E	285	VAL	2.1
1	B	285	VAL	2.1
1	D	307	GLN	2.0
1	E	282	ASP	2.0
1	F	84	ARG	2.0
1	C	272	ASP	2.0
1	C	159	VAL	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.