



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:00 PM UTC

PDB ID : 3F4L / pdb_00003f4l
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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
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

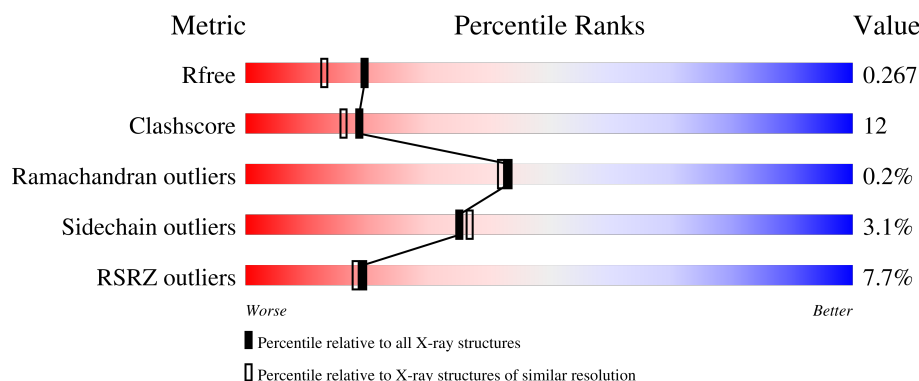
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	345	3% 80% 17% .
1	B	345	6% 75% 23% .
1	C	345	8% 74% 24% .
1	D	345	8% 77% 21% .

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Mol	Chain	Length	Quality of chain
1	E	345	9% 75% 22%
1	F	345	11% 75% 22%

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			

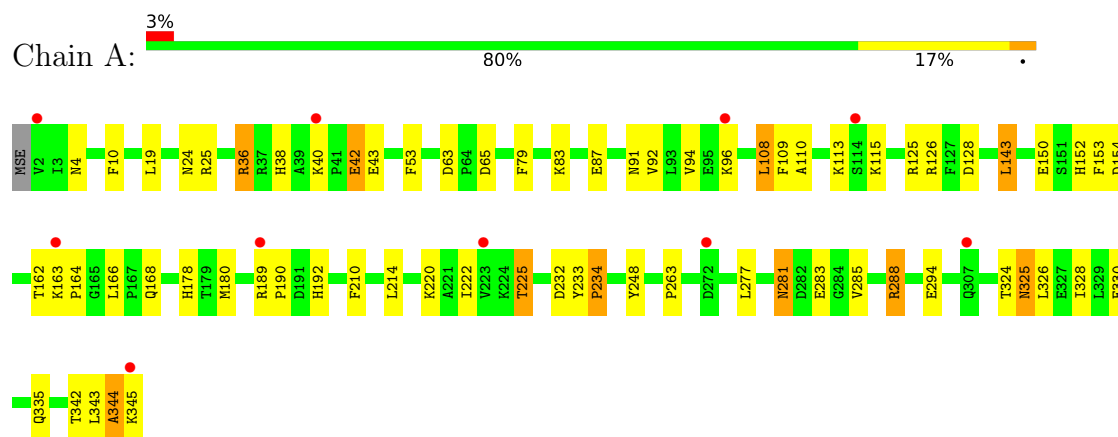
- Molecule 2 is water.

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		

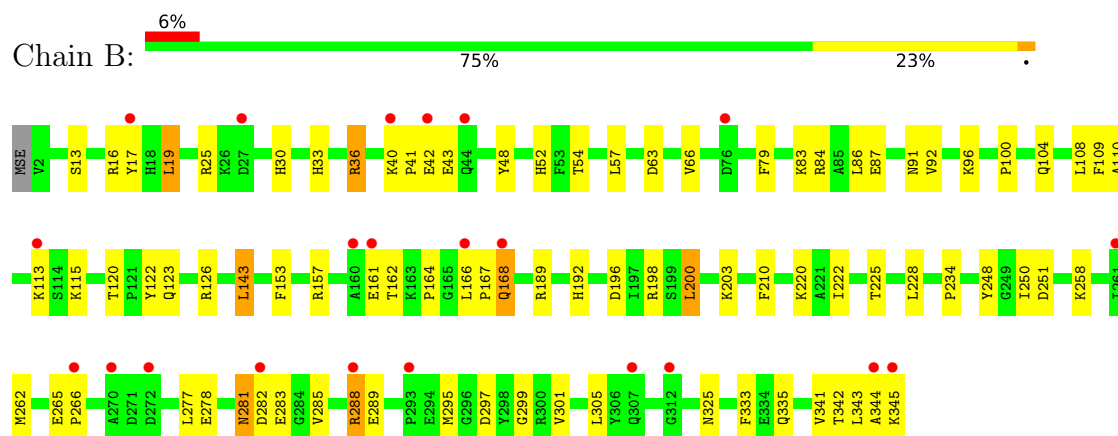
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, 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.

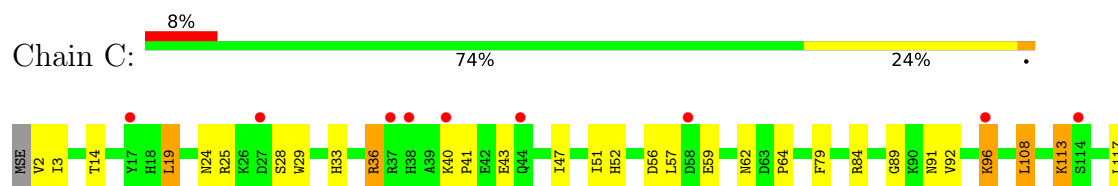
• Molecule 1: Putative oxidoreductase yhhX

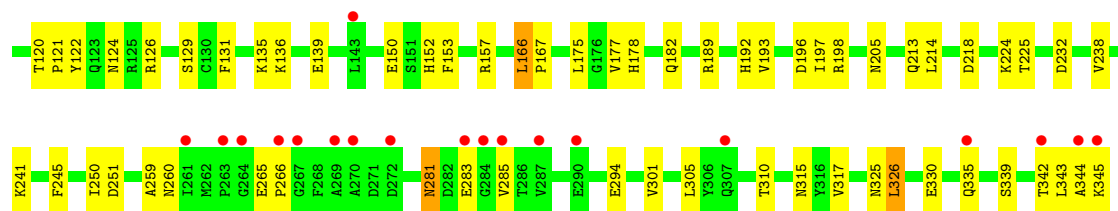


• Molecule 1: Putative oxidoreductase yhhX

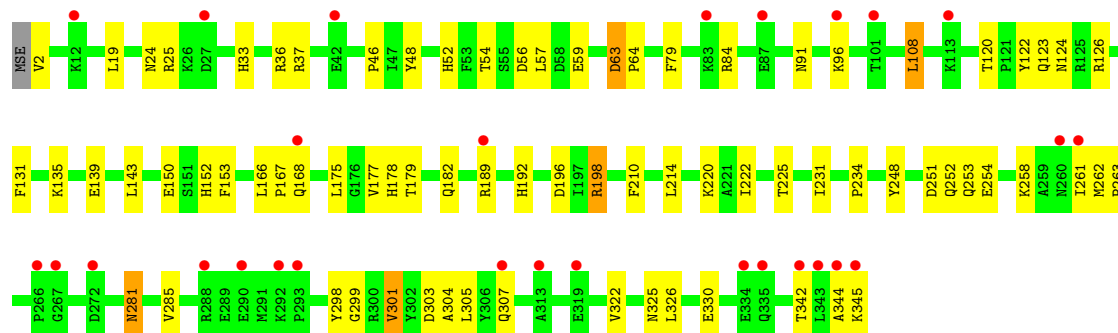
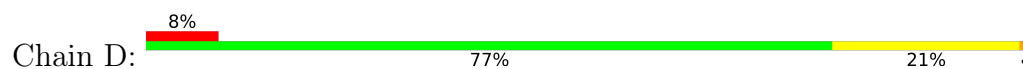


• Molecule 1: Putative oxidoreductase yhhX

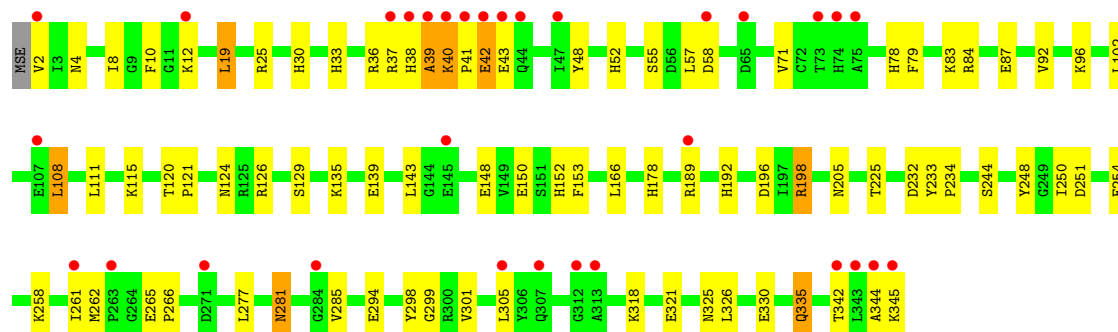
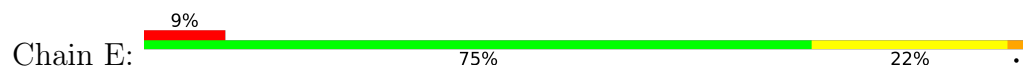




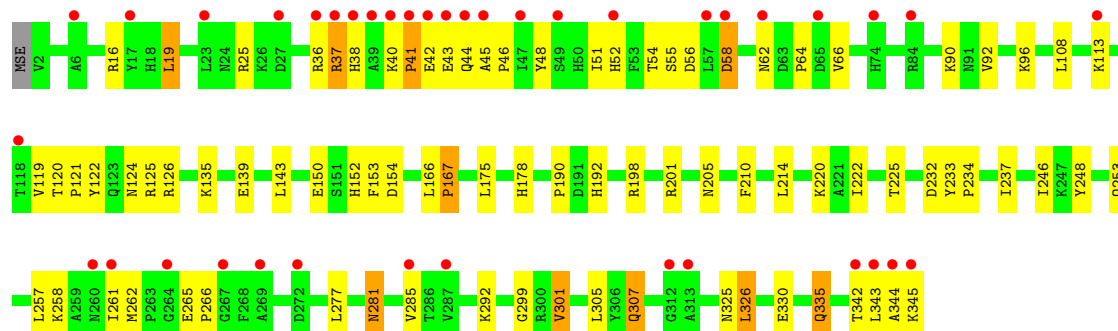
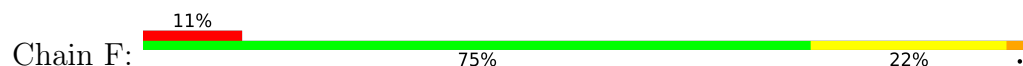
• Molecule 1: Putative oxidoreductase yhhX



• 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, α , β , γ	66.19Å 100.78Å 102.17Å 114.40° 103.31° 101.62°	Depositor
Resolution (Å)	42.60 – 2.00 42.60 – 2.00	Depositor EDS
% Data completeness (in resolution range)	86.4 (42.60-2.00) 89.5 (42.60-2.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	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 ²)	22.9	Xtriage
Anisotropy	0.210	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 33.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	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 ²)	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.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.41	0/2795	0.89	11/3780 (0.3%)
1	B	0.39	0/2795	0.88	8/3780 (0.2%)
1	C	0.39	0/2795	0.89	11/3780 (0.3%)
1	D	0.40	0/2795	0.89	7/3780 (0.2%)
1	E	0.38	0/2795	0.88	7/3780 (0.2%)
1	F	0.38	0/2795	0.90	10/3780 (0.3%)
All	All	0.39	0/16770	0.89	54/22680 (0.2%)

There are no bond length outliers.

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	344	ALA	N-CA-C	-7.69	104.42	113.88
1	D	344	ALA	N-CA-C	-7.67	101.79	112.45
1	C	126	ARG	N-CA-C	-7.22	104.35	113.01
1	B	344	ALA	N-CA-C	-7.17	104.00	112.89
1	E	126	ARG	N-CA-C	-7.09	104.60	113.18
1	D	126	ARG	N-CA-C	-6.75	105.01	113.18
1	A	126	ARG	N-CA-C	-6.68	105.09	113.18
1	F	126	ARG	N-CA-C	-6.58	105.24	113.20
1	B	120	THR	N-CA-C	6.57	114.13	108.22
1	B	92	VAL	N-CA-C	6.38	117.49	108.17
1	E	232	ASP	N-CA-C	6.33	118.96	110.35
1	F	292	LYS	N-CA-C	-6.30	101.71	109.65
1	A	344	ALA	N-CA-C	-6.28	104.53	111.82
1	F	92	VAL	N-CA-C	6.26	117.71	108.45
1	B	126	ARG	N-CA-C	-6.18	105.59	113.01
1	E	233	TYR	N-CA-C	-6.11	103.11	110.13
1	D	91	ASN	N-CA-C	-6.07	101.49	110.48
1	E	205	ASN	N-CA-C	5.95	118.33	110.36
1	A	234	PRO	N-CA-C	-5.94	102.50	111.41
1	A	233	TYR	N-CA-C	-5.86	103.39	110.13
1	A	232	ASP	N-CA-C	5.75	118.16	110.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	92	VAL	N-CA-C	5.70	116.33	108.12
1	C	205	ASN	N-CA-C	5.63	117.90	110.36
1	C	157	ARG	CA-C-N	5.54	125.38	119.28
1	C	157	ARG	C-N-CA	5.54	125.38	119.28
1	F	301	VAL	N-CA-C	-5.51	105.13	110.42
1	B	157	ARG	CA-C-N	5.46	125.29	119.28
1	B	157	ARG	C-N-CA	5.46	125.29	119.28
1	C	301	VAL	N-CA-C	-5.46	105.21	110.72
1	C	344	ALA	N-CA-C	-5.41	104.25	112.04
1	F	232	ASP	N-CA-C	5.41	118.07	110.23
1	F	198	ARG	N-CA-C	5.35	116.94	108.96
1	A	63	ASP	CA-C-N	5.33	125.14	119.28
1	A	63	ASP	C-N-CA	5.33	125.14	119.28
1	C	224	LYS	N-CA-C	5.33	117.18	108.76
1	C	91	ASN	N-CA-C	-5.32	102.61	110.48
1	B	200	LEU	N-CA-C	5.27	119.71	113.12
1	E	92	VAL	N-CA-C	5.24	116.26	108.46
1	E	78	HIS	N-CA-C	5.22	116.66	111.07
1	A	94	VAL	N-CA-C	5.20	115.40	108.11
1	F	51	ILE	N-CA-C	5.20	116.78	108.87
1	F	201	ARG	N-CA-C	5.20	116.76	111.14
1	F	233	TYR	N-CA-C	-5.20	101.46	109.41
1	C	92	VAL	N-CA-C	5.18	116.18	108.46
1	F	205	ASN	N-CA-C	5.18	116.09	110.13
1	A	91	ASN	N-CA-C	-5.17	102.83	110.48
1	A	294	GLU	N-CA-C	-5.16	103.23	110.50
1	D	252	GLN	N-CA-C	5.09	119.18	112.92
1	C	232	ASP	N-CA-C	5.08	117.27	110.35
1	C	51	ILE	N-CA-C	5.03	116.51	108.87
1	B	91	ASN	N-CA-C	-5.01	103.92	110.53
1	D	301	VAL	N-CA-C	-5.01	105.82	110.53
1	D	63	ASP	CA-C-N	5.00	124.49	119.19
1	D	63	ASP	C-N-CA	5.00	124.49	119.19

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	2731	0	2684	54	0
1	B	2731	0	2684	77	0
1	C	2731	0	2684	68	0
1	D	2731	0	2684	61	0
1	E	2731	0	2684	70	0
1	F	2731	0	2684	63	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
2	E	88	0	0	2	0
2	F	84	0	0	3	0
All	All	16970	0	16104	388	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 12.

All (388) 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	1.17	0.92
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:B:189:ARG:HD3	1:B:345:LYS:C	2.02	0.84
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:E:12:LYS:H	1:E:12:LYS:CD	1.95	0.80
1:D:96:LYS:H	1:D:96:LYS:HD3	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:B:40:LYS:HZ1	1:B:42:GLU:HB2	1.49	0.77
1:D:189:ARG:HD3	1:D:345:LYS:C	2.09	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:37:ARG:HG2	1:F:38:HIS:H	1.51	0.76
1:E:12:LYS:HD2	1:E:12:LYS:N	1.98	0.75
1:C:281:ASN:HD22	1:C:281:ASN:C	1.95	0.74
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: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:36:ARG:HD3	1:A:43:GLU:OE1	1.90	0.72
1:A:189:ARG:CZ	1:A:345:LYS:HB3	2.20	0.72
1:B:96:LYS:HE2	1:B:96:LYS:H	1.56	0.70
1:B:96:LYS:HD3	2:B:526:HOH:O	1.91	0.70
1:E:281:ASN:C	1:E:281:ASN:HD22	2.00	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:F:281:ASN:C	1:F:281:ASN:HD22	2.02	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:HD12	1:D:143:LEU:N	2.09	0.67
1:F:40:LYS:C	1:F:42:GLU:H	2.02	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:ARG:HG3	1:A:324:THR:OG1	1.98	0.63
1:F:325:ASN:HD22	1:F:325:ASN:C	2.07	0.63
1:D:281:ASN:C	1:D:281:ASN:HD22	2.06	0.63
1:C:28:SER:HG	1:C:29:TRP:CD1	2.17	0.62
1:E:57:LEU:HD23	1:E:84:ARG:CZ	2.29	0.62
1:D:96:LYS:H	1:D:96:LYS:CD	2.12	0.62
1:D:33:HIS:ND1	1:D:52:HIS:HD2	1.97	0.62
1:F:143:LEU:N	1:F:143:LEU:HD12	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
1:D:150:GLU:OE2	1:D:152:HIS:HE1	1.83	0.61
1:A:281:ASN:C	1:A:281:ASN:HD22	2.08	0.61
1:B:281:ASN:HD21	1:B:285:VAL:H	1.49	0.61
1:B:281:ASN:C	1:B:281:ASN:HD22	2.09	0.60
1:D:220:LYS:HE3	1:D:222:ILE:HD11	1.82	0.60
1:D:96:LYS:HD3	1:D:96:LYS:N	2.15	0.60
1:A:83:LYS:O	1:A:87:GLU:HG3	2.01	0.60
1:A:96:LYS:HE3	1:A:178:HIS:CE1	2.37	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:HD3	1:A:96:LYS:N	2.15	0.59
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:B:40:LYS:NZ	1:B:42:GLU:HB2	2.19	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:D:168:GLN:H	1:D:168:GLN:CD	2.12	0.56
1:C:281:ASN:ND2	1:C:283:GLU:H	2.02	0.56
1:A:234:PRO:HG3	1:A:248:TYR:CZ	2.40	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:C:325:ASN:C	1:C:325:ASN:HD22	2.14	0.55
1:E:96:LYS:N	1:E:96:LYS:HD3	2.22	0.55
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:E:39:ALA:O	1:E:40:LYS:HB2	2.05	0.54
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:F:40:LYS:C	1:F:42:GLU:N	2.66	0.54
1:A:143:LEU:N	1:A:143:LEU:HD12	2.23	0.54
1:D:131:PHE:CE1	1:D:182:GLN:HB3	2.43	0.54
1:C:136:LYS:HE3	1:C:294:GLU:CD	2.32	0.54
1:E:42:GLU:HB3	1:E:48:TYR:CE2	2.43	0.54
1:E:261:ILE:C	1:E:262:MSE:HE2	2.32	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:HG3	1:F:125:ARG:HH11	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:42:GLU:O	1:E:48:TYR:HD2	1.91	0.53
1:E:318:LYS:HB2	1:E:321:GLU:HG3	1.89	0.53
1:F:153:PHE:HB3	1:F:225:THR:HG22	1.91	0.53
1:B:86:LEU:O	1:B:115:LYS:HE2	2.09	0.53
1:B:109:PHE:O	1:B:113:LYS:HE2	2.08	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:288:ARG:HG3	1:A:288:ARG:HH11	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:HD2	1:F:37:ARG:H	1.76	0.51
1:A:189:ARG:HG2	1:A:345:LYS:OXT	2.11	0.51
1:B:343:LEU:HB3	1:B:345:LYS:OXT	2.10	0.51
1:C:96:LYS:N	1:C:96:LYS:CD	2.72	0.51
1:C:113:LYS:HB2	1:C:113:LYS:NZ	2.26	0.51
1:B:96:LYS:HZ2	1:B:122:TYR:HB3	1.75	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:E:281:ASN:C	1:E:281:ASN:ND2	2.69	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:C:57:LEU:HD23	1:C:84:ARG:CZ	2.41	0.50
1:B:281:ASN:ND2	1:B:281:ASN:C	2.69	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: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: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:C:259:ALA:O	1:C:260:ASN:HB2	2.13	0.49
1:D:143:LEU:N	1:D:143:LEU:CD1	2.74	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:16:ARG:HH12	1:B:258:LYS:HZ1	1.60	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:203:LYS:HZ1	1:C:218:ASP:CG	2.21	0.49
1:B:203:LYS:NZ	1:C:218:ASP:OD2	2.43	0.49
1:B:295:MSE:HE2	2:B:365:HOH:O	2.12	0.49
1:F:56:ASP:CG	1:F:58:ASP:HB2	2.38	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
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:B:168:GLN:H	1:B:168:GLN:CD	2.21	0.48
1:C:281:ASN:C	1:C:281:ASN:ND2	2.64	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:H	1:B:96:LYS:CD	2.26	0.48
1:C:2:VAL:HG23	1:C:28:SER:C	2.39	0.48
1:F:36:ARG:C	1:F:36:ARG:CD	2.87	0.48
1:A:109:PHE:C	1:A:113:LYS:HE2	2.39	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:N	1:E:166:LEU:HD22	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:325:ASN:C	1:B:325:ASN:HD22	2.22	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:D:281:ASN:ND2	1:D:281:ASN:C	2.73	0.47
1:E:265:GLU:HB3	1:E:266:PRO:HD2	1.97	0.47
1:F:143:LEU:N	1:F:143:LEU:CD1	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:N	1:B:166:LEU:HD22	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:F:261:ILE:C	1:F:262:MSE:HE2	2.40	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
1:E:178:HIS:HD2	2:E:352:HOH:O	1.98	0.47
1:F:96:LYS:HE2	1:F:96:LYS:H	1.79	0.47
1:B:200:LEU:CD2	1:C:218:ASP:HB3	2.45	0.46
1:C:113:LYS:HB2	1:C:113:LYS:HZ2	1.81	0.46
1:A:189:ARG:HD3	1:A:345:LYS:C	2.40	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:B:109:PHE:C	1:B:113:LYS:HE2	2.41	0.46
1:D:251:ASP:OD2	1:D:253:GLN:HG3	2.16	0.46
1:F:96:LYS:H	1:F:96:LYS:CE	2.29	0.46
1:D:166:LEU:N	1:D:166:LEU:HD22	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:175:LEU:N	1:F:175:LEU:HD22	2.32	0.45
1:A:344:ALA:O	1:A:345:LYS:HG3	2.17	0.45
1:E:96:LYS:H	1:E:96:LYS:CD	2.29	0.45
1:E:325:ASN:C	1:E:325:ASN:HD22	2.25	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: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:N	1:C:175:LEU:HD22	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:N	1:E:96:LYS:CD	2.80	0.45
1:D:254:GLU:HG2	1:D:258:LYS:HE2	1.99	0.45
1:D:304:ALA:HA	1:D:307:GLN:CD	2.42	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:40:LYS:HD2	1:B:41:PRO:CD	2.46	0.44
1:B:109:PHE:O	1:B:113:LYS:HG3	2.18	0.44
1:F:119:VAL:HG13	1:F:119:VAL:O	2.16	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
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:HD22	1:A:166:LEU:N	2.33	0.44
1:E:36:ARG:HG2	1:E:36:ARG:HH11	1.83	0.44
1:E:42:GLU:O	1:E:43:GLU:C	2.60	0.44
1:C:250:ILE:HG12	1:C:251:ASP:N	2.33	0.44
1:D:2:VAL:O	1:D:2:VAL:HG13	2.17	0.44
1:D:261:ILE:C	1:D:262:MSE:HE2	2.43	0.44
1:B:96:LYS:HD3	1:B:96:LYS:N	2.33	0.44
1:E:36:ARG:NE	1:E:43:GLU:OE2	2.42	0.44
1:E:124:ASN:HB2	1:E:298:TYR:CD1	2.53	0.44
1:B:33:HIS:ND1	1:B:52:HIS:HD2	2.15	0.44
1:C:214:LEU:N	1:C:214:LEU:HD12	2.33	0.44
1:D:46:PRO:HD2	2:D:468:HOH:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:124:ASN:HB2	1:D:298:TYR:CD1	2.53	0.44
1:E:8:ILE:HB	1:E:71:VAL:HA	1.99	0.44
1:A:189:ARG:CD	1:A:345:LYS:C	2.91	0.43
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:40:LYS:CD	1:C:41:PRO:HD2	2.49	0.43
1:C:122:TYR:CZ	1:C:124:ASN:HB3	2.52	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: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:D:325:ASN:C	1:D:325:ASN:HD22	2.26	0.43
1:E:153:PHE:HB3	1:E:225:THR:HG22	1.99	0.43
1:F:210:PHE:CZ	1:F:225:THR:HG23	2.53	0.43
1:C:197:ILE:C	1:C:198:ARG:HG2	2.44	0.43
1:E:40:LYS:C	1:E:42:GLU:N	2.76	0.43
1:F:96:LYS:H	1:F:96:LYS:CD	2.32	0.43
1:C:166:LEU:HB3	1:C:167:PRO:CD	2.49	0.42
1:E:10:PHE:HE2	1:E:42:GLU:HB2	1.83	0.42
1:E:281:ASN:ND2	1:E:285:VAL:H	2.17	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
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:HG2	1:C:25:ARG:HH11	1.85	0.42
1:C:335:GLN:HG3	1:C:339:SER:HB3	2.02	0.42
1:D:214:LEU:N	1:D:214:LEU:HD12	2.35	0.42
1:B:40:LYS:CD	1:B:41:PRO:HD2	2.45	0.41
1:B:100:PRO:HD2	1:B:104:GLN:OE1	2.20	0.41
1:F:96:LYS:HD3	1:F:96:LYS:N	2.34	0.41
1:B:166:LEU:HB3	1:B:167:PRO:HD2	2.01	0.41
1:B:167:PRO:HB3	1:B:333:PHE:CE1	2.54	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:325:ASN:HD22	1:A:325:ASN:C	2.27	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:F:37:ARG:O	1:F:38:HIS:C	2.62	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:A:283:GLU:N	1:A:283:GLU:CD	2.79	0.41
1:E:40:LYS:C	1:E:42:GLU:H	2.29	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: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: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:63:ASP:HA	1:D:64:PRO:HD3	1.87	0.40
1:D:192:HIS:ND1	1:D:342:THR:HG22	2.36	0.40
1:E:38:HIS:O	1:E:39:ALA:HB2	2.21	0.40
1:C:36:ARG:HD2	1:C:43:GLU:OE1	2.21	0.40
1:D:168:GLN:CD	1:D:168:GLN:N	2.78	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:HD13	1:C:305:LEU:HA	1.96	0.40
1:C:335:GLN:HG3	1:C:339:SER:CB	2.52	0.40
1:E:2:VAL:HG13	1:E:2:VAL:O	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:F:16:ARG:NH1	1:F:258:LYS:NZ	2.69	0.40
1:F:43:GLU:C	1:F:45:ALA:H	2.30	0.40
1:A:288:ARG:HG3	1:A:288:ARG:NH1	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%)	21	17
1	F	342/345 (99%)	323 (94%)	17 (5%)	2 (1%)	21	17
All	All	2052/2070 (99%)	1993 (97%)	55 (3%)	4 (0%)	43	42

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 [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	294/290 (101%)	283 (96%)	11 (4%)	30	30

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	294/290 (101%)	286 (97%)	8 (3%)	39	42
1	C	294/290 (101%)	284 (97%)	10 (3%)	32	33
1	D	294/290 (101%)	288 (98%)	6 (2%)	48	54
1	E	294/290 (101%)	286 (97%)	8 (3%)	39	42
1	F	294/290 (101%)	283 (96%)	11 (4%)	30	30
All	All	1764/1740 (101%)	1710 (97%)	54 (3%)	35	37

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LEU
1	A	36	ARG
1	A	42	GLU
1	A	108	LEU
1	A	143	LEU
1	A	163	LYS
1	A	168	GLN
1	A	225	THR
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	143	LEU
1	B	168	GLN
1	B	281	ASN
1	B	288	ARG
1	B	289	GLU
1	C	19	LEU
1	C	24	ASN
1	C	36	ARG
1	C	96	LYS
1	C	108	LEU
1	C	113	LYS
1	C	166	LEU
1	C	281	ASN
1	C	326	LEU
1	C	343	LEU
1	D	36	ARG

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Mol	Chain	Res	Type
1	D	54	THR
1	D	108	LEU
1	D	120	THR
1	D	198	ARG
1	D	281	ASN
1	E	19	LEU
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	335	GLN
1	F	19	LEU
1	F	37	ARG
1	F	44	GLN
1	F	55	SER
1	F	58	ASP
1	F	108	LEU
1	F	167	PRO
1	F	281	ASN
1	F	307	GLN
1	F	326	LEU
1	F	335	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (58) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	ASN
1	A	24	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

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Mol	Chain	Res	Type
1	B	213	GLN
1	B	252	GLN
1	B	281	ASN
1	B	311	HIS
1	B	315	ASN
1	B	325	ASN
1	B	335	GLN
1	C	4	ASN
1	C	30	HIS
1	C	44	GLN
1	C	52	HIS
1	C	104	GLN
1	C	152	HIS
1	C	213	GLN
1	C	260	ASN
1	C	281	ASN
1	C	307	GLN
1	C	325	ASN
1	D	4	ASN
1	D	24	ASN
1	D	44	GLN
1	D	52	HIS
1	D	152	HIS
1	D	168	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	252	GLN
1	E	281	ASN
1	E	307	GLN
1	E	311	HIS
1	E	325	ASN
1	F	24	ASN
1	F	62	ASN
1	F	152	HIS
1	F	260	ASN
1	F	281	ASN
1	F	307	GLN

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Mol	Chain	Res	Type
1	F	311	HIS
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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
1	A	340/345 (98%)	0.32	10 (2%)	53	52	13, 22, 36, 52	0
1	B	340/345 (98%)	0.61	22 (6%)	25	23	12, 25, 43, 52	0
1	C	340/345 (98%)	0.74	28 (8%)	17	16	14, 27, 46, 55	0
1	D	340/345 (98%)	0.62	28 (8%)	17	16	14, 27, 41, 58	0
1	E	340/345 (98%)	0.78	31 (9%)	15	13	15, 29, 48, 67	0
1	F	340/345 (98%)	0.87	39 (11%)	9	8	15, 29, 48, 65	0
All	All	2040/2070 (98%)	0.66	158 (7%)	19	18	12, 26, 44, 67	0

All (158) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	41	PRO	8.7
1	F	41	PRO	6.6
1	F	39	ALA	6.5
1	E	38	HIS	6.1
1	A	345	LYS	6.0
1	B	345	LYS	6.0
1	F	42	GLU	5.8
1	C	345	LYS	5.2
1	E	40	LYS	5.2
1	F	37	ARG	5.2
1	F	345	LYS	5.1
1	F	344	ALA	4.9
1	E	39	ALA	4.8
1	E	43	GLU	4.7
1	F	38	HIS	4.6
1	D	288	ARG	4.4
1	E	345	LYS	4.2
1	F	40	LYS	4.2
1	E	37	ARG	4.2

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Mol	Chain	Res	Type	RSRZ
1	F	44	GLN	4.1
1	C	285	VAL	4.1
1	D	345	LYS	3.9
1	D	307	GLN	3.9
1	B	261	ILE	3.8
1	E	307	GLN	3.8
1	F	17	TYR	3.8
1	E	261	ILE	3.8
1	F	58	ASP	3.7
1	E	42	GLU	3.7
1	C	344	ALA	3.6
1	E	44	GLN	3.6
1	F	57	LEU	3.5
1	F	43	GLU	3.5
1	B	272	ASP	3.5
1	B	266	PRO	3.4
1	E	312	GLY	3.4
1	C	44	GLN	3.4
1	D	96	LYS	3.4
1	C	143	LEU	3.4
1	E	2	VAL	3.3
1	E	344	ALA	3.3
1	C	40	LYS	3.3
1	C	27	ASP	3.3
1	F	264	GLY	3.3
1	D	344	ALA	3.3
1	C	272	ASP	3.2
1	D	335	GLN	3.2
1	F	261	ILE	3.1
1	F	343	LEU	3.1
1	C	58	ASP	3.1
1	F	272	ASP	3.1
1	B	168	GLN	3.1
1	D	334	GLU	3.1
1	C	37	ARG	3.0
1	D	342	THR	3.0
1	B	42	GLU	3.0
1	C	267	GLY	3.0
1	B	166	LEU	2.9
1	E	12	LYS	2.9
1	E	342	THR	2.8
1	F	49	SER	2.8

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Mol	Chain	Res	Type	RSRZ
1	E	313	ALA	2.8
1	C	261	ILE	2.8
1	B	282	ASP	2.8
1	C	264	GLY	2.8
1	F	342	THR	2.8
1	C	287	VAL	2.8
1	D	260	ASN	2.7
1	E	58	ASP	2.7
1	D	83	LYS	2.7
1	B	288	ARG	2.7
1	E	284	GLY	2.7
1	F	267	GLY	2.7
1	E	107	GLU	2.6
1	A	189	ARG	2.6
1	B	161	GLU	2.6
1	C	270	ALA	2.6
1	E	263	PRO	2.6
1	A	307	GLN	2.6
1	F	6	ALA	2.6
1	A	2	VAL	2.6
1	B	344	ALA	2.6
1	C	266	PRO	2.6
1	B	307	GLN	2.5
1	F	52	HIS	2.5
1	C	307	GLN	2.5
1	D	313	ALA	2.5
1	A	96	LYS	2.5
1	E	343	LEU	2.5
1	D	12	LYS	2.5
1	E	74	HIS	2.5
1	C	284	GLY	2.5
1	D	267	GLY	2.5
1	D	261	ILE	2.5
1	A	272	ASP	2.4
1	D	87	GLU	2.4
1	C	342	THR	2.4
1	B	40	LYS	2.4
1	A	163	LYS	2.4
1	E	189	ARG	2.4
1	D	343	LEU	2.4
1	F	36	ARG	2.4
1	C	17	TYR	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	319	GLU	2.3
1	B	312	GLY	2.3
1	C	96	LYS	2.3
1	B	160	ALA	2.3
1	F	62	ASN	2.3
1	F	27	ASP	2.3
1	F	84	ARG	2.3
1	F	313	ALA	2.3
1	B	17	TYR	2.3
1	D	189	ARG	2.3
1	F	312	GLY	2.3
1	D	42	GLU	2.3
1	D	293	PRO	2.3
1	B	27	ASP	2.3
1	F	23	LEU	2.3
1	E	145	GLU	2.2
1	E	75	ALA	2.2
1	D	27	ASP	2.2
1	E	271	ASP	2.2
1	D	290	GLU	2.2
1	F	45	ALA	2.2
1	C	263	PRO	2.2
1	D	101	THR	2.2
1	E	73	THR	2.2
1	D	272	ASP	2.2
1	C	38	HIS	2.2
1	E	305	LEU	2.2
1	A	40	LYS	2.2
1	A	223	VAL	2.1
1	F	287	VAL	2.1
1	F	74	HIS	2.1
1	B	44	GLN	2.1
1	D	168	GLN	2.1
1	E	47	ILE	2.1
1	C	283	GLU	2.1
1	F	285	VAL	2.1
1	A	114	SER	2.1
1	E	65	ASP	2.1
1	D	113	LYS	2.1
1	B	270	ALA	2.1
1	C	114	SER	2.1
1	C	269	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	76	ASP	2.0
1	C	335	GLN	2.0
1	D	266	PRO	2.0
1	F	118	THR	2.0
1	F	47	ILE	2.0
1	C	290	GLU	2.0
1	F	269	ALA	2.0
1	F	260	ASN	2.0
1	F	65	ASP	2.0
1	B	113	LYS	2.0
1	D	292	LYS	2.0
1	F	113	LYS	2.0
1	B	293	PRO	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.