



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:51 PM UTC

PDB ID : 5YIK / pdb_00005yik
Title : Structure of a Legionella effector with its substrate
Authors : Feng, Y.; Dong, Y.; Liu, Z.
Deposited on : 2017-10-05
Resolution : 3.10 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

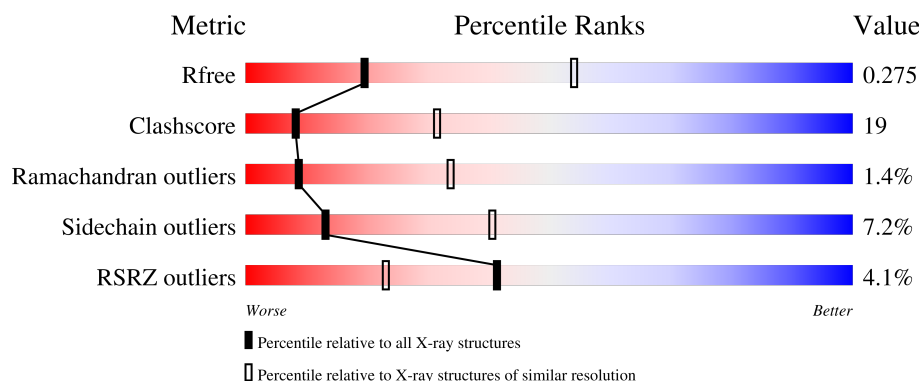
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

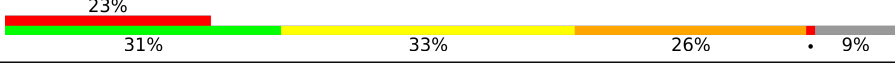
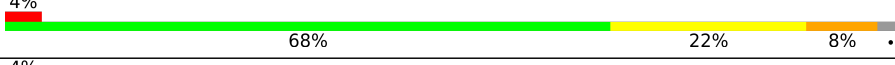
The reported resolution of this entry is 3.10 Å.

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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	968	
2	C	78	
2	D	78	
2	F	78	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9381 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 SdeA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	959	7601	4767	1330	1477	27	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1191	LEU	-	expression tag	UNP Q6RCR0
A	1192	GLU	-	expression tag	UNP Q6RCR0
A	1193	HIS	-	expression tag	UNP Q6RCR0
A	1194	HIS	-	expression tag	UNP Q6RCR0
A	1195	HIS	-	expression tag	UNP Q6RCR0
A	1196	HIS	-	expression tag	UNP Q6RCR0
A	1197	HIS	-	expression tag	UNP Q6RCR0
A	1198	HIS	-	expression tag	UNP Q6RCR0

- Molecule 2 is a protein called ubiquitin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	71	563	356	94	112	1	0	0	0
2	D	76	601	378	105	117	1	0	0	0
2	F	78	616	387	109	119	1	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	SER	-	expression tag	UNP P0CG47
C	0	HIS	-	expression tag	UNP P0CG47
D	-1	SER	-	expression tag	UNP P0CG47
D	0	HIS	-	expression tag	UNP P0CG47

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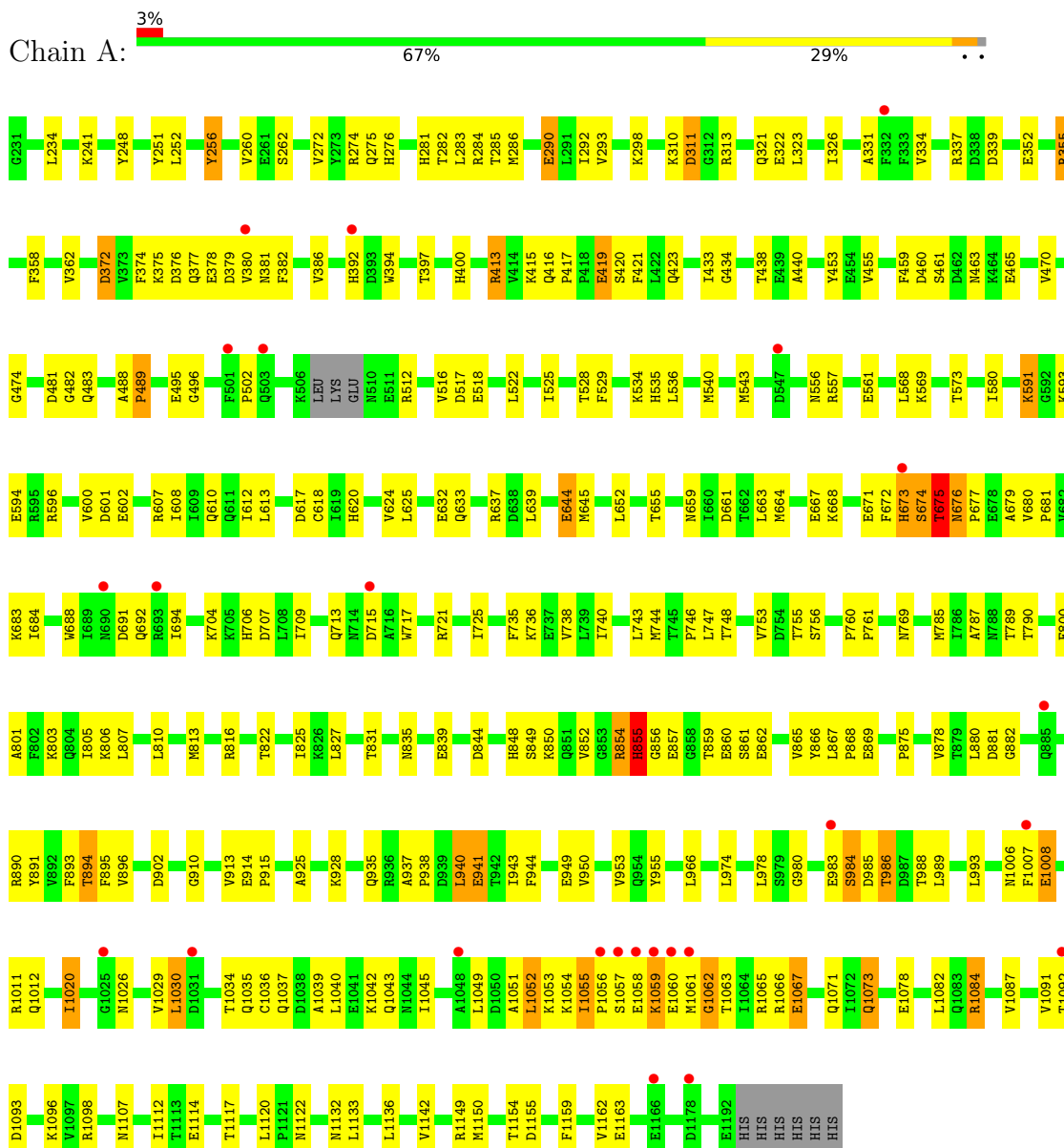
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Chain	Residue	Modelled	Actual	Comment	Reference
F	-1	SER	-	expression tag	UNP P0CG47
F	0	HIS	-	expression tag	UNP P0CG47

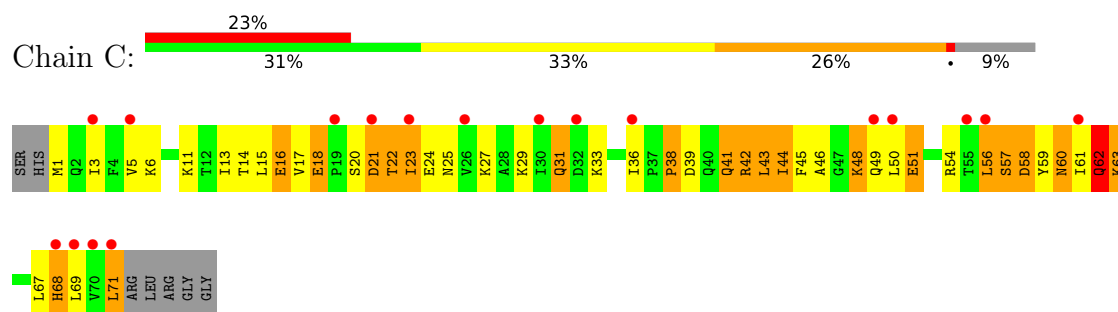
3 Residue-property plots

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.

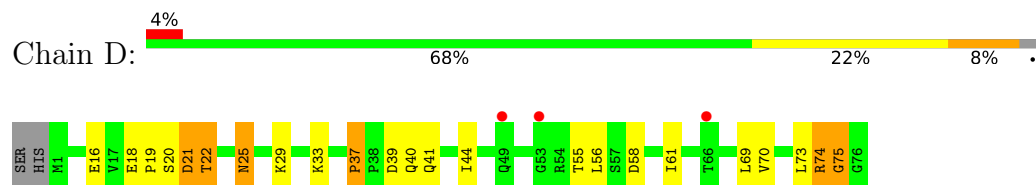
• Molecule 1: SdeA



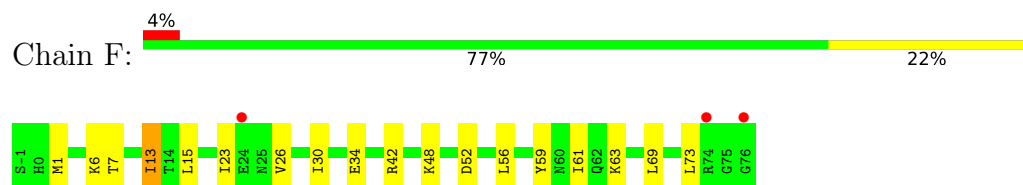
• Molecule 2: ubiquitin



- Molecule 2: ubiquitin



- Molecule 2: ubiquitin



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	108.85Å 145.89Å 104.09Å 90.00° 104.46° 90.00°	Depositor
Resolution (Å)	42.54 – 3.10 42.54 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.7 (42.54-3.10) 88.5 (42.54-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.53 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.223 , 0.279 0.234 , 0.275	Depositor DCC
R_{free} test set	1290 reflections (4.52%)	wwPDB-VP
Wilson B-factor (Å ²)	54.3	Xtriage
Anisotropy	0.205	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 36.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	9381	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.86% 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 [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	0.61	7/7734 (0.1%)	0.82	14/10439 (0.1%)
2	C	0.71	0/569	1.10	3/767 (0.4%)
2	D	0.55	0/607	0.93	3/816 (0.4%)
2	F	0.34	0/623	0.83	0/838
All	All	0.60	7/9533 (0.1%)	0.85	20/12860 (0.2%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	893	PHE	C-O	-5.73	1.17	1.24
1	A	676	ASN	CA-C	-5.67	1.49	1.53
1	A	894	THR	C-O	-5.49	1.17	1.24
1	A	676	ASN	C-N	5.42	1.40	1.33
1	A	895	PHE	CA-C	-5.27	1.46	1.52
1	A	1039	ALA	C-O	-5.09	1.18	1.24
1	A	677	PRO	N-CD	5.04	1.54	1.47

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	854	ARG	N-CA-C	9.28	121.48	111.36
2	D	75	GLY	N-CA-C	9.24	123.82	112.73
1	A	938	PRO	N-CA-CB	7.63	109.69	103.36
1	A	855	HIS	N-CA-C	7.50	124.43	113.40
1	A	984	SER	CB-CA-C	-7.46	105.90	116.34
1	A	676	ASN	CA-C-N	-6.61	112.82	119.56
1	A	676	ASN	C-N-CA	-6.61	112.82	119.56
2	C	62	GLN	N-CA-C	6.55	118.16	108.60
1	A	941	GLU	N-CA-C	-6.26	104.54	111.36
1	A	529	PHE	CA-C-N	6.15	125.84	119.56
1	A	529	PHE	C-N-CA	6.15	125.84	119.56
1	A	1067	GLU	N-CA-C	-6.00	105.61	113.17
1	A	675	THR	N-CA-C	-5.98	100.46	109.79

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	591	LYS	N-CA-C	5.86	117.47	111.14
1	A	674	SER	N-CA-C	-5.83	104.83	111.07
2	C	23	ILE	N-CA-C	-5.79	104.71	110.62
2	C	33	LYS	N-CA-C	5.69	117.17	110.97
2	D	37	PRO	CA-C-N	5.39	125.05	119.56
2	D	37	PRO	C-N-CA	5.39	125.05	119.56
1	A	1063	THR	N-CA-C	-5.28	106.43	112.92

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	7601	0	7518	263	0
2	C	563	0	586	79	0
2	D	601	0	629	21	0
2	F	616	0	638	11	0
All	All	9381	0	9371	362	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 19.

All (362) 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:822:THR:HG21	1:A:827:LEU:CD2	1.33	1.57
2:C:42:ARG:HG2	2:C:49:GLN:NE2	1.54	1.20
1:A:419:GLU:HG3	1:A:516:VAL:HG11	1.21	1.12
1:A:822:THR:CG2	1:A:827:LEU:CD2	2.29	1.10
1:A:822:THR:CG2	1:A:827:LEU:HD23	1.80	1.10
1:A:1026:ASN:HD22	1:A:1061:MET:HE1	1.15	1.09
2:C:48:LYS:NZ	2:C:49:GLN:O	1.88	1.06
1:A:822:THR:HG21	1:A:827:LEU:HD23	1.03	1.01
2:C:20:SER:O	2:C:56:LEU:HD22	1.59	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:1:MET:HB3	2:C:17:VAL:O	1.61	1.00
1:A:822:THR:HG21	1:A:827:LEU:HD22	1.43	1.00
1:A:1058:GLU:HG2	1:A:1059:LYS:HD3	1.45	0.98
1:A:940:LEU:HD12	1:A:940:LEU:H	1.29	0.97
1:A:375:LYS:HB3	1:A:379:ASP:OD2	1.65	0.97
1:A:709:ILE:CG2	1:A:852:VAL:HG23	1.95	0.96
1:A:1154:THR:O	1:A:1155:ASP:OD1	1.82	0.96
1:A:993:LEU:HD21	1:A:1060:GLU:HB3	1.45	0.94
1:A:415:LYS:HD3	1:A:421:PHE:CE2	2.02	0.94
2:C:42:ARG:HG2	2:C:49:GLN:HE22	1.22	0.93
1:A:1055:ILE:HG12	1:A:1071:GLN:OE1	1.69	0.92
2:C:21:ASP:O	2:C:22:THR:OG1	1.89	0.90
1:A:993:LEU:CD2	1:A:1060:GLU:HB3	2.01	0.90
2:C:42:ARG:CG	2:C:49:GLN:NE2	2.35	0.89
1:A:1093:ASP:HB3	1:A:1096:LYS:HB3	1.54	0.89
1:A:928:LYS:HE2	1:A:1034:THR:HG22	1.55	0.89
1:A:415:LYS:CD	1:A:421:PHE:CE2	2.56	0.88
2:C:58:ASP:O	2:C:59:TYR:CD2	2.27	0.88
1:A:1058:GLU:N	1:A:1058:GLU:OE2	2.07	0.87
2:C:43:LEU:C	2:C:44:ILE:HD12	1.98	0.87
1:A:709:ILE:HG22	1:A:852:VAL:HG23	1.57	0.86
1:A:940:LEU:O	1:A:944:PHE:HD1	1.57	0.86
2:C:20:SER:O	2:C:56:LEU:CD2	2.23	0.86
2:C:17:VAL:HG22	2:C:18:GLU:H	1.39	0.86
1:A:397:THR:HG1	1:A:400:HIS:HD1	1.24	0.85
2:C:51:GLU:OE2	2:C:51:GLU:N	2.09	0.85
1:A:1056:PRO:O	1:A:1057:SER:OG	1.91	0.85
2:C:20:SER:O	2:C:56:LEU:HD13	1.76	0.84
1:A:706:HIS:CD2	2:D:75:GLY:HA3	2.13	0.83
1:A:855:HIS:HA	1:A:859:THR:HG21	1.59	0.82
2:D:40:GLN:HG2	2:D:73:LEU:HD13	1.58	0.82
2:D:41:GLN:HB2	2:D:69:LEU:HD11	1.60	0.81
1:A:1035:GLN:OE1	1:A:1054:LYS:NZ	2.14	0.81
2:C:24:GLU:HG3	2:C:25:ASN:N	1.96	0.80
2:C:42:ARG:CG	2:C:49:GLN:HE22	1.93	0.80
1:A:985:ASP:O	1:A:986:THR:OG1	1.98	0.79
1:A:1026:ASN:ND2	1:A:1061:MET:HE1	1.95	0.79
2:C:56:LEU:HB3	2:C:61:ILE:HD13	1.65	0.78
2:C:56:LEU:HB3	2:C:61:ILE:CD1	2.14	0.78
1:A:744:MET:O	1:A:748:THR:HG23	1.85	0.77
1:A:1036:CYS:SG	1:A:1052:LEU:CD2	2.73	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:803:LYS:O	1:A:807:LEU:HG	1.86	0.76
2:C:44:ILE:HA	2:C:48:LYS:O	1.84	0.76
1:A:928:LYS:CE	1:A:1034:THR:HG22	2.16	0.75
2:C:20:SER:O	2:C:56:LEU:CD1	2.35	0.75
2:C:43:LEU:HD22	2:C:43:LEU:H	1.51	0.75
1:A:986:THR:OG1	1:A:988:THR:HG22	1.87	0.74
2:C:43:LEU:HB2	2:C:50:LEU:HD12	1.66	0.74
1:A:1036:CYS:SG	1:A:1052:LEU:HD23	2.29	0.73
1:A:859:THR:HG23	1:A:861:SER:H	1.53	0.73
2:C:17:VAL:HG22	2:C:18:GLU:N	2.03	0.73
1:A:419:GLU:HG3	1:A:516:VAL:CG1	2.10	0.72
2:C:43:LEU:HD22	2:C:43:LEU:N	2.04	0.72
1:A:940:LEU:HD12	1:A:940:LEU:N	2.03	0.72
1:A:633:GLN:OE1	1:A:637:ARG:NH1	2.22	0.72
2:C:58:ASP:O	2:C:59:TYR:HD2	1.70	0.71
1:A:854:ARG:O	1:A:855:HIS:HB2	1.89	0.71
1:A:822:THR:HG21	1:A:827:LEU:HD21	1.62	0.71
1:A:481:ASP:O	1:A:483:GLN:N	2.23	0.71
1:A:415:LYS:HD2	1:A:421:PHE:CE2	2.24	0.71
2:C:62:GLN:N	2:C:62:GLN:OE1	2.24	0.70
2:C:1:MET:HE2	2:C:18:GLU:HA	1.74	0.69
1:A:674:SER:HA	1:A:721:ARG:NH2	2.08	0.69
1:A:1091:VAL:O	1:A:1093:ASP:N	2.26	0.69
1:A:415:LYS:CD	1:A:421:PHE:HE2	2.07	0.68
1:A:831:THR:CG2	2:D:44:ILE:HD12	2.23	0.68
1:A:663:LEU:O	1:A:667:GLU:HG2	1.93	0.68
1:A:707:ASP:OD1	2:D:74:ARG:HD3	1.94	0.68
1:A:910:GLY:HA2	1:A:913:VAL:HG12	1.74	0.67
1:A:1051:ALA:O	1:A:1054:LYS:HB2	1.94	0.67
1:A:1055:ILE:CG1	1:A:1071:GLN:OE1	2.43	0.67
2:C:22:THR:OG1	2:C:25:ASN:HB2	1.94	0.67
1:A:1059:LYS:O	1:A:1061:MET:N	2.27	0.67
2:C:1:MET:HG3	2:C:63:LYS:HB3	1.77	0.66
1:A:857:GLU:O	1:A:857:GLU:HG2	1.95	0.66
2:C:45:PHE:HD2	2:C:67:LEU:CD2	2.09	0.66
2:C:44:ILE:HG23	2:C:48:LYS:N	2.12	0.65
2:C:43:LEU:HA	2:C:69:LEU:HD22	1.79	0.65
1:A:953:VAL:HG21	1:A:966:LEU:HD12	1.79	0.64
2:C:45:PHE:HD2	2:C:67:LEU:HD23	1.62	0.64
1:A:985:ASP:HA	1:A:989:LEU:HD13	1.80	0.64
1:A:593:LYS:HE3	1:A:844:ASP:O	1.98	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1036:CYS:SG	1:A:1052:LEU:HD21	2.38	0.63
1:A:940:LEU:O	1:A:944:PHE:CD1	2.48	0.63
1:A:1059:LYS:HD2	1:A:1059:LYS:N	2.12	0.63
1:A:707:ASP:OD1	2:D:74:ARG:NE	2.30	0.63
1:A:902:ASP:HA	1:A:1149:ARG:HH22	1.63	0.63
2:C:22:THR:OG1	2:C:25:ASN:CB	2.48	0.62
1:A:417:PRO:HG2	1:A:420:SER:HB2	1.82	0.62
1:A:880:LEU:HD13	1:A:891:TYR:CE1	2.35	0.62
1:A:1029:VAL:HG22	1:A:1056:PRO:HD2	1.82	0.62
1:A:1049:LEU:O	1:A:1053:LYS:HG3	2.00	0.61
1:A:709:ILE:HG22	1:A:852:VAL:CG2	2.29	0.61
2:C:56:LEU:O	2:C:61:ILE:HG23	2.00	0.61
1:A:512:ARG:NH2	1:A:518:GLU:OE2	2.34	0.60
1:A:709:ILE:HG21	1:A:852:VAL:HG23	1.80	0.60
1:A:607:ARG:NH1	1:A:610:GLN:OE1	2.34	0.60
1:A:1058:GLU:HG2	1:A:1059:LYS:CD	2.26	0.60
1:A:1056:PRO:C	1:A:1057:SER:HG	1.99	0.59
1:A:596:ARG:NH1	1:A:869:GLU:OE2	2.35	0.59
1:A:620:HIS:O	1:A:736:LYS:NZ	2.35	0.59
1:A:692:GLN:NE2	1:A:707:ASP:OD2	2.36	0.59
1:A:707:ASP:OD1	2:D:74:ARG:CD	2.50	0.59
1:A:983:GLU:O	1:A:984:SER:OG	2.11	0.59
1:A:1159:PHE:O	1:A:1163:GLU:HG2	2.03	0.59
2:C:60:ASN:O	2:C:62:GLN:NE2	2.35	0.59
1:A:282:THR:CG2	1:A:334:VAL:HG12	2.33	0.58
1:A:362:VAL:HG11	1:A:380:VAL:HG11	1.85	0.58
2:D:16:GLU:O	2:D:29:LYS:NZ	2.33	0.58
2:D:20:SER:O	2:D:21:ASP:HB2	2.02	0.58
1:A:377:GLN:O	1:A:381:ASN:ND2	2.35	0.58
2:F:13:ILE:HD13	2:F:34:GLU:HG3	1.85	0.58
1:A:416:GLN:HA	1:A:417:PRO:C	2.28	0.58
1:A:415:LYS:HD3	1:A:421:PHE:CZ	2.39	0.58
2:C:17:VAL:CG2	2:C:18:GLU:H	2.13	0.58
2:C:58:ASP:O	2:C:59:TYR:CB	2.52	0.58
1:A:806:LYS:NZ	1:A:902:ASP:O	2.36	0.58
2:D:56:LEU:HB3	2:D:61:ILE:HB	1.85	0.57
1:A:596:ARG:NH2	1:A:602:GLU:OE1	2.37	0.57
2:C:44:ILE:HD12	2:C:44:ILE:N	2.18	0.57
1:A:256:TYR:HB3	1:A:262:SER:HB3	1.87	0.57
1:A:676:ASN:O	1:A:679:ALA:HB3	2.05	0.57
1:A:377:GLN:HG3	1:A:381:ASN:ND2	2.19	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:248:TYR:HA	1:A:252:LEU:HB2	1.87	0.57
1:A:260:VAL:N	1:A:339:ASP:OD2	2.37	0.56
1:A:241:LYS:HD3	1:A:568:LEU:HD12	1.86	0.56
1:A:474:GLY:HA2	1:A:502:PRO:HD3	1.87	0.56
1:A:593:LYS:CE	1:A:844:ASP:O	2.54	0.56
1:A:1112:ILE:HG12	1:A:1136:LEU:HD13	1.87	0.56
2:C:60:ASN:O	2:C:62:GLN:CD	2.48	0.56
1:A:925:ALA:HB2	1:A:1037:GLN:HE22	1.70	0.56
2:C:31:GLN:HB2	2:C:38:PRO:HG3	1.89	0.55
2:C:45:PHE:O	2:C:46:ALA:HB3	2.05	0.55
1:A:673:HIS:HB3	1:A:675:THR:O	2.07	0.55
1:A:715:ASP:HB3	1:A:717:TRP:NE1	2.22	0.55
1:A:985:ASP:CG	1:A:986:THR:H	2.13	0.55
1:A:1154:THR:O	1:A:1155:ASP:CG	2.48	0.55
1:A:849:SER:HB3	1:A:865:VAL:HG22	1.89	0.55
2:C:56:LEU:N	2:C:56:LEU:HD12	2.22	0.55
1:A:517:ASP:OD1	1:A:517:ASP:N	2.39	0.54
1:A:613:LEU:HD11	1:A:747:LEU:HD11	1.88	0.54
2:F:1:MET:HB2	2:F:63:LYS:HG2	1.89	0.54
1:A:928:LYS:NZ	1:A:1034:THR:CG2	2.70	0.54
1:A:311:ASP:OD1	1:A:311:ASP:N	2.36	0.54
1:A:1007:PHE:CZ	1:A:1011:ARG:HD3	2.43	0.54
2:C:20:SER:H	2:C:56:LEU:HD22	1.73	0.54
1:A:935:GLN:OE1	1:A:1026:ASN:ND2	2.41	0.53
1:A:706:HIS:HD2	2:D:75:GLY:HA3	1.70	0.53
1:A:655:THR:O	1:A:659:ASN:ND2	2.31	0.53
2:C:16:GLU:H	2:C:16:GLU:CD	2.16	0.53
2:C:56:LEU:HB3	2:C:61:ILE:HD11	1.89	0.53
1:A:859:THR:OG1	1:A:860:GLU:N	2.41	0.53
1:A:949:GLU:OE2	2:F:6:LYS:NZ	2.37	0.53
1:A:1058:GLU:O	1:A:1059:LYS:HG2	2.09	0.53
1:A:372:ASP:OD1	1:A:372:ASP:N	2.41	0.53
2:C:56:LEU:CD1	2:C:56:LEU:H	2.22	0.53
2:C:56:LEU:CD1	2:C:56:LEU:N	2.72	0.52
1:A:955:TYR:HB3	2:F:42:ARG:NH1	2.24	0.52
1:A:600:VAL:HG23	1:A:601:ASP:N	2.22	0.52
1:A:859:THR:HG23	1:A:861:SER:N	2.24	0.52
1:A:789:THR:HG22	1:A:801:ALA:HB1	1.91	0.52
2:C:44:ILE:N	2:C:44:ILE:CD1	2.73	0.52
1:A:753:VAL:HA	1:A:850:LYS:HE3	1.92	0.52
1:A:1059:LYS:N	1:A:1061:MET:O	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1065:ARG:O	1:A:1067:GLU:N	2.40	0.52
1:A:854:ARG:O	1:A:855:HIS:CB	2.53	0.52
2:F:23:ILE:HB	2:F:52:ASP:HA	1.91	0.52
1:A:985:ASP:CB	1:A:989:LEU:HD22	2.40	0.52
2:C:17:VAL:HG22	2:C:18:GLU:O	2.10	0.52
1:A:674:SER:OG	1:A:675:THR:N	2.43	0.51
1:A:691:ASP:OD1	2:D:74:ARG:NH2	2.43	0.51
1:A:985:ASP:CA	1:A:989:LEU:HD13	2.39	0.51
1:A:460:ASP:HB3	1:A:463:ASN:HB2	1.92	0.51
1:A:1059:LYS:N	1:A:1059:LYS:CD	2.73	0.50
1:A:283:LEU:HD22	1:A:568:LEU:HD21	1.94	0.50
2:C:21:ASP:C	2:C:22:THR:HG1	2.08	0.50
2:C:58:ASP:O	2:C:59:TYR:CG	2.63	0.50
1:A:375:LYS:CB	1:A:379:ASP:OD2	2.50	0.50
1:A:860:GLU:HG2	2:D:70:VAL:HG11	1.93	0.50
1:A:675:THR:HB	1:A:676:ASN:HD22	1.75	0.50
1:A:668:LYS:HG2	1:A:683:LYS:HD2	1.93	0.50
1:A:985:ASP:HB2	1:A:989:LEU:HD22	1.93	0.49
2:C:71:LEU:H	2:C:71:LEU:HD22	1.77	0.49
2:D:25:ASN:O	2:D:29:LYS:HG3	2.11	0.49
1:A:419:GLU:OE1	1:A:460:ASP:O	2.30	0.49
2:C:43:LEU:CB	2:C:50:LEU:HD12	2.41	0.49
1:A:852:VAL:HG12	1:A:862:GLU:O	2.11	0.49
1:A:988:THR:HG23	1:A:989:LEU:HD12	1.94	0.49
1:A:928:LYS:CE	1:A:1034:THR:CG2	2.90	0.49
1:A:358:PHE:O	1:A:362:VAL:HG23	2.12	0.49
1:A:453:TYR:OH	1:A:556:ASN:ND2	2.36	0.49
1:A:580:ILE:HG12	1:A:896:VAL:HG11	1.94	0.49
1:A:608:ILE:O	1:A:612:ILE:HG12	2.13	0.49
2:C:1:MET:HE2	2:C:18:GLU:CA	2.40	0.49
1:A:313:ARG:NH2	1:A:322:GLU:OE1	2.45	0.49
1:A:985:ASP:HA	1:A:989:LEU:CD1	2.41	0.49
1:A:993:LEU:HD21	1:A:1060:GLU:CB	2.32	0.49
1:A:955:TYR:HB3	2:F:42:ARG:HH11	1.78	0.49
1:A:983:GLU:C	1:A:984:SER:HG	2.11	0.49
2:D:18:GLU:C	2:D:20:SER:H	2.21	0.49
1:A:292:ILE:HG23	1:A:433:ILE:HD12	1.94	0.49
1:A:1008:GLU:O	1:A:1011:ARG:NH2	2.45	0.49
2:C:24:GLU:O	2:C:27:LYS:HB2	2.13	0.48
2:C:20:SER:O	2:C:56:LEU:CG	2.61	0.48
1:A:323:LEU:HA	1:A:326:ILE:HD12	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:557:ARG:NH2	1:A:561:GLU:OE2	2.47	0.48
1:A:902:ASP:HA	1:A:1149:ARG:NH2	2.29	0.48
1:A:624:VAL:HG23	1:A:735:PHE:HZ	1.79	0.48
1:A:785:MET:HG2	1:A:805:ILE:HG12	1.96	0.48
1:A:1030:LEU:O	1:A:1034:THR:HG23	2.14	0.48
1:A:1059:LYS:C	1:A:1061:MET:N	2.72	0.48
1:A:334:VAL:HG22	1:A:337:ARG:HG3	1.96	0.47
1:A:800:GLU:OE1	1:A:803:LYS:NZ	2.39	0.47
2:C:45:PHE:CD2	2:C:67:LEU:CD2	2.94	0.47
1:A:807:LEU:HD21	1:A:1142:VAL:CG2	2.44	0.47
1:A:852:VAL:O	1:A:852:VAL:HG13	2.13	0.47
1:A:459:PHE:HB3	1:A:536:LEU:HB2	1.96	0.47
1:A:785:MET:O	1:A:789:THR:HG23	2.15	0.47
1:A:1049:LEU:HD23	1:A:1049:LEU:HA	1.73	0.47
2:C:6:LYS:O	2:C:68:HIS:HA	2.15	0.47
2:C:58:ASP:O	2:C:59:TYR:HB2	2.13	0.47
1:A:256:TYR:HD2	1:A:262:SER:HA	1.80	0.47
1:A:950:VAL:HG22	1:A:966:LEU:HB3	1.97	0.47
1:A:1084:ARG:NE	1:A:1084:ARG:O	2.48	0.47
1:A:283:LEU:HD23	1:A:286:MET:HE3	1.97	0.47
2:C:22:THR:HB	2:C:24:GLU:HG2	1.96	0.47
1:A:807:LEU:HD21	1:A:1142:VAL:HG21	1.97	0.46
1:A:382:PHE:O	1:A:386:VAL:HG23	2.16	0.46
2:C:45:PHE:CD1	2:C:46:ALA:N	2.84	0.46
2:C:57:SER:O	2:C:58:ASP:C	2.58	0.46
2:F:7:THR:HG22	2:F:69:LEU:HD23	1.98	0.46
1:A:854:ARG:HH12	1:A:856:GLY:HA3	1.81	0.46
2:C:6:LYS:N	2:C:67:LEU:O	2.38	0.46
1:A:1061:MET:HG3	1:A:1062:GLY:N	2.31	0.46
2:C:27:LYS:HG2	2:C:41:GLN:CD	2.41	0.46
1:A:286:MET:HE2	1:A:331:ALA:HA	1.96	0.46
1:A:596:ARG:HD2	1:A:1087:VAL:HG21	1.98	0.46
1:A:831:THR:HG22	2:D:44:ILE:HD12	1.95	0.46
1:A:671:GLU:HG3	1:A:684:ILE:HB	1.98	0.45
1:A:800:GLU:OE2	1:A:1107:ASN:ND2	2.49	0.45
1:A:839:GLU:HB3	1:A:894:THR:HA	1.98	0.45
2:C:43:LEU:O	2:C:44:ILE:HD12	2.14	0.45
1:A:463:ASN:OD1	1:A:465:GLU:HG2	2.16	0.45
1:A:1011:ARG:HG3	1:A:1012:GLN:N	2.32	0.45
1:A:1058:GLU:C	1:A:1059:LYS:CG	2.88	0.45
2:C:56:LEU:O	2:C:61:ILE:CG2	2.64	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:41:GLN:HB2	2:D:69:LEU:CD1	2.41	0.45
1:A:397:THR:OG1	1:A:400:HIS:ND1	2.25	0.45
1:A:461:SER:HB3	1:A:535:HIS:CE1	2.52	0.45
1:A:352:GLU:O	1:A:355:ARG:HG3	2.16	0.45
1:A:419:GLU:OE1	1:A:460:ASP:N	2.44	0.45
1:A:1040:LEU:HA	1:A:1040:LEU:HD23	1.76	0.45
2:F:26:VAL:O	2:F:30:ILE:HG13	2.17	0.45
1:A:725:ILE:HG21	1:A:738:VAL:HA	1.99	0.45
2:C:62:GLN:N	2:C:62:GLN:CD	2.73	0.45
1:A:358:PHE:CE2	1:A:380:VAL:HG13	2.52	0.45
1:A:822:THR:CB	1:A:827:LEU:HD23	2.46	0.45
1:A:1059:LYS:C	1:A:1060:GLU:HG3	2.41	0.45
1:A:1122:ASN:OD1	1:A:1122:ASN:N	2.49	0.45
2:C:22:THR:OG1	2:C:25:ASN:HB3	2.17	0.45
1:A:1042:LYS:O	1:A:1043:GLN:HB2	2.16	0.45
2:C:60:ASN:O	2:C:62:GLN:HG3	2.17	0.45
1:A:251:TYR:O	1:A:274:ARG:NH2	2.49	0.44
1:A:540:MET:HG2	1:A:543:MET:HE3	1.99	0.44
1:A:293:VAL:HG21	1:A:323:LEU:HD13	1.99	0.44
1:A:980:GLY:HA3	1:A:989:LEU:CD1	2.47	0.44
2:D:22:THR:OG1	2:D:25:ASN:HB2	2.16	0.44
1:A:612:ILE:HD12	1:A:624:VAL:HG11	1.99	0.44
1:A:807:LEU:CD2	1:A:1142:VAL:HG21	2.48	0.44
1:A:481:ASP:N	1:A:481:ASP:OD1	2.49	0.44
1:A:645:MET:SD	1:A:746:PRO:HG3	2.58	0.44
1:A:652:LEU:HD11	1:A:746:PRO:HB2	2.00	0.44
1:A:807:LEU:CD2	1:A:1142:VAL:CG2	2.96	0.44
1:A:392:HIS:CG	1:A:392:HIS:O	2.70	0.44
1:A:760:PRO:HA	1:A:761:PRO:HD2	1.83	0.44
1:A:1082:LEU:HD23	1:A:1082:LEU:HA	1.90	0.44
2:C:16:GLU:CD	2:C:16:GLU:N	2.76	0.44
1:A:618:CYS:HB2	1:A:625:LEU:O	2.18	0.44
2:C:3:ILE:HD13	2:C:17:VAL:HG12	2.00	0.44
2:C:24:GLU:HG3	2:C:25:ASN:H	1.80	0.44
1:A:275:GLN:HG3	1:A:276:HIS:ND1	2.33	0.44
1:A:453:TYR:CE1	1:A:556:ASN:HB3	2.53	0.44
2:F:56:LEU:HD22	2:F:61:ILE:HD12	1.99	0.44
1:A:489:PRO:HD3	1:A:495:GLU:HB2	2.00	0.43
1:A:569:LYS:O	1:A:573:THR:HG23	2.18	0.43
1:A:645:MET:HB2	1:A:866:TYR:CZ	2.53	0.43
1:A:488:ALA:HA	1:A:489:PRO:HD3	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:661:ASP:CG	1:A:704:LYS:HZ1	2.18	0.43
1:A:1006:ASN:HB3	1:A:1007:PHE:H	1.41	0.43
1:A:675:THR:HB	1:A:676:ASN:ND2	2.34	0.43
1:A:1071:GLN:HE21	1:A:1071:GLN:HB2	1.57	0.43
1:A:234:LEU:HB2	1:A:290:GLU:HG3	2.00	0.43
1:A:272:VAL:HG11	1:A:561:GLU:HB2	1.99	0.43
1:A:377:GLN:HG3	1:A:381:ASN:HD21	1.83	0.43
1:A:672:PHE:CD1	1:A:681:PRO:HB3	2.54	0.43
1:A:914:GLU:HB2	1:A:915:PRO:HD3	2.01	0.43
1:A:434:GLY:O	1:A:438:THR:OG1	2.33	0.43
1:A:461:SER:HB2	1:A:534:LYS:O	2.18	0.43
1:A:664:MET:HG3	1:A:688:TRP:CD1	2.52	0.43
1:A:1098:ARG:HA	1:A:1150:MET:HE1	2.01	0.43
1:A:282:THR:O	1:A:286:MET:HG3	2.18	0.43
1:A:525:ILE:HA	1:A:528:THR:HG22	2.00	0.43
2:C:43:LEU:HB2	2:C:50:LEU:CD1	2.45	0.43
2:D:55:THR:OG1	2:D:58:ASP:OD2	2.23	0.42
1:A:1040:LEU:HD21	1:A:1078:GLU:OE2	2.19	0.42
1:A:284:ARG:HH22	1:A:413:ARG:NH2	2.17	0.42
1:A:852:VAL:HG12	1:A:862:GLU:C	2.44	0.42
1:A:1061:MET:HG3	1:A:1062:GLY:H	1.84	0.42
2:F:48:LYS:HE2	2:F:59:TYR:HE1	1.84	0.42
1:A:298:LYS:HG2	1:A:787:ALA:HA	2.00	0.42
1:A:376:ASP:OD1	1:A:378:GLU:HB3	2.20	0.42
2:D:18:GLU:HB3	2:D:19:PRO:HD2	2.01	0.42
2:D:37:PRO:HB2	2:D:39:ASP:OD1	2.20	0.42
2:F:15:LEU:HD21	2:F:30:ILE:HG12	2.01	0.42
1:A:810:LEU:HD22	1:A:813:MET:HE2	2.01	0.42
1:A:1029:VAL:CG2	1:A:1056:PRO:HD2	2.49	0.42
2:C:21:ASP:HA	2:C:56:LEU:HD13	2.00	0.42
2:C:61:ILE:C	2:C:62:GLN:HG3	2.44	0.42
1:A:283:LEU:HD23	1:A:283:LEU:HA	1.91	0.42
1:A:943:ILE:HD13	1:A:1020:ILE:HG12	2.00	0.42
1:A:1073:GLN:HE21	1:A:1073:GLN:HB3	1.61	0.41
1:A:676:ASN:HB3	1:A:679:ALA:HB2	2.02	0.41
1:A:867:LEU:HD12	1:A:868:PRO:HD2	2.01	0.41
1:A:980:GLY:C	1:A:989:LEU:HD11	2.45	0.41
1:A:392:HIS:CE1	1:A:394:TRP:CZ2	3.09	0.41
1:A:852:VAL:HG11	1:A:862:GLU:HG2	2.01	0.41
1:A:281:HIS:O	1:A:285:THR:HG23	2.20	0.41
1:A:322:GLU:O	1:A:326:ILE:HG13	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:980:GLY:HA3	1:A:989:LEU:HD12	2.02	0.41
1:A:374:PHE:CD2	1:A:380:VAL:HG22	2.56	0.41
1:A:612:ILE:HG21	1:A:743:LEU:HD21	2.01	0.41
1:A:756:SER:HB3	1:A:850:LYS:HD2	2.02	0.41
1:A:769:ASN:HA	1:A:835:ASN:OD1	2.21	0.41
1:A:1026:ASN:HD22	1:A:1061:MET:CE	2.05	0.41
2:C:14:THR:O	2:C:15:LEU:HD23	2.21	0.41
1:A:881:ASP:O	1:A:890:ARG:HB3	2.21	0.41
1:A:984:SER:OG	1:A:985:ASP:N	2.51	0.41
1:A:362:VAL:HG11	1:A:380:VAL:HG21	2.03	0.41
1:A:470:VAL:O	1:A:470:VAL:HG12	2.20	0.41
2:C:23:ILE:HG22	2:C:27:LYS:HD2	2.03	0.41
1:A:785:MET:HE3	1:A:875:PRO:HG3	2.02	0.41
1:A:925:ALA:HB2	1:A:1037:GLN:NE2	2.34	0.41
2:C:43:LEU:O	2:C:49:GLN:HA	2.21	0.41
1:A:321:GLN:CD	1:A:321:GLN:H	2.28	0.41
1:A:1056:PRO:HB2	1:A:1061:MET:HB3	2.03	0.40
1:A:282:THR:CG2	1:A:334:VAL:CG1	2.99	0.40
1:A:440:ALA:HB2	1:A:790:THR:HB	2.04	0.40
2:C:56:LEU:O	2:C:57:SER:C	2.65	0.40
1:A:740:ILE:O	1:A:744:MET:HG2	2.22	0.40
1:A:949:GLU:O	1:A:953:VAL:HG13	2.21	0.40
1:A:988:THR:HG23	1:A:989:LEU:N	2.35	0.40
1:A:825:ILE:H	1:A:825:ILE:HG13	1.66	0.40
1:A:1058:GLU:O	1:A:1059:LYS:CB	2.70	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	955/968 (99%)	891 (93%)	51 (5%)	13 (1%)	9 34

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	C	69/78 (88%)	64 (93%)	3 (4%)	2 (3%)	3	19
2	D	74/78 (95%)	71 (96%)	2 (3%)	1 (1%)	9	34
2	F	76/78 (97%)	75 (99%)	1 (1%)	0	100	100
All	All	1174/1202 (98%)	1101 (94%)	57 (5%)	16 (1%)	9	34

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	489	PRO
1	A	1092	THR
2	C	38	PRO
1	A	482	GLY
1	A	496	GLY
1	A	848	HIS
1	A	882	GLY
1	A	1008	GLU
1	A	644	GLU
1	A	1066	ARG
2	C	22	THR
1	A	632	GLU
1	A	986	THR
2	D	21	ASP
1	A	937	ALA
1	A	1062	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	827/844 (98%)	784 (95%)	43 (5%)	21	51
2	C	65/70 (93%)	40 (62%)	25 (38%)	0	0
2	D	68/70 (97%)	64 (94%)	4 (6%)	18	48
2	F	69/70 (99%)	67 (97%)	2 (3%)	37	66

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	1029/1054 (98%)	955 (93%)	74 (7%)	13	40

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	256	TYR
1	A	290	GLU
1	A	310	LYS
1	A	311	ASP
1	A	355	ARG
1	A	372	ASP
1	A	413	ARG
1	A	419	GLU
1	A	423	GLN
1	A	455	VAL
1	A	522	LEU
1	A	591	LYS
1	A	594	GLU
1	A	617	ASP
1	A	639	LEU
1	A	644	GLU
1	A	673	HIS
1	A	675	THR
1	A	680	VAL
1	A	694	ILE
1	A	713	GLN
1	A	755	THR
1	A	816	ARG
1	A	855	HIS
1	A	878	VAL
1	A	940	LEU
1	A	941	GLU
1	A	974	LEU
1	A	978	LEU
1	A	1020	ILE
1	A	1030	LEU
1	A	1045	ILE
1	A	1052	LEU
1	A	1055	ILE
1	A	1059	LYS
1	A	1073	GLN
1	A	1084	ARG

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Mol	Chain	Res	Type
1	A	1114	GLU
1	A	1117	THR
1	A	1120	LEU
1	A	1132	ASN
1	A	1133	LEU
1	A	1162	VAL
2	C	5	VAL
2	C	11	LYS
2	C	13	ILE
2	C	16	GLU
2	C	18	GLU
2	C	21	ASP
2	C	29	LYS
2	C	31	GLN
2	C	36	ILE
2	C	39	ASP
2	C	41	GLN
2	C	42	ARG
2	C	43	LEU
2	C	44	ILE
2	C	48	LYS
2	C	51	GLU
2	C	54	ARG
2	C	56	LEU
2	C	57	SER
2	C	58	ASP
2	C	60	ASN
2	C	62	GLN
2	C	63	LYS
2	C	68	HIS
2	C	71	LEU
2	D	22	THR
2	D	25	ASN
2	D	33	LYS
2	D	74	ARG
2	F	13	ILE
2	F	73	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	392	HIS

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Mol	Chain	Res	Type
1	A	405	GLN
1	A	620	HIS
1	A	676	ASN
1	A	706	HIS
1	A	713	GLN
1	A	788	ASN
1	A	797	HIS
1	A	818	ASN
1	A	1016	GLN
1	A	1026	ASN
1	A	1037	GLN
2	C	49	GLN
2	F	31	GLN
2	F	60	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	959/968 (99%)	0.15	25 (2%) 57 36	13, 47, 78, 99	0
2	C	71/78 (91%)	1.43	18 (25%) 1 1	26, 118, 141, 152	0
2	D	76/78 (97%)	0.69	3 (3%) 43 24	54, 83, 103, 117	0
2	F	78/78 (100%)	0.32	3 (3%) 44 25	37, 63, 92, 104	0
All	All	1184/1202 (98%)	0.27	49 (4%) 41 23	13, 51, 103, 152	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1060	GLU	4.9
1	A	1092	THR	4.7
1	A	690	ASN	3.8
2	D	53	GLY	3.6
2	C	55	THR	3.4
1	A	1058	GLU	3.2
1	A	1059	LYS	3.2
2	C	32	ASP	3.1
2	C	19	PRO	3.0
2	C	21	ASP	3.0
1	A	673	HIS	2.9
2	D	66	THR	2.9
1	A	501	PHE	2.8
2	C	49	GLN	2.8
2	C	71	LEU	2.8
1	A	1057	SER	2.7
1	A	392	HIS	2.7
1	A	332	PHE	2.7
2	C	68	HIS	2.6
1	A	715	ASP	2.6
2	C	70	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
2	C	30	ILE	2.5
2	F	76	GLY	2.4
2	C	36	ILE	2.4
2	C	50	LEU	2.4
1	A	547	ASP	2.4
1	A	1031	ASP	2.3
2	D	49	GLN	2.3
2	C	56	LEU	2.3
2	F	24	GLU	2.3
1	A	380	VAL	2.2
1	A	885	GLN	2.2
1	A	1178	ASP	2.2
1	A	983	GLU	2.2
2	C	23	ILE	2.2
1	A	1061	MET	2.2
2	F	74	ARG	2.2
2	C	5	VAL	2.2
1	A	503	GLN	2.2
2	C	69	LEU	2.2
1	A	693	ARG	2.2
2	C	61	ILE	2.1
1	A	1007	PHE	2.1
2	C	26	VAL	2.1
2	C	3	ILE	2.1
1	A	1025	GLY	2.0
1	A	1166	GLU	2.0
1	A	1048	ALA	2.0
1	A	1056	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.