



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:55 AM UTC

PDB ID : 3WTW / pdb_00003wtw
Title : Crystal structure of the complex comprised of ETS1(K167A), RUNX1, CBF-BETA, and the tcralpha gene enhancer DNA
Authors : Shiina, M.; Hamada, K.; Ogata, K.
Deposited on : 2014-04-21
Resolution : 2.90 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

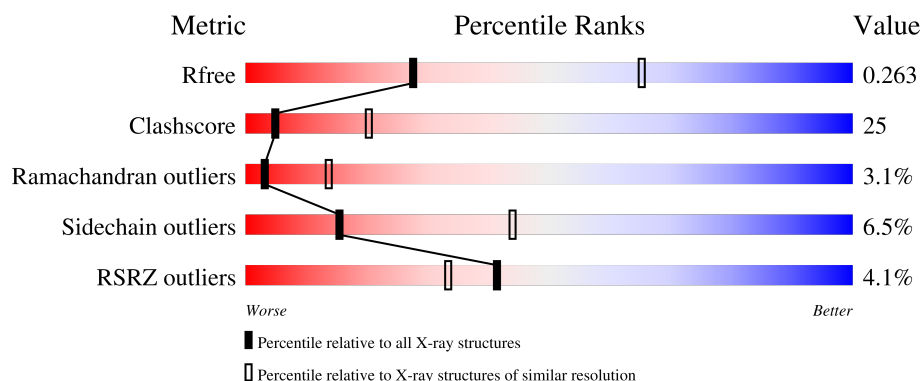
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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.



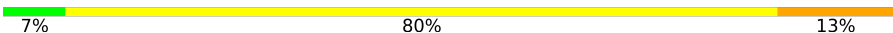
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	204	
1	F	204	
2	B	142	
2	G	142	
3	C	166	

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Mol	Chain	Length	Quality of chain
3	H	166	
4	D	15	
4	I	15	
5	E	15	
5	J	15	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7014 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 Runt-related transcription factor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	118	Total	C	N	O	S	0	0	0
			910	571	168	167	4			
1	F	119	Total	C	N	O	S	0	0	0
			921	577	172	168	4			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	94	LYS	LEU	engineered mutation	UNP Q03347
A	167	ALA	LYS	engineered mutation	UNP Q03347
F	94	LYS	LEU	engineered mutation	UNP Q03347
F	167	ALA	LYS	engineered mutation	UNP Q03347

- Molecule 2 is a protein called Core-binding factor subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	130	Total	C	N	O	S	0	0	0
			1071	671	195	199	6			
2	G	129	Total	C	N	O	S	0	0	0
			1062	666	193	197	6			

- Molecule 3 is a protein called Protein C-ets-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	118	Total	C	N	O	S	0	0	0
			967	627	165	171	4			
3	H	101	Total	C	N	O	S	0	0	0
			853	553	147	149	4			

- Molecule 4 is a DNA chain called DNA (5'-D(*GP*AP*AP*GP*CP*CP*AP*CP*AP*TP*CP*CP*TP*CP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	15	Total 299	C 144	N 54	O 87	P 14	0	0	0
4	I	15	Total 299	C 144	N 54	O 87	P 14	0	0	0

- Molecule 5 is a DNA chain called DNA (5'-D(*AP*GP*AP*GP*GP*AP*TP*GP*TP*GP*GP*CP*TP*TP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	15	Total 310	C 148	N 59	O 89	P 14	0	0	0
5	J	15	Total 310	C 148	N 59	O 89	P 14	0	0	0

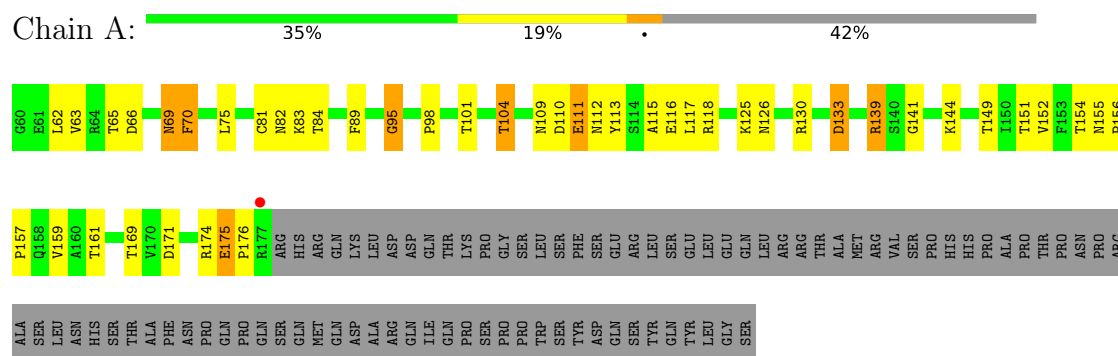
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total 1	O 1	0	0
6	B	1	Total 1	O 1	0	0
6	C	3	Total 3	O 3	0	0
6	F	5	Total 5	O 5	0	0
6	G	1	Total 1	O 1	0	0
6	E	1	Total 1	O 1	0	0

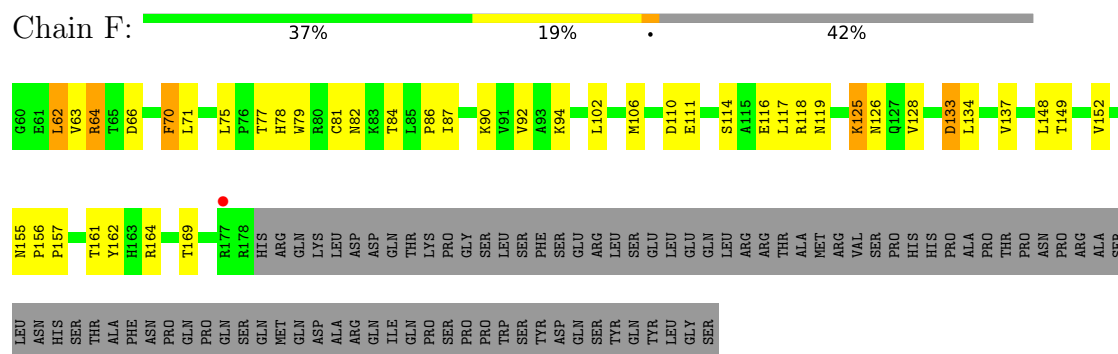
3 Residue-property plots

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.

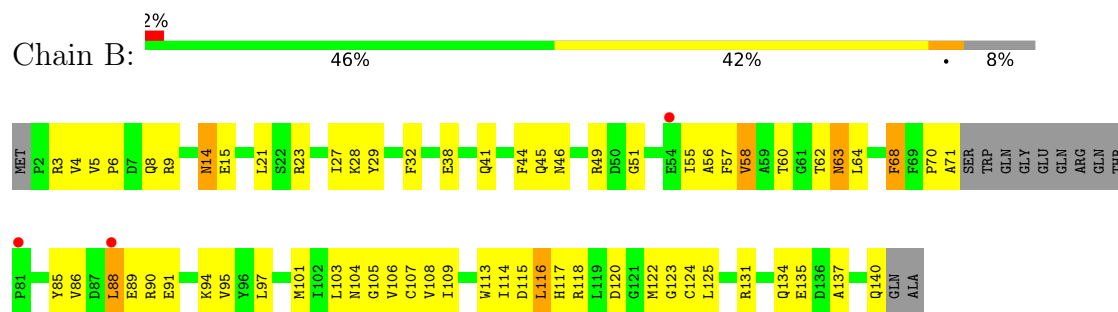
• Molecule 1: Runt-related transcription factor 1



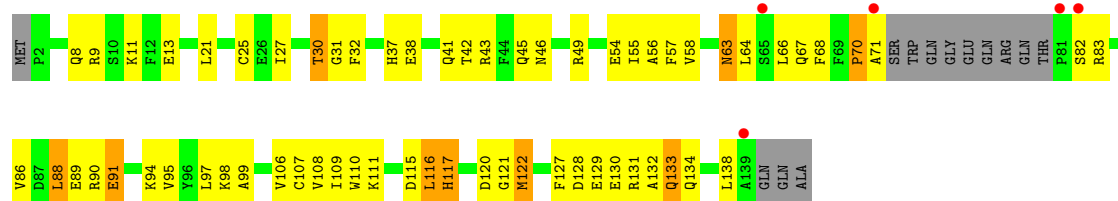
• Molecule 1: Runt-related transcription factor 1



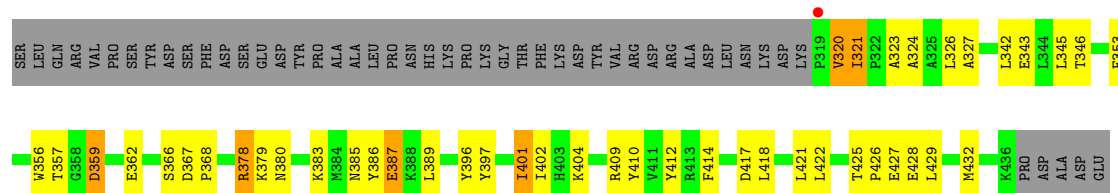
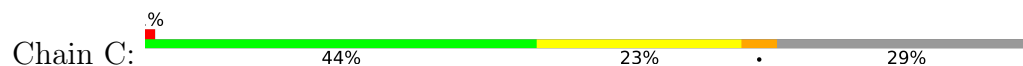
• Molecule 2: Core-binding factor subunit beta



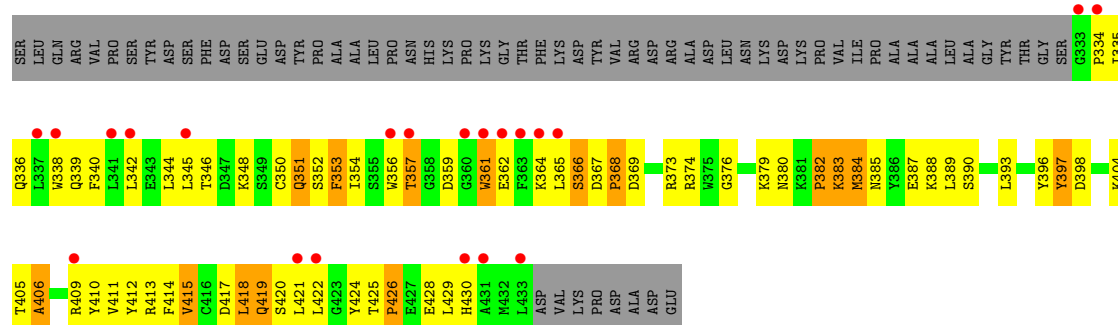
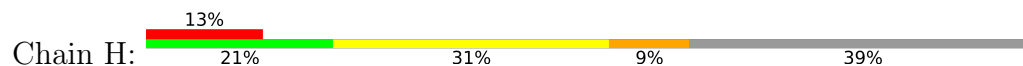
• Molecule 2: Core-binding factor subunit beta



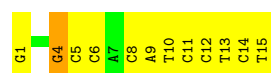
• Molecule 3: Protein C-ets-1



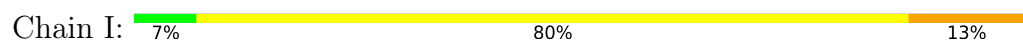
• Molecule 3: Protein C-ets-1



• Molecule 4: DNA (5'-D(*GP*AP*AP*GP*CP*CP*AP*CP*AP*TP*CP*CP*TP*CP*T)-3')



• Molecule 4: DNA (5'-D(*GP*AP*AP*GP*CP*CP*AP*CP*AP*TP*CP*CP*TP*CP*T)-3')



- Molecule 5: DNA (5'-D(*AP*GP*AP*GP*GP*AP*TP*GP*TP*GP*GP*CP*TP*TP*C)-3')

Chain E:  87% 13%



- Molecule 5: DNA (5'-D(*AP*GP*AP*GP*GP*AP*TP*GP*TP*GP*GP*CP*TP*TP*C)-3')

Chain J:  93% 7%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.50Å 100.61Å 194.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.36 – 2.90 42.36 – 2.90	Depositor EDS
% Data completeness (in resolution range)	97.9 (42.36-2.90) 97.9 (42.36-2.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.60 (at 2.90Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.219 , 0.264 0.218 , 0.263	Depositor DCC
R_{free} test set	3414 reflections (9.99%)	wwPDB-VP
Wilson B-factor (Å ²)	77.0	Xtriage
Anisotropy	0.301	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 52.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7014	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% 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	0.68	0/929	1.02	3/1264 (0.2%)
1	F	0.73	0/940	1.09	2/1278 (0.2%)
2	B	0.49	0/1093	0.80	1/1466 (0.1%)
2	G	0.59	1/1084 (0.1%)	0.86	0/1454
3	C	0.66	0/994	1.03	1/1342 (0.1%)
3	H	0.47	0/877	0.82	3/1181 (0.3%)
4	D	0.45	0/334	0.98	1/512 (0.2%)
4	I	0.47	0/334	0.99	1/512 (0.2%)
5	E	0.47	0/348	0.80	0/537
5	J	0.44	0/348	0.85	0/537
All	All	0.59	1/7281 (0.0%)	0.93	12/10083 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
4	D	0	3
4	I	0	2
5	E	0	1
All	All	0	6

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	122	MET	SD-CE	5.82	1.94	1.79

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	4	DG	N9-C1'-C2'	-6.85	103.23	113.50
1	F	70	PHE	N-CA-C	6.41	119.16	109.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	111	GLU	N-CA-C	-6.40	107.32	114.62
1	F	133	ASP	N-CA-C	6.40	120.22	111.39
4	I	4	DG	N9-C1'-C2'	-6.25	104.12	113.50
1	A	133	ASP	N-CA-C	5.79	119.23	111.24
3	H	361	TRP	N-CA-C	-5.45	107.89	114.75
3	C	401	ILE	N-CA-C	-5.40	107.35	111.62
3	H	357	THR	N-CA-C	-5.39	107.36	114.31
3	H	415	VAL	N-CA-C	-5.22	108.20	113.47
2	B	68	PHE	N-CA-C	5.06	117.21	109.62
1	A	70	PHE	N-CA-C	5.05	117.06	109.23

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	D	4	DG	Sidechain
4	D	5	DC	Sidechain
4	D	6	DC	Sidechain
5	E	13	DT	Sidechain
4	I	4	DG	Sidechain
4	I	5	DC	Sidechain

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	910	0	914	38	0
1	F	921	0	927	38	0
2	B	1071	0	1034	71	0
2	G	1062	0	1026	56	0
3	C	967	0	966	36	0
3	H	853	0	847	82	0
4	D	299	0	170	14	0
4	I	299	0	170	17	0
5	E	310	0	171	1	0
5	J	310	0	171	1	0
6	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	1	0	0	1	0
6	C	3	0	0	0	0
6	E	1	0	0	0	0
6	F	5	0	0	1	0
6	G	1	0	0	0	0
All	All	7014	0	6396	339	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 25.

All (339) 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)
3:C:385:ASN:HD21	3:C:387:GLU:HB2	1.22	1.01
1:A:82:ASN:HD21	1:A:118:ARG:HH11	1.04	0.89
2:B:8:GLN:HE21	2:B:107:CYS:H	1.16	0.88
1:A:161:THR:H	2:B:104:ASN:HD21	1.20	0.87
1:A:109:ASN:HD21	1:A:111:GLU:HB2	1.39	0.86
1:F:82:ASN:HD21	1:F:118:ARG:HH11	1.18	0.86
4:I:11:DC:H1'	4:I:12:DC:H5''	1.58	0.84
3:H:345:LEU:HB3	3:H:356:TRP:HE1	1.39	0.84
1:A:111:GLU:HG3	1:A:144:LYS:HZ3	1.43	0.83
2:B:41:GLN:HE21	2:B:117:HIS:HA	1.45	0.81
1:F:87:ILE:HD12	1:F:87:ILE:O	1.81	0.81
1:A:111:GLU:HG3	1:A:144:LYS:NZ	1.95	0.81
2:B:60:THR:HG23	2:B:62:THR:H	1.48	0.79
1:A:82:ASN:HD21	1:A:118:ARG:NH1	1.80	0.78
3:C:385:ASN:ND2	3:C:387:GLU:HB2	1.99	0.78
3:C:425:THR:HB	3:C:426:PRO:HD2	1.67	0.77
3:H:409:ARG:O	3:H:411:VAL:HG23	1.85	0.77
3:H:424:TYR:HB3	3:H:429:LEU:HD13	1.66	0.76
1:F:63:VAL:HG22	1:F:64:ARG:N	2.01	0.75
2:B:108:VAL:HG12	2:B:109:ILE:H	1.53	0.74
2:B:57:PHE:HB2	2:B:60:THR:HG22	1.69	0.74
3:C:327:ALA:HA	3:C:421:LEU:HD23	1.69	0.73
3:H:365:LEU:HD21	3:H:368:PRO:HA	1.69	0.73
1:A:69:ASN:ND2	1:A:95:GLY:H	1.88	0.72
3:C:409:ARG:HG2	3:C:410:TYR:CD2	2.26	0.71
4:D:10:DT:H2''	4:D:11:DC:H5'	1.72	0.71
2:B:71:ALA:HB3	2:B:131:ARG:HD3	1.72	0.70
3:H:345:LEU:HB3	3:H:356:TRP:NE1	2.06	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:326:LEU:HD13	3:C:422:LEU:HA	1.74	0.70
3:H:335:ILE:HD12	3:H:339:GLN:NE2	2.07	0.70
3:H:340:PHE:HZ	3:H:374:ARG:HB3	1.57	0.70
3:H:359:ASP:HB2	3:H:362:GLU:HB3	1.74	0.69
2:B:45:GLN:HE21	2:B:49:ARG:HH12	1.41	0.69
1:F:78:HIS:HB2	6:F:301:HOH:O	1.91	0.69
4:I:14:DC:H1'	4:I:15:DT:H5'	1.74	0.69
2:B:108:VAL:HG12	2:B:109:ILE:N	2.05	0.69
3:C:357:THR:CG2	3:C:362:GLU:HG2	2.23	0.69
1:F:63:VAL:HG22	1:F:64:ARG:H	1.59	0.68
3:H:384:MET:HE1	3:H:388:LYS:HB3	1.76	0.67
1:F:62:LEU:HD21	1:F:92:VAL:HG21	1.74	0.67
1:F:66:ASP:OD2	1:F:161:THR:HB	1.94	0.67
2:G:106:VAL:HG12	2:G:108:VAL:HG23	1.77	0.66
2:B:88:LEU:N	2:B:88:LEU:HD12	2.10	0.66
3:H:365:LEU:HD23	3:H:365:LEU:O	1.96	0.66
3:C:320:VAL:HG11	3:C:432:MET:HE2	1.78	0.65
3:C:359:ASP:OD2	3:C:359:ASP:C	2.38	0.65
3:H:405:THR:HA	3:H:413:ARG:NH2	2.12	0.65
1:F:87:ILE:HD12	1:F:87:ILE:C	2.21	0.65
4:I:1:DG:H8	4:I:1:DG:HO5'	1.45	0.65
4:D:11:DC:H2''	4:D:12:DC:H5'	1.79	0.65
3:C:404:LYS:HD3	3:C:412:TYR:CE2	2.33	0.64
2:B:56:ALA:HB2	2:B:63:ASN:HA	1.80	0.64
1:F:71:LEU:HD11	1:F:94:LYS:HE3	1.79	0.64
2:B:106:VAL:O	2:B:108:VAL:HG23	1.97	0.64
3:H:368:PRO:HG2	3:H:369:ASP:H	1.62	0.64
3:H:348:LYS:HA	3:H:351:GLN:HG3	1.79	0.64
2:B:29:TYR:CE1	2:B:44:PHE:HB2	2.32	0.63
1:A:69:ASN:HD21	1:A:95:GLY:H	1.45	0.63
3:H:379:LYS:O	3:H:380:ASN:HB3	1.98	0.63
2:B:134:GLN:O	2:B:137:ALA:HB3	1.99	0.62
3:H:383:LYS:H	3:H:383:LYS:HD3	1.64	0.62
3:H:350:CYS:C	3:H:352:SER:H	2.08	0.62
2:B:91:GLU:HB3	2:B:94:LYS:HB2	1.79	0.62
2:B:108:VAL:HG11	2:B:125:LEU:HB3	1.81	0.62
2:G:131:ARG:NH1	2:G:131:ARG:HB3	2.13	0.62
3:H:348:LYS:HA	3:H:351:GLN:CD	2.25	0.62
3:H:368:PRO:HG2	3:H:369:ASP:OD1	1.98	0.62
3:H:404:LYS:NZ	3:H:409:ARG:HA	2.14	0.62
2:G:63:ASN:N	2:G:63:ASN:HD22	1.97	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:84:THR:HG23	1:F:133:ASP:OD1	1.99	0.61
3:H:354:ILE:HG13	3:H:365:LEU:HA	1.83	0.61
3:C:417:ASP:O	3:C:421:LEU:HD13	2.00	0.61
1:A:156:PRO:HB2	1:A:157:PRO:HD2	1.82	0.61
3:H:336:GLN:H	3:H:339:GLN:HE21	1.48	0.61
4:I:10:DT:H1'	4:I:11:DC:H5'	1.83	0.60
2:B:63:ASN:N	2:B:63:ASN:HD22	1.98	0.60
4:D:1:DG:O4'	4:I:15:DT:H2'	2.02	0.60
1:A:82:ASN:ND2	1:A:118:ARG:HH11	1.88	0.60
3:H:348:LYS:HA	3:H:351:GLN:CG	2.31	0.60
2:B:115:ASP:OD2	2:B:118:ARG:HD3	2.02	0.60
3:C:357:THR:HG21	3:C:362:GLU:HG2	1.84	0.60
3:H:404:LYS:HG2	3:H:412:TYR:CE2	2.37	0.60
2:G:70:PRO:O	2:G:71:ALA:HB2	2.01	0.60
2:G:82:SER:OG	2:G:83:ARG:N	2.35	0.60
3:H:419:GLN:NE2	3:H:425:THR:HG22	2.17	0.60
3:H:362:GLU:HA	3:H:413:ARG:HB3	1.84	0.60
1:A:104:THR:HG22	1:A:151:THR:HB	1.83	0.59
3:H:389:LEU:O	3:H:389:LEU:HD23	2.03	0.59
2:G:55:ILE:HG22	2:G:56:ALA:N	2.18	0.59
3:H:389:LEU:HD23	3:H:389:LEU:C	2.28	0.59
3:C:323:ALA:HB2	3:C:346:THR:HG21	1.84	0.58
3:H:420:SER:O	3:H:421:LEU:HD12	2.03	0.58
3:H:418:LEU:HD22	3:H:422:LEU:HD22	1.86	0.58
3:C:320:VAL:HG12	3:C:321:ILE:H	1.69	0.58
1:F:161:THR:HG22	1:F:162:TYR:N	2.19	0.58
3:H:351:GLN:HA	3:H:354:ILE:O	2.04	0.58
2:G:134:GLN:O	2:G:138:LEU:HG	2.04	0.57
3:H:357:THR:HA	3:H:364:LYS:NZ	2.19	0.57
2:B:45:GLN:NE2	2:B:49:ARG:HH12	2.02	0.57
3:H:419:GLN:CD	3:H:425:THR:HG22	2.28	0.57
4:I:12:DC:C2'	4:I:13:DT:H72	2.34	0.57
3:C:378:ARG:HG3	3:C:378:ARG:O	2.04	0.57
1:A:159:VAL:HG13	2:B:103:LEU:HA	1.86	0.57
1:F:82:ASN:HD21	1:F:118:ARG:NH1	1.95	0.57
2:B:95:VAL:HG23	2:B:116:LEU:HG	1.88	0.56
2:B:115:ASP:C	2:B:117:HIS:H	2.14	0.55
3:H:348:LYS:HG2	3:H:351:GLN:NE2	2.21	0.55
2:G:94:LYS:HA	2:G:116:LEU:HG	1.87	0.55
3:H:404:LYS:HZ3	3:H:409:ARG:HA	1.72	0.55
4:I:9:DA:H1'	4:I:10:DT:H5''	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:335:ILE:HD11	3:H:339:GLN:HB3	1.88	0.55
1:A:65:THR:HG23	1:A:70:PHE:O	2.07	0.55
4:D:14:DC:H2''	4:D:15:DT:H5'	1.88	0.55
3:C:397:TYR:OH	3:C:404:LYS:HB2	2.07	0.55
2:G:95:VAL:HG13	2:G:116:LEU:HD11	1.89	0.54
2:G:37:HIS:O	2:G:41:GLN:HG3	2.08	0.54
1:F:114:SER:HB2	2:G:30:THR:HG21	1.90	0.54
3:H:336:GLN:H	3:H:339:GLN:NE2	2.06	0.54
2:B:91:GLU:HG2	2:B:94:LYS:HD2	1.89	0.53
1:F:63:VAL:CG2	1:F:64:ARG:N	2.70	0.53
1:F:82:ASN:CG	1:F:82:ASN:O	2.50	0.53
3:H:353:PHE:C	3:H:366:SER:HB3	2.33	0.53
2:G:55:ILE:HG22	2:G:56:ALA:H	1.74	0.53
3:H:419:GLN:HG2	3:H:425:THR:HA	1.91	0.53
2:B:91:GLU:HG2	2:B:94:LYS:CG	2.38	0.53
1:A:70:PHE:CE2	1:A:152:VAL:HG21	2.44	0.53
2:B:60:THR:HG23	2:B:62:THR:N	2.19	0.53
1:A:171:ASP:HB3	1:A:174:ARG:HD2	1.90	0.53
2:B:56:ALA:CB	2:B:63:ASN:HA	2.38	0.52
2:B:94:LYS:HD3	2:B:113:TRP:CE3	2.43	0.52
3:C:320:VAL:HG12	3:C:321:ILE:N	2.24	0.52
2:G:131:ARG:HB3	2:G:131:ARG:CZ	2.40	0.52
3:H:368:PRO:HB3	3:H:410:TYR:CE1	2.44	0.52
1:A:112:ASN:C	1:A:112:ASN:HD22	2.18	0.52
3:H:345:LEU:HD13	3:H:356:TRP:CE2	2.45	0.51
2:B:41:GLN:NE2	2:B:117:HIS:HA	2.20	0.51
1:A:75:LEU:HD21	1:A:89:PHE:CD2	2.45	0.51
4:D:8:DC:H2''	4:D:9:DA:C8	2.45	0.51
2:B:38:GLU:O	2:B:41:GLN:HB2	2.11	0.51
1:F:161:THR:HG22	1:F:162:TYR:H	1.75	0.50
3:H:398:ASP:OD1	3:H:398:ASP:N	2.44	0.50
1:A:149:THR:HG21	2:B:63:ASN:O	2.12	0.50
2:G:21:LEU:HD22	2:G:57:PHE:CD1	2.46	0.50
3:H:361:TRP:CZ2	3:H:426:PRO:HG3	2.47	0.50
1:A:98:PRO:O	1:A:101:THR:OG1	2.27	0.50
1:A:109:ASN:ND2	1:A:111:GLU:HB2	2.17	0.50
3:H:405:THR:HA	3:H:413:ARG:CZ	2.42	0.50
3:H:418:LEU:O	3:H:422:LEU:HB3	2.12	0.50
1:F:149:THR:HG21	2:G:63:ASN:O	2.12	0.49
1:A:175:GLU:O	1:A:176:PRO:C	2.55	0.49
2:B:137:ALA:O	2:B:140:GLN:HG3	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:426:PRO:HG2	3:C:427:GLU:H	1.78	0.49
3:H:342:LEU:HD23	3:H:342:LEU:O	2.12	0.49
2:G:31:GLY:O	2:G:32:PHE:C	2.55	0.49
4:D:9:DA:H1'	4:D:10:DT:C5'	2.43	0.49
2:B:68:PHE:CZ	2:B:97:LEU:HD13	2.47	0.49
2:G:49:ARG:HG2	2:G:49:ARG:HH11	1.78	0.49
1:A:69:ASN:HD21	1:A:95:GLY:N	2.10	0.49
2:G:38:GLU:O	2:G:41:GLN:HB2	2.12	0.49
3:H:406:ALA:H	3:H:413:ARG:NH2	2.11	0.49
3:C:425:THR:CB	3:C:426:PRO:HD2	2.41	0.49
3:H:383:LYS:H	3:H:383:LYS:CD	2.24	0.49
1:F:87:ILE:C	1:F:87:ILE:CD1	2.86	0.48
3:H:406:ALA:H	3:H:413:ARG:HH22	1.61	0.48
1:A:159:VAL:HG21	2:B:64:LEU:HD21	1.95	0.48
2:B:8:GLN:HB3	6:B:201:HOH:O	2.13	0.48
2:B:108:VAL:CG1	2:B:109:ILE:N	2.74	0.48
2:G:98:LYS:HE3	2:G:111:LYS:HE2	1.95	0.48
1:F:63:VAL:CG2	1:F:64:ARG:H	2.23	0.48
2:G:121:GLY:O	2:G:122:MET:HE2	2.14	0.48
3:H:354:ILE:HA	3:H:366:SER:H	1.78	0.48
2:B:108:VAL:CG1	2:B:109:ILE:H	2.22	0.48
3:C:396:TYR:O	3:C:397:TYR:C	2.56	0.48
1:F:155:ASN:HD22	1:F:155:ASN:C	2.21	0.48
4:D:13:DT:H2''	4:D:14:DC:H5'	1.95	0.48
3:C:386:TYR:O	3:C:387:GLU:C	2.57	0.48
2:G:55:ILE:HG12	2:G:66:LEU:HD12	1.96	0.48
3:H:409:ARG:HG3	3:H:410:TYR:HD2	1.79	0.48
2:G:88:LEU:N	2:G:88:LEU:HD12	2.29	0.48
2:G:89:GLU:O	2:G:91:GLU:N	2.47	0.48
3:H:338:TRP:HB3	3:H:396:TYR:CZ	2.48	0.48
4:I:11:DC:C1'	4:I:12:DC:H5''	2.37	0.48
2:G:25:CYS:O	2:G:122:MET:HA	2.14	0.47
2:G:63:ASN:N	2:G:63:ASN:ND2	2.63	0.47
3:H:390:SER:O	3:H:393:LEU:HB2	2.13	0.47
4:I:8:DC:H2''	4:I:9:DA:C8	2.49	0.47
3:H:376:GLY:HA2	3:H:384:MET:HG3	1.97	0.47
3:H:409:ARG:HG3	3:H:410:TYR:CD2	2.50	0.47
2:B:14:ASN:HD22	2:B:14:ASN:N	2.12	0.47
2:B:91:GLU:HG2	2:B:94:LYS:CD	2.44	0.47
2:G:66:LEU:HD22	2:G:99:ALA:HB3	1.97	0.47
1:A:84:THR:HG23	1:A:133:ASP:OD1	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:57:PHE:O	2:B:58:VAL:C	2.57	0.47
3:H:350:CYS:HB3	3:H:353:PHE:HE1	1.80	0.47
2:B:14:ASN:N	2:B:14:ASN:ND2	2.58	0.47
2:G:49:ARG:HH12	2:G:88:LEU:HD22	1.80	0.47
3:H:338:TRP:CE2	3:H:339:GLN:HG3	2.50	0.47
2:G:109:ILE:HG13	2:G:128:ASP:HB2	1.98	0.46
2:B:9:ARG:HG3	2:B:9:ARG:HH11	1.80	0.46
2:B:45:GLN:OE1	2:B:116:LEU:HD22	2.14	0.46
4:I:13:DT:H1'	4:I:14:DC:H5''	1.97	0.46
1:F:79:TRP:CZ2	1:F:86:PRO:HD3	2.51	0.46
3:H:357:THR:HB	3:H:359:ASP:OD2	2.15	0.46
3:H:419:GLN:HG3	3:H:426:PRO:HD3	1.96	0.46
4:I:12:DC:H2'	4:I:13:DT:H72	1.97	0.45
1:A:109:ASN:C	1:A:111:GLU:H	2.25	0.45
3:H:396:TYR:O	3:H:397:TYR:C	2.59	0.45
3:H:414:PHE:N	3:H:414:PHE:CD2	2.84	0.45
3:H:417:ASP:C	3:H:419:GLN:H	2.23	0.45
2:G:95:VAL:HG13	2:G:116:LEU:HD21	1.98	0.45
3:H:366:SER:O	3:H:367:ASP:C	2.58	0.45
2:B:8:GLN:NE2	2:B:107:CYS:H	1.97	0.45
2:B:23:ARG:O	2:B:124:CYS:HB3	2.16	0.45
1:F:70:PHE:CE2	1:F:152:VAL:HG21	2.52	0.45
1:F:148:LEU:HB2	1:F:162:TYR:HB3	1.98	0.45
2:G:9:ARG:O	2:G:13:GLU:HG3	2.16	0.45
2:G:115:ASP:C	2:G:117:HIS:H	2.23	0.45
4:I:9:DA:H1'	4:I:10:DT:C5'	2.46	0.45
2:G:129:GLU:O	2:G:132:ALA:HB3	2.17	0.45
2:G:89:GLU:N	2:G:89:GLU:CD	2.74	0.45
1:A:113:TYR:CZ	2:B:28:LYS:HD2	2.52	0.45
3:H:369:ASP:O	3:H:373:ARG:HG2	2.17	0.45
1:F:110:ASP:CG	1:F:110:ASP:O	2.58	0.45
1:F:102:LEU:N	1:F:102:LEU:HD22	2.32	0.45
2:B:115:ASP:C	2:B:117:HIS:N	2.75	0.44
2:B:123:GLY:O	2:B:124:CYS:HB3	2.17	0.44
2:B:70:PRO:HD3	2:B:85:TYR:CE2	2.53	0.44
3:C:353:PHE:CD2	3:C:353:PHE:N	2.85	0.44
2:G:54:GLU:C	2:G:55:ILE:HD13	2.43	0.44
1:F:75:LEU:HD11	1:F:148:LEU:HD11	2.00	0.44
1:F:81:CYS:HB2	1:F:169:THR:O	2.17	0.44
2:G:11:LYS:HD3	2:G:106:VAL:CG2	2.47	0.44
3:H:382:PRO:HB2	3:H:383:LYS:HD3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:5:VAL:HG12	2:B:105:GLY:O	2.18	0.44
3:C:379:LYS:O	3:C:380:ASN:HB3	2.18	0.44
3:H:342:LEU:HD23	3:H:342:LEU:C	2.43	0.44
4:I:12:DC:H2''	4:I:13:DT:C7	2.48	0.44
1:A:111:GLU:OE2	1:A:141:GLY:HA3	2.18	0.44
2:G:42:THR:OG1	2:G:43:ARG:N	2.51	0.44
2:G:68:PHE:CD1	2:G:99:ALA:HB2	2.53	0.44
1:A:125:LYS:O	1:A:126:ASN:HB2	2.18	0.44
2:B:55:ILE:CD1	2:B:114:ILE:HD11	2.48	0.44
3:H:418:LEU:O	3:H:422:LEU:HD23	2.17	0.44
2:G:63:ASN:HD22	2:G:63:ASN:H	1.64	0.43
2:G:108:VAL:HG12	2:G:109:ILE:N	2.32	0.43
4:D:9:DA:H1'	4:D:10:DT:H5''	2.00	0.43
2:G:70:PRO:O	2:G:71:ALA:CB	2.63	0.43
3:H:417:ASP:OD2	3:H:421:LEU:HD13	2.18	0.43
1:A:155:ASN:HA	1:A:156:PRO:HA	1.89	0.43
2:B:27:ILE:HG22	2:B:57:PHE:CD2	2.53	0.43
2:G:45:GLN:O	2:G:46:ASN:C	2.61	0.43
3:H:335:ILE:HG13	3:H:336:GLN:N	2.33	0.43
2:B:21:LEU:HD21	2:B:60:THR:HG21	2.01	0.43
1:F:71:LEU:CD1	1:F:94:LYS:HE3	2.47	0.43
1:F:157:PRO:CG	2:G:67:GLN:HG3	2.48	0.43
2:G:8:GLN:NE2	2:G:107:CYS:H	2.16	0.43
2:B:115:ASP:O	2:B:117:HIS:N	2.52	0.43
2:G:9:ARG:HA	2:G:127:PHE:CE2	2.54	0.43
1:F:118:ARG:O	1:F:119:ASN:HB2	2.18	0.42
4:I:5:DC:H1'	4:I:6:DC:H5''	2.01	0.42
1:A:81:CYS:C	1:A:83:LYS:H	2.26	0.42
2:G:54:GLU:O	2:G:55:ILE:HD13	2.19	0.42
3:H:340:PHE:O	3:H:344:LEU:HD13	2.19	0.42
2:B:55:ILE:HD13	2:B:114:ILE:HD11	2.01	0.42
3:C:345:LEU:HB3	3:C:356:TRP:CZ2	2.55	0.42
3:H:348:LYS:O	3:H:351:GLN:HB2	2.19	0.42
4:D:11:DC:H1'	4:D:12:DC:C5'	2.49	0.42
3:C:383:LYS:H	3:C:383:LYS:HG2	1.62	0.42
2:B:90:ARG:O	2:B:90:ARG:HG2	2.19	0.42
1:F:86:PRO:O	1:F:87:ILE:HG23	2.20	0.42
4:D:13:DT:H1'	4:D:14:DC:H5''	2.01	0.42
4:I:8:DC:H4'	4:I:8:DC:OP1	2.19	0.42
1:A:115:ALA:O	1:A:117:LEU:HD13	2.19	0.42
2:B:28:LYS:HB2	2:B:28:LYS:HE3	1.88	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:45:GLN:O	2:B:46:ASN:C	2.62	0.42
2:B:89:GLU:C	2:B:91:GLU:H	2.27	0.42
1:A:116:GLU:O	1:A:117:LEU:HD12	2.20	0.42
2:G:27:ILE:HB	2:G:55:ILE:CG2	2.49	0.42
2:B:114:ILE:HG23	2:B:120:ASP:N	2.34	0.42
3:C:418:LEU:HA	3:C:418:LEU:HD23	1.79	0.42
2:G:133:GLN:HE21	2:G:133:GLN:HB2	1.62	0.42
3:H:385:ASN:OD1	3:H:387:GLU:HB2	2.20	0.42
1:A:154:THR:O	1:A:156:PRO:C	2.63	0.41
3:H:335:ILE:HD12	3:H:339:GLN:HE21	1.79	0.41
4:D:9:DA:H2''	4:D:10:DT:H5'	2.01	0.41
4:D:11:DC:C2'	4:D:12:DC:H5'	2.47	0.41
4:I:2:DA:H1'	4:I:3:DA:H5'	2.01	0.41
1:A:66:ASP:O	2:B:104:ASN:HB3	2.20	0.41
3:C:342:LEU:HD13	3:C:414:PHE:CZ	2.54	0.41
1:F:106:MET:HE1	2:G:64:LEU:C	2.46	0.41
3:H:419:GLN:CG	3:H:426:PRO:HD3	2.50	0.41
2:B:6:PRO:C	2:B:8:GLN:H	2.27	0.41
2:B:32:PHE:CD2	2:B:32:PHE:N	2.87	0.41
2:G:68:PHE:CE2	2:G:97:LEU:HD22	2.56	0.41
3:H:345:LEU:HD22	3:H:356:TRP:CD1	2.56	0.41
3:H:383:LYS:HD3	3:H:383:LYS:N	2.33	0.41
4:D:10:DT:H1'	4:D:11:DC:H5''	2.03	0.41
1:A:62:LEU:C	1:A:63:VAL:HG23	2.46	0.41
3:C:320:VAL:CG1	3:C:321:ILE:H	2.31	0.41
3:C:326:LEU:HB3	3:C:421:LEU:O	2.20	0.41
1:F:126:ASN:O	1:F:128:VAL:HG23	2.20	0.41
1:F:157:PRO:HG2	2:G:67:GLN:HG3	2.02	0.41
3:H:420:SER:C	3:H:421:LEU:HD12	2.45	0.41
2:B:3:ARG:HH11	2:B:135:GLU:CD	2.28	0.41
2:B:28:LYS:O	2:B:55:ILE:HG23	2.21	0.41
2:B:91:GLU:HG2	2:B:94:LYS:HG3	2.02	0.41
3:C:379:LYS:O	3:C:380:ASN:CB	2.68	0.41
2:G:89:GLU:C	2:G:91:GLU:N	2.79	0.41
1:A:139:ARG:HD3	1:A:139:ARG:N	2.36	0.41
2:B:55:ILE:CG2	2:B:56:ALA:N	2.84	0.41
2:G:108:VAL:CG1	2:G:109:ILE:N	2.84	0.41
3:H:415:VAL:HG12	3:H:415:VAL:O	2.21	0.41
5:E:15:DC:H2'	5:J:1:DA:O4'	2.21	0.41
3:C:368:PRO:HB3	3:C:386:TYR:CD1	2.55	0.41
3:C:389:LEU:C	3:C:389:LEU:HD23	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:155:ASN:HA	1:F:156:PRO:HA	1.97	0.41
3:H:362:GLU:HG3	3:H:412:TYR:O	2.20	0.41
3:H:419:GLN:C	3:H:421:LEU:H	2.29	0.41
2:B:51:GLY:HA2	2:B:68:PHE:HD2	1.85	0.41
1:F:116:GLU:HG2	1:F:137:VAL:HB	2.03	0.41
3:H:414:PHE:H	3:H:414:PHE:HD2	1.68	0.41
1:F:114:SER:CB	2:G:30:THR:HG21	2.51	0.40
2:G:98:LYS:HA	2:G:110:TRP:O	2.21	0.40
2:G:127:PHE:CE1	2:G:129:GLU:HA	2.56	0.40
4:D:9:DA:H1'	4:D:10:DT:H5'	2.03	0.40
4:I:9:DA:H2''	4:I:10:DT:H5'	2.03	0.40
2:B:21:LEU:HD22	2:B:57:PHE:CD1	2.56	0.40
3:C:324:ALA:HB2	3:C:343:GLU:OE1	2.21	0.40
3:C:353:PHE:HB3	3:C:367:ASP:HB3	2.03	0.40
1:A:159:VAL:HG21	2:B:64:LEU:CD2	2.51	0.40
3:C:353:PHE:O	3:C:366:SER:N	2.54	0.40
2:G:115:ASP:O	2:G:117:HIS:N	2.55	0.40
3:H:357:THR:HA	3:H:364:LYS:HZ3	1.86	0.40
1:A:81:CYS:C	1:A:83:LYS:N	2.80	0.40
1:F:125:LYS:O	1:F:126:ASN:HB2	2.21	0.40
2:B:28:LYS:HE3	2:B:58:VAL:HG21	2.03	0.40
2:B:113:TRP:CZ2	2:B:122:MET:HE2	2.57	0.40
3:C:401:ILE:HG22	3:C:402:ILE:HG13	2.03	0.40
3:H:368:PRO:HG2	3:H:369:ASP:N	2.32	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	116/204 (57%)	108 (93%)	6 (5%)	2 (2%)	7 26

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	117/204 (57%)	109 (93%)	7 (6%)	1 (1%)	14	41
2	B	126/142 (89%)	99 (79%)	24 (19%)	3 (2%)	4	18
2	G	125/142 (88%)	107 (86%)	13 (10%)	5 (4%)	2	9
3	C	116/166 (70%)	109 (94%)	6 (5%)	1 (1%)	14	41
3	H	99/166 (60%)	68 (69%)	21 (21%)	10 (10%)	0	1
All	All	699/1024 (68%)	600 (86%)	77 (11%)	22 (3%)	3	14

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	320	VAL
3	H	366	SER
3	H	382	PRO
2	B	15	GLU
2	B	116	LEU
1	F	111	GLU
2	G	70	PRO
2	G	90	ARG
2	G	116	LEU
3	H	406	ALA
3	H	426	PRO
1	A	95	GLY
3	H	428	GLU
1	A	110	ASP
2	B	58	VAL
3	H	334	PRO
3	H	351	GLN
3	H	418	LEU
2	G	130	GLU
3	H	397	TYR
2	G	58	VAL
3	H	368	PRO

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	99/178 (56%)	93 (94%)	6 (6%)	17	46
1	F	100/178 (56%)	92 (92%)	8 (8%)	11	34
2	B	113/123 (92%)	107 (95%)	6 (5%)	20	52
2	G	112/123 (91%)	104 (93%)	8 (7%)	13	40
3	C	102/145 (70%)	96 (94%)	6 (6%)	18	48
3	H	91/145 (63%)	85 (93%)	6 (7%)	15	43
All	All	617/892 (69%)	577 (94%)	40 (6%)	15	44

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	69	ASN
1	A	104	THR
1	A	130	ARG
1	A	139	ARG
1	A	169	THR
1	A	175	GLU
2	B	4	VAL
2	B	14	ASN
2	B	63	ASN
2	B	86	VAL
2	B	88	LEU
2	B	101	MET
3	C	321	ILE
3	C	359	ASP
3	C	378	ARG
3	C	387	GLU
3	C	428	GLU
3	C	429	LEU
1	F	62	LEU
1	F	64	ARG
1	F	77	THR
1	F	90	LYS
1	F	117	LEU
1	F	125	LYS
1	F	134	LEU
1	F	164	ARG
2	G	30	THR
2	G	63	ASN

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Mol	Chain	Res	Type
2	G	86	VAL
2	G	88	LEU
2	G	91	GLU
2	G	117	HIS
2	G	120	ASP
2	G	133	GLN
3	H	346	THR
3	H	353	PHE
3	H	383	LYS
3	H	384	MET
3	H	419	GLN
3	H	430	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	ASN
1	A	82	ASN
1	A	109	ASN
1	A	112	ASN
1	A	119	ASN
1	A	126	ASN
1	A	132	ASN
2	B	8	GLN
2	B	14	ASN
2	B	41	GLN
2	B	104	ASN
2	B	117	HIS
2	B	134	GLN
3	C	380	ASN
3	C	385	ASN
3	C	419	GLN
1	F	82	ASN
1	F	112	ASN
1	F	119	ASN
1	F	127	GLN
1	F	155	ASN
1	F	163	HIS
2	G	63	ASN
2	G	133	GLN
2	G	134	GLN
3	H	336	GLN

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Mol	Chain	Res	Type
3	H	339	GLN
3	H	380	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	118/204 (57%)	-0.00	1 (0%) 82 77	57, 77, 94, 111	0
1	F	119/204 (58%)	-0.18	1 (0%) 82 77	51, 66, 86, 132	0
2	B	130/142 (91%)	0.61	3 (2%) 61 52	80, 105, 131, 136	0
2	G	129/142 (90%)	0.41	5 (3%) 43 35	68, 89, 117, 122	0
3	C	118/166 (71%)	-0.13	1 (0%) 82 77	52, 72, 104, 115	0
3	H	101/166 (60%)	1.25	21 (20%) 2 2	77, 131, 172, 180	0
4	D	15/15 (100%)	-0.46	0 100 100	59, 67, 79, 80	0
4	I	15/15 (100%)	-0.32	0 100 100	61, 72, 90, 92	0
5	E	15/15 (100%)	-0.64	0 100 100	55, 67, 81, 85	0
5	J	15/15 (100%)	-0.23	0 100 100	59, 69, 92, 98	0
All	All	775/1084 (71%)	0.25	32 (4%) 41 33	51, 82, 143, 180	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	G	81	PRO	6.0
2	G	139	ALA	4.5
3	H	360	GLY	4.0
2	B	81	PRO	4.0
3	H	422	LEU	3.8
3	H	433	LEU	3.7
3	H	356	TRP	3.6
3	H	421	LEU	3.4
3	C	319	PRO	3.3
3	H	338	TRP	3.2
3	H	345	LEU	3.2
3	H	342	LEU	2.8
1	A	177	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
2	B	54	GLU	2.6
3	H	364	LYS	2.6
2	G	71	ALA	2.5
3	H	362	GLU	2.4
3	H	333	GLY	2.4
1	F	177	ARG	2.4
3	H	361	TRP	2.3
3	H	357	THR	2.3
3	H	341	LEU	2.2
2	B	88	LEU	2.2
3	H	431	ALA	2.2
3	H	409	ARG	2.2
3	H	430	HIS	2.2
2	G	65	SER	2.1
3	H	363	PHE	2.1
3	H	334	PRO	2.1
3	H	337	LEU	2.1
3	H	365	LEU	2.1
2	G	82	SER	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.