



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 13, 2026 – 11:45 PM UTC

PDB ID : 8Z92 / pdb_00008z92
Title : Crystal structure of CrtAgo/TIR-APAZ in complex with guide DNA and 16-nt target DNA
Authors : Hu, R.; Chen, J.; Liu, L.
Deposited on : 2024-04-22
Resolution : 3.85 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:

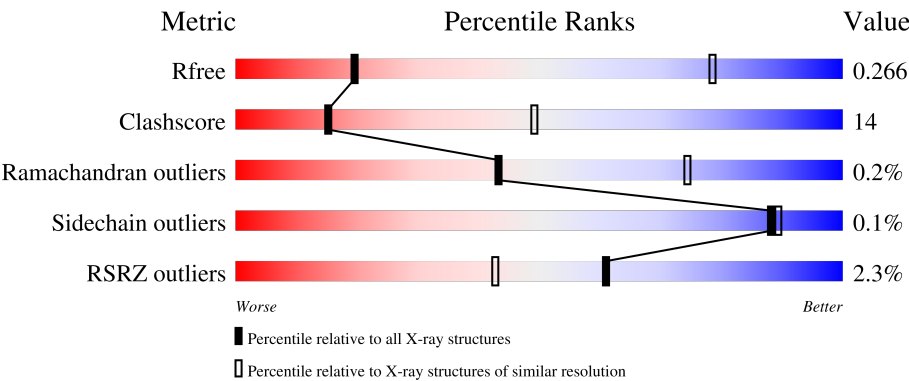
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.85 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	1165 (4.02-3.70)
Clashscore	190562	1207 (4.02-3.70)
Ramachandran outliers	187476	1149 (4.02-3.70)
Sidechain outliers	187428	1142 (4.02-3.70)
RSRZ outliers	180081	1164 (4.02-3.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	507	62%29%8%
1	B	507	5% 65%28%6%
2	E	421	% 69%30%
2	G	421	% 69%30%
3	D	16	25%19%81%

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Mol	Chain	Length	Quality of chain
3	H	16	
4	F	21	
4	I	21	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 16050 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 Piwi domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	476	Total	C	N	O	S	0	0	0
			3842	2488	639	703	12			
1	A	464	Total	C	N	O	S	0	0	0
			3749	2431	623	683	12			

- Molecule 2 is a protein called TIR domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	419	Total	C	N	O	S	0	0	0
			3465	2245	586	623	11			
2	G	419	Total	C	N	O	S	0	0	0
			3480	2255	589	625	11			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	16	Total	C	N	O	P	0	0	0
			315	153	54	93	15			
3	H	16	Total	C	N	O	P	0	0	0
			315	153	54	93	15			

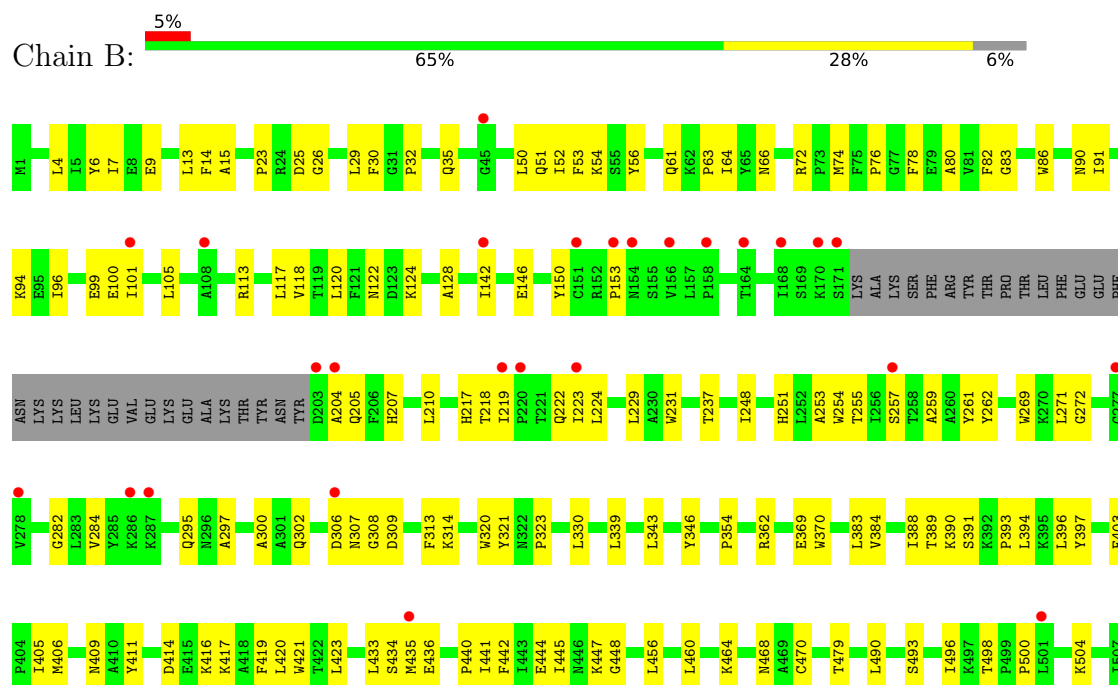
- Molecule 4 is a DNA chain called DNA (5'-D(P*TP*GP*AP*GP*GP*TP*AP*GP*TP*AP*GP*GP*TP*TP*GP*TP*AP*TP*AP*GP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	21	Total	C	N	O	P	0	0	0
			442	210	81	130	21			
4	I	21	Total	C	N	O	P	0	0	0
			442	210	81	130	21			

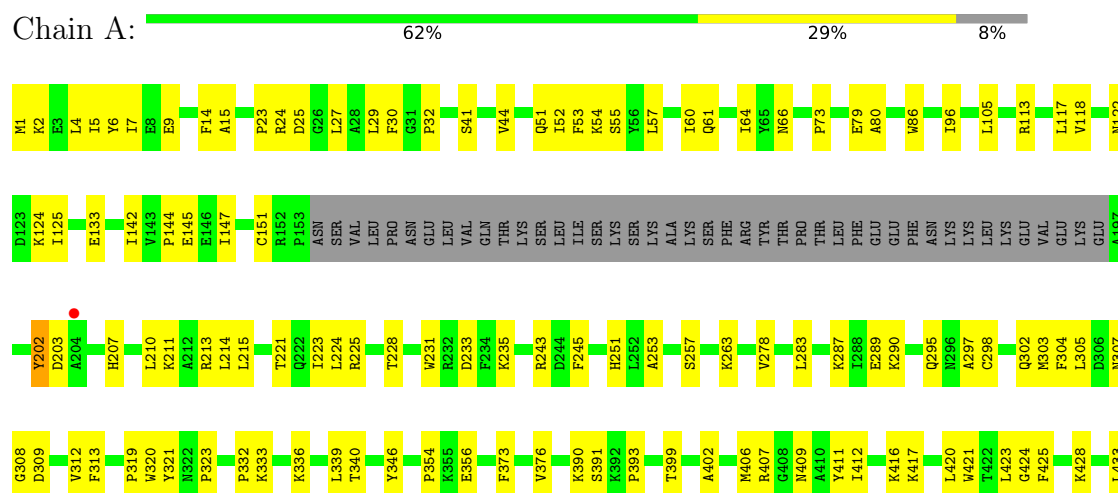
3 Residue-property plots

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: Piwi domain-containing protein

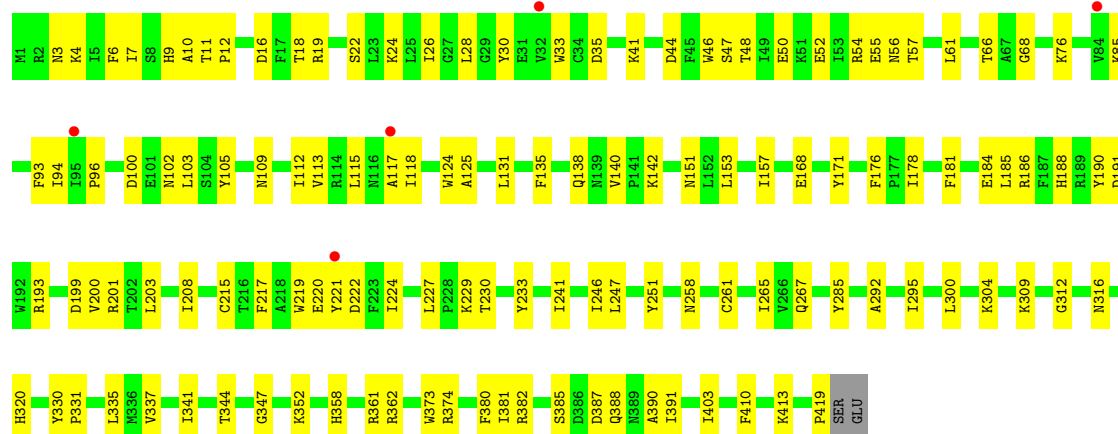


• Molecule 1: Piwi domain-containing protein

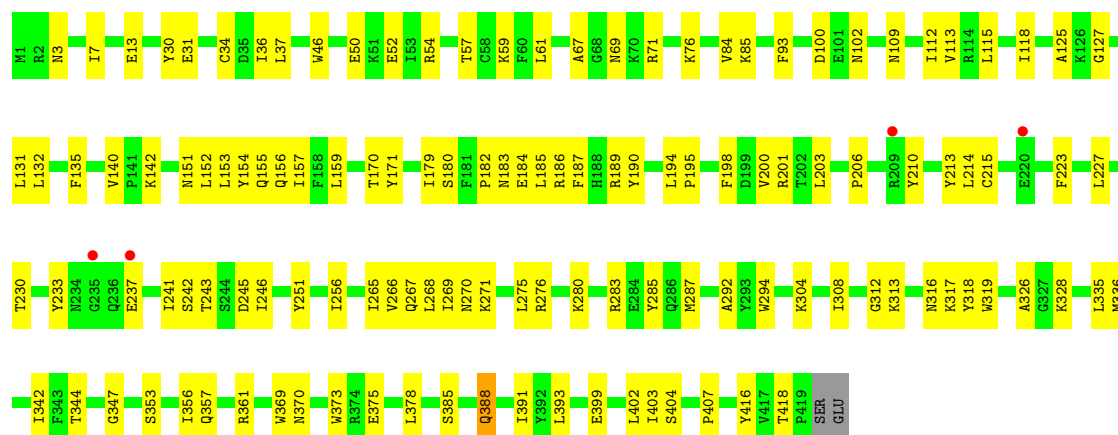




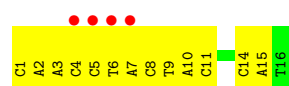
- Molecule 2: TIR domain-containing protein



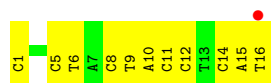
- Molecule 2: TIR domain-containing protein



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



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



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	201.18Å 287.27Å 111.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.25 – 3.85 46.25 – 3.85	Depositor EDS
% Data completeness (in resolution range)	64.8 (46.25-3.85) 77.5 (46.25-3.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.24 (at 3.88Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.247 , 0.265 0.250 , 0.266	Depositor DCC
R_{free} test set	2000 reflections (6.44%)	wwPDB-VP
Wilson B-factor (Å ²)	48.1	Xtriage
Anisotropy	0.234	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 13.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.79	EDS
Total number of atoms	16050	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.58% 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	0.21	0/3845	0.48	0/5206
1	B	0.18	0/3938	0.43	0/5332
2	E	0.18	0/3552	0.43	0/4794
2	G	0.19	0/3568	0.44	0/4814
3	D	0.29	0/351	0.55	0/537
3	H	0.26	0/351	0.54	0/537
4	F	0.28	0/496	0.57	0/765
4	I	0.27	0/496	0.52	0/765
All	All	0.20	0/16597	0.46	0/22750

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
1	A	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	203	ASP	Peptide

5.2 Too-close contacts [i](#)

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

the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3749	0	3752	110	0
1	B	3842	0	3862	107	0
2	E	3465	0	3440	97	0
2	G	3480	0	3467	94	0
3	D	315	0	182	14	0
3	H	315	0	182	12	0
4	F	442	0	240	17	0
4	I	442	0	240	19	0
All	All	16050	0	15365	422	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 14.

All (422) 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:B:210:LEU:HD22	1:B:223:ILE:HD11	1.57	0.86
1:A:9:GLU:OE1	1:A:464:LYS:NZ	2.11	0.83
2:G:30:TYR:HA	2:G:142:LYS:HD2	1.59	0.82
1:B:423:LEU:HD12	1:B:434:SER:HB2	1.59	0.82
2:G:385:SER:HA	2:G:391:ILE:HG12	1.61	0.80
2:G:112:ILE:HG23	2:G:115:LEU:HD12	1.64	0.79
1:B:362:ARG:HH21	1:B:389:THR:HA	1.51	0.75
1:B:118:VAL:O	1:B:122:ASN:ND2	2.18	0.75
2:E:46:TRP:HH2	2:E:76:LYS:HB3	1.51	0.75
1:A:61:GLN:NE2	1:A:86:TRP:O	2.20	0.75
2:G:170:THR:O	1:A:399:THR:OG1	2.05	0.74
1:A:41:SER:OG	1:A:86:TRP:NE1	2.14	0.74
1:B:9:GLU:OE1	1:B:464:LYS:NZ	2.21	0.74
2:G:155:GLN:HA	2:G:159:LEU:HD12	1.69	0.74
1:B:64:ILE:HD11	1:B:253:ALA:HB2	1.69	0.74
2:E:337:VAL:HG11	2:E:381:ILE:HD11	1.68	0.73
1:A:303:MET:HE3	1:A:305:LEU:HD11	1.70	0.73
1:A:423:LEU:HD12	1:A:434:SER:HB2	1.69	0.73
1:B:297:ALA:HB3	1:B:320:TRP:HD1	1.53	0.73
2:G:195:PRO:HG2	2:G:198:PHE:HB2	1.72	0.72
2:G:151:ASN:HB3	1:A:30:PHE:HZ	1.55	0.72
2:E:44:ASP:HB2	2:E:47:SER:HB3	1.71	0.70
2:E:52:GLU:O	2:E:57:THR:OG1	2.09	0.70
1:B:61:GLN:NE2	1:B:86:TRP:O	2.24	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:389:THR:HG23	1:B:442:PHE:HB3	1.74	0.70
1:B:417:LYS:HG2	1:B:444:GLU:HG3	1.72	0.70
1:B:330:LEU:N	1:B:369:GLU:OE1	2.22	0.69
2:G:125:ALA:HB2	1:A:80:ALA:HA	1.74	0.69
1:B:105:LEU:HB2	1:B:117:LEU:HD11	1.75	0.69
1:A:7:ILE:HG23	1:A:457:LYS:HG2	1.75	0.68
1:A:228:THR:OG1	1:A:243:ARG:NH1	2.25	0.68
2:G:266:VAL:O	2:G:270:ASN:ND2	2.27	0.68
1:A:25:ASP:OD2	1:A:428:LYS:NZ	2.23	0.68
2:E:385:SER:HA	2:E:391:ILE:HG12	1.75	0.67
2:E:208:ILE:HD11	2:E:267:GLN:HB2	1.77	0.67
2:E:220:GLU:OE1	2:E:221:TYR:CE2	2.48	0.66
2:G:69:ASN:OD1	2:G:109:ASN:N	2.25	0.66
2:E:200:VAL:HA	2:E:203:LEU:HD23	1.76	0.66
1:B:362:ARG:NH2	1:B:389:THR:HA	2.10	0.65
2:E:219:TRP:N	2:E:222:ASP:OD2	2.29	0.65
3:H:5:DC:O2	4:I:11:DG:N2	2.30	0.65
2:G:190:TYR:HD2	2:G:214:LEU:HD21	1.62	0.65
1:A:278:VAL:HG22	1:A:356:GLU:HB2	1.79	0.65
1:B:51:GLN:HA	1:B:54:LYS:HB2	1.79	0.65
1:A:287:LYS:HE2	1:A:289:GLU:HG2	1.79	0.64
2:E:85:LYS:HG3	2:E:93:PHE:HB3	1.80	0.64
1:A:478:VAL:HA	1:A:481:ARG:HB2	1.79	0.64
2:E:112:ILE:HG22	2:E:115:LEU:HD12	1.80	0.63
1:A:210:LEU:HD23	1:A:223:ILE:HD11	1.80	0.63
2:G:52:GLU:O	2:G:57:THR:OG1	2.16	0.63
1:A:308:GLY:HA3	1:A:465:LEU:HD21	1.81	0.63
1:B:217:HIS:O	1:B:219:ILE:N	2.31	0.63
1:A:15:ALA:HB2	1:A:32:PRO:O	1.98	0.63
2:E:220:GLU:CD	2:E:220:GLU:H	2.07	0.62
2:E:66:THR:OG1	2:E:100:ASP:OD2	2.15	0.62
1:A:51:GLN:HA	1:A:54:LYS:HB2	1.81	0.62
1:B:53:PHE:HD2	1:B:142:ILE:HD11	1.64	0.61
1:A:435:MET:SD	1:A:435:MET:N	2.74	0.61
3:H:9:DT:O2	4:I:7:DG:N2	2.34	0.61
1:A:66:ASN:OD1	1:A:251:HIS:N	2.34	0.61
1:A:215:LEU:HD23	1:A:498:THR:HG21	1.81	0.61
2:G:285:TYR:HB3	2:G:292:ALA:HB3	1.83	0.61
2:G:54:ARG:HA	2:G:84:VAL:HG21	1.82	0.61
1:B:66:ASN:OD1	1:B:251:HIS:N	2.33	0.60
2:G:335:LEU:HD22	2:G:403:ILE:HD13	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:276:ARG:HG2	2:G:393:LEU:HD22	1.83	0.60
2:G:153:LEU:O	2:G:157:ILE:N	2.31	0.60
2:G:283:ARG:HG2	2:G:294:TRP:CZ2	2.36	0.60
4:I:10:DG:H2'	4:I:11:DG:H8	1.66	0.60
1:B:72:ARG:NH1	4:F:14:DG:OP2	2.35	0.60
1:B:391:SER:C	1:B:393:PRO:HD3	2.28	0.59
1:A:308:GLY:HA3	1:A:465:LEU:CD2	2.33	0.59
2:E:382:ARG:O	2:E:385:SER:OG	2.20	0.59
2:G:201:ARG:HH12	4:I:2:DA:P	2.26	0.59
1:A:298:CYS:SG	1:A:490:LEU:HB3	2.42	0.59
1:B:255:THR:HG21	3:D:2:DA:H2''	1.85	0.59
2:E:186:ARG:NH1	2:E:217:PHE:O	2.36	0.59
1:B:435:MET:HG2	4:F:14:DG:O6	2.02	0.58
4:F:12:DT:O3'	4:F:14:DG:N2	2.36	0.58
2:E:335:LEU:HD22	2:E:403:ILE:HD13	1.85	0.58
3:D:5:DC:H2''	3:D:6:DT:H5'	1.86	0.58
1:A:231:TRP:H	1:A:231:TRP:CD1	2.21	0.58
2:G:34:CYS:SG	2:G:36:ILE:HG22	2.43	0.58
1:B:13:LEU:HD23	1:B:272:GLY:HA2	1.86	0.57
2:E:295:ILE:O	2:E:320:HIS:ND1	2.31	0.57
2:E:358:HIS:HB3	2:E:362:ARG:HH22	1.68	0.57
2:G:50:GLU:OE1	2:G:54:ARG:NH2	2.37	0.57
1:B:63:PRO:HG2	2:E:124:TRP:CE2	2.39	0.57
1:B:35:GLN:NE2	1:B:83:GLY:HA3	2.20	0.57
2:G:316:ASN:OD1	2:G:316:ASN:N	2.37	0.57
1:A:14:PHE:HE1	1:A:27:LEU:HD22	1.68	0.57
1:A:390:LYS:O	1:A:390:LYS:HG2	2.04	0.56
1:B:262:TYR:OH	1:B:309:ASP:OD1	2.20	0.56
1:B:297:ALA:HB3	1:B:320:TRP:CD1	2.39	0.56
1:B:370:TRP:CH2	1:B:383:LEU:HG	2.40	0.56
2:G:151:ASN:HB3	1:A:30:PHE:CZ	2.39	0.56
2:G:313:LYS:HA	2:G:318:TYR:HA	1.87	0.56
1:B:99:GLU:OE1	1:B:99:GLU:N	2.36	0.56
1:A:211:LYS:NZ	1:A:221:THR:O	2.38	0.56
2:G:59:LYS:HG3	2:G:135:PHE:HE1	1.71	0.55
2:G:312:GLY:O	2:G:319:TRP:N	2.29	0.55
2:G:182:PRO:HB3	2:G:399:GLU:HB3	1.87	0.55
2:E:247:LEU:H	2:E:247:LEU:HD23	1.70	0.55
2:E:66:THR:HA	2:E:103:LEU:HD11	1.89	0.55
2:E:185:LEU:HD11	2:E:215:CYS:SG	2.47	0.55
1:B:153:PRO:HG3	1:B:204:ALA:HB3	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:9:DT:H72	3:D:10:DA:H62	1.70	0.55
2:E:330:TYR:CD2	2:E:331:PRO:HA	2.42	0.55
2:G:198:PHE:HE1	2:G:227:LEU:HD11	1.72	0.55
1:B:403:PHE:CZ	2:E:419:PRO:HG2	2.41	0.55
2:G:241:ILE:HG21	2:G:246:ILE:HD11	1.89	0.55
4:F:13:DT:O4'	4:F:14:DG:N2	2.40	0.55
1:B:80:ALA:HA	2:E:125:ALA:HB2	1.89	0.54
2:E:190:TYR:O	2:E:193:ARG:HG2	2.07	0.54
2:G:313:LYS:HG2	4:F:18:DA:H5'	1.89	0.54
4:I:10:DG:H2'	4:I:11:DG:C8	2.43	0.54
3:H:5:DC:H2''	3:H:6:DT:H5'	1.89	0.54
3:H:14:DC:H2''	3:H:15:DA:C8	2.42	0.54
1:B:120:LEU:O	1:B:124:LYS:HG2	2.08	0.54
1:A:346:TYR:HE2	1:A:354:PRO:HB3	1.72	0.54
2:E:341:ILE:HB	2:E:361:ARG:HD3	1.90	0.53
1:A:145:GLU:OE2	1:A:225:ARG:NH2	2.33	0.53
1:B:423:LEU:HD13	1:B:433:LEU:HB2	1.90	0.53
1:B:370:TRP:CD1	1:B:447:LYS:HG3	2.44	0.53
1:A:263:LYS:HA	1:A:504:LYS:HD2	1.90	0.53
4:F:0:DT:H2''	4:F:1:DG:C8	2.43	0.53
1:A:117:LEU:HD12	1:A:151:CYS:SG	2.49	0.53
2:G:189:ARG:N	2:G:233:TYR:OH	2.35	0.53
2:G:206:PRO:HB2	2:G:268:LEU:HD22	1.90	0.53
2:E:41:LYS:HG3	2:G:388:GLN:HG3	1.90	0.53
1:B:254:TRP:HB3	1:B:470:CYS:SG	2.48	0.53
1:A:57:LEU:HD21	1:A:86:TRP:CD2	2.43	0.53
2:E:201:ARG:HH21	4:F:1:DG:H4'	1.73	0.53
1:B:406:MET:O	1:B:409:ASN:ND2	2.42	0.53
3:D:3:DA:H2'	3:D:4:DC:H6	1.74	0.52
1:B:390:LYS:O	1:B:390:LYS:HG2	2.09	0.52
2:E:241:ILE:HG21	2:E:246:ILE:HD11	1.91	0.52
1:A:96:ILE:HG12	1:A:124:LYS:HG3	1.91	0.52
1:B:384:VAL:HG22	1:B:448:GLY:HA3	1.90	0.52
1:A:416:LYS:HG2	1:A:445:ILE:HB	1.91	0.52
3:H:11:DC:H42	4:I:4:DG:H1	1.57	0.52
2:E:105:TYR:CE1	2:E:113:VAL:HG13	2.45	0.52
2:E:24:LYS:O	2:E:28:LEU:HG	2.09	0.52
1:B:35:GLN:HA	1:B:261:TYR:OH	2.09	0.52
2:G:198:PHE:CE1	2:G:227:LEU:HD11	2.45	0.52
1:A:231:TRP:CD2	1:A:245:PHE:HB2	2.45	0.52
2:E:109:ASN:O	2:E:113:VAL:HG23	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:417:LYS:HG2	1:A:444:GLU:HG3	1.92	0.51
2:E:9:HIS:NE2	2:E:35:ASP:OD1	2.43	0.51
2:G:85:LYS:HG3	2:G:93:PHE:HB3	1.92	0.51
2:G:267:GLN:HA	2:G:270:ASN:HD22	1.75	0.51
1:B:262:TYR:CE1	1:B:504:LYS:HB3	2.46	0.51
1:A:339:LEU:HD22	1:A:373:PHE:HD1	1.75	0.51
2:E:3:ASN:OD1	2:E:3:ASN:N	2.43	0.51
2:G:233:TYR:CE2	2:G:237:GLU:HB3	2.46	0.51
2:G:265:ILE:O	2:G:269:ILE:HG13	2.10	0.51
1:B:231:TRP:H	1:B:231:TRP:CD1	2.29	0.51
4:I:17:DT:H72	4:I:18:DA:N1	2.26	0.51
2:G:31:GLU:H	2:G:142:LYS:HZ2	1.56	0.51
2:G:7:ILE:HG12	2:G:61:LEU:HB2	1.93	0.50
1:A:51:GLN:HG3	1:A:52:ILE:N	2.26	0.50
2:E:153:LEU:O	2:E:157:ILE:N	2.36	0.50
1:A:2:LYS:HG3	1:A:411:TYR:OH	2.11	0.50
1:A:460:LEU:O	1:A:463:THR:OG1	2.27	0.50
1:A:57:LEU:HD21	1:A:86:TRP:CE2	2.47	0.50
3:D:14:DC:H2''	3:D:15:DA:H8	1.77	0.50
2:E:46:TRP:CH2	2:E:76:LYS:HB3	2.40	0.50
1:B:113:ARG:NE	1:B:150:TYR:O	2.42	0.50
2:E:48:THR:O	2:E:52:GLU:HG2	2.11	0.50
2:E:224:ILE:HG13	2:E:230:THR:HB	1.92	0.50
1:A:2:LYS:HG3	1:A:411:TYR:CZ	2.47	0.50
1:B:7:ILE:HD13	1:B:456:LEU:HB3	1.93	0.50
2:E:184:GLU:HG2	2:E:186:ARG:HH21	1.77	0.50
2:E:199:ASP:OD1	2:E:201:ARG:NH1	2.45	0.49
3:D:10:DA:H4'	3:D:11:DC:OP1	2.11	0.49
1:B:405:ILE:HD11	1:B:421:TRP:CD1	2.48	0.49
2:E:118:ILE:HD12	2:E:131:LEU:HD12	1.93	0.49
2:E:168:GLU:HA	2:E:413:LYS:HA	1.94	0.49
4:I:18:DA:H2''	4:I:19:DG:O4'	2.12	0.49
1:A:253:ALA:O	1:A:257:SER:OG	2.21	0.49
1:A:297:ALA:HB3	1:A:320:TRP:HD1	1.77	0.49
4:I:7:DG:H2'	4:I:8:DT:C6	2.46	0.49
1:B:397:TYR:HE1	2:E:373:TRP:HB3	1.77	0.49
2:G:187:PHE:HE2	2:G:241:ILE:HD12	1.77	0.49
1:A:118:VAL:HG12	1:A:213:ARG:HH21	1.78	0.49
3:H:6:DT:H3	4:I:10:DG:H22	1.60	0.49
1:B:435:MET:O	1:B:436:GLU:HG3	2.12	0.49
1:A:443:ILE:HD13	1:A:459:ILE:HD12	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:295:GLN:HA	1:B:321:TYR:HB3	1.95	0.49
2:E:4:LYS:O	2:E:57:THR:HA	2.12	0.49
2:E:16:ASP:OD1	2:E:19:ARG:NH1	2.45	0.49
1:A:453:GLN:O	1:A:457:LYS:HG3	2.13	0.49
1:B:339:LEU:O	1:B:343:LEU:N	2.39	0.49
1:A:9:GLU:OE1	1:A:407:ARG:NE	2.42	0.49
1:B:416:LYS:HG2	1:B:445:ILE:HB	1.94	0.48
1:B:396:LEU:HD13	2:E:171:TYR:HD2	1.78	0.48
2:E:387:ASP:OD1	2:E:388:GLN:N	2.27	0.48
1:A:340:THR:OG1	1:A:376:VAL:HG11	2.13	0.48
3:D:3:DA:H2'	3:D:4:DC:C6	2.48	0.48
1:A:486:ILE:O	1:A:490:LEU:HD22	2.14	0.48
2:G:180:SER:HB3	2:G:402:LEU:HD12	1.96	0.48
2:E:221:TYR:HA	2:E:224:ILE:HD11	1.95	0.48
1:A:202:TYR:O	1:A:290:LYS:NZ	2.42	0.48
1:A:295:GLN:O	1:A:319:PRO:HA	2.14	0.48
2:E:10:ALA:HB2	2:E:68:GLY:HA2	1.96	0.48
2:E:190:TYR:HE2	2:E:230:THR:HG23	1.78	0.48
1:A:7:ILE:HD13	1:A:456:LEU:HB3	1.95	0.48
1:A:207:HIS:HD1	1:A:223:ILE:HD12	1.78	0.48
1:A:390:LYS:H	1:A:390:LYS:HD3	1.78	0.48
2:E:300:LEU:HD23	2:E:300:LEU:H	1.79	0.48
2:G:287:MET:HE3	3:H:8:DC:H3'	1.95	0.48
1:B:25:ASP:O	1:B:29:LEU:HG	2.12	0.48
1:B:96:ILE:HD13	1:B:120:LEU:HG	1.95	0.48
2:E:93:PHE:HD1	2:E:94:ILE:HG13	1.78	0.47
2:G:125:ALA:HB3	1:A:79:GLU:HG2	1.96	0.47
1:B:50:LEU:O	1:B:54:LYS:N	2.36	0.47
1:A:307:ASN:OD1	1:A:308:GLY:N	2.47	0.47
1:B:396:LEU:HD23	1:B:396:LEU:HA	1.71	0.47
1:B:54:LYS:HE2	1:B:91:ILE:HD12	1.96	0.47
2:E:316:ASN:OD1	2:E:316:ASN:N	2.47	0.47
2:G:34:CYS:SG	2:G:37:LEU:HD23	2.54	0.47
1:B:346:TYR:CE2	1:B:354:PRO:HB3	2.50	0.47
1:A:23:PRO:O	1:A:27:LEU:HD23	2.15	0.47
1:B:51:GLN:HG3	1:B:52:ILE:N	2.30	0.47
1:B:78:PHE:CD1	1:B:82:PHE:HD2	2.33	0.47
1:B:101:ILE:O	1:B:105:LEU:HB3	2.15	0.47
2:G:132:LEU:HA	2:G:135:PHE:HB2	1.96	0.47
1:B:146:GLU:H	1:B:146:GLU:CD	2.23	0.47
1:B:302:GLN:NE2	1:B:479:THR:O	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:6:PHE:HA	2:E:33:TRP:O	2.14	0.47
2:E:30:TYR:HA	2:E:142:LYS:HD2	1.97	0.47
2:G:184:GLU:O	2:G:186:ARG:NH1	2.47	0.47
1:A:346:TYR:CE2	1:A:354:PRO:HB3	2.50	0.47
2:G:171:TYR:CE2	1:A:4:LEU:HD12	2.50	0.47
1:B:282:GLY:O	1:B:302:GLN:HG2	2.15	0.47
2:G:3:ASN:OD1	2:G:3:ASN:N	2.47	0.47
1:A:487:GLY:HA2	1:A:490:LEU:HD23	1.97	0.47
1:B:493:SER:O	1:B:496:ILE:HG12	2.14	0.46
2:E:46:TRP:CD1	2:E:50:GLU:HB2	2.50	0.46
2:G:294:TRP:CD1	2:G:342:ILE:HG13	2.50	0.46
4:F:17:DT:H73	4:I:18:DA:C2	2.50	0.46
1:B:259:ALA:HB2	1:B:468:ASN:HA	1.97	0.46
2:E:285:TYR:HB3	2:E:292:ALA:HB3	1.97	0.46
2:G:46:TRP:HH2	2:G:76:LYS:HB3	1.79	0.46
2:G:179:ILE:HD11	2:G:404:SER:HB2	1.97	0.46
2:G:416:TYR:CE1	2:G:418:THR:HG22	2.50	0.46
3:H:10:DA:H4'	3:H:11:DC:OP1	2.15	0.46
2:E:135:PHE:HD1	2:E:140:VAL:HG21	1.81	0.46
2:E:309:LYS:HZ3	2:E:312:GLY:HA2	1.80	0.46
2:G:67:ALA:O	2:G:71:ARG:HB2	2.16	0.46
2:G:127:GLY:O	2:G:131:LEU:N	2.49	0.46
2:G:369:TRP:O	2:G:373:TRP:HD1	1.98	0.46
3:D:4:DC:H2'	3:D:5:DC:O4'	2.15	0.46
1:B:6:TYR:HD2	1:B:7:ILE:N	2.14	0.46
1:B:314:LYS:HG3	1:B:490:LEU:HD22	1.98	0.46
2:E:135:PHE:CD1	2:E:140:VAL:HG21	2.51	0.46
2:G:190:TYR:HB3	2:G:194:LEU:HG	1.97	0.46
1:B:15:ALA:HB2	1:B:32:PRO:O	2.15	0.46
1:A:211:LYS:HE2	1:A:507:ILE:C	2.41	0.46
1:B:4:LEU:HD11	2:E:410:PHE:HB3	1.97	0.45
1:B:14:PHE:HZ	1:B:23:PRO:HA	1.81	0.45
1:B:90:ASN:C	1:B:90:ASN:HD22	2.22	0.45
2:G:344:THR:HG21	2:G:347:GLY:C	2.41	0.45
1:A:423:LEU:HD13	1:A:433:LEU:HB2	1.98	0.45
1:B:394:LEU:O	1:B:440:PRO:HD2	2.16	0.45
2:G:151:ASN:HA	2:G:154:TYR:HB3	1.98	0.45
1:A:57:LEU:HD23	1:A:57:LEU:HA	1.75	0.45
1:B:14:PHE:CZ	1:B:23:PRO:HA	2.51	0.45
1:B:207:HIS:HB2	1:B:223:ILE:HD12	1.97	0.45
1:B:271:LEU:HD21	1:B:306:ASP:HA	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:308:ILE:HD12	2:G:375:GLU:HB3	1.98	0.45
1:B:7:ILE:CD1	1:B:456:LEU:HB3	2.46	0.45
1:B:30:PHE:HZ	2:E:151:ASN:HB3	1.81	0.45
1:B:100:GLU:HB3	1:B:120:LEU:HD21	1.99	0.45
1:B:7:ILE:HD12	1:B:420:LEU:HD13	1.98	0.45
2:G:242:SER:HB3	2:G:245:ASP:OD2	2.17	0.45
2:G:326:ALA:O	2:G:335:LEU:HD12	2.16	0.45
3:D:15:DA:C2	4:F:1:DG:N1	2.85	0.45
4:F:16:DA:C6	4:F:17:DT:H72	2.51	0.45
1:A:53:PHE:CD2	1:A:142:ILE:HD11	2.51	0.45
2:G:194:LEU:HD12	2:G:194:LEU:O	2.17	0.45
1:A:57:LEU:HD23	1:A:60:ILE:HD11	1.98	0.45
1:A:411:TYR:HD1	1:A:412:ILE:N	2.14	0.45
1:A:214:LEU:HD23	1:A:214:LEU:HA	1.73	0.45
1:B:29:LEU:HB2	1:B:30:PHE:CD2	2.51	0.45
1:B:63:PRO:HG3	1:B:76:PRO:O	2.16	0.45
2:E:295:ILE:HG13	2:E:380:PHE:CE2	2.52	0.45
2:G:152:LEU:O	2:G:156:GLN:HB2	2.17	0.45
1:B:72:ARG:NH2	1:B:248:ILE:HG22	2.32	0.44
1:A:122:ASN:OD1	1:A:214:LEU:HD21	2.17	0.44
1:A:406:MET:HE3	1:A:406:MET:HB2	1.79	0.44
1:B:496:ILE:HD12	1:B:498:THR:HG23	1.98	0.44
2:G:328:LYS:HB2	2:G:336:MET:HE3	1.99	0.44
2:E:96:PRO:HG2	2:E:117:ALA:HA	1.98	0.44
2:G:210:TYR:N	2:G:213:TYR:O	2.33	0.44
3:H:12:DC:H42	4:I:3:DG:H1	1.65	0.44
2:E:7:ILE:HG12	2:E:61:LEU:HB2	1.99	0.44
2:E:100:ASP:OD1	2:E:102:ASN:N	2.49	0.44
1:B:390:LYS:H	1:B:390:LYS:HD3	1.82	0.44
1:A:29:LEU:HB2	1:A:30:PHE:CD2	2.53	0.44
1:A:144:PRO:HD2	1:A:147:ILE:HD12	2.00	0.44
1:A:302:GLN:HA	1:A:312:VAL:HA	1.99	0.44
1:A:420:LEU:HB2	1:A:456:LEU:HD22	1.99	0.44
1:A:435:MET:HE1	3:H:5:DC:H1'	1.99	0.44
2:G:210:TYR:CE1	2:G:256:ILE:HD12	2.53	0.44
2:G:378:LEU:HD21	2:G:407:PRO:CD	2.47	0.44
4:I:7:DG:H2'	4:I:8:DT:H6	1.83	0.44
1:A:223:ILE:O	3:H:1:DC:H4'	2.17	0.44
1:A:313:PHE:HD1	1:A:500:PRO:HG3	1.82	0.44
1:B:56:TYR:CD1	1:B:229:LEU:HD23	2.52	0.44
2:G:200:VAL:HA	2:G:203:LEU:HD23	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:12:DT:C6	4:F:13:DT:H72	2.53	0.44
1:B:222:GLN:NE2	1:B:224:LEU:HD21	2.32	0.44
2:E:304:LYS:HE3	4:I:17:DT:OP1	2.18	0.44
2:G:135:PHE:CD1	2:G:140:VAL:HG21	2.53	0.44
1:A:51:GLN:O	1:A:55:SER:N	2.37	0.44
1:A:421:TRP:CD2	1:A:440:PRO:HB3	2.52	0.44
2:E:229:LYS:HA	2:E:229:LYS:HD3	1.46	0.43
2:E:387:ASP:CG	2:E:388:GLN:H	2.17	0.43
2:G:171:TYR:HE2	1:A:4:LEU:HD12	1.81	0.43
4:F:19:DG:H21	4:I:16:DA:H62	1.66	0.43
2:E:261:CYS:O	2:E:265:ILE:HG13	2.18	0.43
1:A:309:ASP:O	1:A:503:PHE:HB2	2.19	0.43
1:B:74:MET:HE1	2:E:124:TRP:CH2	2.53	0.43
2:E:220:GLU:OE1	2:E:221:TYR:CZ	2.70	0.43
2:G:280:LYS:HD3	2:G:280:LYS:HA	1.86	0.43
1:A:5:ILE:O	1:A:5:ILE:HG13	2.19	0.43
1:A:333:LYS:HE2	1:A:333:LYS:HB2	1.85	0.43
2:G:353:SER:HB3	2:G:356:ILE:HD12	2.01	0.43
2:G:357:GLN:O	2:G:361:ARG:HB2	2.18	0.43
1:B:388:ILE:HG23	1:B:441:ILE:HD12	2.00	0.43
1:B:253:ALA:O	1:B:257:SER:OG	2.24	0.43
1:B:396:LEU:HD21	2:E:410:PHE:CD2	2.54	0.43
1:A:339:LEU:HD22	1:A:373:PHE:CD1	2.54	0.43
2:G:100:ASP:OD1	2:G:102:ASN:N	2.50	0.43
1:A:60:ILE:HA	1:A:64:ILE:HD11	2.01	0.43
2:E:41:LYS:CG	2:G:388:GLN:HG3	2.48	0.43
1:A:6:TYR:HD2	1:A:7:ILE:N	2.17	0.43
1:A:493:SER:HB2	1:A:496:ILE:HG12	1.99	0.43
1:B:321:TYR:O	1:B:323:PRO:HD3	2.19	0.42
2:E:188:HIS:HB3	2:E:233:TYR:HE1	1.84	0.42
2:G:317:LYS:H	2:G:317:LYS:HG3	1.63	0.42
1:A:14:PHE:CE1	1:A:27:LEU:HD22	2.53	0.42
1:B:205:GLN:CG	3:D:1:DC:H41	2.32	0.42
2:E:178:ILE:HG21	2:E:181:PHE:CE1	2.54	0.42
2:E:193:ARG:O	2:E:227:LEU:HD21	2.18	0.42
1:A:133:GLU:OE2	1:A:133:GLU:N	2.44	0.42
1:B:313:PHE:HD1	1:B:500:PRO:HG3	1.84	0.42
2:G:13:GLU:OE2	2:G:13:GLU:N	2.46	0.42
2:G:271:LYS:HE3	2:G:275:LEU:HD11	2.00	0.42
2:G:378:LEU:HD23	2:G:378:LEU:HA	1.79	0.42
2:G:109:ASN:O	2:G:113:VAL:HG23	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:283:LEU:HD23	1:A:283:LEU:HA	1.85	0.42
1:A:321:TYR:O	1:A:323:PRO:HD3	2.20	0.42
1:A:406:MET:HA	1:A:425:PHE:HB3	2.01	0.42
1:B:411:TYR:HB3	1:B:419:PHE:HB2	2.01	0.42
2:E:142:LYS:HA	2:E:142:LYS:HD3	1.86	0.42
2:E:190:TYR:CE2	2:E:230:THR:HG23	2.54	0.42
2:G:201:ARG:NH1	4:I:1:DG:H4'	2.35	0.42
4:F:16:DA:N1	4:F:17:DT:H72	2.34	0.42
4:F:18:DA:N1	4:I:17:DT:C7	2.82	0.42
1:B:94:LYS:HG3	1:B:128:ALA:HB2	2.01	0.42
1:B:307:ASN:OD1	1:B:308:GLY:N	2.53	0.42
1:B:405:ILE:HD11	1:B:421:TRP:CG	2.54	0.42
2:G:183:ASN:O	2:G:243:THR:OG1	2.33	0.42
1:A:1:MET:HE3	1:A:1:MET:HB2	1.85	0.42
1:A:118:VAL:CG1	1:A:213:ARG:HH21	2.32	0.42
1:A:233:ASP:O	1:A:235:LYS:HG3	2.20	0.42
2:G:185:LEU:HD11	2:G:215:CYS:HB2	2.01	0.42
1:A:105:LEU:HD12	1:A:113:ARG:HD2	2.02	0.42
1:A:503:PHE:HE1	1:A:507:ILE:HD11	1.85	0.42
1:B:26:GLY:O	1:B:30:PHE:N	2.48	0.42
2:G:46:TRP:CD1	2:G:50:GLU:HB2	2.55	0.42
3:D:9:DT:H6	3:D:9:DT:H2'	1.70	0.42
1:B:460:LEU:HD12	1:B:460:LEU:HA	1.90	0.41
2:E:193:ARG:HD2	2:E:229:LYS:O	2.20	0.41
2:E:358:HIS:CD2	3:D:8:DC:H4'	2.55	0.41
2:G:46:TRP:HH2	2:G:76:LYS:CB	2.33	0.41
2:G:304:LYS:HE3	4:F:17:DT:OP1	2.20	0.41
1:A:406:MET:O	1:A:409:ASN:ND2	2.53	0.41
1:B:53:PHE:CD2	1:B:142:ILE:HD11	2.50	0.41
2:E:11:THR:HA	2:E:12:PRO:HA	1.80	0.41
2:G:246:ILE:HD13	2:G:251:TYR:HB3	2.02	0.41
2:G:313:LYS:HB3	2:G:313:LYS:HE3	1.82	0.41
2:E:138:GLN:O	2:E:140:VAL:HG23	2.20	0.41
1:A:313:PHE:CD1	1:A:500:PRO:HG3	2.55	0.41
1:A:24:ARG:HD2	1:A:73:PRO:HD2	2.02	0.41
3:H:15:DA:H2''	3:H:16:DT:C6	2.56	0.41
1:A:424:GLY:HA2	1:A:438:PRO:HG3	2.03	0.41
1:B:205:GLN:HG2	3:D:1:DC:H41	1.86	0.41
2:E:50:GLU:O	2:E:54:ARG:NH1	2.54	0.41
2:E:344:THR:HG21	2:E:347:GLY:C	2.45	0.41
2:G:276:ARG:HG2	2:G:393:LEU:CD2	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:370:ASN:HB3	1:A:402:ALA:C	2.46	0.41
4:F:17:DT:H73	4:I:18:DA:H2	1.86	0.41
2:E:176:PHE:CE2	2:E:381:ILE:HG21	2.56	0.41
2:G:223:PHE:O	2:G:230:THR:HG21	2.20	0.41
1:A:391:SER:C	1:A:393:PRO:HD3	2.45	0.41
1:B:396:LEU:HD11	2:E:410:PHE:HD2	1.86	0.41
1:B:414:ASP:OD1	1:B:417:LYS:N	2.52	0.41
2:E:251:TYR:O	2:E:258:ASN:HB2	2.20	0.41
1:B:14:PHE:CE1	1:B:269:TRP:HB3	2.56	0.41
1:B:397:TYR:CD1	2:E:374:ARG:HB2	2.56	0.41
2:E:55:GLU:HG2	2:E:56:ASN:N	2.35	0.41
1:A:447:LYS:HA	1:A:447:LYS:HD3	1.85	0.41
3:D:7:DA:N1	4:F:9:DA:H2	2.19	0.41
1:B:66:ASN:ND2	1:B:72:ARG:O	2.45	0.40
2:E:22:SER:O	2:E:26:ILE:HG13	2.21	0.40
2:E:387:ASP:OD2	2:E:390:ALA:HB3	2.20	0.40
2:G:118:ILE:HD12	2:G:131:LEU:HD12	2.03	0.40
2:E:105:TYR:HD1	2:E:105:TYR:HA	1.77	0.40
2:E:352:LYS:HD3	2:E:352:LYS:HA	1.89	0.40
1:A:304:PHE:HD1	1:A:479:THR:HG22	1.85	0.40
2:G:190:TYR:CD2	2:G:214:LEU:HD21	2.50	0.40
1:A:24:ARG:HD3	1:A:470:CYS:O	2.21	0.40
1:A:44:VAL:HG21	1:A:125:ILE:HG12	2.03	0.40
1:A:332:PRO:O	1:A:336:LYS:HG3	2.21	0.40
1:B:284:VAL:HB	1:B:300:ALA:HB3	2.04	0.40
1:B:346:TYR:HE2	1:B:354:PRO:HB3	1.85	0.40
2:E:18:THR:O	2:E:22:SER:OG	2.35	0.40
2:E:309:LYS:NZ	4:I:17:DT:H3'	2.36	0.40
1:A:223:ILE:O	1:A:224:LEU:HD23	2.22	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	460/507 (91%)	441 (96%)	18 (4%)	1 (0%)	43	74
1	B	472/507 (93%)	454 (96%)	17 (4%)	1 (0%)	43	74
2	E	417/421 (99%)	400 (96%)	16 (4%)	1 (0%)	43	74
2	G	417/421 (99%)	399 (96%)	17 (4%)	1 (0%)	43	74
All	All	1766/1856 (95%)	1694 (96%)	68 (4%)	4 (0%)	43	74

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	218	THR
2	E	191	ASP
2	G	388	GLN
1	A	202	TYR

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	401/446 (90%)	401 (100%)	0	100	100
1	B	416/446 (93%)	415 (100%)	1 (0%)	87	87
2	E	374/387 (97%)	374 (100%)	0	100	100
2	G	378/387 (98%)	378 (100%)	0	100	100
All	All	1569/1666 (94%)	1568 (100%)	1 (0%)	88	89

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	237	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	205	GLN

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Mol	Chain	Res	Type
1	B	302	GLN
2	E	129	GLN
2	E	226	GLN
2	E	262	GLN
2	E	349	ASN
2	G	226	GLN
2	G	270	ASN
1	A	107	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	464/507 (91%)	0.09	1 (0%) 91 81	26, 33, 69, 114	0
1	B	476/507 (93%)	0.64	26 (5%) 30 24	31, 83, 132, 164	0
2	E	419/421 (99%)	0.29	5 (1%) 76 55	27, 55, 104, 126	0
2	G	419/421 (99%)	0.17	4 (0%) 79 59	26, 42, 86, 108	0
3	D	16/16 (100%)	1.52	4 (25%) 2 4	86, 122, 164, 172	0
3	H	16/16 (100%)	0.91	1 (6%) 26 22	31, 62, 171, 174	0
4	F	21/21 (100%)	0.97	0 100 100	44, 102, 135, 153	0
4	I	21/21 (100%)	0.89	2 (9%) 14 16	34, 71, 142, 155	0
All	All	1852/1930 (95%)	0.33	43 (2%) 61 43	26, 51, 117, 174	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	153	PRO	4.0
1	B	204	ALA	3.7
1	B	287	LYS	3.3
3	D	5	DC	3.3
1	B	151	CYS	3.1
1	B	435	MET	3.1
2	G	235	GLY	3.1
1	B	219	ILE	3.1
1	B	171	SER	3.0
4	I	0	DT	2.9
1	B	45	GLY	2.9
1	B	278	VAL	2.9
3	D	6	DT	2.7
1	B	108	ALA	2.7
1	B	257	SER	2.6
1	A	204	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
2	G	237	GLU	2.6
1	B	277	GLY	2.5
2	E	117	ALA	2.5
1	B	154	ASN	2.5
3	D	7	DA	2.4
1	B	156	VAL	2.4
1	B	164	THR	2.4
1	B	223	ILE	2.4
2	G	209	ARG	2.4
1	B	170	LYS	2.3
3	H	16	DT	2.2
1	B	286	LYS	2.2
2	E	221	TYR	2.2
1	B	158	PRO	2.1
1	B	203	ASP	2.1
1	B	168	ILE	2.1
1	B	306	ASP	2.1
2	G	220	GLU	2.1
4	I	6	DA	2.0
2	E	84	VAL	2.0
1	B	220	PRO	2.0
2	E	95	ILE	2.0
3	D	4	DC	2.0
1	B	101	ILE	2.0
1	B	142	ILE	2.0
1	B	501	LEU	2.0
2	E	32	VAL	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.