



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:22 AM UTC

PDB ID : 3JV4 / pdb_00003jv4
Title : Crystal structure of the dimerization domains p50 and RelB
Authors : Vu, D.; Huang, D.B.; Ghosh, G.
Deposited on : 2009-09-15
Resolution : 3.15 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

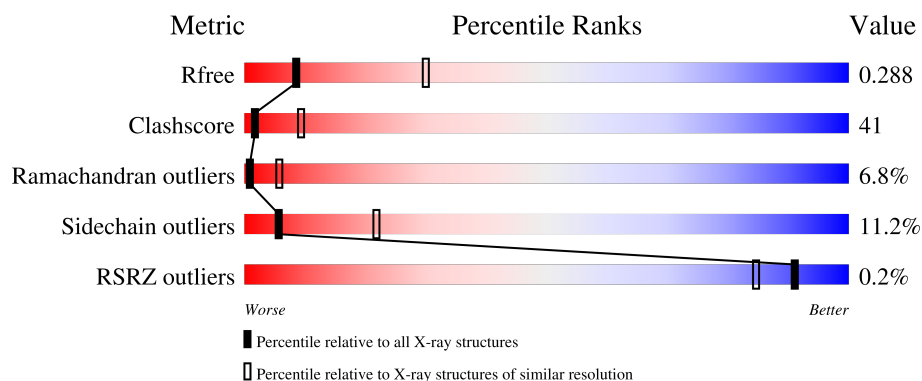
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.15 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	2361 (3.20-3.12)
Clashscore	190562	2486 (3.20-3.12)
Ramachandran outliers	187476	2405 (3.20-3.12)
Sidechain outliers	187428	2404 (3.20-3.12)
RSRZ outliers	180081	2361 (3.20-3.12)

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	101	35% 51% 13% .
1	C	101	37% 50% 13% .
1	E	101	36% 54% 10%
2	B	115	30% 56% 11% .
2	D	115	40% 44% 15% .

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Mol	Chain	Length	Quality of chain
2	F	115	34% 56% 7% . .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5185 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 Transcription factor RelB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	101	Total	C	N	O	S	0	0	0
			800	507	132	157	4			
1	C	101	Total	C	N	O	S	0	0	0
			800	507	132	157	4			
1	E	101	Total	C	N	O	S	0	0	0
			800	507	132	157	4			

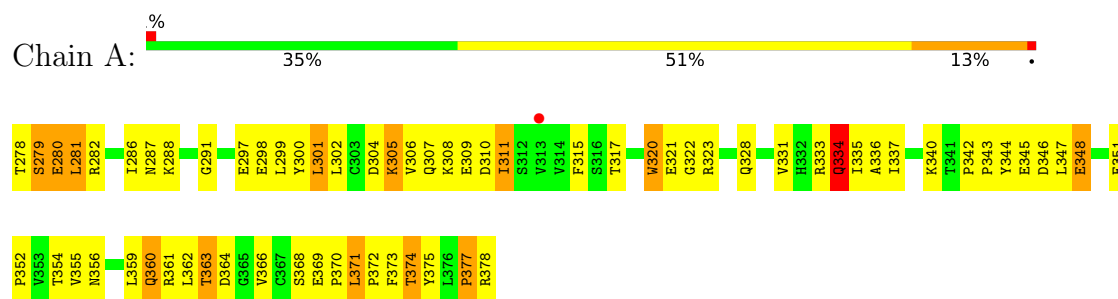
- Molecule 2 is a protein called Nuclear factor NF-kappa-B p105 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	115	Total	C	N	O	S	0	0	0
			932	594	154	181	3			
2	D	115	Total	C	N	O	S	0	0	0
			942	600	158	181	3			
2	F	112	Total	C	N	O	S	0	0	0
			911	581	150	177	3			

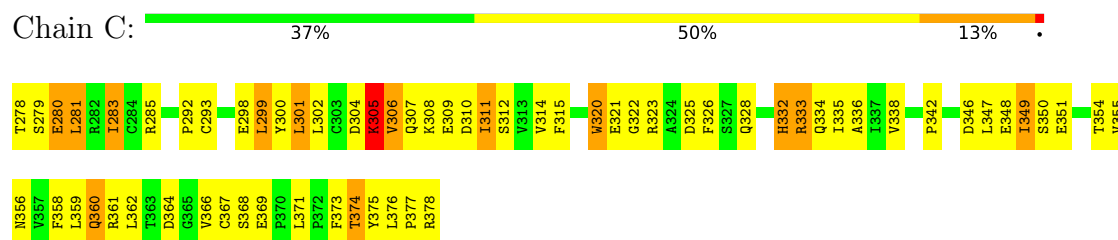
3 Residue-property plots

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.

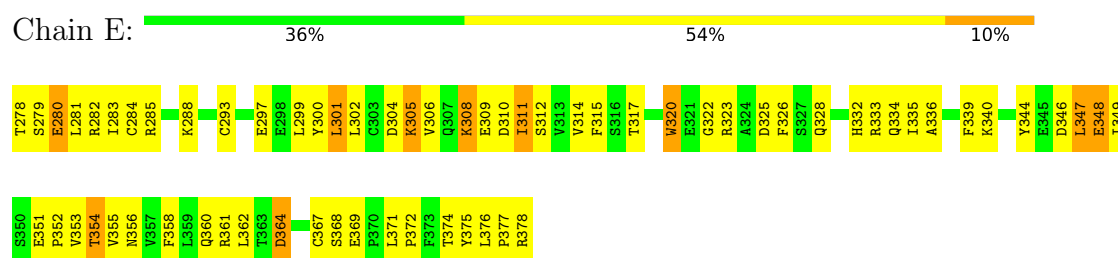
• Molecule 1: Transcription factor RelB



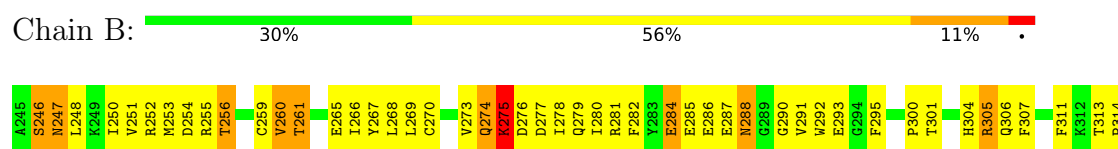
• Molecule 1: Transcription factor RelB



• Molecule 1: Transcription factor RelB

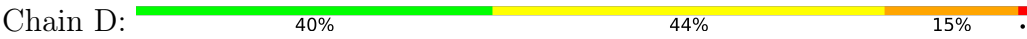


• Molecule 2: Nuclear factor NF-kappa-B p105 subunit

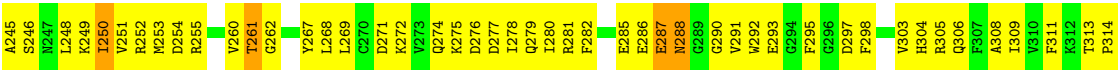
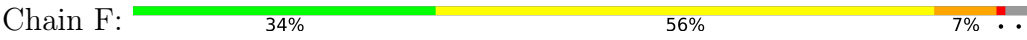




● Molecule 2: Nuclear factor NF-kappa-B p105 subunit



● Molecule 2: Nuclear factor NF-kappa-B p105 subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	106.17Å 106.17Å 143.78Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.76 – 3.15 29.76 – 3.15	Depositor EDS
% Data completeness (in resolution range)	81.6 (29.76-3.15) 90.7 (29.76-3.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 3.12Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.225 , 0.280 0.231 , 0.288	Depositor DCC
R_{free} test set	792 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	83.3	Xtriage
Anisotropy	0.297	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 44.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.035 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5185	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% 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	0.47	0/817	0.91	1/1111 (0.1%)
1	C	0.52	0/817	1.06	3/1111 (0.3%)
1	E	0.43	0/817	0.94	1/1111 (0.1%)
2	B	0.59	0/952	1.10	6/1285 (0.5%)
2	D	0.55	0/962	1.08	4/1296 (0.3%)
2	F	0.49	0/931	1.02	4/1256 (0.3%)
All	All	0.51	0/5296	1.02	19/7170 (0.3%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	280	GLU	N-CA-C	8.38	121.17	111.11
2	F	246	SER	N-CA-C	6.25	117.89	111.14
2	D	341	GLU	N-CA-C	-6.18	101.79	109.64
2	D	248	LEU	N-CA-C	-5.96	99.98	109.76
2	B	348	TYR	CA-C-N	5.80	125.81	119.89
2	B	348	TYR	C-N-CA	5.80	125.81	119.89
1	C	332	HIS	N-CA-C	5.67	117.97	107.99
1	C	320	TRP	N-CA-C	5.56	118.29	109.50
2	F	346	LEU	N-CA-C	5.51	117.94	109.07
2	B	356	GLU	N-CA-C	-5.48	107.16	112.97
2	B	246	SER	N-CA-C	5.47	117.11	108.96
2	F	348	TYR	CA-C-N	5.44	126.64	119.84
2	F	348	TYR	C-N-CA	5.44	126.64	119.84
2	B	338	GLU	N-CA-C	-5.40	102.48	110.48
2	D	321	ILE	CB-CA-C	-5.33	105.47	111.45
1	A	320	TRP	N-CA-C	5.24	117.78	109.50
2	B	339	THR	N-CA-C	5.13	117.44	108.76
2	D	249	LYS	N-CA-C	5.10	117.66	108.48
1	E	320	TRP	N-CA-C	5.08	117.53	109.50

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	800	0	785	71	0
1	C	800	0	785	75	0
1	E	800	0	785	67	0
2	B	932	0	904	106	0
2	D	942	0	926	77	0
2	F	911	0	885	62	0
All	All	5185	0	5070	417	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 41.

All (417) 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:C:306:VAL:HG11	1:C:311:ILE:HG12	1.41	1.02
1:E:347:LEU:HD12	1:E:347:LEU:H	1.30	0.96
2:B:317:LYS:HG3	2:B:318:ASP:H	1.31	0.94
1:C:293:CYS:HA	1:C:375:TYR:HD2	1.31	0.94
1:C:283:ILE:HG12	1:C:359:LEU:CD1	2.00	0.92
2:F:317:LYS:HG3	2:F:318:ASP:H	1.34	0.91
2:F:287:GLU:HG2	2:F:288:ASN:H	1.36	0.90
1:A:311:ILE:HD11	1:A:359:LEU:HD22	1.52	0.90
1:E:293:CYS:HA	1:E:375:TYR:HD2	1.38	0.87
1:C:377:PRO:O	1:C:378:ARG:HB2	1.74	0.86
2:F:278:ILE:HD11	2:F:331:LEU:HB3	1.58	0.85
2:B:336:ASP:O	2:B:337:LEU:HD12	1.75	0.85
2:F:260:VAL:HG21	2:F:321:ILE:HD11	1.60	0.83
1:C:283:ILE:HG12	1:C:359:LEU:HD11	1.58	0.83
1:C:301:LEU:HD12	1:C:301:LEU:C	2.02	0.83
2:D:300:PRO:O	2:D:303:VAL:HG23	1.77	0.83
2:B:250:ILE:HG21	2:B:253:MET:HE2	1.61	0.81
1:C:338:VAL:HG11	2:D:252:ARG:HH11	1.44	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:260:VAL:HG22	2:B:316:TYR:HB3	1.63	0.81
1:A:286:ILE:HG22	1:A:288:LYS:H	1.46	0.80
1:A:301:LEU:HD23	1:A:301:LEU:O	1.81	0.80
2:D:260:VAL:HG13	2:D:319:VAL:HG23	1.63	0.80
2:F:317:LYS:CG	2:F:318:ASP:H	1.95	0.79
2:F:317:LYS:HG3	2:F:318:ASP:N	1.97	0.79
2:B:322:THR:HG22	2:D:255:ARG:NH1	1.98	0.79
2:D:317:LYS:HG3	2:D:318:ASP:H	1.48	0.79
1:E:377:PRO:O	1:E:378:ARG:HB2	1.83	0.78
2:F:324:PRO:HB2	2:F:346:LEU:HD11	1.65	0.78
2:F:250:ILE:HG21	2:F:343:LYS:HG3	1.63	0.78
2:B:317:LYS:HG3	2:B:318:ASP:N	1.99	0.78
1:C:283:ILE:HG23	1:C:301:LEU:HD13	1.66	0.77
1:C:293:CYS:HA	1:C:375:TYR:CD2	2.18	0.77
2:F:333:ARG:HH21	2:F:338:GLU:HG2	1.49	0.77
2:B:357:VAL:O	2:B:358:GLN:HB2	1.86	0.76
1:A:282:ARG:HA	1:A:368:SER:OG	1.86	0.75
1:C:298:GLU:C	1:C:299:LEU:HD23	2.12	0.75
2:D:250:ILE:HG21	2:D:343:LYS:HG3	1.67	0.74
1:A:335:ILE:CD1	2:B:307:PHE:HB2	2.17	0.74
1:A:306:VAL:O	1:A:334:GLN:HB3	1.88	0.74
2:F:278:ILE:HD12	2:F:279:GLN:H	1.53	0.73
2:D:277:ASP:OD1	2:D:334:LYS:HB2	1.87	0.73
1:A:311:ILE:N	1:A:362:LEU:HD13	2.04	0.71
2:F:303:VAL:HG13	2:F:309:ILE:HG12	1.73	0.71
2:D:338:GLU:OE1	2:D:338:GLU:HA	1.90	0.70
2:D:355:GLU:O	2:D:357:VAL:N	2.24	0.70
1:E:349:ILE:HD12	1:E:351:GLU:O	1.92	0.70
1:C:302:LEU:HB3	2:D:304:HIS:CD2	2.27	0.69
1:A:335:ILE:HD13	2:B:307:PHE:HB2	1.74	0.69
2:B:256:THR:OG1	2:D:323:LYS:HD2	1.93	0.69
2:B:274:GLN:O	2:B:276:ASP:N	2.26	0.69
1:A:351:GLU:OE2	1:A:352:PRO:HD2	1.93	0.68
2:D:275:LYS:HG2	2:D:300:PRO:HB3	1.75	0.68
2:D:303:VAL:HG13	2:D:309:ILE:HG12	1.73	0.68
2:F:287:GLU:HG2	2:F:288:ASN:N	2.09	0.68
1:A:320:TRP:CD1	1:A:345:GLU:HB2	2.29	0.67
2:B:268:LEU:C	2:B:268:LEU:HD23	2.20	0.67
1:C:283:ILE:CG2	1:C:301:LEU:HD13	2.23	0.67
1:E:317:THR:HG23	1:E:320:TRP:HE3	1.60	0.67
2:F:250:ILE:CG2	2:F:343:LYS:HG3	2.26	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:341:GLU:HG3	2:B:342:PRO:HD2	1.77	0.66
1:A:335:ILE:HB	2:B:304:HIS:NE2	2.11	0.66
2:B:279:GLN:OE1	2:B:332:ARG:HD3	1.96	0.66
1:C:279:SER:H	1:E:308:LYS:HZ2	1.43	0.66
2:B:322:THR:HG22	2:D:255:ARG:HH12	1.59	0.65
2:B:261:THR:O	2:B:315:LYS:HB2	1.96	0.65
2:D:305:ARG:O	2:D:306:GLN:HB2	1.95	0.65
1:A:363:THR:OG1	1:A:364:ASP:N	2.29	0.65
2:D:259:CYS:SG	2:D:350:GLU:HG2	2.36	0.65
1:C:304:ASP:CG	2:D:305:ARG:NH1	2.54	0.64
2:F:250:ILE:HG12	2:F:268:LEU:HD21	1.79	0.64
1:A:287:ASN:HD22	1:A:300:TYR:HB2	1.63	0.63
1:A:354:THR:HA	1:A:374:THR:HA	1.80	0.63
2:B:254:ASP:O	2:B:255:ARG:HD2	1.98	0.63
1:E:304:ASP:O	1:E:306:VAL:HG13	1.99	0.63
1:E:301:LEU:C	1:E:301:LEU:HD12	2.24	0.62
2:F:333:ARG:HG2	2:F:335:SER:OG	1.99	0.62
1:A:331:VAL:HG22	1:A:337:ILE:HG12	1.80	0.62
2:D:292:TRP:CD1	2:D:317:LYS:HB3	2.35	0.62
1:A:311:ILE:H	1:A:362:LEU:HD13	1.61	0.61
1:E:293:CYS:HA	1:E:375:TYR:CD2	2.28	0.61
1:E:283:ILE:HG23	1:E:301:LEU:HD13	1.83	0.61
1:C:307:GLN:HG3	1:C:309:GLU:OE1	2.01	0.61
1:A:288:LYS:HE3	1:A:297:GLU:OE1	2.00	0.61
2:D:277:ASP:HA	2:D:334:LYS:HD2	1.83	0.61
1:A:287:ASN:ND2	1:A:300:TYR:HB2	2.15	0.60
1:A:361:ARG:HH12	1:A:366:VAL:HG21	1.65	0.60
2:B:247:ASN:HD22	2:B:333:ARG:HH22	1.48	0.60
2:F:252:ARG:O	2:F:253:MET:HG3	2.02	0.60
2:B:250:ILE:HG12	2:B:331:LEU:HG	1.84	0.60
2:B:354:LYS:C	2:B:356:GLU:H	2.08	0.60
1:C:306:VAL:HG11	1:C:311:ILE:CG1	2.26	0.59
2:D:316:TYR:O	2:D:317:LYS:C	2.43	0.59
2:B:273:VAL:HG22	2:B:306:GLN:O	2.01	0.59
1:C:308:LYS:O	1:C:328:GLN:HG2	2.02	0.59
1:C:311:ILE:HD11	1:C:359:LEU:HB3	1.84	0.59
2:D:333:ARG:HH21	2:D:338:GLU:HG3	1.68	0.59
1:E:299:LEU:HB2	1:E:339:PHE:CE1	2.38	0.59
1:A:361:ARG:NH1	1:A:366:VAL:HG21	2.18	0.59
2:B:321:ILE:O	2:B:349:PRO:HG3	2.02	0.59
1:C:346:ASP:O	1:C:349:ILE:HG23	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:344:TYR:HB2	1:E:375:TYR:CE2	2.38	0.58
2:F:249:LYS:HA	2:F:340:SER:OG	2.03	0.58
2:B:277:ASP:O	2:B:333:ARG:HA	2.04	0.58
1:C:285:ARG:CZ	2:D:310:VAL:HG21	2.34	0.58
2:B:286:GLU:O	2:B:287:GLU:HB3	2.03	0.58
1:C:298:GLU:O	1:C:299:LEU:HD23	2.04	0.58
1:E:288:LYS:HB2	1:E:299:LEU:CD2	2.34	0.58
1:E:308:LYS:O	1:E:328:GLN:HG2	2.03	0.58
2:D:250:ILE:CG2	2:D:343:LYS:HG3	2.34	0.57
1:C:364:ASP:CG	1:C:366:VAL:HG23	2.29	0.57
2:B:352:LYS:HE3	2:D:255:ARG:HH21	1.70	0.57
2:B:265:GLU:O	2:B:266:ILE:HD12	2.04	0.57
2:B:292:TRP:NE1	2:B:317:LYS:HB3	2.20	0.57
1:C:361:ARG:CD	1:C:364:ASP:OD2	2.52	0.57
1:E:317:THR:HG22	1:E:355:VAL:HG12	1.86	0.57
2:D:322:THR:O	2:D:323:LYS:HD3	2.04	0.57
1:E:278:THR:HG23	1:E:279:SER:H	1.68	0.57
2:D:354:LYS:O	2:D:355:GLU:C	2.47	0.57
1:A:304:ASP:CG	2:B:305:ARG:HH22	2.13	0.57
1:C:338:VAL:HG11	2:D:252:ARG:NH1	2.17	0.57
2:D:272:LYS:HD3	2:D:306:GLN:HE22	1.67	0.57
1:C:283:ILE:HB	1:C:371:LEU:HD22	1.87	0.57
2:D:250:ILE:HG13	2:D:268:LEU:HD21	1.87	0.57
2:F:295:PHE:O	2:F:314:PRO:HB3	2.05	0.56
1:A:333:ARG:C	1:A:334:GLN:HG2	2.29	0.56
1:A:306:VAL:H	1:A:334:GLN:HB2	1.71	0.56
1:A:317:THR:HG22	1:A:355:VAL:HG12	1.87	0.56
2:D:331:LEU:H	2:D:340:SER:HB3	1.70	0.56
1:E:354:THR:HA	1:E:374:THR:HA	1.87	0.56
2:F:274:GLN:O	2:F:276:ASP:N	2.38	0.56
2:D:272:LYS:HD3	2:D:306:GLN:NE2	2.21	0.56
1:E:333:ARG:O	1:E:335:ILE:HG23	2.05	0.56
1:E:351:GLU:OE2	1:E:352:PRO:HD2	2.06	0.56
2:F:321:ILE:O	2:F:349:PRO:HG3	2.05	0.56
1:A:278:THR:O	1:A:279:SER:HB2	2.06	0.56
2:D:317:LYS:HG3	2:D:318:ASP:N	2.20	0.56
2:F:277:ASP:O	2:F:333:ARG:HA	2.05	0.56
1:E:300:TYR:CE2	2:F:254:ASP:HA	2.41	0.56
2:B:275:LYS:HB2	2:B:300:PRO:HB3	1.86	0.55
2:B:354:LYS:O	2:B:356:GLU:N	2.36	0.55
2:B:278:ILE:HD11	2:B:331:LEU:HB3	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:359:LEU:HD12	1:C:359:LEU:H	1.71	0.55
1:A:300:TYR:CE1	2:B:269:LEU:HD11	2.41	0.55
1:C:281:LEU:HG	1:C:367:CYS:O	2.05	0.55
1:E:377:PRO:O	1:E:378:ARG:CB	2.53	0.55
2:B:322:THR:HA	2:D:255:ARG:NH1	2.21	0.55
1:C:361:ARG:HD2	1:C:366:VAL:HB	1.88	0.55
1:E:282:ARG:NH1	1:E:284:CYS:SG	2.79	0.55
1:A:287:ASN:ND2	1:A:300:TYR:H	2.05	0.55
1:A:302:LEU:HB3	2:B:304:HIS:ND1	2.22	0.55
1:A:335:ILE:HD11	2:B:307:PHE:HB2	1.88	0.55
2:B:304:HIS:CD2	2:B:305:ARG:HB2	2.42	0.55
1:C:278:THR:HG23	1:E:308:LYS:NZ	2.22	0.55
2:D:354:LYS:NZ	2:D:354:LYS:HB3	2.21	0.55
2:F:346:LEU:HD12	2:F:347:TYR:N	2.22	0.55
2:D:356:GLU:O	2:D:356:GLU:HG2	2.05	0.55
2:F:308:ALA:C	2:F:309:ILE:HG13	2.32	0.55
1:A:298:GLU:C	1:A:299:LEU:HD23	2.33	0.54
2:B:280:ILE:HG21	2:B:311:PHE:CZ	2.42	0.54
2:B:352:LYS:NZ	2:D:265:GLU:O	2.40	0.54
1:E:344:TYR:HB2	1:E:375:TYR:CD2	2.42	0.54
1:E:315:PHE:HA	1:E:356:ASN:O	2.07	0.54
2:B:333:ARG:O	2:B:337:LEU:HA	2.07	0.54
2:D:333:ARG:NH2	2:D:338:GLU:HG3	2.23	0.54
2:D:356:GLU:O	2:D:357:VAL:HG12	2.07	0.54
1:C:281:LEU:HD12	1:C:359:LEU:HB2	1.90	0.54
1:C:325:ASP:OD2	1:C:325:ASP:C	2.51	0.54
2:B:252:ARG:O	2:B:253:MET:HG3	2.08	0.54
2:B:280:ILE:HG13	2:B:311:PHE:CD2	2.43	0.53
2:B:286:GLU:HG2	2:B:291:VAL:HA	1.89	0.53
2:B:277:ASP:OD2	2:B:333:ARG:HG3	2.08	0.53
2:D:350:GLU:OE2	2:D:351:ILE:N	2.42	0.53
1:E:301:LEU:HD23	1:E:339:PHE:CE2	2.44	0.53
2:B:317:LYS:CG	2:B:318:ASP:H	2.12	0.53
1:E:344:TYR:OH	1:E:353:VAL:HG11	2.09	0.53
2:D:250:ILE:HD12	2:D:331:LEU:HG	1.90	0.53
1:E:378:ARG:HE	1:E:378:ARG:HA	1.73	0.53
2:F:250:ILE:HD11	2:F:268:LEU:HD11	1.90	0.53
2:D:268:LEU:HD23	2:D:268:LEU:C	2.34	0.53
2:F:314:PRO:O	2:F:315:LYS:C	2.52	0.53
1:E:308:LYS:HE3	1:E:328:GLN:OE1	2.09	0.53
1:C:279:SER:H	1:E:308:LYS:NZ	2.06	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:308:LYS:HG2	1:E:328:GLN:HG2	1.90	0.52
1:C:311:ILE:H	1:C:362:LEU:CD2	2.22	0.52
1:C:354:THR:HA	1:C:374:THR:HA	1.90	0.52
2:F:280:ILE:HD12	2:F:311:PHE:CG	2.45	0.52
2:B:274:GLN:H	2:B:274:GLN:NE2	2.06	0.52
2:F:332:ARG:HA	2:F:338:GLU:O	2.10	0.52
1:E:301:LEU:HB3	1:E:339:PHE:HE2	1.75	0.52
2:F:248:LEU:HB3	2:F:340:SER:HB2	1.91	0.52
2:F:336:ASP:OD2	2:F:338:GLU:HB2	2.10	0.52
2:D:346:LEU:C	2:D:346:LEU:HD23	2.34	0.52
2:F:308:ALA:O	2:F:309:ILE:HG13	2.09	0.52
2:B:247:ASN:HD22	2:B:333:ARG:NH2	2.08	0.52
1:C:350:SER:O	1:C:351:GLU:HG2	2.10	0.52
1:A:287:ASN:ND2	1:A:300:TYR:N	2.57	0.51
2:F:333:ARG:NH2	2:F:338:GLU:HG2	2.23	0.51
2:B:268:LEU:HD21	2:B:270:CYS:HB3	1.92	0.51
1:A:344:TYR:HB2	1:A:375:TYR:CD1	2.46	0.51
1:C:359:LEU:HD12	1:C:359:LEU:N	2.25	0.51
2:F:278:ILE:CD1	2:F:331:LEU:HB3	2.38	0.51
2:F:245:ALA:HB3	2:F:339:THR:O	2.10	0.51
1:A:302:LEU:CD2	2:B:267:TYR:CE2	2.94	0.50
2:B:274:GLN:C	2:B:276:ASP:H	2.19	0.50
1:E:317:THR:HG22	1:E:355:VAL:CG1	2.41	0.50
1:E:360:GLN:HG3	1:E:367:CYS:SG	2.50	0.50
1:A:364:ASP:OD2	1:A:366:VAL:HG22	2.10	0.50
1:A:361:ARG:NH1	1:A:366:VAL:CG2	2.74	0.50
2:B:274:GLN:N	2:B:274:GLN:CD	2.68	0.50
1:C:309:GLU:HG2	1:C:310:ASP:H	1.76	0.50
2:B:314:PRO:O	2:B:315:LYS:C	2.55	0.50
1:E:314:VAL:HB	1:E:358:PHE:CE1	2.46	0.50
2:B:356:GLU:C	2:B:358:GLN:H	2.19	0.50
1:C:281:LEU:O	1:C:368:SER:HB2	2.11	0.50
1:C:368:SER:O	1:C:369:GLU:C	2.51	0.50
2:B:250:ILE:HG21	2:B:253:MET:CE	2.36	0.50
1:C:306:VAL:H	1:C:334:GLN:HB3	1.77	0.50
2:B:247:ASN:ND2	2:B:333:ARG:HH22	2.10	0.49
1:E:304:ASP:O	1:E:305:LYS:C	2.55	0.49
2:D:336:ASP:OD2	2:D:338:GLU:HG2	2.11	0.49
1:E:281:LEU:O	1:E:368:SER:HB2	2.11	0.49
2:B:284:GLU:HB2	2:B:327:VAL:HG12	1.94	0.49
2:D:251:VAL:O	2:D:252:ARG:HG2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:355:GLU:O	2:F:356:GLU:HG2	2.13	0.49
2:B:357:VAL:O	2:B:358:GLN:CB	2.59	0.49
1:C:355:VAL:HG22	1:C:373:PHE:O	2.13	0.49
2:B:322:THR:HA	2:D:255:ARG:HH11	1.78	0.49
1:C:310:ASP:O	1:C:311:ILE:HB	2.13	0.49
2:F:285:GLU:OE2	2:F:291:VAL:HG22	2.12	0.49
2:B:350:GLU:O	2:B:351:ILE:C	2.55	0.49
1:A:297:GLU:O	1:A:340:LYS:HA	2.13	0.48
1:E:285:ARG:HG3	2:F:267:TYR:OH	2.14	0.48
1:E:346:ASP:O	1:E:349:ILE:HG23	2.14	0.48
2:F:274:GLN:C	2:F:276:ASP:H	2.21	0.48
2:B:285:GLU:O	2:B:286:GLU:C	2.56	0.48
1:C:320:TRP:CG	1:C:321:GLU:N	2.81	0.48
2:F:279:GLN:OE1	2:F:332:ARG:HD2	2.13	0.48
1:A:333:ARG:O	1:A:335:ILE:HG23	2.14	0.48
2:B:356:GLU:OE1	2:B:356:GLU:HA	2.12	0.48
1:C:314:VAL:HG13	1:C:322:GLY:O	2.14	0.48
2:B:253:MET:HE1	2:B:329:VAL:CG1	2.44	0.48
1:E:278:THR:HG23	1:E:279:SER:N	2.28	0.48
2:B:259:CYS:SG	2:B:261:THR:HB	2.53	0.48
2:F:304:HIS:O	2:F:306:GLN:N	2.46	0.48
1:A:368:SER:O	1:A:369:GLU:C	2.57	0.48
1:E:361:ARG:HD2	1:E:364:ASP:OD1	2.12	0.48
1:A:307:GLN:HB2	1:A:310:ASP:OD2	2.13	0.48
2:B:252:ARG:O	2:B:253:MET:CG	2.61	0.48
2:F:260:VAL:CG2	2:F:316:TYR:HB3	2.43	0.48
1:A:304:ASP:CB	2:B:305:ARG:HH12	2.26	0.48
2:B:320:ASN:HD21	1:C:292:PRO:HG3	1.79	0.48
2:B:304:HIS:CD2	2:B:304:HIS:C	2.92	0.47
2:B:305:ARG:O	2:B:306:GLN:HB2	2.13	0.47
1:C:335:ILE:HG12	2:D:307:PHE:HB2	1.95	0.47
2:B:295:PHE:N	2:B:295:PHE:CD1	2.81	0.47
1:C:361:ARG:HD2	1:C:364:ASP:OD2	2.14	0.47
2:D:351:ILE:HG22	2:D:354:LYS:H	1.78	0.47
1:E:314:VAL:HG22	1:E:323:ARG:HA	1.96	0.47
1:C:354:THR:OG1	1:C:374:THR:HB	2.14	0.47
1:C:285:ARG:NH1	2:D:310:VAL:HG21	2.29	0.47
1:C:285:ARG:NH2	2:D:310:VAL:HG21	2.28	0.47
2:D:279:GLN:HG2	2:D:332:ARG:O	2.15	0.47
1:E:306:VAL:H	1:E:334:GLN:HB3	1.79	0.47
1:C:346:ASP:C	1:C:348:GLU:H	2.23	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:353:ASP:O	2:D:354:LYS:C	2.58	0.47
1:E:368:SER:O	1:E:369:GLU:C	2.58	0.47
2:F:345:PHE:CD1	2:F:345:PHE:C	2.91	0.47
1:A:360:GLN:HA	1:A:366:VAL:O	2.15	0.47
2:B:275:LYS:O	2:B:300:PRO:HB3	2.14	0.47
2:D:307:PHE:O	2:D:308:ALA:HB2	2.15	0.47
1:E:346:ASP:C	1:E:348:GLU:H	2.23	0.47
2:F:261:THR:CG2	2:F:262:GLY:N	2.77	0.47
2:D:353:ASP:O	2:D:354:LYS:O	2.33	0.47
1:C:305:LYS:HB2	1:C:305:LYS:NZ	2.31	0.46
2:F:260:VAL:HG22	2:F:316:TYR:HB3	1.97	0.46
1:C:360:GLN:HA	1:C:366:VAL:O	2.14	0.46
2:D:259:CYS:HA	2:D:348:TYR:CE2	2.50	0.46
2:D:331:LEU:HB2	2:D:340:SER:CB	2.45	0.46
1:E:310:ASP:OD2	1:E:361:ARG:HD3	2.16	0.46
2:D:341:GLU:HA	2:D:341:GLU:OE1	2.15	0.46
2:F:317:LYS:CG	2:F:318:ASP:N	2.59	0.46
1:A:377:PRO:O	1:A:378:ARG:C	2.58	0.46
2:B:250:ILE:CG1	2:B:331:LEU:HG	2.44	0.46
2:D:332:ARG:HA	2:D:338:GLU:O	2.16	0.46
2:B:248:LEU:HB3	2:B:340:SER:HB2	1.97	0.46
2:B:331:LEU:H	2:B:340:SER:HB3	1.80	0.46
2:D:327:VAL:HG22	2:D:345:PHE:O	2.15	0.46
2:D:355:GLU:O	2:D:355:GLU:HG3	2.16	0.46
1:E:314:VAL:HG22	1:E:323:ARG:HG3	1.97	0.46
1:A:373:PHE:HE1	1:A:375:TYR:CD2	2.33	0.46
2:B:253:MET:HE1	2:B:329:VAL:HG11	1.98	0.46
1:C:300:TYR:CD2	1:C:338:VAL:HG22	2.51	0.46
2:B:292:TRP:CD1	2:B:317:LYS:HB3	2.50	0.46
1:C:364:ASP:OD2	1:C:366:VAL:HG23	2.14	0.46
1:E:297:GLU:O	1:E:340:LYS:HA	2.16	0.46
1:A:279:SER:O	1:A:281:LEU:HD13	2.15	0.46
2:F:286:GLU:O	2:F:287:GLU:C	2.58	0.46
2:B:277:ASP:CG	2:B:333:ARG:HG3	2.40	0.46
1:C:304:ASP:O	1:C:305:LYS:C	2.59	0.46
1:A:311:ILE:H	1:A:362:LEU:CD1	2.29	0.45
2:B:254:ASP:OD2	2:B:267:TYR:N	2.45	0.45
1:C:349:ILE:HD12	1:C:351:GLU:O	2.17	0.45
1:A:342:PRO:HA	1:A:343:PRO:HD3	1.88	0.45
1:A:308:LYS:O	1:A:328:GLN:HG2	2.17	0.45
2:B:351:ILE:O	2:B:352:LYS:C	2.58	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:361:ARG:CD	1:C:366:VAL:HB	2.46	0.45
2:D:345:PHE:CD1	2:D:345:PHE:C	2.95	0.45
1:E:280:GLU:C	1:E:281:LEU:HD12	2.42	0.45
1:A:301:LEU:HD23	1:A:301:LEU:C	2.41	0.45
2:B:251:VAL:HB	2:B:269:LEU:O	2.16	0.45
1:C:301:LEU:HD12	1:C:302:LEU:N	2.31	0.45
1:E:293:CYS:HB3	1:E:376:LEU:O	2.17	0.45
1:E:299:LEU:HD12	1:E:339:PHE:HE1	1.81	0.45
1:A:300:TYR:CE2	2:B:254:ASP:HA	2.51	0.45
1:E:347:LEU:H	1:E:347:LEU:CD1	2.02	0.45
1:E:371:LEU:HD22	1:E:372:PRO:HD2	1.98	0.45
2:D:305:ARG:NH2	2:D:305:ARG:HG3	2.31	0.45
1:A:356:ASN:HB3	1:A:370:PRO:HB3	1.99	0.45
2:D:286:GLU:C	2:D:288:ASN:H	2.24	0.45
2:F:281:ARG:HG3	2:F:295:PHE:CE2	2.52	0.45
2:B:259:CYS:SG	2:B:260:VAL:N	2.90	0.44
2:B:354:LYS:C	2:B:356:GLU:N	2.75	0.44
1:C:314:VAL:HB	1:C:358:PHE:CZ	2.52	0.44
1:A:287:ASN:HD21	1:A:300:TYR:N	2.14	0.44
1:A:371:LEU:HD12	1:A:372:PRO:HD2	1.99	0.44
2:D:260:VAL:HG13	2:D:319:VAL:CG2	2.42	0.44
1:E:312:SER:HB2	1:E:362:LEU:HD21	1.99	0.44
1:A:291:GLY:HA3	1:A:297:GLU:OE2	2.17	0.44
1:C:301:LEU:O	1:C:336:ALA:HA	2.18	0.44
1:C:332:HIS:HB3	1:C:336:ALA:HB3	2.00	0.44
1:A:335:ILE:HB	2:B:304:HIS:CD2	2.51	0.44
2:B:260:VAL:HG22	2:B:316:TYR:CB	2.40	0.44
1:E:310:ASP:O	1:E:311:ILE:HB	2.18	0.44
1:A:301:LEU:HD22	1:A:337:ILE:HB	1.99	0.44
1:A:300:TYR:CE2	2:B:254:ASP:HB3	2.53	0.44
1:A:302:LEU:HD21	2:B:267:TYR:CE2	2.53	0.44
1:A:304:ASP:O	1:A:305:LYS:C	2.60	0.44
2:D:275:LYS:NZ	2:D:275:LYS:HB3	2.33	0.44
2:D:286:GLU:HB2	2:D:290:GLY:O	2.17	0.44
2:D:351:ILE:HD13	2:D:351:ILE:HA	1.79	0.44
1:A:305:LYS:HD3	1:A:334:GLN:HG3	1.99	0.43
2:B:254:ASP:O	2:B:255:ARG:CD	2.63	0.43
1:C:306:VAL:CG1	1:C:311:ILE:HG12	2.30	0.43
2:D:252:ARG:HB2	2:D:269:LEU:HB2	1.99	0.43
2:D:323:LYS:HB3	2:D:324:PRO:HD2	1.99	0.43
2:F:297:ASP:O	2:F:298:PHE:HB3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:282:PHE:O	2:B:293:GLU:HA	2.18	0.43
1:E:282:ARG:CZ	1:E:284:CYS:SG	3.06	0.43
2:F:303:VAL:HG13	2:F:309:ILE:CG1	2.47	0.43
1:A:302:LEU:HB3	2:B:304:HIS:CE1	2.53	0.43
2:B:268:LEU:CD2	2:B:270:CYS:HB3	2.48	0.43
2:B:295:PHE:O	2:B:314:PRO:HB3	2.18	0.43
1:C:315:PHE:HA	1:C:356:ASN:O	2.18	0.43
1:A:298:GLU:O	1:A:299:LEU:HD23	2.18	0.43
2:D:331:LEU:HB2	2:D:340:SER:HB3	1.98	0.43
1:A:301:LEU:O	1:A:336:ALA:HA	2.17	0.43
1:A:346:ASP:C	1:A:348:GLU:H	2.26	0.43
2:B:254:ASP:OD2	2:B:254:ASP:N	2.51	0.43
2:D:346:LEU:HD23	2:D:347:TYR:C	2.44	0.43
1:A:300:TYR:CD2	2:B:254:ASP:HB3	2.54	0.43
1:A:315:PHE:HA	1:A:356:ASN:O	2.19	0.43
1:E:306:VAL:O	1:E:334:GLN:HG2	2.19	0.43
2:B:315:LYS:HG2	2:B:316:TYR:O	2.18	0.43
2:B:333:ARG:HG2	2:B:335:SER:OG	2.18	0.43
1:E:302:LEU:HG	2:F:304:HIS:CD2	2.54	0.43
1:E:306:VAL:HG22	1:E:334:GLN:O	2.19	0.43
2:F:282:PHE:O	2:F:293:GLU:HA	2.19	0.43
2:F:248:LEU:CB	2:F:340:SER:HB2	2.48	0.43
2:F:345:PHE:CD1	2:F:346:LEU:N	2.87	0.43
1:C:325:ASP:OD2	1:C:326:PHE:N	2.52	0.42
2:F:278:ILE:HD12	2:F:279:GLN:N	2.27	0.42
1:A:355:VAL:HG22	1:A:373:PHE:O	2.20	0.42
2:B:304:HIS:O	2:B:306:GLN:N	2.52	0.42
1:C:310:ASP:OD2	1:C:362:LEU:N	2.52	0.42
1:E:301:LEU:HD23	1:E:339:PHE:HE2	1.81	0.42
1:A:279:SER:O	1:A:280:GLU:C	2.62	0.42
1:C:311:ILE:HD12	1:C:312:SER:H	1.85	0.42
1:E:332:HIS:HB2	2:F:269:LEU:HD13	2.01	0.42
2:D:271:ASP:O	2:D:272:LYS:C	2.62	0.42
1:E:301:LEU:O	1:E:336:ALA:HA	2.19	0.42
2:B:281:ARG:HD2	2:B:295:PHE:CE2	2.55	0.42
2:D:274:GLN:HB2	2:D:277:ASP:HB3	2.01	0.42
1:E:349:ILE:CD1	1:E:353:VAL:HG23	2.50	0.42
2:F:268:LEU:HD23	2:F:268:LEU:C	2.45	0.42
2:F:292:TRP:CG	2:F:293:GLU:N	2.87	0.42
2:B:287:GLU:C	2:B:288:ASN:CG	2.86	0.42
1:C:335:ILE:HD12	1:C:335:ILE:C	2.45	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:250:ILE:HG23	2:B:268:LEU:HD11	2.01	0.41
1:C:283:ILE:HB	1:C:371:LEU:CD2	2.51	0.41
2:B:274:GLN:C	2:B:276:ASP:N	2.78	0.41
1:A:311:ILE:CD1	1:A:359:LEU:HD22	2.36	0.41
1:A:320:TRP:CZ2	1:A:322:GLY:HA3	2.56	0.41
2:B:260:VAL:HG13	2:B:319:VAL:HA	2.02	0.41
2:D:356:GLU:C	2:D:358:GLN:N	2.78	0.41
2:F:253:MET:HB3	2:F:255:ARG:O	2.21	0.41
2:F:274:GLN:C	2:F:276:ASP:N	2.79	0.41
1:C:285:ARG:NH2	2:D:310:VAL:CG2	2.83	0.41
1:C:293:CYS:HB3	1:C:376:LEU:O	2.20	0.41
2:B:246:SER:OG	2:B:247:ASN:N	2.53	0.41
2:B:277:ASP:OD2	2:B:333:ARG:CG	2.68	0.41
2:B:279:GLN:HG3	2:B:332:ARG:HB3	2.01	0.41
1:C:285:ARG:HD3	1:C:302:LEU:HD12	2.03	0.41
2:D:292:TRP:CG	2:D:293:GLU:N	2.88	0.41
1:E:344:TYR:CE2	1:E:346:ASP:HB3	2.56	0.41
1:E:314:VAL:HG13	1:E:322:GLY:O	2.21	0.41
1:A:344:TYR:HB2	1:A:375:TYR:CE1	2.56	0.41
2:B:335:SER:OG	2:B:336:ASP:N	2.54	0.41
1:A:335:ILE:HD11	2:B:307:PHE:CB	2.51	0.41
1:E:354:THR:HB	1:E:374:THR:HB	2.03	0.41
1:C:300:TYR:CE2	1:C:338:VAL:HG22	2.56	0.40
2:D:355:GLU:C	2:D:357:VAL:N	2.78	0.40
1:E:325:ASP:OD2	1:E:326:PHE:N	2.54	0.40
1:C:333:ARG:HG3	2:D:307:PHE:CD2	2.56	0.40
2:F:271:ASP:HB3	2:F:272:LYS:H	1.76	0.40
2:F:334:LYS:HD2	2:F:334:LYS:HA	1.94	0.40
1:C:375:TYR:O	1:C:376:LEU:HD12	2.22	0.40
2:F:251:VAL:HB	2:F:269:LEU:O	2.21	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	99/101 (98%)	79 (80%)	11 (11%)	9 (9%)	0	2
1	C	99/101 (98%)	84 (85%)	10 (10%)	5 (5%)	1	10
1	E	99/101 (98%)	81 (82%)	15 (15%)	3 (3%)	3	19
2	B	113/115 (98%)	82 (73%)	22 (20%)	9 (8%)	1	4
2	D	113/115 (98%)	92 (81%)	11 (10%)	10 (9%)	0	3
2	F	110/115 (96%)	91 (83%)	12 (11%)	7 (6%)	1	6
All	All	633/648 (98%)	509 (80%)	81 (13%)	43 (7%)	1	6

All (43) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	279	SER
1	A	377	PRO
2	B	275	LYS
2	B	288	ASN
2	B	305	ARG
2	B	355	GLU
2	B	358	GLN
2	D	354	LYS
2	D	356	GLU
2	F	275	LYS
2	F	288	ASN
2	F	305	ARG
1	A	280	GLU
1	C	311	ILE
2	D	272	LYS
2	D	288	ASN
2	D	317	LYS
2	D	335	SER
2	D	352	LYS
2	D	355	GLU
1	E	308	LYS
2	F	287	GLU
2	F	317	LYS
2	F	355	GLU
1	A	305	LYS
1	C	305	LYS
1	E	305	LYS

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Mol	Chain	Res	Type
1	A	334	GLN
1	A	360	GLN
1	A	363	THR
2	B	317	LYS
1	C	347	LEU
2	D	287	GLU
2	D	289	GLY
1	E	311	ILE
2	B	353	ASP
1	A	347	LEU
2	B	351	ILE
1	C	360	GLN
2	F	290	GLY
1	A	311	ILE
1	C	349	ILE
2	B	290	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	92/92 (100%)	83 (90%)	9 (10%)	7	28
1	C	92/92 (100%)	81 (88%)	11 (12%)	5	20
1	E	92/92 (100%)	85 (92%)	7 (8%)	12	37
2	B	102/104 (98%)	89 (87%)	13 (13%)	4	18
2	D	104/104 (100%)	88 (85%)	16 (15%)	2	12
2	F	100/104 (96%)	91 (91%)	9 (9%)	9	30
All	All	582/588 (99%)	517 (89%)	65 (11%)	6	22

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	281	LEU
1	A	301	LEU

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Mol	Chain	Res	Type
1	A	309	GLU
1	A	321	GLU
1	A	323	ARG
1	A	334	GLN
1	A	348	GLU
1	A	371	LEU
1	A	374	THR
2	B	247	ASN
2	B	256	THR
2	B	260	VAL
2	B	261	THR
2	B	274	GLN
2	B	275	LYS
2	B	284	GLU
2	B	301	THR
2	B	313	THR
2	B	315	LYS
2	B	337	LEU
2	B	351	ILE
2	B	356	GLU
1	C	280	GLU
1	C	281	LEU
1	C	283	ILE
1	C	299	LEU
1	C	301	LEU
1	C	305	LYS
1	C	306	VAL
1	C	323	ARG
1	C	333	ARG
1	C	342	PRO
1	C	374	THR
2	D	250	ILE
2	D	260	VAL
2	D	275	LYS
2	D	286	GLU
2	D	287	GLU
2	D	301	THR
2	D	303	VAL
2	D	321	ILE
2	D	329	VAL
2	D	337	LEU
2	D	338	GLU

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Mol	Chain	Res	Type
2	D	340	SER
2	D	351	ILE
2	D	354	LYS
2	D	358	GLN
2	D	359	ARG
1	E	280	GLU
1	E	301	LEU
1	E	309	GLU
1	E	347	LEU
1	E	348	GLU
1	E	354	THR
1	E	364	ASP
2	F	250	ILE
2	F	261	THR
2	F	313	THR
2	F	319	VAL
2	F	321	ILE
2	F	330	GLN
2	F	349	PRO
2	F	352	LYS
2	F	355	GLU

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	334	GLN
1	A	360	GLN
2	B	247	ASN
2	B	274	GLN
2	B	306	GLN
2	B	320	ASN
1	C	328	GLN
2	D	247	ASN
2	D	306	GLN
1	E	307	GLN
1	E	334	GLN
1	E	356	ASN
1	E	360	GLN
2	F	304	HIS
2	F	306	GLN

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	101/101 (100%)	-0.20	1 (0%) 79 62	66, 102, 124, 140	0
1	C	101/101 (100%)	-0.47	0 100 100	56, 83, 111, 131	0
1	E	101/101 (100%)	-0.36	0 100 100	79, 115, 138, 140	0
2	B	115/115 (100%)	-0.41	0 100 100	49, 70, 115, 136	0
2	D	115/115 (100%)	-0.59	0 100 100	51, 71, 116, 140	0
2	F	112/115 (97%)	-0.46	0 100 100	57, 88, 115, 135	0
All	All	645/648 (99%)	-0.42	1 (0%) 91 84	49, 87, 131, 140	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	313	VAL	2.2

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.