



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:03 AM UTC

PDB ID : 1D00 / pdb_00001d00
Title : STRUCTURE OF TNF RECEPTOR ASSOCIATED FACTOR 2 IN COM-
PLEX WITH A 5-RESIDUE CD40 PEPTIDE
Authors : Ye, H.; Park, Y.C.; Kreishman, M.; Kieff, E.; Wu, H.
Deposited on : 1999-09-07
Resolution : 2.00 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

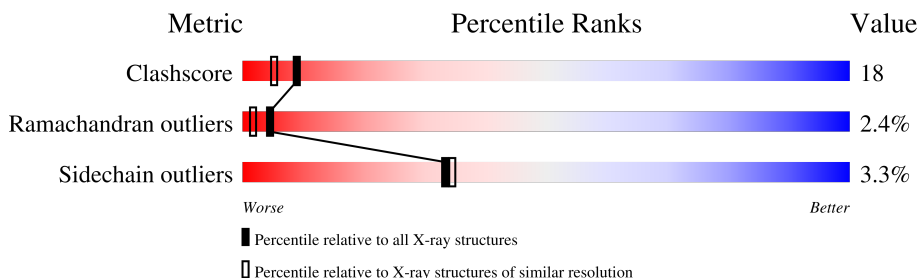
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	168	69% 26% 5%
1	B	168	64% 30% 6%
1	C	168	66% 29% 5%
1	D	168	65% 30% . .
1	E	168	70% 25% 5% .
1	F	168	67% 29% .
1	G	168	66% 30% .
1	H	168	64% 32% 5%

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Mol	Chain	Length	Quality of chain
2	I	7	57% 29% 14%
2	J	7	29% 57% 14%
2	K	7	43% 43% 14%
2	L	7	57% 43%
2	M	7	57% 29% 14%
2	N	7	57% 43%
2	O	7	57% 29% 14%
2	P	7	43% 43% 14%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10670 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 TUMOR NECROSIS FACTOR RECEPTOR ASSOCIATED PROTEIN 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	168	Total	C	N	O	S	0	0	0
			1282	825	217	231	9			
1	B	168	Total	C	N	O	S	0	0	0
			1282	825	217	231	9			
1	C	168	Total	C	N	O	S	0	0	0
			1282	825	217	231	9			
1	D	168	Total	C	N	O	S	0	0	0
			1282	825	217	231	9			
1	E	168	Total	C	N	O	S	0	0	0
			1282	825	217	231	9			
1	F	168	Total	C	N	O	S	0	0	0
			1282	825	217	231	9			
1	G	168	Total	C	N	O	S	0	0	0
			1282	825	217	231	9			
1	H	168	Total	C	N	O	S	0	0	0
			1282	825	217	231	9			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	362	ALA	PRO	conflict	GB 1363002
A	365	ARG	LEU	conflict	GB 1363002
B	362	ALA	PRO	conflict	GB 1363002
B	365	ARG	LEU	conflict	GB 1363002
C	362	ALA	PRO	conflict	GB 1363002
C	365	ARG	LEU	conflict	GB 1363002
D	362	ALA	PRO	conflict	GB 1363002
D	365	ARG	LEU	conflict	GB 1363002
E	362	ALA	PRO	conflict	GB 1363002
E	365	ARG	LEU	conflict	GB 1363002
F	362	ALA	PRO	conflict	GB 1363002
F	365	ARG	LEU	conflict	GB 1363002

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Chain	Residue	Modelled	Actual	Comment	Reference
G	362	ALA	PRO	conflict	GB 1363002
G	365	ARG	LEU	conflict	GB 1363002
H	362	ALA	PRO	conflict	GB 1363002
H	365	ARG	LEU	conflict	GB 1363002

- Molecule 2 is a protein called B-CELL SURFACE ANTIGEN CD40.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	I	7	Total	C	N	O	0	0	1
			43	26	7	10			
2	J	7	Total	C	N	O	0	0	1
			43	26	7	10			
2	K	7	Total	C	N	O	0	0	1
			43	26	7	10			
2	L	7	Total	C	N	O	0	0	1
			43	26	7	10			
2	M	7	Total	C	N	O	0	0	1
			43	26	7	10			
2	N	7	Total	C	N	O	0	0	1
			43	26	7	10			
2	O	7	Total	C	N	O	0	0	1
			43	26	7	10			
2	P	7	Total	C	N	O	0	0	1
			43	26	7	10			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	12	Total	O	0	0
			12	12		
3	B	3	Total	O	0	0
			3	3		
3	C	11	Total	O	0	0
			11	11		
3	D	7	Total	O	0	0
			7	7		
3	E	5	Total	O	0	0
			5	5		
3	F	9	Total	O	0	0
			9	9		
3	G	8	Total	O	0	0
			8	8		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	13	Total 13	O 13	0	0
3	K	1	Total 1	O 1	0	0
3	L	1	Total 1	O 1	0	0

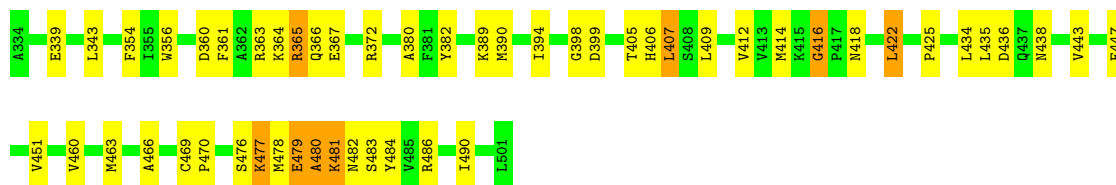
3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

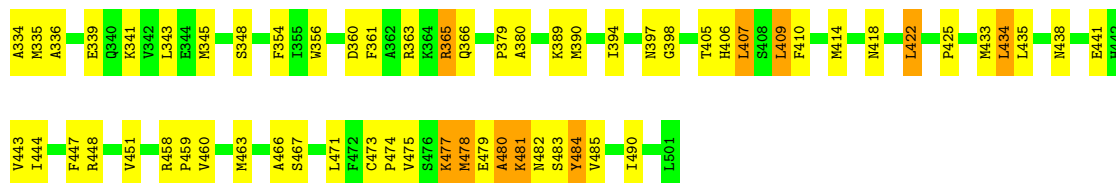
• Molecule 1: TUMOR NECROSIS FACTOR RECEPTOR ASSOCIATED PROTEIN 2

Chain A: 



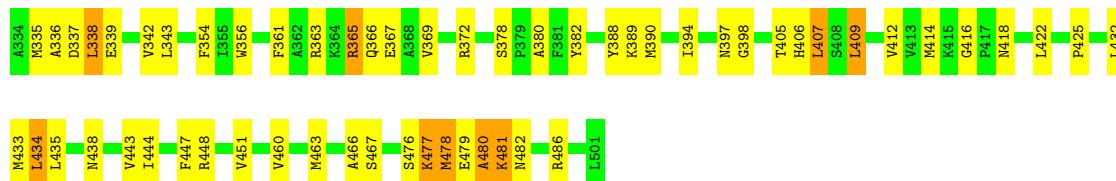
• Molecule 1: TUMOR NECROSIS FACTOR RECEPTOR ASSOCIATED PROTEIN 2

Chain B: 



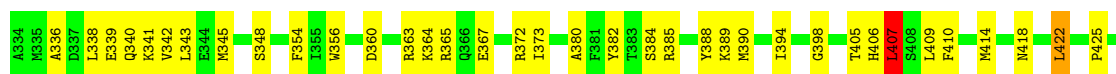
• Molecule 1: TUMOR NECROSIS FACTOR RECEPTOR ASSOCIATED PROTEIN 2

Chain C: 



• Molecule 1: TUMOR NECROSIS FACTOR RECEPTOR ASSOCIATED PROTEIN 2

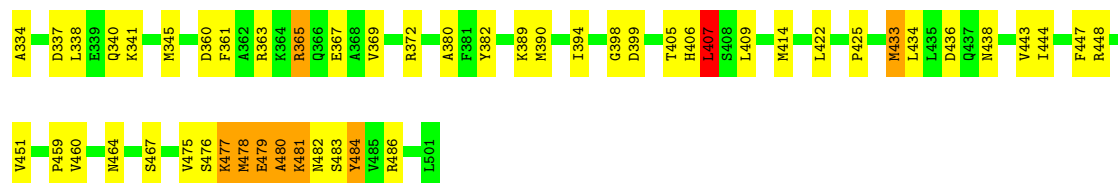
Chain D: 





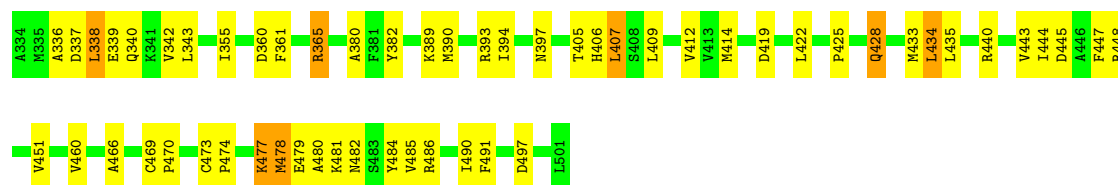
- Molecule 1: TUMOR NECROSIS FACTOR RECEPTOR ASSOCIATED PROTEIN 2

Chain E: 70% 25% 5% •



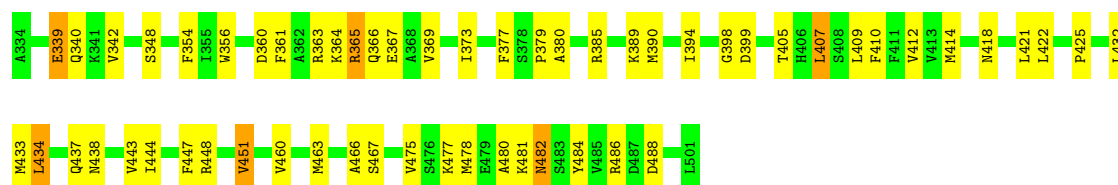
- Molecule 1: TUMOR NECROSIS FACTOR RECEPTOR ASSOCIATED PROTEIN 2

Chain F: 67% 29% •



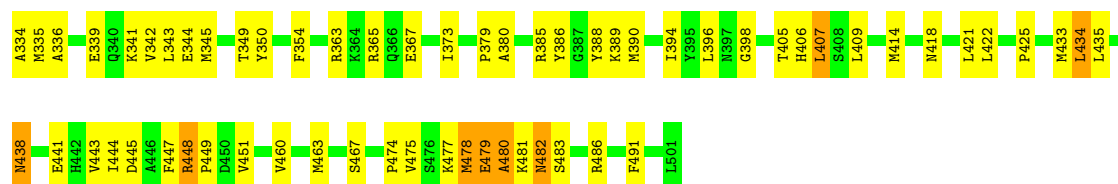
- Molecule 1: TUMOR NECROSIS FACTOR RECEPTOR ASSOCIATED PROTEIN 2

Chain G: 66% 30% •



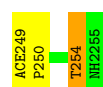
- Molecule 1: TUMOR NECROSIS FACTOR RECEPTOR ASSOCIATED PROTEIN 2

Chain H: 64% 32% 5% •

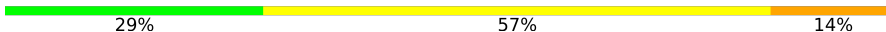


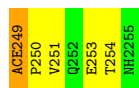
- Molecule 2: B-CELL SURFACE ANTIGEN CD40

Chain I: 57% 29% 14%

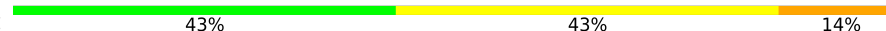


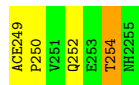
- Molecule 2: B-CELL SURFACE ANTIGEN CD40

Chain J:  29% 57% 14%



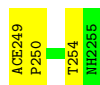
- Molecule 2: B-CELL SURFACE ANTIGEN CD40

Chain K:  43% 43% 14%



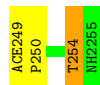
- Molecule 2: B-CELL SURFACE ANTIGEN CD40

Chain L:  57% 43%



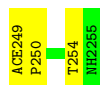
- Molecule 2: B-CELL SURFACE ANTIGEN CD40

Chain M:  57% 29% 14%



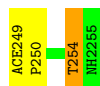
- Molecule 2: B-CELL SURFACE ANTIGEN CD40

Chain N:  57% 43%

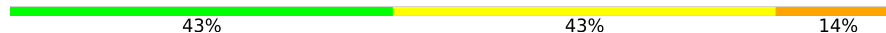


- Molecule 2: B-CELL SURFACE ANTIGEN CD40

Chain O:  57% 29% 14%



- Molecule 2: B-CELL SURFACE ANTIGEN CD40

Chain P:  43% 43% 14%



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	R 3	Depositor
Cell constants a, b, c, α , β , γ	111.40Å 111.40Å 111.40Å 103.70° 103.70° 103.70°	Depositor
Resolution (Å)	20.00 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.00)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.219 , 0.254	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	10670	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ACE, NH2

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	0.50	0/1312	1.08	7/1781 (0.4%)
1	B	0.49	0/1312	1.07	9/1781 (0.5%)
1	C	0.48	0/1312	1.11	13/1781 (0.7%)
1	D	0.48	0/1312	1.08	10/1781 (0.6%)
1	E	0.51	0/1312	1.06	6/1781 (0.3%)
1	F	0.48	0/1312	1.06	8/1781 (0.4%)
1	G	0.49	0/1312	1.04	10/1781 (0.6%)
1	H	0.47	0/1312	1.04	9/1781 (0.5%)
2	I	1.49	1/40 (2.5%)	0.86	0/55
2	J	1.47	1/40 (2.5%)	0.87	0/55
2	K	1.51	1/40 (2.5%)	0.88	0/55
2	L	1.44	1/40 (2.5%)	1.04	0/55
2	M	1.54	1/40 (2.5%)	0.87	0/55
2	N	1.50	1/40 (2.5%)	0.87	0/55
2	O	1.45	1/40 (2.5%)	0.90	0/55
2	P	1.49	1/40 (2.5%)	0.91	0/55
All	All	0.55	8/10816 (0.1%)	1.06	72/14688 (0.5%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	M	249	ACE	C-N	-9.34	1.34	1.43
2	K	249	ACE	C-N	-9.06	1.34	1.43
2	N	249	ACE	C-N	-9.05	1.34	1.43
2	P	249	ACE	C-N	-8.98	1.34	1.43
2	I	249	ACE	C-N	-8.93	1.34	1.43
2	L	249	ACE	C-N	-8.80	1.34	1.43
2	O	249	ACE	C-N	-8.71	1.34	1.43
2	J	249	ACE	C-N	-8.71	1.34	1.43

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	335	MET	N-CA-C	-10.20	100.96	113.50
1	A	398	GLY	N-CA-C	8.29	122.51	112.48
1	F	460	VAL	N-CA-C	-8.27	104.69	112.96
1	B	335	MET	N-CA-C	-8.14	103.10	112.87
1	B	460	VAL	N-CA-C	-8.12	104.84	112.96
1	A	460	VAL	N-CA-C	-7.92	105.04	112.96
1	C	460	VAL	N-CA-C	-7.69	105.27	112.96
1	E	460	VAL	N-CA-C	-7.67	105.29	112.96
1	D	336	ALA	N-CA-C	-7.66	103.66	113.16
1	E	407	LEU	N-CA-C	-7.55	96.96	108.96
1	F	407	LEU	N-CA-C	-7.45	97.09	109.46
1	B	398	GLY	N-CA-C	7.40	122.27	112.65
1	E	398	GLY	N-CA-C	7.31	121.89	112.68
1	A	438	ASN	N-CA-C	-7.23	104.49	113.38
1	C	416	GLY	CA-C-N	7.17	126.80	119.56
1	C	416	GLY	C-N-CA	7.17	126.80	119.56
1	D	373	ILE	CA-C-N	7.14	126.97	119.05
1	D	373	ILE	C-N-CA	7.14	126.97	119.05
1	H	460	VAL	N-CA-C	-7.07	105.89	112.96
1	C	407	LEU	N-CA-C	-6.77	98.22	109.46
1	A	407	LEU	N-CA-C	-6.75	98.24	109.24
1	G	460	VAL	N-CA-C	-6.68	106.28	112.96
1	D	460	VAL	N-CA-C	-6.65	106.31	112.96
1	G	398	GLY	N-CA-C	6.63	121.03	112.68
1	D	407	LEU	N-CA-C	-6.58	98.53	109.46
1	H	407	LEU	N-CA-C	-6.53	98.09	108.73
1	G	407	LEU	N-CA-C	-6.51	98.61	108.96
1	B	435	LEU	N-CA-C	6.50	119.20	109.25
1	B	407	LEU	N-CA-C	-6.50	98.63	108.96
1	G	448	ARG	N-CA-C	-6.40	100.33	109.62
1	B	348	SER	N-CA-C	6.37	119.01	110.35
1	C	438	ASN	N-CA-C	-6.35	105.56	113.38
1	C	398	GLY	N-CA-C	6.32	120.86	112.65
1	H	398	GLY	N-CA-C	6.31	120.63	112.68
1	C	338	LEU	N-CA-C	-6.25	106.19	113.88
1	F	435	LEU	N-CA-C	6.07	119.42	109.76
1	F	337	ASP	N-CA-C	-6.07	106.23	113.21
1	C	435	LEU	N-CA-C	5.94	119.20	109.76
1	E	484	TYR	N-CA-C	-5.79	106.05	113.23
1	D	448	ARG	N-CA-C	-5.72	102.41	110.31
1	E	448	ARG	N-CA-C	-5.65	101.43	109.62
1	D	398	GLY	N-CA-C	5.58	119.71	112.68
1	A	416	GLY	CA-C-N	5.54	125.15	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	416	GLY	C-N-CA	5.54	125.15	119.56
1	D	348	SER	N-CA-C	5.52	117.42	110.24
1	C	448	ARG	N-CA-C	-5.46	102.77	110.31
1	D	444	ILE	N-CA-C	5.39	115.32	107.88
1	B	448	ARG	N-CA-C	-5.36	101.27	109.64
1	H	373	ILE	CA-C-N	5.32	124.95	119.05
1	H	373	ILE	C-N-CA	5.32	124.95	119.05
1	F	419	ASP	N-CA-C	5.26	118.48	111.75
1	H	434	LEU	N-CA-C	-5.24	98.85	108.02
1	G	373	ILE	CA-C-N	5.19	124.81	119.05
1	G	373	ILE	C-N-CA	5.19	124.81	119.05
1	C	361	PHE	N-CA-C	5.18	117.34	111.02
1	E	433	MET	N-CA-C	5.17	117.86	109.06
1	F	382	TYR	N-CA-C	5.17	117.54	109.52
1	C	434	LEU	N-CA-C	-5.17	98.83	107.99
1	A	435	LEU	N-CA-C	5.17	117.98	109.76
1	F	338	LEU	N-CA-C	-5.16	106.32	113.18
1	G	438	ASN	N-CA-C	-5.15	107.00	113.28
1	B	484	TYR	N-CA-C	-5.13	105.77	112.23
1	G	339	GLU	N-CA-C	-5.12	105.61	111.14
1	G	348	SER	N-CA-C	5.11	117.85	110.28
1	H	448	ARG	N-CA-C	-5.11	101.67	109.64
1	C	378	SER	N-CA-C	-5.09	104.20	110.31
1	F	434	LEU	N-CA-C	-5.04	99.20	108.02
1	G	434	LEU	N-CA-C	-5.04	99.20	108.02
1	D	438	ASN	N-CA-C	-5.03	107.14	113.28
1	B	434	LEU	N-CA-C	-5.02	99.10	107.99
1	H	438	ASN	N-CA-C	-5.01	106.75	112.92
1	H	435	LEU	N-CA-C	5.00	117.71	109.76

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	1282	0	1242	41	0
1	B	1282	0	1242	52	0
1	C	1282	0	1242	44	0
1	D	1282	0	1242	51	0
1	E	1282	0	1242	46	0
1	F	1282	0	1242	47	0
1	G	1282	0	1242	52	0
1	H	1282	0	1242	57	0
2	I	43	0	40	3	0
2	J	43	0	40	8	0
2	K	43	0	40	6	0
2	L	43	0	40	1	0
2	M	43	0	40	3	0
2	N	43	0	40	1	0
2	O	43	0	40	4	0
2	P	43	0	40	9	0
3	A	12	0	0	1	0
3	B	3	0	0	0	0
3	C	11	0	0	0	0
3	D	7	0	0	3	0
3	E	5	0	0	1	0
3	F	9	0	0	4	0
3	G	8	0	0	0	0
3	H	13	0	0	5	0
3	K	1	0	0	0	0
3	L	1	0	0	0	0
All	All	10670	0	10256	385	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 18.

All (385) 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:G:389:LYS:HG2	1:G:414:MET:HE3	1.29	1.09
1:A:389:LYS:HG2	1:A:414:MET:HE3	1.36	1.06
1:F:389:LYS:HG2	1:F:414:MET:HE3	1.37	1.06
1:D:389:LYS:HG2	1:D:414:MET:HE3	1.32	1.06
1:C:389:LYS:HG2	1:C:414:MET:HE3	1.40	1.02
1:H:425:PRO:HG3	1:H:451:VAL:HG13	1.41	1.00
1:A:425:PRO:HG3	1:A:451:VAL:HG13	1.43	0.99
1:H:394:ILE:HD11	1:H:407:LEU:HD21	1.41	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:389:LYS:HG2	1:H:414:MET:HE3	1.45	0.97
1:E:389:LYS:HG2	1:E:414:MET:HE3	1.47	0.95
1:E:425:PRO:HG3	1:E:451:VAL:HG13	1.51	0.93
3:F:1052:HOH:O	1:H:421:LEU:HD21	1.69	0.92
1:A:447:PHE:CD1	2:I:250:PRO:HG3	2.09	0.87
1:F:339:GLU:O	1:F:343:LEU:HD13	1.73	0.87
1:F:394:ILE:HD11	1:F:407:LEU:HD21	1.56	0.86
1:B:389:LYS:HG2	1:B:414:MET:HE3	1.57	0.86
1:C:447:PHE:CD1	2:K:250:PRO:HG3	2.10	0.86
1:D:425:PRO:HB2	3:D:1021:HOH:O	1.77	0.85
1:D:409:LEU:HD22	1:D:432:LEU:HD13	1.57	0.84
1:F:428:GLN:HG3	1:F:497:ASP:O	1.78	0.84
1:H:447:PHE:CD1	2:P:250:PRO:HG3	2.11	0.84
1:D:390:MET:C	1:D:414:MET:HE2	2.04	0.83
1:H:380:ALA:CB	1:H:414:MET:HE1	2.10	0.82
1:F:433:MET:HE2	1:F:444:ILE:HD12	1.61	0.82
1:E:380:ALA:CB	1:E:414:MET:HE1	2.10	0.82
1:D:447:PHE:CD1	2:L:250:PRO:HG3	2.14	0.82
1:A:425:PRO:CG	1:A:451:VAL:HG13	2.10	0.81
1:B:425:PRO:HG3	1:B:451:VAL:HG13	1.60	0.81
1:H:451:VAL:HA	3:H:1032:HOH:O	1.78	0.81
1:E:367:GLU:HG2	1:E:372:ARG:NH2	1.96	0.81
1:G:389:LYS:HG2	1:G:414:MET:CE	2.11	0.80
1:E:394:ILE:HD11	1:E:407:LEU:HD21	1.64	0.80
1:B:380:ALA:CB	1:B:414:MET:HE1	2.11	0.80
1:F:447:PHE:CD1	2:N:250:PRO:HG3	2.17	0.80
2:J:254:THR:HG22	2:J:254:THR:O	1.81	0.80
1:A:380:ALA:CB	1:A:414:MET:HE1	2.12	0.80
1:E:447:PHE:CD1	2:M:250:PRO:HG3	2.18	0.79
1:C:434:LEU:HB3	1:C:443:VAL:HB	1.63	0.78
1:G:380:ALA:CB	1:G:414:MET:HE1	2.13	0.78
1:G:390:MET:C	1:G:414:MET:HE2	2.10	0.76
1:A:394:ILE:HD11	1:A:407:LEU:HD21	1.66	0.76
1:C:339:GLU:O	1:C:343:LEU:HD13	1.86	0.75
1:C:390:MET:C	1:C:414:MET:HE2	2.12	0.75
1:D:380:ALA:CB	1:D:414:MET:HE1	2.17	0.74
1:G:380:ALA:HB1	1:G:414:MET:HE1	1.68	0.74
1:F:380:ALA:CB	1:F:414:MET:HE1	2.17	0.74
1:E:425:PRO:CG	1:E:451:VAL:HG13	2.17	0.73
1:A:380:ALA:HB2	1:A:414:MET:HE1	1.71	0.73
1:C:433:MET:HE2	1:C:444:ILE:HD12	1.70	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:447:PHE:CD1	2:O:250:PRO:HG3	2.24	0.72
1:H:425:PRO:CG	1:H:451:VAL:HG13	2.19	0.72
1:F:425:PRO:HG3	1:F:451:VAL:HG13	1.70	0.72
1:A:434:LEU:HB3	1:A:443:VAL:HB	1.69	0.71
1:B:447:PHE:CD1	2:J:250:PRO:HG3	2.25	0.71
1:A:425:PRO:HG3	1:A:451:VAL:CG1	2.19	0.71
1:E:380:ALA:HB2	1:E:414:MET:HE1	1.70	0.71
1:H:481:LYS:O	1:H:482:ASN:HB2	1.91	0.71
1:H:394:ILE:CD1	1:H:407:LEU:HD21	2.20	0.71
1:A:339:GLU:O	1:A:343:LEU:HD13	1.91	0.71
1:H:434:LEU:HB3	1:H:443:VAL:HB	1.73	0.71
1:F:448:ARG:HG3	2:J:249:ACE:H1	1.72	0.70
1:H:433:MET:HE2	1:H:444:ILE:HD12	1.74	0.70
1:C:380:ALA:CB	1:C:414:MET:HE1	2.21	0.70
1:B:390:MET:C	1:B:414:MET:HE2	2.16	0.70
1:D:448:ARG:NH2	2:P:251:VAL:HG22	2.06	0.70
1:H:380:ALA:HB2	1:H:414:MET:HE1	1.74	0.70
1:E:390:MET:C	1:E:414:MET:HE2	2.17	0.70
1:H:478:MET:SD	1:H:479:GLU:N	2.65	0.70
1:G:365:ARG:O	1:G:369:VAL:HG23	1.92	0.69
1:D:456:PHE:HB2	3:D:1021:HOH:O	1.92	0.69
1:G:481:LYS:O	1:G:482:ASN:HB2	1.92	0.69
1:E:433:MET:HE2	1:E:444:ILE:HD12	1.76	0.68
1:F:434:LEU:HB3	1:F:443:VAL:HB	1.73	0.68
1:H:390:MET:C	1:H:414:MET:HE2	2.18	0.68
2:K:254:THR:HG22	2:K:254:THR:O	1.92	0.68
1:A:390:MET:C	1:A:414:MET:HE2	2.18	0.68
1:F:389:LYS:HG2	1:F:414:MET:CE	2.20	0.67
1:G:434:LEU:HB3	1:G:443:VAL:HB	1.75	0.67
1:E:367:GLU:HG2	1:E:372:ARG:HH21	1.60	0.67
1:C:480:ALA:O	1:C:481:LYS:O	2.14	0.67
1:F:380:ALA:HB2	1:F:414:MET:HE1	1.78	0.66
1:A:478:MET:HE1	1:A:490:ILE:HG21	1.77	0.66
1:H:334:ALA:O	1:H:336:ALA:N	2.24	0.66
1:B:380:ALA:HB2	1:B:414:MET:HE1	1.76	0.66
1:D:363:ARG:NH1	1:D:367:GLU:OE1	2.28	0.66
1:A:405:THR:OG1	1:A:406:HIS:HD2	1.80	0.66
1:C:409:LEU:HD22	1:C:432:LEU:HD13	1.77	0.65
1:G:418:ASN:O	1:G:422:LEU:HD13	1.95	0.65
1:G:394:ILE:HD11	1:G:407:LEU:HD21	1.77	0.65
1:B:434:LEU:HB3	1:B:443:VAL:HB	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:448:ARG:HG3	2:P:249:ACE:H1	1.77	0.65
1:D:363:ARG:HG2	1:D:363:ARG:HH11	1.63	0.64
1:D:389:LYS:HG2	1:D:414:MET:CE	2.20	0.64
1:E:478:MET:SD	1:E:479:GLU:N	2.71	0.64
1:B:405:THR:C	1:B:475:VAL:HG23	2.22	0.64
1:E:478:MET:HB3	1:E:484:TYR:CD1	2.33	0.64
1:H:380:ALA:HB1	1:H:414:MET:HE1	1.78	0.64
1:C:363:ARG:HG2	1:C:363:ARG:HH11	1.63	0.64
1:A:486:ARG:HH11	1:A:486:ARG:HG3	1.63	0.63
1:E:380:ALA:HB1	1:E:414:MET:HE1	1.79	0.63
1:C:418:ASN:O	1:C:422:LEU:HD13	1.98	0.63
1:D:338:LEU:O	1:D:342:VAL:HG23	1.98	0.63
1:G:478:MET:O	1:G:480:ALA:N	2.30	0.63
1:D:380:ALA:HB1	1:D:414:MET:HE1	1.81	0.62
1:H:425:PRO:HG3	1:H:451:VAL:CG1	2.23	0.62
1:F:365:ARG:NH2	1:F:397:ASN:OD1	2.33	0.62
1:B:433:MET:HE2	1:B:444:ILE:HD12	1.81	0.62
1:B:380:ALA:HB1	1:B:414:MET:HE1	1.80	0.61
1:F:433:MET:CE	1:F:444:ILE:HD12	2.29	0.61
1:G:385:ARG:HG2	1:H:349:THR:HG22	1.82	0.61
1:H:341:LYS:O	1:H:345:MET:HG3	2.00	0.61
1:A:389:LYS:HG2	1:A:414:MET:CE	2.23	0.61
1:A:418:ASN:O	1:A:422:LEU:HD13	2.01	0.61
1:F:390:MET:C	1:F:414:MET:HE2	2.26	0.61
1:C:478:MET:SD	1:C:479:GLU:N	2.74	0.61
1:C:342:VAL:HG13	1:E:345:MET:SD	2.41	0.61
1:F:478:MET:SD	1:F:479:GLU:N	2.73	0.61
1:C:380:ALA:HB2	1:C:414:MET:HE1	1.83	0.60
1:C:394:ILE:HD11	1:C:407:LEU:HD21	1.82	0.60
1:E:341:LYS:O	1:E:345:MET:HG3	2.01	0.60
1:B:394:ILE:HD11	1:B:407:LEU:HD21	1.83	0.60
1:E:434:LEU:HB3	1:E:443:VAL:HB	1.83	0.59
1:H:481:LYS:O	1:H:482:ASN:CB	2.49	0.59
1:B:405:THR:O	1:B:475:VAL:HG23	2.02	0.59
1:F:380:ALA:HB1	1:F:414:MET:HE1	1.83	0.59
1:H:334:ALA:C	1:H:336:ALA:H	2.09	0.59
1:G:409:LEU:HD22	1:G:432:LEU:HD13	1.83	0.59
1:C:414:MET:HG2	1:C:463:MET:HG2	1.85	0.59
1:E:360:ASP:OD1	1:E:363:ARG:NE	2.28	0.59
1:H:365:ARG:HG2	3:H:1001:HOH:O	2.03	0.58
1:A:394:ILE:CD1	1:A:407:LEU:HD21	2.32	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:485:VAL:HG22	1:B:490:ILE:CG2	2.32	0.58
1:F:491:PHE:HB2	3:F:1052:HOH:O	2.03	0.58
1:D:405:THR:OG1	1:D:406:HIS:HD2	1.87	0.58
1:B:334:ALA:C	1:B:336:ALA:N	2.59	0.58
1:H:486:ARG:HG3	1:H:486:ARG:HH11	1.68	0.58
1:D:434:LEU:HB3	1:D:443:VAL:HB	1.86	0.58
1:G:361:PHE:O	1:G:365:ARG:HB2	2.03	0.57
1:B:467:SER:OG	2:J:250:PRO:HB2	2.04	0.57
1:H:363:ARG:O	1:H:367:GLU:HG3	2.05	0.57
1:B:354:PHE:CE2	1:B:356:TRP:HB2	2.40	0.57
1:E:478:MET:O	1:E:480:ALA:N	2.37	0.56
1:F:338:LEU:O	1:F:342:VAL:HG23	2.05	0.56
1:A:360:ASP:OD1	1:A:363:ARG:NE	2.33	0.56
1:D:380:ALA:HB2	1:D:414:MET:HE1	1.88	0.56
1:F:394:ILE:CD1	1:F:407:LEU:HD21	2.31	0.56
1:F:412:VAL:HG22	1:F:466:ALA:HB2	1.87	0.56
1:D:367:GLU:HG2	1:D:372:ARG:HH21	1.71	0.56
1:B:425:PRO:CG	1:B:451:VAL:HG13	2.35	0.56
1:H:438:ASN:ND2	1:H:483:SER:HB3	2.20	0.56
1:B:478:MET:HB3	1:B:484:TYR:CD1	2.41	0.56
1:A:380:ALA:HB1	1:A:414:MET:HE1	1.87	0.56
1:E:337:ASP:O	1:E:340:GLN:N	2.30	0.55
1:E:405:THR:OG1	1:E:406:HIS:HD2	1.89	0.55
1:D:448:ARG:HH21	2:P:251:VAL:HG22	1.71	0.55
1:G:481:LYS:O	1:G:482:ASN:CB	2.53	0.55
1:H:451:VAL:HG12	1:H:451:VAL:O	2.06	0.55
2:J:254:THR:O	2:J:254:THR:CG2	2.53	0.55
1:C:365:ARG:HH21	1:C:397:ASN:HA	1.72	0.55
1:G:405:THR:O	1:G:475:VAL:HG23	2.07	0.55
1:G:425:PRO:HG2	1:G:451:VAL:HG12	1.89	0.55
1:C:481:LYS:O	1:C:482:ASN:HB2	2.06	0.54
1:F:425:PRO:CG	1:F:451:VAL:HG13	2.36	0.54
1:B:478:MET:SD	1:B:479:GLU:N	2.80	0.54
1:B:481:LYS:O	1:B:482:ASN:HB2	2.06	0.54
1:D:367:GLU:HG2	1:D:372:ARG:NH2	2.23	0.54
1:H:418:ASN:O	1:H:422:LEU:HD13	2.07	0.54
1:G:360:ASP:HB2	1:G:364:LYS:HE3	1.89	0.54
1:H:396:LEU:HD22	3:H:1001:HOH:O	2.07	0.54
1:E:486:ARG:HG3	1:E:486:ARG:HH11	1.73	0.54
1:G:412:VAL:HG22	1:G:466:ALA:HB2	1.89	0.53
1:G:425:PRO:CG	1:G:451:VAL:HG12	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:478:MET:SD	1:A:479:GLU:N	2.81	0.53
1:E:436:ASP:OD1	1:E:483:SER:OG	2.22	0.53
1:G:380:ALA:HB2	1:G:414:MET:HE1	1.88	0.53
1:H:389:LYS:HG2	1:H:414:MET:CE	2.29	0.53
1:F:434:LEU:HD21	1:F:478:MET:HG2	1.90	0.53
1:E:425:PRO:HG3	1:E:451:VAL:CG1	2.33	0.53
1:E:334:ALA:HB2	3:E:1061:HOH:O	2.09	0.53
1:E:451:VAL:HG12	1:E:451:VAL:O	2.09	0.53
1:E:467:SER:OG	2:M:250:PRO:HB2	2.09	0.53
1:D:418:ASN:O	1:D:422:LEU:HD13	2.09	0.53
1:E:405:THR:O	1:E:475:VAL:HG23	2.08	0.53
1:G:354:PHE:HZ	1:G:379:PRO:HD2	1.72	0.53
1:H:339:GLU:O	1:H:343:LEU:HD13	2.09	0.53
1:A:476:SER:O	1:A:477:LYS:C	2.52	0.52
1:A:481:LYS:O	1:A:482:ASN:HB2	2.09	0.52
1:C:380:ALA:HB1	1:C:414:MET:HE1	1.88	0.52
1:E:365:ARG:O	1:E:369:VAL:HG23	2.09	0.52
1:B:441:GLU:HA	1:F:440:ARG:HA	1.91	0.52
1:A:478:MET:HB2	1:A:484:TYR:HB2	1.90	0.52
1:D:476:SER:O	1:D:477:LYS:C	2.52	0.52
1:B:389:LYS:C	1:B:390:MET:HG3	2.33	0.52
1:B:477:LYS:C	1:B:478:MET:O	2.50	0.52
1:A:479:GLU:O	1:A:480:ALA:O	2.28	0.52
1:C:476:SER:O	1:C:477:LYS:O	2.27	0.51
1:F:393:ARG:NH1	3:F:1012:HOH:O	2.43	0.51
1:D:448:ARG:HG3	2:P:249:ACE:CH3	2.39	0.51
1:E:394:ILE:CD1	1:E:407:LEU:HD21	2.38	0.51
1:B:334:ALA:C	1:B:336:ALA:H	2.16	0.51
1:G:409:LEU:HD23	1:G:410:PHE:N	2.26	0.51
1:G:478:MET:HB3	1:G:484:TYR:CD1	2.46	0.51
1:B:414:MET:HG2	1:B:463:MET:HG2	1.93	0.51
1:E:478:MET:HB2	1:E:484:TYR:HB2	1.93	0.51
1:B:365:ARG:HH21	1:B:397:ASN:HA	1.76	0.51
1:G:478:MET:C	1:G:480:ALA:H	2.18	0.51
1:D:409:LEU:HD23	1:D:410:PHE:N	2.26	0.51
1:E:407:LEU:HD23	1:E:407:LEU:C	2.35	0.51
1:H:354:PHE:HZ	1:H:379:PRO:HD2	1.75	0.51
1:H:478:MET:O	1:H:480:ALA:N	2.44	0.51
1:G:467:SER:OG	2:O:250:PRO:HB2	2.11	0.51
1:F:336:ALA:C	1:F:338:LEU:H	2.18	0.50
1:H:480:ALA:C	1:H:481:LYS:O	2.53	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:405:THR:OG1	1:F:406:HIS:HD2	1.93	0.50
1:A:367:GLU:HG2	1:A:372:ARG:HH21	1.76	0.50
1:B:354:PHE:HZ	1:B:379:PRO:HD2	1.76	0.50
1:C:336:ALA:C	1:C:338:LEU:H	2.20	0.50
1:D:480:ALA:O	1:D:481:LYS:O	2.30	0.50
1:E:477:LYS:O	1:E:478:MET:O	2.29	0.50
1:F:478:MET:HB2	1:F:484:TYR:HB2	1.92	0.50
1:C:365:ARG:O	1:C:369:VAL:HG23	2.12	0.50
1:D:394:ILE:HD11	1:D:407:LEU:HD21	1.93	0.50
1:G:390:MET:CA	1:G:414:MET:HE2	2.41	0.50
1:H:344:GLU:O	1:H:345:MET:C	2.54	0.50
1:G:339:GLU:O	1:G:340:GLN:C	2.54	0.49
1:H:405:THR:O	1:H:475:VAL:HG23	2.13	0.49
1:D:339:GLU:O	1:D:340:GLN:C	2.56	0.49
1:H:388:TYR:CD2	1:H:422:LEU:HD23	2.48	0.49
1:B:478:MET:O	1:B:479:GLU:CB	2.60	0.49
1:D:360:ASP:HB3	1:D:364:LYS:HE3	1.93	0.49
1:H:451:VAL:CG1	1:H:451:VAL:O	2.61	0.49
1:D:477:LYS:O	1:D:482:ASN:OD1	2.32	0.48
1:H:444:ILE:HG12	1:H:445:ASP:N	2.27	0.48
1:H:467:SER:OG	2:P:250:PRO:HB2	2.13	0.48
1:A:354:PHE:CE2	1:A:356:TRP:HB2	2.49	0.48
1:D:388:TYR:CD2	1:D:422:LEU:HD23	2.48	0.48
1:B:485:VAL:HG22	1:B:490:ILE:HG22	1.95	0.48
1:C:338:LEU:HG	1:E:338:LEU:HD21	1.94	0.48
1:C:476:SER:O	1:C:477:LYS:C	2.57	0.48
1:F:477:LYS:C	1:F:478:MET:O	2.55	0.48
1:F:486:ARG:HG3	1:F:486:ARG:HH11	1.79	0.48
1:A:416:GLY:HA3	3:A:1067:HOH:O	2.13	0.48
1:C:363:ARG:HG2	1:C:363:ARG:NH1	2.28	0.48
1:H:447:PHE:CG	2:P:250:PRO:HG3	2.49	0.48
1:G:363:ARG:O	1:G:367:GLU:HG3	2.14	0.48
1:F:338:LEU:HD11	1:G:342:VAL:HG21	1.96	0.48
1:F:448:ARG:HG3	2:J:249:ACE:CH3	2.43	0.48
1:B:343:LEU:N	1:B:343:LEU:HD12	2.29	0.47
1:G:385:ARG:HG3	1:G:385:ARG:HH11	1.78	0.47
1:A:414:MET:HG2	1:A:463:MET:HG2	1.95	0.47
1:D:414:MET:HG2	1:D:463:MET:HG2	1.97	0.47
1:B:405:THR:OG1	1:B:406:HIS:HD2	1.97	0.47
1:G:478:MET:HE2	1:G:484:TYR:HB2	1.96	0.47
1:H:396:LEU:HB3	3:H:1001:HOH:O	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:339:GLU:O	1:B:343:LEU:CD1	2.63	0.47
1:D:382:TYR:N	1:D:382:TYR:CD1	2.82	0.47
1:D:478:MET:SD	1:D:479:GLU:N	2.87	0.47
1:G:407:LEU:C	1:G:407:LEU:HD23	2.40	0.47
1:H:334:ALA:C	1:H:336:ALA:N	2.69	0.47
1:C:447:PHE:CG	2:K:250:PRO:HG3	2.50	0.47
1:B:363:ARG:HG2	1:B:363:ARG:HH11	1.80	0.47
1:A:382:TYR:CD1	1:A:382:TYR:N	2.82	0.47
1:B:479:GLU:O	1:B:480:ALA:O	2.33	0.47
1:C:365:ARG:NH2	1:C:397:ASN:OD1	2.48	0.47
1:D:448:ARG:HH21	2:P:251:VAL:CG2	2.29	0.46
1:H:339:GLU:C	1:H:343:LEU:HD13	2.40	0.46
1:C:405:THR:OG1	1:C:406:HIS:HD2	1.99	0.46
1:E:438:ASN:ND2	1:E:483:SER:CB	2.77	0.46
1:E:438:ASN:ND2	1:E:483:SER:HB3	2.31	0.46
1:C:367:GLU:HG2	1:C:372:ARG:NH2	2.31	0.46
1:F:355:ILE:HD12	1:H:386:TYR:CG	2.51	0.46
1:C:425:PRO:HG3	1:C:451:VAL:HG13	1.97	0.46
1:E:382:TYR:N	1:E:382:TYR:CD1	2.84	0.46
1:G:385:ARG:HG3	1:G:385:ARG:NH1	2.31	0.46
1:B:409:LEU:HD23	1:B:410:PHE:N	2.31	0.46
1:C:339:GLU:O	1:C:343:LEU:CD1	2.61	0.46
1:G:399:ASP:OD1	2:O:254:THR:HB	2.17	0.45
1:G:433:MET:HE3	1:G:444:ILE:HD12	1.98	0.45
1:B:478:MET:HE2	1:B:484:TYR:HB2	1.98	0.45
1:H:422:LEU:CD1	1:H:422:LEU:N	2.79	0.45
1:B:478:MET:CB	1:B:484:TYR:CD1	3.00	0.45
1:C:354:PHE:CE2	1:C:356:TRP:HB2	2.52	0.45
1:F:448:ARG:NH2	2:J:251:VAL:HG22	2.31	0.45
2:K:254:THR:O	2:K:254:THR:CG2	2.62	0.45
1:D:407:LEU:C	1:D:407:LEU:HD23	2.42	0.45
1:F:478:MET:HB3	1:F:484:TYR:CD1	2.52	0.44
1:A:360:ASP:HB3	1:A:364:LYS:HE3	2.00	0.44
1:D:390:MET:CA	1:D:414:MET:HE2	2.47	0.44
1:F:444:ILE:CG1	1:F:445:ASP:N	2.81	0.44
1:A:451:VAL:HG12	1:A:451:VAL:O	2.17	0.44
1:C:389:LYS:HG2	1:C:414:MET:CE	2.29	0.44
1:B:361:PHE:O	1:B:365:ARG:HB2	2.16	0.44
1:H:438:ASN:ND2	1:H:483:SER:CB	2.80	0.44
1:F:355:ILE:HB	1:H:386:TYR:CE2	2.52	0.44
1:G:354:PHE:CE2	1:G:356:TRP:HB2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:388:TYR:CD2	1:C:422:LEU:HD23	2.53	0.44
1:C:412:VAL:HG22	1:C:466:ALA:HB2	2.00	0.44
1:H:339:GLU:O	1:H:342:VAL:N	2.51	0.44
1:A:407:LEU:C	1:A:407:LEU:HD23	2.43	0.44
1:C:382:TYR:CD1	1:C:382:TYR:N	2.86	0.44
1:D:448:ARG:NH2	2:P:251:VAL:CG2	2.80	0.44
1:H:350:TYR:CD2	1:H:385:ARG:HA	2.53	0.44
1:F:481:LYS:O	1:F:482:ASN:HB2	2.18	0.43
1:H:414:MET:HG2	1:H:463:MET:HG2	1.99	0.43
1:D:363:ARG:NH1	1:D:363:ARG:HG2	2.28	0.43
1:C:479:GLU:O	1:C:480:ALA:O	2.36	0.43
1:A:447:PHE:CG	2:I:250:PRO:HG3	2.51	0.43
1:B:406:HIS:CD2	1:B:471:LEU:HD22	2.54	0.43
1:E:405:THR:C	1:E:475:VAL:HG23	2.43	0.43
1:B:438:ASN:ND2	1:B:483:SER:CB	2.82	0.43
1:H:406:HIS:CE1	1:H:474:PRO:HG3	2.53	0.43
1:E:476:SER:O	1:E:477:LYS:O	2.36	0.43
1:B:478:MET:HB2	1:B:484:TYR:HB2	2.01	0.43
1:C:486:ARG:HG3	1:C:486:ARG:HH11	1.84	0.43
1:D:338:LEU:HD12	1:D:338:LEU:HA	1.82	0.43
1:D:354:PHE:CE2	1:D:356:TRP:HB2	2.54	0.43
1:E:451:VAL:CG1	1:E:451:VAL:O	2.66	0.43
1:B:341:LYS:O	1:B:345:MET:HG3	2.19	0.43
1:H:425:PRO:HB2	3:H:1032:HOH:O	2.19	0.43
1:A:476:SER:O	1:A:477:LYS:O	2.36	0.43
1:F:473:CYS:HA	1:F:474:PRO:HD3	1.87	0.43
1:C:447:PHE:CD1	1:C:447:PHE:C	2.97	0.43
1:D:394:ILE:CD1	1:D:407:LEU:HD21	2.49	0.43
1:A:469:CYS:HA	1:A:470:PRO:HD2	1.95	0.42
1:E:486:ARG:HG3	1:E:486:ARG:NH1	2.34	0.42
1:G:481:LYS:C	1:G:482:ASN:OD1	2.61	0.42
1:B:365:ARG:NH2	1:B:397:ASN:OD1	2.52	0.42
1:B:466:ALA:HB3	2:J:253:GLU:CD	2.44	0.42
1:C:425:PRO:CG	1:C:451:VAL:HG13	2.49	0.42
1:G:414:MET:HG2	1:G:463:MET:HG2	2.00	0.42
1:A:486:ARG:HG3	1:A:486:ARG:NH1	2.32	0.42
1:A:478:MET:HB2	1:A:478:MET:HE2	1.97	0.42
1:F:361:PHE:O	1:F:365:ARG:HB2	2.20	0.42
1:D:478:MET:O	1:D:479:GLU:CB	2.68	0.42
1:F:447:PHE:CD1	1:F:447:PHE:C	2.97	0.42
1:F:491:PHE:CB	3:F:1052:HOH:O	2.64	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:412:VAL:HG22	1:G:466:ALA:CB	2.50	0.42
1:E:434:LEU:HD21	1:E:478:MET:HG2	2.00	0.42
1:F:390:MET:O	1:F:414:MET:HE2	2.20	0.42
1:G:389:LYS:C	1:G:390:MET:HG3	2.43	0.42
1:G:434:LEU:HD21	1:G:478:MET:HE3	2.00	0.42
1:B:447:PHE:CD1	1:B:447:PHE:C	2.98	0.42
1:G:339:GLU:O	1:G:342:VAL:N	2.53	0.42
1:A:361:PHE:O	1:A:365:ARG:HB2	2.20	0.41
1:A:412:VAL:HG22	1:A:466:ALA:HB2	2.03	0.41
1:E:480:ALA:O	1:E:481:LYS:O	2.37	0.41
1:D:389:LYS:C	1:D:390:MET:HG3	2.46	0.41
1:D:440:ARG:HA	1:H:441:GLU:HA	2.02	0.41
1:B:354:PHE:HE2	1:B:356:TRP:HB2	1.84	0.41
1:B:407:LEU:C	1:B:407:LEU:HD23	2.46	0.41
1:D:486:ARG:O	1:D:487:ASP:C	2.63	0.41
1:E:361:PHE:O	1:E:365:ARG:HB2	2.20	0.41
1:F:425:PRO:HG3	1:F:451:VAL:CG1	2.46	0.41
1:C:477:LYS:C	1:C:478:MET:O	2.63	0.41
1:G:405:THR:C	1:G:475:VAL:HG23	2.45	0.41
1:F:469:CYS:HA	1:F:470:PRO:HD2	1.96	0.41
1:G:478:MET:C	1:G:480:ALA:N	2.78	0.41
1:A:436:ASP:OD1	1:A:483:SER:OG	2.37	0.41
1:D:339:GLU:O	1:D:343:LEU:HD12	2.21	0.41
1:D:363:ARG:O	1:D:367:GLU:HG3	2.21	0.41
1:A:399:ASP:OD1	2:I:254:THR:HB	2.21	0.41
1:C:467:SER:OG	2:K:250:PRO:HB2	2.21	0.41
1:E:459:PRO:HG3	1:E:464:ASN:HD21	1.86	0.41
1:B:434:LEU:HD21	1:B:478:MET:HG2	2.02	0.41
1:D:384:SER:O	1:D:385:ARG:C	2.64	0.41
1:D:451:VAL:HA	3:D:1021:HOH:O	2.21	0.41
1:E:399:ASP:OD1	2:M:254:THR:HB	2.21	0.40
1:G:394:ILE:CD1	1:G:407:LEU:HD21	2.48	0.40
1:B:473:CYS:HA	1:B:474:PRO:HD3	1.91	0.40
1:G:488:ASP:O	1:G:488:ASP:CG	2.63	0.40
1:H:448:ARG:HA	1:H:449:PRO:HD3	1.91	0.40
1:H:486:ARG:HG3	1:H:486:ARG:NH1	2.33	0.40
1:B:418:ASN:O	1:B:422:LEU:HD13	2.22	0.40
1:C:336:ALA:C	1:C:338:LEU:N	2.80	0.40
1:C:467:SER:HA	2:K:252:GLN:HA	2.03	0.40
1:F:485:VAL:HG22	1:F:490:ILE:CG2	2.51	0.40
1:G:377:PHE:CE2	1:G:463:MET:SD	3.14	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:447:PHE:CG	2:O:250:PRO:HG3	2.57	0.40
1:B:394:ILE:CD1	1:B:407:LEU:HD21	2.49	0.40
1:D:341:LYS:O	1:D:345:MET:HG3	2.22	0.40
1:F:340:GLN:HA	1:F:343:LEU:HD22	2.02	0.40
1:B:458:ARG:HA	1:B:459:PRO:HD3	1.97	0.40
1:D:405:THR:C	1:D:475:VAL:HG23	2.46	0.40
1:G:421:LEU:HD11	1:H:491:PHE:CD2	2.56	0.40
1:G:437:GLN:NE2	1:G:486:ARG:HB3	2.35	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	166/168 (99%)	156 (94%)	6 (4%)	4 (2%)	4	2
1	B	166/168 (99%)	156 (94%)	6 (4%)	4 (2%)	4	2
1	C	166/168 (99%)	154 (93%)	7 (4%)	5 (3%)	3	1
1	D	166/168 (99%)	156 (94%)	7 (4%)	3 (2%)	6	3
1	E	166/168 (99%)	155 (93%)	5 (3%)	6 (4%)	2	1
1	F	166/168 (99%)	158 (95%)	5 (3%)	3 (2%)	6	3
1	G	166/168 (99%)	155 (93%)	9 (5%)	2 (1%)	10	6
1	H	166/168 (99%)	152 (92%)	8 (5%)	6 (4%)	2	1
2	I	5/7 (71%)	5 (100%)	0	0	100	100
2	J	5/7 (71%)	4 (80%)	1 (20%)	0	100	100
2	K	5/7 (71%)	5 (100%)	0	0	100	100
2	L	5/7 (71%)	5 (100%)	0	0	100	100
2	M	5/7 (71%)	5 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	N	5/7 (71%)	5 (100%)	0	0	100	100
2	O	5/7 (71%)	5 (100%)	0	0	100	100
2	P	5/7 (71%)	5 (100%)	0	0	100	100
All	All	1368/1400 (98%)	1281 (94%)	54 (4%)	33 (2%)	4	2

All (33) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	477	LYS
1	A	480	ALA
1	B	477	LYS
1	B	480	ALA
1	C	477	LYS
1	C	480	ALA
1	C	481	LYS
1	D	481	LYS
1	E	477	LYS
1	E	478	MET
1	E	480	ALA
1	E	481	LYS
1	F	477	LYS
1	F	480	ALA
1	H	480	ALA
1	H	482	ASN
1	D	480	ALA
1	E	479	GLU
1	G	482	ASN
1	H	335	MET
1	H	477	LYS
1	A	481	LYS
1	B	478	MET
1	E	482	ASN
1	F	478	MET
1	G	477	LYS
1	H	478	MET
1	A	479	GLU
1	B	481	LYS
1	C	478	MET
1	H	479	GLU
1	D	477	LYS
1	C	337	ASP

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	132/144 (92%)	128 (97%)	4 (3%)	36	38
1	B	132/144 (92%)	127 (96%)	5 (4%)	29	29
1	C	132/144 (92%)	129 (98%)	3 (2%)	44	49
1	D	132/144 (92%)	128 (97%)	4 (3%)	36	38
1	E	132/144 (92%)	128 (97%)	4 (3%)	36	38
1	F	132/144 (92%)	127 (96%)	5 (4%)	29	29
1	G	132/144 (92%)	129 (98%)	3 (2%)	44	49
1	H	132/144 (92%)	131 (99%)	1 (1%)	73	80
2	I	5/5 (100%)	4 (80%)	1 (20%)	1	0
2	J	5/5 (100%)	5 (100%)	0	100	100
2	K	5/5 (100%)	4 (80%)	1 (20%)	1	0
2	L	5/5 (100%)	4 (80%)	1 (20%)	1	0
2	M	5/5 (100%)	4 (80%)	1 (20%)	1	0
2	N	5/5 (100%)	4 (80%)	1 (20%)	1	0
2	O	5/5 (100%)	4 (80%)	1 (20%)	1	0
2	P	5/5 (100%)	4 (80%)	1 (20%)	1	0
All	All	1096/1192 (92%)	1060 (97%)	36 (3%)	33	34

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	365	ARG
1	A	366	GLN
1	A	409	LEU
1	A	422	LEU
1	B	360	ASP
1	B	365	ARG
1	B	366	GLN
1	B	409	LEU

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Mol	Chain	Res	Type
1	B	422	LEU
1	C	365	ARG
1	C	366	GLN
1	C	409	LEU
1	D	365	ARG
1	D	407	LEU
1	D	422	LEU
1	D	451	VAL
1	E	365	ARG
1	E	407	LEU
1	E	409	LEU
1	E	422	LEU
1	F	360	ASP
1	F	365	ARG
1	F	409	LEU
1	F	422	LEU
1	F	428	GLN
1	G	365	ARG
1	G	366	GLN
1	G	451	VAL
1	H	409	LEU
2	I	254	THR
2	K	254	THR
2	L	254	THR
2	M	254	THR
2	N	254	THR
2	O	254	THR
2	P	254	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	366	GLN
1	A	406	HIS
1	A	438	ASN
1	A	439	ASN
1	A	461	ASN
1	B	406	HIS
1	B	438	ASN
1	B	461	ASN
1	C	406	HIS
1	C	428	GLN

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Mol	Chain	Res	Type
1	C	461	ASN
1	D	406	HIS
1	D	439	ASN
1	D	457	GLN
1	D	461	ASN
1	E	366	GLN
1	E	406	HIS
1	E	438	ASN
1	E	461	ASN
1	F	366	GLN
1	F	406	HIS
1	F	461	ASN
1	G	406	HIS
1	G	461	ASN
1	H	366	GLN
1	H	406	HIS
1	H	438	ASN
1	H	461	ASN
2	J	252	GLN
2	K	252	GLN
2	M	252	GLN
2	N	252	GLN
2	O	252	GLN
2	P	252	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.