



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:02 PM UTC

PDB ID : 1JYM / pdb_00001jym
Title : Crystals of Peptide Deformylase from Plasmodium falciparum with Ten Subunits per Asymmetric Unit Reveal Critical Characteristics of the Active Site for Drug Design
Authors : Kumar, A.; Nguyen, K.T.; Srivathsan, S.; Ornstein, B.; Turley, S.; Hirsh, I.; Pei, D.; Hol, W.G.J.
Deposited on : 2001-09-12
Resolution : 2.80 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)	:	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

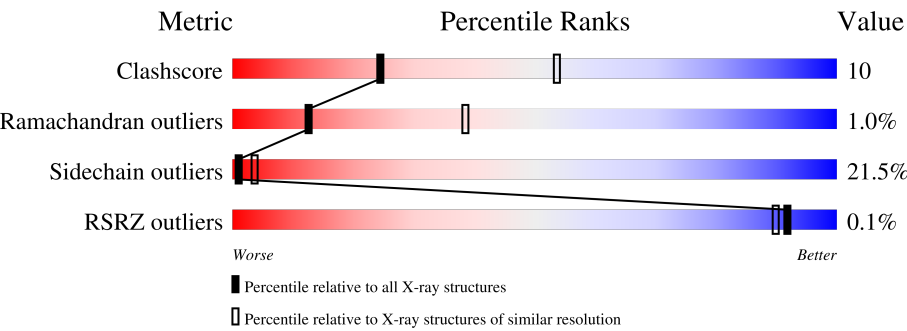
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	183	
1	B	183	
1	C	183	
1	D	183	
1	E	183	
1	F	183	

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Mol	Chain	Length	Quality of chain
1	G	183	58%27%9%6%
1	H	183	54%33%7%5%
1	I	183	53%34%5%6%
1	J	183	%56%27%11%6%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 14584 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 Peptide Deformylase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	179	Total	C	N	O	S	Se	0	0	0
			1487	953	263	267	1	3			
1	B	179	Total	C	N	O	S	Se	0	0	0
			1487	953	263	267	1	3			
1	C	168	Total	C	N	O	S	Se	0	0	0
			1394	898	246	246	1	3			
1	D	173	Total	C	N	O	S	Se	0	0	0
			1435	921	253	257	1	3			
1	E	176	Total	C	N	O	S	Se	0	0	0
			1458	936	256	262	1	3			
1	F	172	Total	C	N	O	S	Se	0	0	0
			1426	916	252	254	1	3			
1	G	172	Total	C	N	O	S	Se	0	0	0
			1425	915	252	254	1	3			
1	H	173	Total	C	N	O	S	Se	0	0	0
			1435	921	253	257	1	3			
1	I	172	Total	C	N	O	S	Se	0	0	0
			1425	915	252	254	1	3			
1	J	172	Total	C	N	O	S	Se	0	0	0
			1425	915	252	254	1	3			

- Molecule 2 is COBALT (II) ION (CCD ID: CO) (formula: Co).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Co	0	0
			1	1		
2	B	1	Total	Co	0	0
			1	1		
2	C	1	Total	Co	0	0
			1	1		
2	D	1	Total	Co	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	E	1	Total 1	Co 1	0	0
2	F	1	Total 1	Co 1	0	0
2	G	1	Total 1	Co 1	0	0
2	H	1	Total 1	Co 1	0	0
2	I	1	Total 1	Co 1	0	0
2	J	1	Total 1	Co 1	0	0

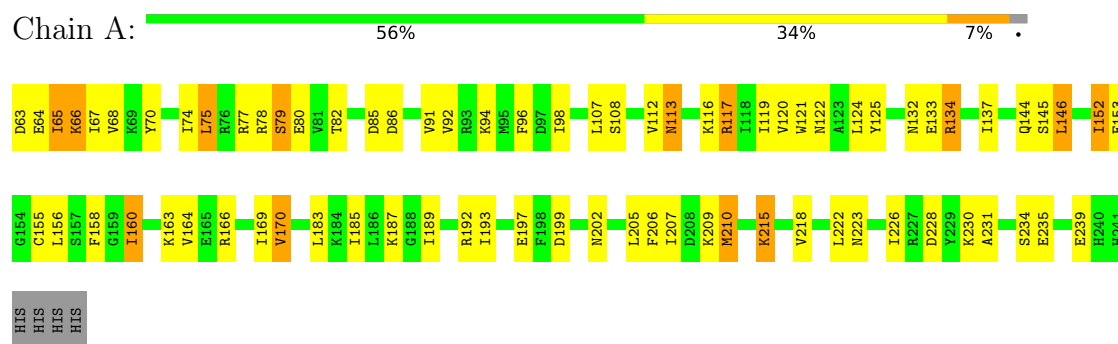
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	22	Total 22	O 22	0	0
3	B	17	Total 17	O 17	0	0
3	C	17	Total 17	O 17	0	0
3	D	16	Total 16	O 16	0	0
3	E	19	Total 19	O 19	0	0
3	F	26	Total 26	O 26	0	0
3	G	13	Total 13	O 13	0	0
3	H	15	Total 15	O 15	0	0
3	I	16	Total 16	O 16	0	0
3	J	16	Total 16	O 16	0	0

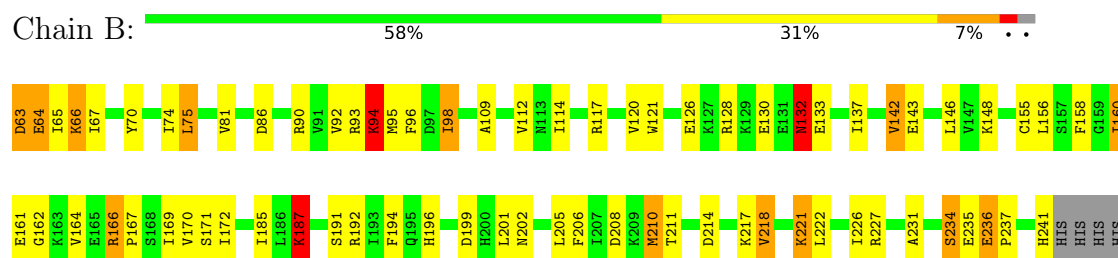
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

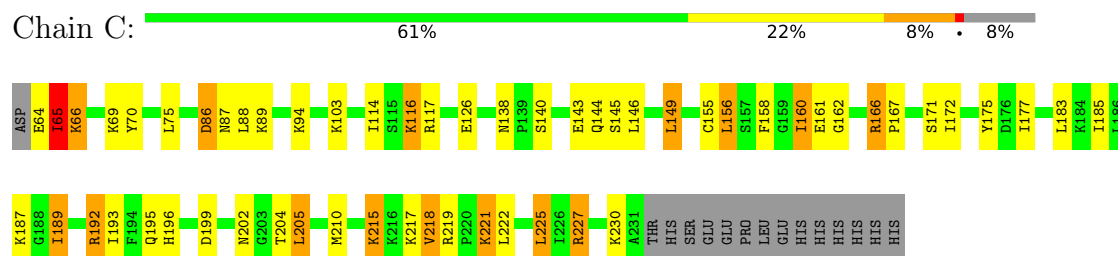
• Molecule 1: Peptide Deformylase



• Molecule 1: Peptide Deformylase

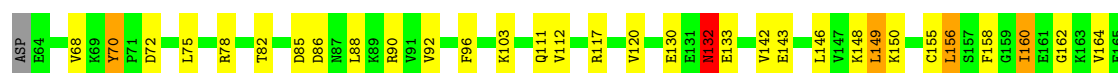


• Molecule 1: Peptide Deformylase

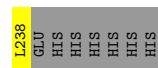


• Molecule 1: Peptide Deformylase

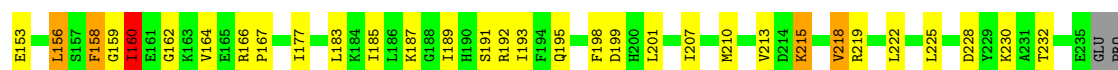




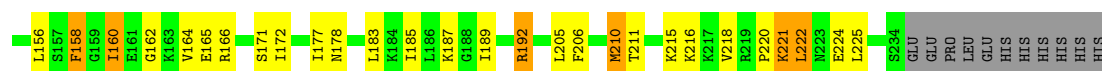
• Molecule 1: Peptide Deformylase



• Molecule 1: Peptide Deformylase

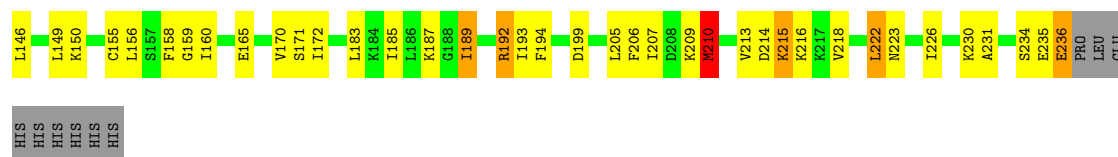


• Molecule 1: Peptide Deformylase



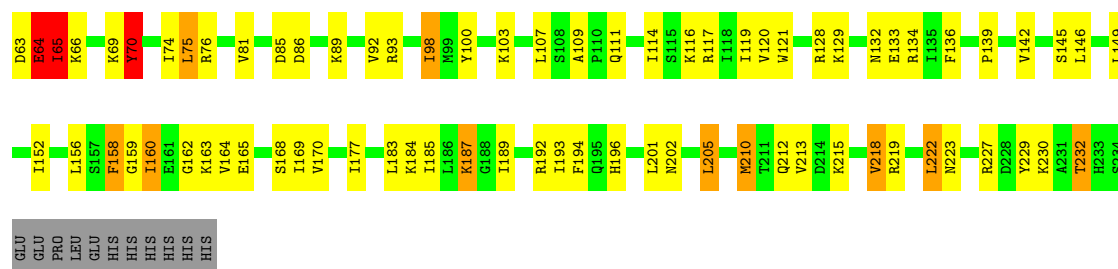
• Molecule 1: Peptide Deformylase





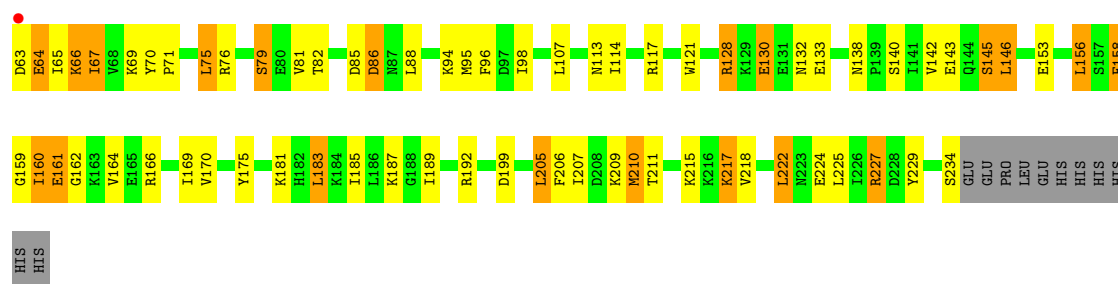
• Molecule 1: Peptide Deformylase

Chain I: 53% 34% 5% 6%



• Molecule 1: Peptide Deformylase

Chain J: 56% 27% 11% 6%



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	121.26Å 121.26Å 177.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.80 30.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	82.5 (30.00-2.80) 82.5 (30.00-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.89 (at 2.80Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.231 , 0.296 (Not available) , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	43.5	Xtriage
Anisotropy	0.414	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 45.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.052 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	14584	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.78% 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 ⓘ

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

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	0.64	0/1513	1.11	5/2026 (0.2%)
1	B	0.73	1/1513 (0.1%)	1.15	8/2026 (0.4%)
1	C	0.81	0/1416	1.20	6/1893 (0.3%)
1	D	0.74	0/1458	1.12	5/1950 (0.3%)
1	E	0.74	0/1482	1.16	7/1984 (0.4%)
1	F	0.96	4/1449 (0.3%)	1.25	8/1938 (0.4%)
1	G	0.78	0/1448	1.17	7/1937 (0.4%)
1	H	0.85	1/1458 (0.1%)	1.27	9/1950 (0.5%)
1	I	0.71	0/1448	1.17	10/1937 (0.5%)
1	J	0.81	0/1448	1.23	8/1937 (0.4%)
All	All	0.78	6/14633 (0.0%)	1.18	73/19578 (0.4%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	138	ASN	C-O	-6.73	1.20	1.23
1	B	210	MSE	SE-CE	-6.33	1.76	1.95
1	H	210	MSE	SE-CE	-6.19	1.76	1.95
1	F	138	ASN	CA-C	-5.42	1.49	1.53
1	F	137	ILE	CA-CB	-5.39	1.47	1.54
1	F	65	ILE	CA-CB	5.17	1.61	1.54

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	189	ILE	N-CA-C	-15.75	96.27	113.43
1	I	132	ASN	N-CA-C	-10.45	97.94	113.61
1	A	132	ASN	N-CA-C	-9.38	101.36	113.17
1	B	235	GLU	N-CA-C	-9.34	101.30	112.89
1	F	134	ARG	N-CA-C	9.02	123.14	108.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	132	ASN	N-CA-C	-8.99	99.68	113.89
1	H	132	ASN	N-CA-C	-8.87	99.87	113.89
1	C	89	LYS	N-CA-C	-8.84	102.04	113.17
1	D	88	LEU	N-CA-C	-8.46	100.42	110.41
1	F	89	LYS	N-CA-C	-7.91	103.36	113.16
1	A	121	TRP	N-CA-C	7.82	118.94	108.38
1	I	64	GLU	N-CA-C	7.80	122.17	109.85
1	G	132	ASN	N-CA-C	-7.64	103.06	113.30
1	E	132	ASN	N-CA-C	-7.60	104.04	113.38
1	E	232	THR	N-CA-C	-7.57	101.93	112.45
1	I	65	ILE	N-CA-C	-7.38	104.31	112.80
1	H	121	TRP	N-CA-C	7.27	118.19	108.38
1	D	112	VAL	N-CA-C	-7.24	105.77	112.43
1	H	67	ILE	N-CA-C	-7.17	100.34	109.58
1	C	70	TYR	N-CA-C	6.94	125.14	109.81
1	F	160	ILE	N-CA-CB	-6.85	104.26	112.07
1	I	89	LYS	N-CA-C	-6.78	105.01	113.28
1	E	160	ILE	N-CA-CB	-6.76	103.88	112.60
1	J	158	PHE	CB-CA-C	-6.60	102.78	111.42
1	F	121	TRP	N-CA-C	6.52	117.18	108.38
1	I	158	PHE	CB-CA-C	-6.40	102.74	110.94
1	B	132	ASN	N-CA-C	-6.35	105.47	113.72
1	G	222	LEU	N-CA-C	-6.32	104.47	111.36
1	F	158	PHE	CB-CA-C	-6.29	101.05	110.24
1	A	235	GLU	N-CA-C	-6.28	105.11	112.89
1	J	88	LEU	N-CA-C	-6.19	103.02	110.44
1	B	121	TRP	N-CA-C	6.14	118.18	108.79
1	C	87	ASN	N-CA-C	-6.03	105.78	113.02
1	D	132	ASN	N-CA-C	-6.01	105.24	113.30
1	J	121	TRP	N-CA-C	5.98	116.31	107.88
1	B	126	GLU	N-CA-C	-5.94	106.00	113.18
1	G	67	ILE	N-CA-C	-5.91	101.39	109.55
1	H	87	ASN	N-CA-C	5.82	119.60	111.74
1	E	163	LYS	N-CA-C	5.82	117.69	108.67
1	A	170	VAL	N-CA-C	5.79	116.63	108.17
1	E	228	ASP	N-CA-C	-5.72	105.12	111.36
1	F	70	TYR	N-CA-C	5.68	122.37	109.81
1	I	163	LYS	N-CA-C	5.62	118.60	110.28
1	D	162	GLY	N-CA-C	5.61	118.93	110.18
1	F	159	GLY	CA-C-N	5.61	128.27	122.27
1	F	159	GLY	C-N-CA	5.61	128.27	122.27
1	H	88	LEU	N-CA-C	-5.55	105.49	114.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	64	GLU	N-CA-C	5.54	119.12	110.32
1	B	94	LYS	N-CA-C	-5.51	105.17	111.07
1	J	128	ARG	N-CA-C	5.50	117.27	108.41
1	H	86	ASP	N-CA-C	5.41	116.75	108.52
1	E	196	HIS	N-CA-C	-5.41	105.30	111.14
1	C	158	PHE	CB-CA-C	-5.40	100.76	109.89
1	C	66	LYS	N-CA-C	5.39	118.13	110.10
1	D	190	HIS	N-CA-C	-5.38	105.45	112.23
1	H	235	GLU	N-CA-C	-5.38	106.75	113.15
1	I	70	TYR	N-CA-C	5.37	121.68	109.81
1	C	65	ILE	N-CA-C	5.37	120.50	109.34
1	J	86	ASP	N-CA-C	5.36	116.99	108.79
1	I	121	TRP	N-CA-C	5.34	115.60	108.38
1	G	89	LYS	N-CA-C	-5.32	106.49	112.87
1	B	187	LYS	N-CA-C	5.30	116.96	107.80
1	H	142	VAL	N-CA-C	5.28	116.06	110.72
1	I	184	LYS	N-CA-C	5.26	117.08	108.55
1	A	91	VAL	N-CA-C	-5.24	105.39	110.42
1	G	70	TYR	N-CA-C	5.18	121.26	109.81
1	I	196	HIS	N-CA-C	-5.17	105.54	111.07
1	G	158	PHE	CB-CA-C	-5.15	102.72	110.24
1	J	227	ARG	CA-C-N	5.14	127.43	120.38
1	J	227	ARG	C-N-CA	5.14	127.43	120.38
1	B	234	SER	N-CA-C	5.13	116.96	107.73
1	G	121	TRP	N-CA-C	5.05	115.97	108.60
1	E	158	PHE	CB-CA-C	-5.01	103.31	110.62

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	1487	0	1533	35	0
1	B	1487	0	1533	34	0
1	C	1394	0	1460	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1435	0	1491	23	0
1	E	1458	0	1513	29	0
1	F	1426	0	1485	27	0
1	G	1425	0	1483	27	0
1	H	1435	0	1491	31	0
1	I	1425	0	1483	35	0
1	J	1425	0	1483	37	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
3	A	22	0	0	0	0
3	B	17	0	0	0	0
3	C	17	0	0	0	0
3	D	16	0	0	0	0
3	E	19	0	0	0	0
3	F	26	0	0	1	0
3	G	13	0	0	0	0
3	H	15	0	0	1	0
3	I	16	0	0	1	0
3	J	16	0	0	0	0
All	All	14584	0	14955	290	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 10.

All (290) 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:J:164:VAL:O	1:J:192:ARG:NH2	2.05	0.90
1:I:65:ILE:HA	1:I:98:ILE:HD12	1.57	0.86
1:D:117:ARG:HH22	1:D:202:ASN:HD21	1.26	0.84
1:F:65:ILE:HA	1:F:98:ILE:HD11	1.63	0.80
1:A:231:ALA:HA	1:A:234:SER:HB3	1.64	0.80
1:F:160:ILE:HG23	1:F:218:VAL:HG21	1.66	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:231:ALA:HA	1:E:234:SER:HB3	1.65	0.77
1:B:164:VAL:O	1:B:192:ARG:NH2	2.16	0.76
1:C:117:ARG:HH21	1:C:202:ASN:HD21	1.31	0.76
1:F:167:PRO:HD2	1:F:195:GLN:OE1	1.86	0.75
1:H:231:ALA:HA	1:H:234:SER:HB3	1.68	0.73
1:F:65:ILE:O	1:F:65:ILE:HG13	1.89	0.72
1:C:185:ILE:HG12	1:D:185:ILE:HG12	1.72	0.72
1:A:94:LYS:O	1:A:98:ILE:HG12	1.91	0.71
1:D:164:VAL:O	1:D:192:ARG:NH2	2.23	0.70
1:I:164:VAL:O	1:I:192:ARG:NH2	2.26	0.69
1:J:158:PHE:C	1:J:160:ILE:H	1.99	0.69
1:G:162:GLY:HA3	1:G:210:MSE:HE2	1.76	0.68
1:B:117:ARG:HH22	1:B:202:ASN:ND2	1.92	0.68
1:A:68:VAL:HG21	1:A:75:LEU:HD11	1.75	0.67
1:F:164:VAL:O	1:F:192:ARG:NH2	2.27	0.67
1:H:94:LYS:O	1:H:98:ILE:HG12	1.94	0.67
1:I:119:ILE:HD12	1:I:139:PRO:HG3	1.75	0.67
1:D:117:ARG:NH2	1:D:202:ASN:HD21	1.93	0.66
1:A:155:CYS:HB3	1:A:158:PHE:HD1	1.60	0.66
1:B:117:ARG:HA	1:B:201:LEU:HD22	1.78	0.65
1:F:158:PHE:O	3:F:5305:HOH:O	2.14	0.65
1:B:158:PHE:HB3	1:B:160:ILE:HB	1.79	0.65
1:A:199:ASP:OD2	1:A:209:LYS:NZ	2.25	0.65
1:I:100:TYR:OH	1:I:133:GLU:OE2	2.15	0.65
1:B:65:ILE:HA	1:B:98:ILE:HD12	1.77	0.64
1:I:134:ARG:HH21	1:I:136:PHE:HZ	1.45	0.64
1:D:117:ARG:HA	1:D:201:LEU:HD22	1.80	0.63
1:F:185:ILE:HG12	1:H:185:ILE:HG12	1.79	0.63
1:H:90:ARG:HA	1:H:93:ARG:HB2	1.81	0.63
1:E:166:ARG:NH2	1:E:199:ASP:OD2	2.33	0.62
1:D:158:PHE:HB3	1:D:160:ILE:HB	1.80	0.62
1:J:199:ASP:OD2	1:J:209:LYS:NZ	2.30	0.62
1:F:143:GLU:HG3	1:H:142:VAL:HG11	1.80	0.61
1:I:64:GLU:HB3	1:I:66:LYS:H	1.63	0.61
1:F:66:LYS:O	1:F:112:VAL:HA	2.00	0.60
1:C:117:ARG:NH2	1:C:202:ASN:HD21	1.98	0.59
1:B:231:ALA:HA	1:B:234:SER:HB3	1.83	0.59
1:A:164:VAL:O	1:A:192:ARG:NH2	2.36	0.58
1:E:70:TYR:CD1	1:E:70:TYR:C	2.81	0.58
1:I:158:PHE:HB3	1:I:160:ILE:HB	1.84	0.58
1:G:220:PRO:O	1:G:224:GLU:HG3	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:227:ARG:HB3	1:B:241:HIS:CE1	2.39	0.58
1:A:74:ILE:HD12	1:A:77:ARG:HH11	1.68	0.58
1:E:149:LEU:O	1:E:192:ARG:NH2	2.27	0.58
1:C:117:ARG:HH21	1:C:202:ASN:ND2	2.02	0.57
1:H:199:ASP:OD2	1:H:209:LYS:NZ	2.26	0.57
1:I:64:GLU:HB3	1:I:66:LYS:HB2	1.86	0.57
1:D:68:VAL:HG22	1:D:111:GLN:HA	1.87	0.57
1:I:117:ARG:HH22	1:I:202:ASN:ND2	2.02	0.57
1:H:105:ILE:HD11	1:H:123:ALA:HA	1.87	0.56
1:C:114:ILE:HG22	1:C:116:LYS:HB2	1.86	0.56
1:C:143:GLU:HB2	1:C:171:SER:HB2	1.87	0.56
1:G:164:VAL:O	1:G:192:ARG:NH2	2.39	0.56
1:D:143:GLU:HB2	1:D:171:SER:HB3	1.88	0.55
1:E:148:LYS:HA	1:E:167:PRO:HA	1.89	0.55
1:I:145:SER:HB3	1:I:169:ILE:HB	1.90	0.55
1:J:67:ILE:HD11	1:J:156:LEU:HB3	1.88	0.54
1:E:158:PHE:C	1:E:160:ILE:H	2.15	0.54
1:B:63:ASP:O	1:B:94:LYS:HD3	2.08	0.54
1:A:117:ARG:NH2	1:A:202:ASN:OD1	2.35	0.54
1:J:71:PRO:HG2	1:J:229:TYR:HB2	1.90	0.54
1:B:162:GLY:HA3	1:B:210:MSE:HE3	1.90	0.54
1:I:229:TYR:O	1:I:232:THR:HB	2.07	0.54
1:I:70:TYR:C	1:I:70:TYR:CD2	2.86	0.53
1:D:169:ILE:HG12	1:D:187:LYS:HG3	1.90	0.53
1:J:158:PHE:HB3	1:J:160:ILE:HB	1.90	0.53
1:E:167:PRO:HD2	1:E:195:GLN:OE1	2.09	0.53
1:F:105:ILE:O	1:F:121:TRP:HB2	2.09	0.53
1:I:219:ARG:NH1	1:J:161:GLU:OE2	2.41	0.53
1:E:185:ILE:HG12	1:J:185:ILE:HG12	1.92	0.52
1:J:70:TYR:C	1:J:70:TYR:CD2	2.86	0.52
1:E:158:PHE:O	1:E:160:ILE:N	2.32	0.52
1:G:114:ILE:HG22	1:G:116:LYS:HB2	1.91	0.52
1:I:158:PHE:C	1:I:160:ILE:H	2.16	0.52
1:B:96:PHE:CZ	1:B:133:GLU:HB3	2.45	0.52
1:C:210:MSE:HB3	1:C:215:LYS:HG2	1.92	0.52
1:J:153:GLU:OE2	1:J:166:ARG:NH1	2.43	0.52
1:I:117:ARG:HH22	1:I:202:ASN:HD21	1.58	0.51
1:B:95:MSE:HG2	1:B:112:VAL:HG21	1.92	0.51
1:A:185:ILE:HG12	1:B:185:ILE:HG12	1.93	0.51
1:F:105:ILE:HD11	1:F:193:ILE:HD11	1.92	0.51
1:F:162:GLY:HA3	1:F:210:MSE:HE2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:63:ASP:O	1:G:94:LYS:HD3	2.10	0.51
1:B:155:CYS:HB2	1:B:196:HIS:CE1	2.46	0.51
1:I:117:ARG:HA	1:I:201:LEU:HD22	1.92	0.51
1:J:158:PHE:O	1:J:160:ILE:N	2.34	0.51
1:C:166:ARG:HG3	1:C:192:ARG:CZ	2.41	0.51
1:G:158:PHE:HB3	1:G:160:ILE:HB	1.92	0.50
1:A:74:ILE:HG21	1:A:113:ASN:ND2	2.27	0.50
1:E:149:LEU:HD23	1:E:151:LEU:HD11	1.93	0.50
1:I:165:GLU:OE1	3:I:8304:HOH:O	2.20	0.50
1:H:223:ASN:HA	1:H:226:ILE:HD12	1.93	0.50
1:F:107:LEU:HD13	1:F:156:LEU:HD23	1.93	0.50
1:J:218:VAL:HG13	1:J:222:LEU:HD22	1.94	0.50
1:E:66:LYS:O	1:E:112:VAL:HA	2.12	0.50
1:E:193:ILE:O	1:E:197:GLU:HG2	2.11	0.50
1:F:207:ILE:HB	1:F:215:LYS:HE2	1.94	0.50
1:G:158:PHE:C	1:G:160:ILE:H	2.18	0.50
1:B:70:TYR:HB2	1:B:158:PHE:CE2	2.47	0.49
1:F:64:GLU:C	1:F:66:LYS:H	2.21	0.49
1:B:210:MSE:HE3	1:B:214:ASP:HB3	1.94	0.49
1:D:96:PHE:CZ	1:D:133:GLU:HB3	2.46	0.49
1:G:95:MSE:HE2	1:G:120:VAL:HG22	1.93	0.49
1:A:230:LYS:HG3	1:D:149:LEU:HD11	1.94	0.49
1:J:70:TYR:HB2	1:J:158:PHE:CE2	2.47	0.49
1:B:142:VAL:HG22	1:B:171:SER:HB3	1.94	0.49
1:H:210:MSE:HE3	1:H:214:ASP:HB3	1.95	0.49
1:E:158:PHE:HB3	1:E:160:ILE:HB	1.95	0.49
1:G:153:GLU:OE2	1:G:166:ARG:NH1	2.46	0.49
1:H:143:GLU:HB2	1:H:171:SER:HB3	1.94	0.49
1:E:65:ILE:O	1:E:65:ILE:HG13	2.12	0.49
1:E:160:ILE:HD11	1:E:221:LYS:HB2	1.94	0.48
1:A:65:ILE:HA	1:A:98:ILE:HD11	1.95	0.48
1:G:150:LYS:HG2	1:G:165:GLU:HB2	1.95	0.48
1:J:158:PHE:C	1:J:160:ILE:N	2.71	0.48
1:J:162:GLY:HA3	1:J:210:MSE:HE3	1.96	0.48
1:A:79:SER:HB3	1:A:117:ARG:HG2	1.96	0.48
1:B:148:LYS:HA	1:B:167:PRO:HA	1.94	0.48
1:A:122:ASN:HB3	1:A:125:TYR:HB3	1.95	0.48
1:A:226:ILE:O	1:A:230:LYS:HG2	2.13	0.47
1:B:158:PHE:C	1:B:160:ILE:H	2.21	0.47
1:D:158:PHE:C	1:D:160:ILE:H	2.22	0.47
1:J:95:MSE:HE2	1:J:95:MSE:HB3	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:96:PHE:CZ	1:J:133:GLU:HB3	2.49	0.47
1:D:117:ARG:NH2	1:D:202:ASN:ND2	2.62	0.47
1:H:114:ILE:HG22	1:H:116:LYS:HB2	1.97	0.47
1:E:220:PRO:O	1:E:224:GLU:HG3	2.14	0.47
1:B:96:PHE:CZ	1:B:120:VAL:HB	2.49	0.47
1:B:75:LEU:HA	1:B:75:LEU:HD12	1.59	0.47
1:B:109:ALA:HB1	1:B:114:ILE:HB	1.96	0.47
1:E:96:PHE:CZ	1:E:133:GLU:HB3	2.50	0.47
1:I:223:ASN:O	1:I:227:ARG:HG3	2.15	0.47
1:J:64:GLU:C	1:J:66:LYS:H	2.22	0.47
1:A:96:PHE:CZ	1:A:120:VAL:HB	2.50	0.47
1:C:167:PRO:HD2	1:C:195:GLN:OE1	2.13	0.47
1:C:162:GLY:HA3	1:C:210:MSE:HE2	1.96	0.47
1:J:79:SER:HB3	1:J:117:ARG:HG2	1.97	0.47
1:E:142:VAL:HG11	1:J:143:GLU:HG3	1.97	0.47
1:F:117:ARG:HA	1:F:201:LEU:HD22	1.96	0.47
1:J:107:LEU:HD12	1:J:107:LEU:HA	1.69	0.47
1:A:193:ILE:O	1:A:197:GLU:HG2	2.14	0.46
1:A:234:SER:O	1:A:234:SER:OG	2.31	0.46
1:E:192:ARG:HD3	1:E:192:ARG:HA	1.71	0.46
1:I:107:LEU:HD11	1:I:111:GLN:HB2	1.97	0.46
1:J:192:ARG:HD2	1:J:192:ARG:HA	1.68	0.46
1:D:231:ALA:O	1:D:235:GLU:HG3	2.16	0.46
1:B:132:ASN:HD22	1:B:132:ASN:HA	1.54	0.46
1:B:160:ILE:HG23	1:B:218:VAL:HG21	1.98	0.46
1:H:207:ILE:HB	1:H:215:LYS:HE2	1.96	0.46
1:J:75:LEU:HD12	1:J:75:LEU:HA	1.59	0.46
1:B:64:GLU:HB3	1:B:66:LYS:HB2	1.97	0.46
1:E:119:ILE:HG13	1:E:201:LEU:HD11	1.97	0.46
1:F:119:ILE:HD13	1:F:119:ILE:HG21	1.60	0.46
1:I:66:LYS:HE2	1:I:66:LYS:HB3	1.77	0.46
1:A:153:GLU:OE2	1:A:166:ARG:NH1	2.49	0.46
1:B:187:LYS:HE2	1:B:187:LYS:HB3	1.78	0.46
1:H:95:MSE:HG2	1:H:112:VAL:HG21	1.98	0.46
1:A:183:LEU:HD21	1:B:143:GLU:HB3	1.98	0.46
1:D:206:PHE:H	1:D:206:PHE:HD2	1.64	0.46
1:H:65:ILE:HA	1:H:98:ILE:HD12	1.97	0.45
1:A:74:ILE:HG21	1:A:113:ASN:HD22	1.81	0.45
1:E:183:LEU:HB3	1:J:185:ILE:HD12	1.98	0.45
1:I:218:VAL:HG13	1:I:222:LEU:HD22	1.98	0.45
1:J:160:ILE:HG23	1:J:218:VAL:HG21	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:162:GLY:HA3	1:E:210:MSE:HE2	1.97	0.45
1:E:109:ALA:HB1	1:E:114:ILE:HB	1.98	0.45
1:F:66:LYS:HD2	1:F:66:LYS:HA	1.80	0.45
1:A:107:LEU:HD21	1:A:112:VAL:HG13	1.99	0.45
1:C:86:ASP:HB3	1:C:88:LEU:H	1.81	0.45
1:F:166:ARG:NH2	1:F:199:ASP:OD2	2.50	0.45
1:C:155:CYS:HB2	1:C:196:HIS:CE1	2.50	0.45
1:C:205:LEU:HD12	1:C:205:LEU:HA	1.84	0.45
1:J:205:LEU:HD12	1:J:205:LEU:HA	1.86	0.45
1:A:66:LYS:HD2	1:A:66:LYS:HA	1.79	0.45
1:D:148:LYS:HA	1:D:167:PRO:HA	1.99	0.45
1:A:152:ILE:HG13	1:A:163:LYS:HG2	1.99	0.44
1:A:210:MSE:HB3	1:A:215:LYS:HG2	1.99	0.44
1:H:222:LEU:HD23	1:I:152:ILE:HD13	1.98	0.44
1:E:75:LEU:HD22	1:E:75:LEU:HA	1.73	0.44
1:F:153:GLU:OE1	1:F:166:ARG:HD2	2.16	0.44
1:C:65:ILE:HG22	1:C:66:LYS:N	2.32	0.44
1:F:72:ASP:HA	1:F:73:PRO:HD3	1.84	0.44
1:B:166:ARG:NH2	1:B:199:ASP:OD2	2.50	0.44
1:D:156:LEU:HD12	1:D:156:LEU:HA	1.51	0.44
1:G:143:GLU:HG3	1:I:142:VAL:HG11	1.99	0.44
1:I:76:ARG:CZ	1:I:205:LEU:HD22	2.48	0.44
1:C:189:ILE:HG12	1:C:193:ILE:HD12	2.00	0.44
1:F:105:ILE:HG21	1:F:105:ILE:HD13	1.62	0.44
1:H:79:SER:HB3	1:H:117:ARG:HG2	2.00	0.44
1:H:192:ARG:HH11	1:H:192:ARG:HD3	1.56	0.44
1:J:153:GLU:OE1	1:J:166:ARG:HD2	2.18	0.44
1:H:159:GLY:N	3:H:7303:HOH:O	2.50	0.44
1:G:64:GLU:HB3	1:G:66:LYS:HB2	1.99	0.43
1:H:234:SER:O	1:H:234:SER:OG	2.30	0.43
1:A:207:ILE:O	1:A:215:LYS:HE3	2.17	0.43
1:B:90:ARG:O	1:B:94:LYS:HG2	2.18	0.43
1:F:93:ARG:HD2	1:F:93:ARG:HA	1.69	0.43
1:F:119:ILE:HD11	1:F:198:PHE:HA	1.99	0.43
1:H:192:ARG:HD3	1:H:192:ARG:HA	1.78	0.43
1:I:230:LYS:HA	1:I:230:LYS:HD2	1.83	0.43
1:J:145:SER:HB3	1:J:169:ILE:H	1.83	0.43
1:J:175:TYR:OH	1:J:181:LYS:HD2	2.18	0.43
1:A:145:SER:HB2	1:A:169:ILE:HB	1.99	0.43
1:C:225:LEU:HD12	1:C:225:LEU:HA	1.90	0.43
1:G:171:SER:HA	1:G:185:ILE:HD13	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:155:CYS:HB3	1:H:158:PHE:HD1	1.83	0.43
1:H:193:ILE:O	1:H:194:PHE:C	2.60	0.43
1:J:113:ASN:O	1:J:113:ASN:CG	2.62	0.43
1:A:153:GLU:OE1	1:A:166:ARG:HD2	2.19	0.43
1:G:93:ARG:HD2	1:G:93:ARG:HA	1.82	0.43
1:G:98:ILE:HD13	1:G:98:ILE:N	2.34	0.43
1:B:234:SER:O	1:B:234:SER:OG	2.34	0.43
1:E:185:ILE:HD12	1:J:183:LEU:HB3	2.01	0.43
1:D:70:TYR:CD2	1:D:70:TYR:C	2.96	0.43
1:C:149:LEU:HD13	1:C:149:LEU:HA	1.80	0.43
1:J:217:LYS:HE3	1:J:217:LYS:HB3	1.69	0.43
1:B:221:LYS:H	1:B:221:LYS:HG2	1.59	0.42
1:C:218:VAL:HG22	1:C:221:LYS:HG3	2.01	0.42
1:A:70:TYR:HB2	1:A:158:PHE:CE2	2.54	0.42
1:F:70:TYR:HB2	1:F:158:PHE:CE2	2.53	0.42
1:I:111:GLN:HE21	1:I:111:GLN:HB3	1.62	0.42
1:A:133:GLU:C	1:A:134:ARG:HG2	2.44	0.42
1:G:88:LEU:C	1:G:90:ARG:N	2.76	0.42
1:I:75:LEU:HD12	1:I:75:LEU:HA	1.62	0.42
1:D:155:CYS:HB2	1:D:196:HIS:CE1	2.54	0.42
1:F:137:ILE:HD11	1:F:177:ILE:HD12	2.00	0.42
1:I:162:GLY:HA3	1:I:210:MSE:HE2	2.00	0.42
1:G:185:ILE:HG12	1:I:185:ILE:HG12	2.00	0.42
1:H:88:LEU:O	1:H:90:ARG:N	2.52	0.42
1:H:98:ILE:HG13	1:H:112:VAL:HG11	2.02	0.42
1:J:224:GLU:HA	1:J:227:ARG:HG3	2.02	0.42
1:C:94:LYS:HA	1:C:94:LYS:HD3	1.73	0.42
1:C:160:ILE:HG23	1:C:218:VAL:HG21	2.02	0.42
1:A:192:ARG:HD2	1:A:192:ARG:HA	1.81	0.42
1:B:98:ILE:H	1:B:98:ILE:HG12	1.58	0.42
1:E:95:MSE:HE2	1:E:95:MSE:HB3	1.93	0.42
1:D:202:ASN:HD22	1:D:202:ASN:HA	1.71	0.42
1:H:93:ARG:HD2	1:H:93:ARG:HA	1.84	0.42
1:H:99:MSE:HB2	1:H:107:LEU:HD22	2.01	0.42
1:J:76:ARG:HD3	1:J:205:LEU:HD22	2.01	0.42
1:G:146:LEU:HA	1:G:146:LEU:HD12	1.57	0.42
1:C:138:ASN:HB2	1:C:175:TYR:CD2	2.55	0.41
1:C:192:ARG:HD3	1:C:192:ARG:HA	1.82	0.41
1:D:96:PHE:CZ	1:D:120:VAL:HB	2.54	0.41
1:A:119:ILE:HG21	1:A:119:ILE:HD13	1.81	0.41
1:B:155:CYS:HB2	1:B:196:HIS:NE2	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:132:ASN:HD22	1:D:132:ASN:HA	1.70	0.41
1:I:187:LYS:HE2	1:I:187:LYS:HB3	1.35	0.41
1:J:82:THR:HB	1:J:138:ASN:HD21	1.85	0.41
1:J:130:GLU:H	1:J:130:GLU:HG3	1.54	0.41
1:A:146:LEU:HD12	1:A:146:LEU:HA	1.94	0.41
1:E:199:ASP:OD2	1:E:209:LYS:NZ	2.45	0.41
1:H:109:ALA:HB1	1:H:114:ILE:HB	2.02	0.41
1:I:139:PRO:HB2	1:I:194:PHE:HZ	1.85	0.41
1:A:223:ASN:HA	1:A:226:ILE:HD12	2.03	0.41
1:C:160:ILE:CG2	1:C:161:GLU:N	2.83	0.41
1:D:192:ARG:HD2	1:D:192:ARG:HA	1.49	0.41
1:G:66:LYS:HB3	1:G:66:LYS:HE2	1.64	0.41
1:B:236:GLU:HA	1:B:237:PRO:HD3	1.78	0.41
1:G:119:ILE:HD13	1:G:119:ILE:HG21	1.89	0.41
1:G:221:LYS:H	1:G:221:LYS:HG2	1.40	0.41
1:G:149:LEU:HG	1:G:151:LEU:HD21	2.03	0.41
1:G:160:ILE:HD11	1:G:221:LYS:HB2	2.02	0.41
1:C:156:LEU:HD12	1:C:156:LEU:HA	1.85	0.41
1:I:92:VAL:HG13	1:I:120:VAL:HG11	2.03	0.41
1:G:111:GLN:HE21	1:G:111:GLN:HB3	1.72	0.41
1:H:234:SER:C	1:H:236:GLU:H	2.29	0.41
1:I:63:ASP:HA	1:I:114:ILE:HD11	2.02	0.41
1:I:114:ILE:HG21	1:I:114:ILE:HD13	1.83	0.41
1:E:72:ASP:HA	1:E:73:PRO:HD3	1.85	0.40
1:H:150:LYS:HG2	1:H:165:GLU:HB2	2.01	0.40
1:C:227:ARG:HH11	1:C:227:ARG:HD3	1.72	0.40
1:F:80:GLU:C	1:F:117:ARG:HG3	2.46	0.40
1:G:76:ARG:HH11	1:G:76:ARG:HD3	1.75	0.40
1:H:119:ILE:HD13	1:H:119:ILE:HG21	1.72	0.40
1:E:176:ASP:OD1	1:E:179:GLY:N	2.54	0.40
1:G:75:LEU:HA	1:G:75:LEU:HD12	1.74	0.40
1:H:96:PHE:CZ	1:H:120:VAL:HB	2.56	0.40
1:J:146:LEU:HD13	1:J:146:LEU:HA	1.91	0.40
1:A:158:PHE:HB3	1:A:160:ILE:HB	2.02	0.40
1:B:172:ILE:HD12	1:B:194:PHE:CE2	2.57	0.40
1:G:128:ARG:H	1:G:128:ARG:HG3	1.50	0.40
1:C:199:ASP:HB3	1:C:204:THR:HB	2.02	0.40
1:I:109:ALA:HB1	1:I:114:ILE:HB	2.03	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	177/183 (97%)	164 (93%)	12 (7%)	1 (1%)	21	51
1	B	177/183 (97%)	157 (89%)	20 (11%)	0	100	100
1	C	166/183 (91%)	154 (93%)	11 (7%)	1 (1%)	21	51
1	D	171/183 (93%)	153 (90%)	17 (10%)	1 (1%)	21	51
1	E	174/183 (95%)	160 (92%)	11 (6%)	3 (2%)	7	25
1	F	170/183 (93%)	157 (92%)	8 (5%)	5 (3%)	3	13
1	G	170/183 (93%)	149 (88%)	20 (12%)	1 (1%)	21	51
1	H	171/183 (93%)	154 (90%)	15 (9%)	2 (1%)	10	34
1	I	170/183 (93%)	160 (94%)	8 (5%)	2 (1%)	10	34
1	J	170/183 (93%)	157 (92%)	12 (7%)	1 (1%)	21	51
All	All	1716/1830 (94%)	1565 (91%)	134 (8%)	17 (1%)	12	38

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	89	LYS
1	J	159	GLY
1	A	117	ARG
1	C	65	ILE
1	I	70	TYR
1	D	70	TYR
1	E	217	LYS
1	F	86	ASP
1	E	159	GLY
1	F	128	ARG
1	E	70	TYR
1	F	65	ILE
1	F	230	LYS
1	G	88	LEU

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Mol	Chain	Res	Type
1	F	70	TYR
1	H	65	ILE
1	I	159	GLY

5.3.2 Protein sidechains [i](#)

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	168/169 (99%)	133 (79%)	35 (21%)	1	4
1	B	168/169 (99%)	133 (79%)	35 (21%)	1	4
1	C	157/169 (93%)	126 (80%)	31 (20%)	1	5
1	D	162/169 (96%)	132 (82%)	30 (18%)	1	6
1	E	165/169 (98%)	126 (76%)	39 (24%)	1	3
1	F	161/169 (95%)	126 (78%)	35 (22%)	1	3
1	G	161/169 (95%)	122 (76%)	39 (24%)	1	2
1	H	162/169 (96%)	127 (78%)	35 (22%)	1	3
1	I	161/169 (95%)	128 (80%)	33 (20%)	1	4
1	J	161/169 (95%)	124 (77%)	37 (23%)	1	3
All	All	1626/1690 (96%)	1277 (78%)	349 (22%)	1	4

All (349) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	ASP
1	A	64	GLU
1	A	65	ILE
1	A	66	LYS
1	A	67	ILE
1	A	75	LEU
1	A	78	ARG
1	A	79	SER
1	A	80	GLU
1	A	82	THR

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Mol	Chain	Res	Type
1	A	85	ASP
1	A	86	ASP
1	A	92	VAL
1	A	108	SER
1	A	113	ASN
1	A	116	LYS
1	A	124	LEU
1	A	134	ARG
1	A	137	ILE
1	A	144	GLN
1	A	146	LEU
1	A	152	ILE
1	A	156	LEU
1	A	160	ILE
1	A	170	VAL
1	A	187	LYS
1	A	189	ILE
1	A	205	LEU
1	A	206	PHE
1	A	210	MSE
1	A	215	LYS
1	A	218	VAL
1	A	222	LEU
1	A	228	ASP
1	A	239	GLU
1	B	63	ASP
1	B	66	LYS
1	B	67	ILE
1	B	74	ILE
1	B	75	LEU
1	B	81	VAL
1	B	86	ASP
1	B	92	VAL
1	B	93	ARG
1	B	94	LYS
1	B	98	ILE
1	B	128	ARG
1	B	130	GLU
1	B	132	ASN
1	B	137	ILE
1	B	142	VAL
1	B	146	LEU

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Mol	Chain	Res	Type
1	B	156	LEU
1	B	160	ILE
1	B	161	GLU
1	B	166	ARG
1	B	169	ILE
1	B	170	VAL
1	B	187	LYS
1	B	191	SER
1	B	205	LEU
1	B	206	PHE
1	B	208	ASP
1	B	211	THR
1	B	217	LYS
1	B	218	VAL
1	B	221	LYS
1	B	222	LEU
1	B	226	ILE
1	B	236	GLU
1	C	64	GLU
1	C	69	LYS
1	C	75	LEU
1	C	86	ASP
1	C	103	LYS
1	C	116	LYS
1	C	126	GLU
1	C	140	SER
1	C	144	GLN
1	C	145	SER
1	C	146	LEU
1	C	149	LEU
1	C	156	LEU
1	C	160	ILE
1	C	166	ARG
1	C	172	ILE
1	C	177	ILE
1	C	183	LEU
1	C	187	LYS
1	C	189	ILE
1	C	192	ARG
1	C	205	LEU
1	C	215	LYS
1	C	217	LYS

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Mol	Chain	Res	Type
1	C	218	VAL
1	C	219	ARG
1	C	221	LYS
1	C	222	LEU
1	C	225	LEU
1	C	227	ARG
1	C	230	LYS
1	D	72	ASP
1	D	75	LEU
1	D	78	ARG
1	D	82	THR
1	D	85	ASP
1	D	86	ASP
1	D	90	ARG
1	D	92	VAL
1	D	103	LYS
1	D	130	GLU
1	D	132	ASN
1	D	142	VAL
1	D	146	LEU
1	D	149	LEU
1	D	150	LYS
1	D	156	LEU
1	D	160	ILE
1	D	166	ARG
1	D	177	ILE
1	D	183	LEU
1	D	189	ILE
1	D	192	ARG
1	D	208	ASP
1	D	210	MSE
1	D	215	LYS
1	D	217	LYS
1	D	218	VAL
1	D	222	LEU
1	D	225	LEU
1	D	227	ARG
1	E	63	ASP
1	E	64	GLU
1	E	66	LYS
1	E	67	ILE
1	E	75	LEU

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Mol	Chain	Res	Type
1	E	79	SER
1	E	82	THR
1	E	85	ASP
1	E	86	ASP
1	E	92	VAL
1	E	94	LYS
1	E	98	ILE
1	E	114	ILE
1	E	127	LYS
1	E	128	ARG
1	E	129	LYS
1	E	130	GLU
1	E	137	ILE
1	E	146	LEU
1	E	150	LYS
1	E	156	LEU
1	E	160	ILE
1	E	161	GLU
1	E	169	ILE
1	E	170	VAL
1	E	183	LEU
1	E	187	LYS
1	E	191	SER
1	E	192	ARG
1	E	205	LEU
1	E	206	PHE
1	E	211	THR
1	E	215	LYS
1	E	218	VAL
1	E	221	LYS
1	E	222	LEU
1	E	225	LEU
1	E	236	GLU
1	E	238	LEU
1	F	65	ILE
1	F	66	LYS
1	F	68	VAL
1	F	69	LYS
1	F	75	LEU
1	F	79	SER
1	F	89	LYS
1	F	93	ARG

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Mol	Chain	Res	Type
1	F	94	LYS
1	F	101	GLU
1	F	103	LYS
1	F	105	ILE
1	F	116	LYS
1	F	117	ARG
1	F	119	ILE
1	F	132	ASN
1	F	134	ARG
1	F	142	VAL
1	F	144	GLN
1	F	146	LEU
1	F	149	LEU
1	F	156	LEU
1	F	160	ILE
1	F	183	LEU
1	F	187	LYS
1	F	189	ILE
1	F	191	SER
1	F	213	VAL
1	F	215	LYS
1	F	218	VAL
1	F	219	ARG
1	F	222	LEU
1	F	225	LEU
1	F	228	ASP
1	F	232	THR
1	G	63	ASP
1	G	64	GLU
1	G	75	LEU
1	G	79	SER
1	G	81	VAL
1	G	82	THR
1	G	86	ASP
1	G	89	LYS
1	G	94	LYS
1	G	98	ILE
1	G	113	ASN
1	G	116	LYS
1	G	127	LYS
1	G	130	GLU
1	G	132	ASN

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Mol	Chain	Res	Type
1	G	134	ARG
1	G	137	ILE
1	G	142	VAL
1	G	146	LEU
1	G	149	LEU
1	G	156	LEU
1	G	160	ILE
1	G	172	ILE
1	G	177	ILE
1	G	178	ASN
1	G	183	LEU
1	G	187	LYS
1	G	189	ILE
1	G	192	ARG
1	G	205	LEU
1	G	206	PHE
1	G	210	MSE
1	G	211	THR
1	G	215	LYS
1	G	216	LYS
1	G	218	VAL
1	G	221	LYS
1	G	222	LEU
1	G	225	LEU
1	H	65	ILE
1	H	67	ILE
1	H	68	VAL
1	H	75	LEU
1	H	78	ARG
1	H	82	THR
1	H	85	ASP
1	H	86	ASP
1	H	94	LYS
1	H	116	LYS
1	H	119	ILE
1	H	128	ARG
1	H	130	GLU
1	H	137	ILE
1	H	145	SER
1	H	146	LEU
1	H	149	LEU
1	H	156	LEU

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Mol	Chain	Res	Type
1	H	160	ILE
1	H	170	VAL
1	H	172	ILE
1	H	183	LEU
1	H	187	LYS
1	H	189	ILE
1	H	192	ARG
1	H	205	LEU
1	H	206	PHE
1	H	210	MSE
1	H	213	VAL
1	H	215	LYS
1	H	216	LYS
1	H	218	VAL
1	H	222	LEU
1	H	230	LYS
1	H	236	GLU
1	I	64	GLU
1	I	65	ILE
1	I	69	LYS
1	I	74	ILE
1	I	75	LEU
1	I	81	VAL
1	I	85	ASP
1	I	86	ASP
1	I	93	ARG
1	I	98	ILE
1	I	103	LYS
1	I	116	LYS
1	I	128	ARG
1	I	129	LYS
1	I	146	LEU
1	I	149	LEU
1	I	156	LEU
1	I	160	ILE
1	I	168	SER
1	I	170	VAL
1	I	177	ILE
1	I	183	LEU
1	I	187	LYS
1	I	189	ILE
1	I	193	ILE

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Mol	Chain	Res	Type
1	I	205	LEU
1	I	210	MSE
1	I	212	GLN
1	I	213	VAL
1	I	215	LYS
1	I	218	VAL
1	I	222	LEU
1	I	232	THR
1	J	63	ASP
1	J	64	GLU
1	J	65	ILE
1	J	66	LYS
1	J	67	ILE
1	J	69	LYS
1	J	75	LEU
1	J	79	SER
1	J	81	VAL
1	J	85	ASP
1	J	86	ASP
1	J	94	LYS
1	J	98	ILE
1	J	114	ILE
1	J	128	ARG
1	J	130	GLU
1	J	140	SER
1	J	142	VAL
1	J	145	SER
1	J	146	LEU
1	J	156	LEU
1	J	160	ILE
1	J	161	GLU
1	J	170	VAL
1	J	183	LEU
1	J	187	LYS
1	J	189	ILE
1	J	205	LEU
1	J	206	PHE
1	J	207	ILE
1	J	210	MSE
1	J	211	THR
1	J	215	LYS
1	J	217	LYS

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Mol	Chain	Res	Type
1	J	222	LEU
1	J	225	LEU
1	J	234	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	ASN
1	A	144	GLN
1	B	113	ASN
1	B	144	GLN
1	B	182	HIS
1	B	202	ASN
1	C	87	ASN
1	C	113	ASN
1	C	202	ASN
1	D	111	GLN
1	D	132	ASN
1	D	144	GLN
1	D	202	ASN
1	E	111	GLN
1	E	144	GLN
1	E	202	ASN
1	F	111	GLN
1	F	113	ASN
1	F	132	ASN
1	F	144	GLN
1	F	202	ASN
1	G	113	ASN
1	G	132	ASN
1	G	144	GLN
1	G	195	GLN
1	G	202	ASN
1	G	223	ASN
1	H	144	GLN
1	H	178	ASN
1	I	113	ASN
1	I	144	GLN
1	I	178	ASN
1	I	202	ASN
1	I	233	HIS
1	J	132	ASN

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Mol	Chain	Res	Type
1	J	202	ASN
1	J	233	HIS

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 ⓘ

Of 10 ligands modelled in this entry, 10 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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(Å²)	Q<0.9
1	A	176/183 (96%)	-1.47	0	100	100	32, 54, 96, 104	0
1	B	176/183 (96%)	-1.50	0	100	100	22, 46, 96, 106	0
1	C	165/183 (90%)	-1.58	0	100	100	13, 38, 81, 92	0
1	D	170/183 (92%)	-1.48	0	100	100	21, 47, 92, 102	0
1	E	173/183 (94%)	-1.49	0	100	100	23, 47, 99, 106	0
1	F	169/183 (92%)	-1.60	0	100	100	15, 32, 80, 104	0
1	G	169/183 (92%)	-1.51	0	100	100	18, 43, 88, 104	0
1	H	170/183 (92%)	-1.59	0	100	100	17, 39, 82, 106	0
1	I	169/183 (92%)	-1.51	0	100	100	18, 42, 90, 105	0
1	J	169/183 (92%)	-1.48	1 (0%)	85	80	20, 42, 89, 105	0
All	All	1706/1830 (93%)	-1.52	1 (0%)	92	90	13, 43, 92, 106	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	63	ASP	2.3

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CO	A	300	1/1	1.00	0.02	46,46,46,46	0
2	CO	B	1300	1/1	1.00	0.01	39,39,39,39	0
2	CO	C	2300	1/1	1.00	0.01	35,35,35,35	0
2	CO	D	3300	1/1	1.00	0.02	34,34,34,34	0
2	CO	E	4300	1/1	1.00	0.01	39,39,39,39	0
2	CO	F	5300	1/1	1.00	0.00	24,24,24,24	0
2	CO	G	6300	1/1	1.00	0.01	30,30,30,30	0
2	CO	H	7300	1/1	1.00	0.01	34,34,34,34	0
2	CO	I	8300	1/1	1.00	0.00	45,45,45,45	0
2	CO	J	9300	1/1	1.00	0.01	44,44,44,44	0

6.5 Other polymers

There are no such residues in this entry.