



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:01 AM UTC

PDB ID : 3TDM / pdb_00003tdm
Title : Computationally designed TIM-barrel protein, HalfFLR
Authors : Harp, J.M.; Fortenberry, C.; Bowman, E.; Profitt, W.; Dorr, B.; Mizoue, L.
Deposited on : 2011-08-11
Resolution : 2.40 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

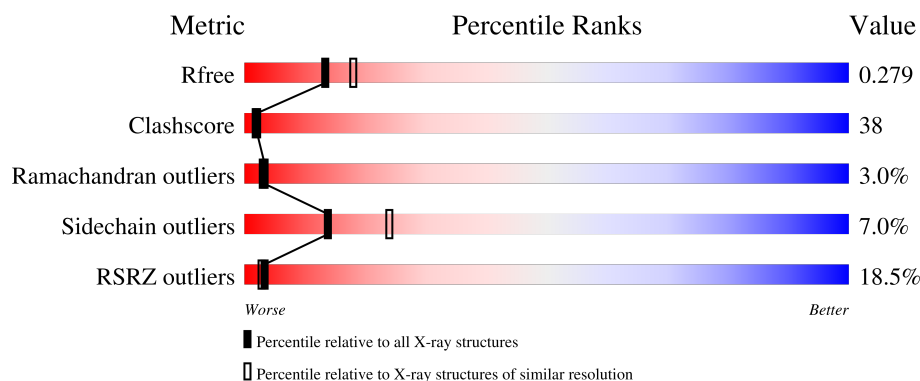
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	126	6% 56% 37% • 5%
1	B	126	11% 52% 31% 12% 5%
1	C	126	21% 48% 29% 10% • 12%
1	D	126	32% 48% 37% 9% • 5%

2 Entry composition [i](#)

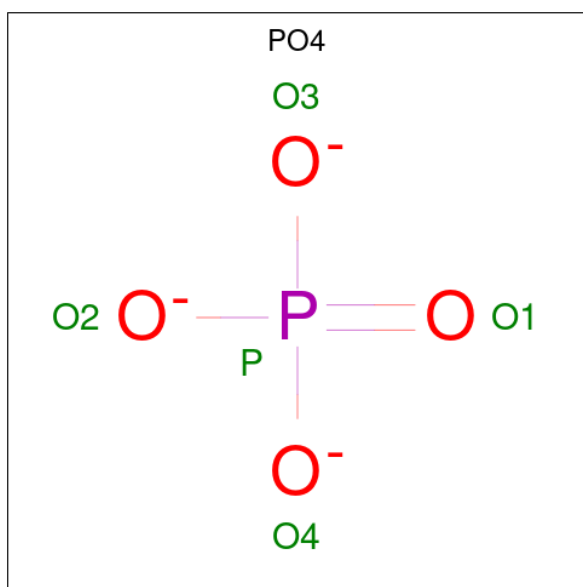
There are 3 unique types of molecules in this entry. The entry contains 3684 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Computationally designed two-fold symmetric TIM-barrel protein, FLR (half molecule).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	120	Total	C	N	O	S	0	0	0
			916	582	157	174	3			
1	B	120	Total	C	N	O	S	0	0	0
			918	583	157	175	3			
1	C	111	Total	C	N	O	S	0	0	0
			846	535	146	162	3			
1	D	120	Total	C	N	O	S	0	0	0
			918	583	157	175	3			

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		

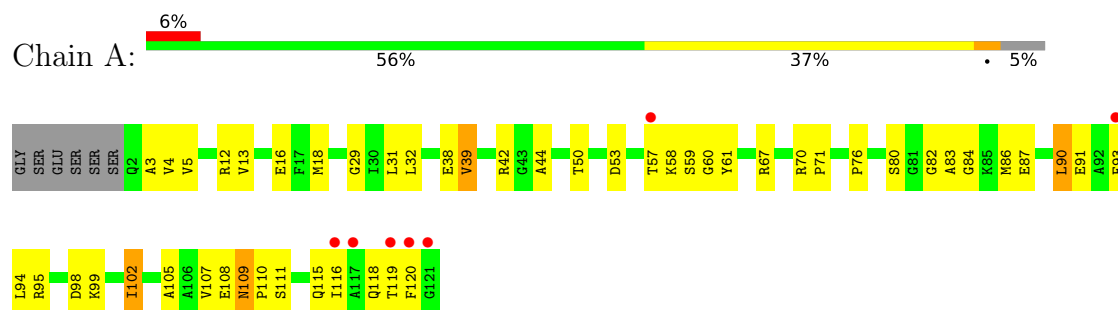
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	27	Total	O	0	0
			27	27		
3	B	23	Total	O	0	0
			23	23		
3	C	7	Total	O	0	0
			7	7		
3	D	9	Total	O	0	0
			9	9		

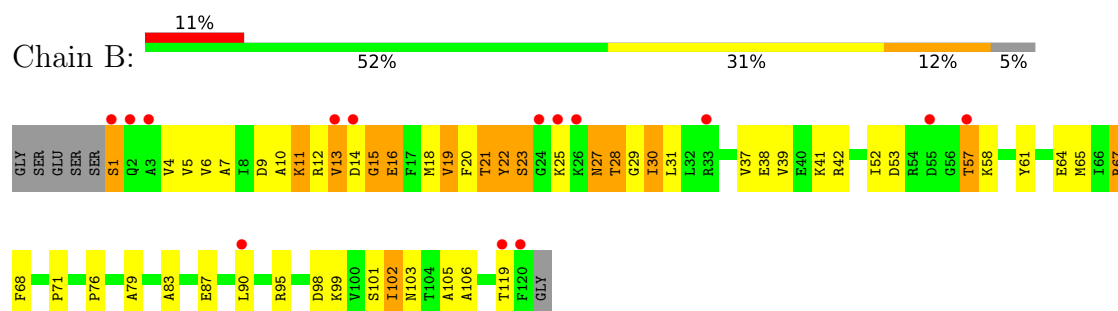
3 Residue-property plots [i](#)

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.

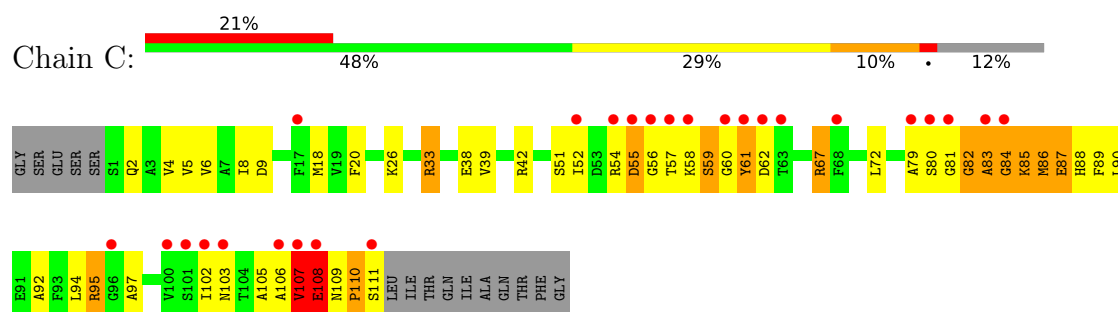
- Molecule 1: Computationally designed two-fold symmetric TIM-barrel protein, FLR (half molecule)



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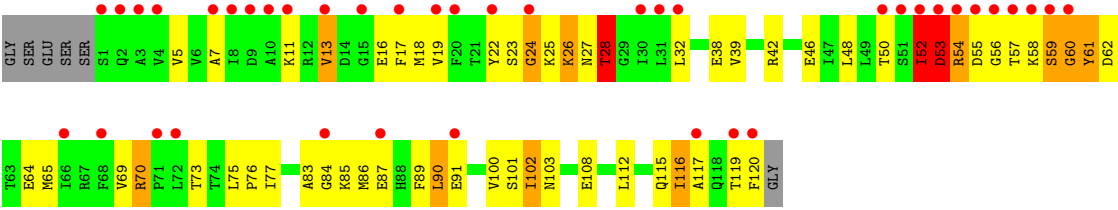


- Molecule 1: Computationally designed two-fold symmetric TIM-barrel protein, FLR (half molecule)



- Molecule 1: Computationally designed two-fold symmetric TIM-barrel protein, FLR (half molecule)





4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	60.49Å 60.49Å 140.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.98 – 2.40 36.98 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.7 (36.98-2.40) 99.8 (36.98-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	12.47 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.242 , 0.288 0.237 , 0.279	Depositor DCC
R_{free} test set	1080 reflections (5.48%)	wwPDB-VP
Wilson B-factor (Å ²)	44.9	Xtriage
Anisotropy	0.194	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 54.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.053 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3684	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 39.18 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.2995e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: PO4

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	1.25	1/929 (0.1%)	1.20	2/1254 (0.2%)
1	B	1.67	14/931 (1.5%)	1.47	10/1257 (0.8%)
1	C	1.68	15/858 (1.7%)	1.42	14/1157 (1.2%)
1	D	1.34	4/931 (0.4%)	1.64	22/1257 (1.8%)
All	All	1.49	34/3649 (0.9%)	1.44	48/4925 (1.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	26	LYS	C-N	18.56	1.57	1.33
1	C	84	GLY	C-N	12.98	1.51	1.33
1	B	79	ALA	CA-CB	-7.87	1.41	1.53
1	B	30	ILE	CA-CB	-7.75	1.45	1.54
1	C	6	VAL	C-O	-7.41	1.16	1.24
1	B	9	ASP	C-O	-7.16	1.15	1.24
1	D	101	SER	C-N	-7.07	1.27	1.33
1	B	7	ALA	CA-CB	-6.87	1.44	1.53
1	B	30	ILE	C-O	-6.80	1.16	1.24
1	B	5	VAL	C-O	-6.64	1.16	1.24
1	C	5	VAL	C-O	-6.46	1.17	1.24
1	C	6	VAL	CA-C	-6.24	1.44	1.52
1	A	13	VAL	CA-CB	-6.21	1.46	1.53
1	D	28	THR	CA-C	-5.96	1.43	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	6	VAL	CA-CB	-5.92	1.47	1.54
1	B	41	LYS	C-O	-5.83	1.17	1.24
1	C	33	ARG	C-O	-5.82	1.17	1.24
1	B	4	VAL	C-O	-5.75	1.18	1.24
1	C	5	VAL	CA-CB	-5.74	1.47	1.54
1	C	8	ILE	C-O	-5.71	1.18	1.24
1	D	103	ASN	C-O	-5.71	1.16	1.24
1	C	4	VAL	C-O	-5.65	1.18	1.24
1	B	37	VAL	CA-CB	-5.63	1.47	1.54
1	B	30	ILE	CA-C	-5.62	1.46	1.52
1	B	7	ALA	C-O	-5.55	1.17	1.24
1	B	21	THR	N-CA	-5.50	1.39	1.46
1	C	79	ALA	CA-C	5.50	1.59	1.52
1	B	39	VAL	C-O	-5.33	1.17	1.24
1	C	4	VAL	CA-CB	-5.32	1.47	1.54
1	C	97	ALA	C-O	-5.29	1.17	1.23
1	B	6	VAL	C-O	-5.25	1.18	1.24
1	C	5	VAL	N-CA	-5.10	1.40	1.46
1	C	92	ALA	CA-C	-5.06	1.46	1.52
1	C	83	ALA	CA-CB	-5.05	1.46	1.53

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1	SER	O-C-N	-15.13	98.79	123.00
1	D	101	SER	O-C-N	-14.68	106.22	123.10
1	D	28	THR	N-CA-C	-13.67	97.20	114.56
1	D	101	SER	CA-C-N	12.02	138.06	121.96
1	D	101	SER	C-N-CA	12.02	138.06	121.96
1	D	26	LYS	O-C-N	-11.56	108.57	123.13
1	C	84	GLY	CA-C-N	11.04	142.64	121.54
1	C	84	GLY	C-N-CA	11.04	142.64	121.54
1	D	55	ASP	CA-C-N	-10.75	106.26	122.28
1	D	55	ASP	C-N-CA	-10.75	106.26	122.28
1	B	1	SER	CA-C-N	10.15	136.84	120.60
1	B	1	SER	C-N-CA	10.15	136.84	120.60
1	D	26	LYS	CA-C-N	10.13	138.97	120.95
1	D	26	LYS	C-N-CA	10.13	138.97	120.95
1	B	19	VAL	O-C-N	-9.91	112.69	123.20
1	C	84	GLY	O-C-N	-9.73	110.05	122.70
1	D	54	ARG	CA-C-N	-9.65	109.39	123.00
1	D	54	ARG	C-N-CA	-9.65	109.39	123.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	95	ARG	N-CA-C	-9.52	101.17	113.17
1	B	57	THR	N-CA-C	-9.45	97.80	110.55
1	C	81	GLY	CA-C-N	-7.92	111.31	122.00
1	C	81	GLY	C-N-CA	-7.92	111.31	122.00
1	C	79	ALA	CA-C-N	7.50	134.72	121.75
1	C	79	ALA	C-N-CA	7.50	134.72	121.75
1	D	55	ASP	N-CA-C	7.40	120.48	108.41
1	B	58	LYS	N-CA-C	-7.08	100.91	111.81
1	B	16	GLU	N-CA-C	7.03	120.86	109.40
1	B	119	THR	N-CA-C	-6.73	105.22	113.50
1	C	79	ALA	N-CA-C	6.66	120.54	109.95
1	C	4	VAL	N-CA-C	6.60	117.39	107.75
1	B	11	LYS	N-CA-C	6.43	119.20	109.23
1	C	82	GLY	N-CA-C	-6.30	106.21	115.72
1	D	53	ASP	CA-C-N	-6.15	113.77	123.12
1	D	53	ASP	C-N-CA	-6.15	113.77	123.12
1	A	4	VAL	N-CA-C	6.11	116.41	107.37
1	A	102	ILE	CB-CA-C	-6.00	100.34	110.71
1	C	81	GLY	N-CA-C	-5.98	99.00	113.18
1	D	116	ILE	CB-CA-C	-5.74	102.96	112.26
1	D	102	ILE	CB-CA-C	-5.67	103.92	111.80
1	D	60	GLY	N-CA-C	5.55	117.22	111.95
1	B	15	GLY	N-CA-C	-5.44	107.02	114.64
1	D	27	ASN	CA-C-N	-5.40	113.82	122.08
1	D	27	ASN	C-N-CA	-5.40	113.82	122.08
1	C	2	GLN	N-CA-C	-5.33	106.00	112.88
1	D	52	ILE	O-C-N	5.24	129.12	122.57
1	C	107	VAL	CB-CA-C	-5.23	102.72	111.29
1	D	70	ARG	CA-C-N	-5.03	115.69	120.83
1	D	70	ARG	C-N-CA	-5.03	115.69	120.83

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	1	SER	Mainchain

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	916	0	935	38	0
1	B	918	0	940	61	2
1	C	846	0	863	87	0
1	D	918	0	940	93	2
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
3	A	27	0	0	3	0
3	B	23	0	0	1	0
3	C	7	0	0	1	0
3	D	9	0	0	3	0
All	All	3684	0	3678	274	2

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 38.

All (274) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:20:PHE:CE1	1:B:27:ASN:HB2	1.34	1.60
1:D:89:PHE:CE1	1:D:116:ILE:HD11	1.35	1.58
1:D:89:PHE:CE1	1:D:116:ILE:CD1	1.90	1.51
1:C:109:ASN:ND2	1:C:111:SER:HB2	1.13	1.42
1:D:52:ILE:CG1	1:D:53:ASP:H	1.34	1.37
1:D:89:PHE:HE1	1:D:116:ILE:CD1	1.24	1.37
1:D:83:ALA:HB1	1:D:89:PHE:CE2	1.58	1.35
1:C:109:ASN:ND2	1:C:111:SER:CB	1.90	1.34
1:B:20:PHE:CD1	1:B:27:ASN:HA	1.62	1.32
1:B:20:PHE:CE1	1:B:27:ASN:CB	2.14	1.29
1:B:18:MET:HE1	1:B:29:GLY:O	1.19	1.27
1:A:93:PHE:CE2	1:A:120:PHE:CZ	2.24	1.24
1:C:109:ASN:O	1:C:111:SER:N	1.74	1.21
1:A:116:ILE:HD11	1:A:120:PHE:CE1	1.74	1.20
1:D:83:ALA:CB	1:D:89:PHE:HE2	1.54	1.20
1:C:85:LYS:HE2	1:C:88:HIS:CE1	1.78	1.19
1:D:89:PHE:CD1	1:D:116:ILE:HG12	1.79	1.18
1:B:20:PHE:HE1	1:B:27:ASN:CB	1.50	1.16
1:A:93:PHE:CD2	1:A:120:PHE:CZ	2.35	1.15

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:PHE:CD2	1:A:120:PHE:HZ	1.64	1.15
1:C:108:GLU:H	1:C:109:ASN:HA	1.08	1.14
1:B:20:PHE:CD1	1:B:27:ASN:CA	2.31	1.12
1:B:67:ARG:NH1	1:B:67:ARG:HG2	1.43	1.10
1:C:109:ASN:HD22	1:C:111:SER:CB	1.53	1.10
1:D:89:PHE:HD1	1:D:116:ILE:HG12	1.13	1.10
1:B:18:MET:CE	1:B:29:GLY:O	1.99	1.09
1:C:107:VAL:O	1:C:108:GLU:HB3	1.45	1.09
1:C:109:ASN:C	1:C:111:SER:H	1.58	1.09
1:D:89:PHE:CD1	1:D:116:ILE:CD1	2.35	1.09
1:D:11:LYS:HD2	1:D:52:ILE:HD13	1.17	1.08
1:D:52:ILE:HG13	1:D:53:ASP:N	1.31	1.08
1:C:107:VAL:CA	1:C:110:PRO:HD3	1.83	1.08
1:C:57:THR:O	1:C:82:GLY:HA2	1.54	1.08
1:D:11:LYS:CD	1:D:52:ILE:HD13	1.87	1.05
1:B:67:ARG:HH11	1:B:67:ARG:CG	1.62	1.05
1:C:108:GLU:N	1:C:109:ASN:HA	1.58	1.04
1:C:108:GLU:H	1:C:110:PRO:CD	1.70	1.03
1:D:89:PHE:CD1	1:D:116:ILE:CG1	2.41	1.03
1:D:87:GLU:O	1:D:91:GLU:HG3	1.57	1.01
1:B:19:VAL:O	1:B:28:THR:HB	1.58	1.01
1:A:116:ILE:HD11	1:A:120:PHE:HE1	1.07	1.01
1:C:85:LYS:O	1:C:89:PHE:CD1	2.13	1.00
1:C:87:GLU:HA	1:C:87:GLU:OE1	1.56	1.00
1:D:19:VAL:O	1:D:28:THR:HG23	1.60	1.00
1:D:89:PHE:CE1	1:D:116:ILE:HD13	1.92	1.00
1:C:107:VAL:HA	1:C:110:PRO:HD3	1.42	0.98
1:C:9:ASP:HB3	1:C:52:ILE:HD11	1.44	0.96
1:D:89:PHE:HD1	1:D:116:ILE:CG1	1.77	0.95
1:A:93:PHE:CE2	1:A:120:PHE:CE1	2.55	0.94
1:A:111:SER:O	1:A:115:GLN:HG3	1.67	0.93
1:A:108:GLU:O	1:A:110:PRO:HD3	1.69	0.93
1:C:108:GLU:N	1:C:110:PRO:HD3	1.85	0.91
1:D:83:ALA:CB	1:D:89:PHE:CE2	2.38	0.90
1:C:106:ALA:O	1:C:107:VAL:HG23	1.69	0.90
1:C:108:GLU:O	1:C:108:GLU:HG2	1.72	0.90
1:C:61:TYR:CZ	1:C:83:ALA:HB2	2.08	0.89
1:C:108:GLU:N	1:C:109:ASN:CA	2.36	0.89
1:B:67:ARG:HG2	1:B:67:ARG:HH11	0.74	0.89
1:B:20:PHE:CE1	1:B:27:ASN:CA	2.52	0.88
1:D:23:SER:O	1:D:25:LYS:HG2	1.74	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:85:LYS:CE	1:C:88:HIS:CE1	2.57	0.88
1:C:109:ASN:C	1:C:111:SER:N	2.19	0.87
1:D:19:VAL:O	1:D:28:THR:CG2	2.23	0.87
1:C:85:LYS:O	1:C:89:PHE:CE1	2.28	0.86
1:C:85:LYS:O	1:C:89:PHE:HD1	1.56	0.86
1:B:28:THR:HG23	1:B:30:ILE:H	1.41	0.86
1:B:57:THR:O	3:B:142:HOH:O	1.93	0.86
1:B:20:PHE:HD1	1:B:27:ASN:CA	1.88	0.85
1:C:108:GLU:H	1:C:109:ASN:CA	1.87	0.85
1:D:19:VAL:HG12	1:D:28:THR:HG21	1.59	0.85
1:B:67:ARG:NH1	1:B:67:ARG:CG	2.30	0.84
1:C:107:VAL:C	1:C:110:PRO:HD3	2.03	0.84
1:C:67:ARG:HG3	1:C:67:ARG:HH11	1.43	0.84
1:D:87:GLU:O	1:D:91:GLU:CG	2.25	0.84
1:C:109:ASN:HD21	1:C:111:SER:CB	1.87	0.83
1:A:99:LYS:NZ	3:A:149:HOH:O	2.12	0.83
1:D:85:LYS:HD3	1:D:87:GLU:H	1.43	0.83
1:D:83:ALA:HB1	1:D:89:PHE:HE2	0.71	0.83
1:D:89:PHE:CE1	1:D:116:ILE:CG1	2.60	0.82
1:B:11:LYS:HB2	1:B:53:ASP:OD2	1.79	0.82
1:B:18:MET:CE	1:B:29:GLY:C	2.53	0.82
1:D:89:PHE:CD1	1:D:116:ILE:HD13	2.10	0.81
1:A:116:ILE:CD1	1:A:120:PHE:CE1	2.61	0.81
1:B:12:ARG:O	1:B:13:VAL:CG2	2.29	0.81
1:D:11:LYS:CE	1:D:52:ILE:CD1	2.59	0.80
1:C:106:ALA:O	1:C:107:VAL:CG2	2.29	0.80
1:B:22:TYR:O	1:B:25:LYS:HB2	1.83	0.79
1:D:19:VAL:C	1:D:28:THR:CG2	2.55	0.79
1:B:20:PHE:CE1	1:B:27:ASN:HA	2.13	0.79
1:C:9:ASP:CB	1:C:52:ILE:HD11	2.13	0.79
1:B:19:VAL:O	1:B:28:THR:CB	2.30	0.78
1:C:108:GLU:N	1:C:110:PRO:CD	2.41	0.78
1:D:85:LYS:HD2	1:D:87:GLU:HB3	1.64	0.78
1:D:11:LYS:NZ	1:D:52:ILE:CD1	2.47	0.77
1:C:57:THR:O	1:C:82:GLY:CA	2.30	0.77
1:D:11:LYS:CE	1:D:52:ILE:HD13	2.15	0.77
1:C:87:GLU:OE1	1:C:87:GLU:CA	2.30	0.76
1:C:58:LYS:HD3	1:C:84:GLY:HA3	1.68	0.76
1:C:102:ILE:HD11	1:C:105:ALA:HB3	1.68	0.76
1:C:106:ALA:O	1:C:107:VAL:CB	2.30	0.75
1:D:13:VAL:N	1:D:16:GLU:O	2.18	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:13:VAL:O	1:B:14:ASP:HB2	1.86	0.74
1:C:106:ALA:C	1:C:107:VAL:HG23	2.10	0.74
1:D:117:ALA:O	1:D:120:PHE:N	2.17	0.73
1:C:61:TYR:CZ	1:C:83:ALA:CB	2.70	0.73
1:C:9:ASP:HB3	1:C:52:ILE:CD1	2.19	0.73
1:C:108:GLU:O	1:C:108:GLU:CG	2.37	0.73
1:C:62:ASP:OD1	1:C:62:ASP:C	2.30	0.73
1:A:119:THR:O	1:C:95:ARG:NH2	2.20	0.72
1:C:107:VAL:HA	1:C:110:PRO:CD	2.19	0.72
1:B:19:VAL:O	1:B:28:THR:CG2	2.38	0.72
1:C:107:VAL:O	1:C:108:GLU:CB	2.30	0.72
1:B:12:ARG:O	1:B:13:VAL:HG23	1.89	0.71
1:C:106:ALA:O	1:C:107:VAL:HB	1.91	0.70
1:D:11:LYS:HB3	1:D:52:ILE:HG12	1.72	0.70
1:B:102:ILE:HG12	1:B:106:ALA:HB2	1.72	0.69
1:A:60:GLY:N	1:A:82:GLY:O	2.23	0.69
1:A:86:MET:HE1	1:A:116:ILE:HA	1.74	0.69
1:A:93:PHE:HE2	1:A:120:PHE:CE1	2.06	0.69
1:A:116:ILE:CD1	1:A:120:PHE:HE1	1.97	0.69
1:B:20:PHE:HD1	1:B:27:ASN:N	1.90	0.68
1:D:11:LYS:HD2	1:D:52:ILE:CD1	2.10	0.68
1:B:20:PHE:HD1	1:B:27:ASN:HA	1.41	0.68
1:B:27:ASN:OD1	1:B:28:THR:N	2.27	0.67
1:D:62:ASP:C	1:D:62:ASP:OD1	2.36	0.67
1:D:11:LYS:HE3	1:D:52:ILE:HD11	1.77	0.66
1:B:102:ILE:HD11	1:B:105:ALA:HB3	1.77	0.66
1:B:12:ARG:C	1:B:13:VAL:HG23	2.21	0.66
1:C:54:ARG:HD2	1:C:62:ASP:HA	1.78	0.66
3:A:149:HOH:O	1:B:99:LYS:NZ	2.27	0.66
1:D:53:ASP:O	1:D:54:ARG:NE	2.30	0.65
1:A:53:ASP:OD2	3:A:133:HOH:O	2.14	0.64
1:C:55:ASP:OD1	1:C:56:GLY:N	2.30	0.64
1:D:85:LYS:CD	1:D:87:GLU:HB3	2.27	0.64
1:D:86:MET:O	1:D:90:LEU:HD13	1.97	0.64
1:D:11:LYS:CD	1:D:52:ILE:CD1	2.69	0.63
1:D:52:ILE:HG13	1:D:53:ASP:CA	2.24	0.63
1:A:3:ALA:HA	1:B:99:LYS:HG2	1.81	0.63
1:D:53:ASP:O	1:D:54:ARG:CZ	2.46	0.63
1:D:19:VAL:C	1:D:28:THR:HG23	2.22	0.63
1:C:61:TYR:CE2	1:C:83:ALA:HB2	2.33	0.62
1:D:17:PHE:CD1	1:D:65:MET:HE2	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:27:ASN:OD1	1:B:27:ASN:C	2.43	0.61
1:B:12:ARG:O	1:B:13:VAL:HG22	1.98	0.61
1:C:109:ASN:ND2	1:C:111:SER:CA	2.63	0.61
1:D:11:LYS:CE	1:D:52:ILE:HD11	2.29	0.60
1:C:51:SER:O	1:C:55:ASP:N	2.34	0.59
1:C:57:THR:HG22	1:C:59:SER:OG	2.03	0.59
1:D:11:LYS:NZ	1:D:52:ILE:HD13	2.17	0.58
1:D:61:TYR:CE2	1:D:83:ALA:HB2	2.38	0.58
1:C:109:ASN:O	1:C:110:PRO:C	2.36	0.58
1:C:85:LYS:H	1:C:89:PHE:HE1	1.51	0.57
1:D:52:ILE:O	1:D:54:ARG:N	2.37	0.57
1:A:58:LYS:HB3	1:A:84:GLY:HA2	1.86	0.57
1:C:55:ASP:OD1	1:C:55:ASP:C	2.45	0.57
1:D:70:ARG:NH2	1:D:75:LEU:O	2.34	0.57
1:D:56:GLY:HA2	3:D:131:HOH:O	2.04	0.57
1:B:18:MET:HE3	1:B:29:GLY:C	2.29	0.57
1:D:86:MET:HE1	1:D:115:GLN:O	2.04	0.57
1:D:89:PHE:N	1:D:89:PHE:CD2	2.73	0.56
1:C:61:TYR:CE2	1:C:83:ALA:CB	2.88	0.56
1:C:54:ARG:HD2	1:C:62:ASP:CA	2.36	0.56
1:A:91:GLU:O	1:A:95:ARG:HG2	2.06	0.56
1:B:20:PHE:CD1	1:B:27:ASN:N	2.71	0.56
1:A:12:ARG:HA	1:A:16:GLU:O	2.05	0.56
1:A:50:THR:HA	1:A:80:SER:O	2.06	0.56
1:B:12:ARG:C	1:B:13:VAL:CG2	2.78	0.55
1:C:57:THR:CG2	1:C:59:SER:OG	2.55	0.55
1:C:86:MET:O	1:C:87:GLU:C	2.49	0.55
1:D:11:LYS:HG3	1:D:13:VAL:HG22	1.88	0.55
1:C:109:ASN:HD21	1:C:111:SER:HB3	1.66	0.54
1:D:61:TYR:CZ	1:D:83:ALA:HB2	2.42	0.54
1:B:76:PRO:HA	1:B:98:ASP:OD2	2.08	0.54
1:C:85:LYS:O	1:C:89:PHE:HE1	1.89	0.54
1:D:7:ALA:HA	1:D:48:LEU:HB3	1.90	0.54
1:C:108:GLU:H	1:C:110:PRO:HD2	1.66	0.54
1:D:56:GLY:CA	3:D:131:HOH:O	2.56	0.54
1:D:60:GLY:C	1:D:61:TYR:O	2.50	0.54
1:D:61:TYR:OH	1:D:100:VAL:HB	2.07	0.54
1:D:13:VAL:HG21	1:D:18:MET:HB2	1.89	0.53
1:C:80:SER:O	3:C:129:HOH:O	2.19	0.53
1:C:107:VAL:CA	1:C:110:PRO:CD	2.72	0.53
1:A:70:ARG:HB3	1:A:71:PRO:HD3	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:86:MET:HE3	1:D:90:LEU:HD11	1.89	0.53
1:D:89:PHE:HE1	1:D:116:ILE:HD11	0.43	0.53
1:D:89:PHE:CZ	1:D:102:ILE:HG21	2.44	0.53
1:B:38:GLU:OE1	1:B:42:ARG:NE	2.34	0.53
1:D:22:TYR:HB2	1:D:26:LYS:HB2	1.91	0.53
1:D:112:LEU:O	1:D:116:ILE:HG13	2.08	0.52
1:B:20:PHE:HE1	1:B:27:ASN:HB2	0.58	0.52
1:A:39:VAL:HG13	1:A:44:ALA:HB2	1.92	0.52
1:D:32:LEU:HD21	1:D:69:VAL:HG23	1.91	0.52
1:D:11:LYS:HZ2	1:D:52:ILE:CD1	2.21	0.52
1:D:52:ILE:HG13	1:D:53:ASP:H	0.46	0.52
1:B:21:THR:O	1:B:22:TYR:C	2.52	0.51
1:B:21:THR:O	1:B:22:TYR:O	2.27	0.51
1:C:67:ARG:HG3	1:C:67:ARG:NH1	2.17	0.51
1:A:67:ARG:O	1:A:67:ARG:HG2	2.10	0.51
1:C:106:ALA:C	1:C:107:VAL:CG2	2.76	0.51
1:D:58:LYS:HB3	1:D:84:GLY:HA3	1.92	0.51
1:B:30:ILE:HG22	1:B:31:LEU:O	2.11	0.50
1:D:11:LYS:HZ2	1:D:52:ILE:HD13	1.77	0.50
1:D:62:ASP:OD1	1:D:62:ASP:O	2.30	0.50
1:C:62:ASP:OD1	1:C:62:ASP:O	2.30	0.50
1:D:19:VAL:C	1:D:28:THR:HG21	2.36	0.50
1:B:13:VAL:C	1:B:15:GLY:H	2.19	0.50
1:C:60:GLY:O	1:C:61:TYR:O	2.30	0.50
1:D:60:GLY:O	1:D:61:TYR:O	2.30	0.50
1:D:73:THR:OG1	1:D:75:LEU:HG	2.12	0.50
1:B:30:ILE:HG22	1:B:31:LEU:N	2.25	0.49
1:C:54:ARG:O	1:C:55:ASP:C	2.56	0.49
1:B:13:VAL:O	1:B:16:GLU:N	2.30	0.48
1:B:19:VAL:HB	1:B:28:THR:HG21	1.94	0.48
1:B:61:TYR:CZ	1:B:83:ALA:HB2	2.49	0.48
1:C:9:ASP:HB3	1:C:52:ILE:CG1	2.43	0.48
1:C:20:PHE:HA	1:C:26:LYS:O	2.13	0.48
1:B:102:ILE:HG13	1:B:103:ASN:N	2.28	0.47
1:D:13:VAL:CG2	1:D:18:MET:HB2	2.44	0.47
1:C:85:LYS:HE2	1:C:88:HIS:ND1	2.24	0.47
1:C:56:GLY:O	1:C:57:THR:C	2.56	0.47
1:D:32:LEU:HD21	1:D:69:VAL:CG2	2.45	0.47
1:C:58:LYS:CD	1:C:84:GLY:HA3	2.43	0.47
1:D:17:PHE:CD1	1:D:65:MET:CE	2.97	0.47
1:B:10:ALA:O	1:B:65:MET:HE1	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:ILE:CG2	1:B:31:LEU:N	2.74	0.46
1:D:58:LYS:NZ	1:D:108:GLU:OE1	2.32	0.46
1:B:13:VAL:C	1:B:15:GLY:N	2.72	0.46
1:D:86:MET:HE1	1:D:116:ILE:HA	1.97	0.46
1:A:5:VAL:O	1:B:101:SER:HA	2.16	0.46
1:A:57:THR:O	1:A:59:SER:N	2.47	0.46
1:B:65:MET:O	1:B:68:PHE:HB3	2.15	0.46
1:A:31:LEU:O	1:A:32:LEU:C	2.59	0.46
1:A:18:MET:HE1	1:A:29:GLY:C	2.42	0.45
1:D:23:SER:O	1:D:24:GLY:C	2.60	0.45
1:C:108:GLU:H	1:C:110:PRO:N	2.14	0.45
1:A:115:GLN:O	1:A:118:GLN:HG2	2.16	0.45
1:A:38:GLU:O	1:A:42:ARG:HG2	2.17	0.44
1:A:105:ALA:O	1:A:109:ASN:N	2.41	0.44
1:C:105:ALA:O	1:C:108:GLU:C	2.61	0.44
1:D:46:GLU:HG3	1:D:76:PRO:HB2	1.98	0.44
1:D:17:PHE:CE1	1:D:65:MET:HE3	2.53	0.44
1:D:38:GLU:O	1:D:42:ARG:HG2	2.17	0.44
1:A:116:ILE:CD1	1:A:120:PHE:CD1	2.99	0.44
1:C:85:LYS:N	1:C:89:PHE:HE1	2.16	0.44
1:C:90:LEU:HD23	1:C:90:LEU:HA	1.85	0.44
1:A:61:TYR:CZ	1:A:83:ALA:HB2	2.53	0.43
1:B:67:ARG:NH2	1:B:95:ARG:NH2	2.66	0.43
1:A:94:LEU:HD23	1:A:94:LEU:HA	1.66	0.43
1:D:54:ARG:HD3	1:D:54:ARG:HA	1.71	0.43
1:C:38:GLU:OE1	1:C:42:ARG:NE	2.44	0.43
1:D:70:ARG:CZ	1:D:77:ILE:HD12	2.48	0.43
1:B:64:GLU:O	1:B:68:PHE:HB2	2.19	0.43
1:D:5:VAL:HA	1:D:46:GLU:O	2.18	0.43
1:C:102:ILE:CD1	1:C:105:ALA:HB3	2.45	0.43
1:A:90:LEU:HD22	1:C:87:GLU:OE1	2.18	0.42
1:D:89:PHE:N	1:D:89:PHE:HD2	2.14	0.42
1:D:52:ILE:CG1	1:D:53:ASP:N	2.15	0.42
1:A:90:LEU:CD2	1:C:87:GLU:OE1	2.67	0.42
1:D:85:LYS:CD	1:D:87:GLU:H	2.23	0.42
1:C:60:GLY:HA2	1:C:88:HIS:HB3	2.02	0.42
1:A:76:PRO:HA	1:A:98:ASP:OD2	2.19	0.41
1:D:11:LYS:HB3	1:D:52:ILE:CG1	2.48	0.41
1:C:33:ARG:HA	1:C:72:LEU:HD13	2.02	0.41
1:C:85:LYS:C	1:C:89:PHE:CE1	2.98	0.41
1:C:103:ASN:OD1	1:C:103:ASN:C	2.61	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:57:THR:O	1:A:57:THR:OG1	2.33	0.41
1:B:11:LYS:HE2	1:B:52:ILE:HG21	2.02	0.41
1:B:28:THR:CG2	1:B:30:ILE:H	2.20	0.41
1:B:95:ARG:HD2	1:B:95:ARG:HA	1.82	0.41
1:B:68:PHE:O	1:B:71:PRO:HD2	2.21	0.40
1:D:59:SER:HB3	3:D:130:HOH:O	2.21	0.40
1:C:108:GLU:H	1:C:109:ASN:C	2.29	0.40
1:D:17:PHE:CE1	1:D:65:MET:CE	3.04	0.40
1:D:83:ALA:HB1	1:D:89:PHE:CD2	2.38	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:GLU:OE2	1:D:85:LYS:NZ[4_465]	1.34	0.86
1:B:95:ARG:NH2	1:D:119:THR:O[4_465]	1.56	0.64

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	118/126 (94%)	112 (95%)	5 (4%)	1 (1%)	16	25
1	B	118/126 (94%)	107 (91%)	8 (7%)	3 (2%)	4	5
1	C	109/126 (86%)	97 (89%)	6 (6%)	6 (6%)	1	1
1	D	118/126 (94%)	103 (87%)	11 (9%)	4 (3%)	3	2
All	All	463/504 (92%)	419 (90%)	30 (6%)	14 (3%)	3	3

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	61	TYR
1	C	85	LYS
1	C	108	GLU
1	C	110	PRO
1	D	52	ILE
1	D	53	ASP
1	D	61	TYR
1	D	24	GLY
1	B	22	TYR
1	B	23	SER
1	B	13	VAL
1	C	55	ASP
1	C	107	VAL
1	A	109	ASN

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	97/102 (95%)	92 (95%)	5 (5%)	21	36
1	B	98/102 (96%)	92 (94%)	6 (6%)	17	30
1	C	90/102 (88%)	82 (91%)	8 (9%)	9	15
1	D	98/102 (96%)	90 (92%)	8 (8%)	10	18
All	All	383/408 (94%)	356 (93%)	27 (7%)	14	24

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	VAL
1	A	87	GLU
1	A	90	LEU
1	A	102	ILE
1	A	107	VAL
1	B	23	SER
1	B	27	ASN
1	B	28	THR

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Mol	Chain	Res	Type
1	B	67	ARG
1	B	90	LEU
1	B	102	ILE
1	C	18	MET
1	C	39	VAL
1	C	59	SER
1	C	67	ARG
1	C	86	MET
1	C	87	GLU
1	C	94	LEU
1	C	108	GLU
1	D	13	VAL
1	D	28	THR
1	D	39	VAL
1	D	50	THR
1	D	57	THR
1	D	59	SER
1	D	64	GLU
1	D	90	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	GLN
1	A	118	GLN
1	C	109	ASN

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 [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PO4	C	122	-	4,4,4	0.87	0	6,6,6	0.30	0
2	PO4	A	122	-	4,4,4	1.07	0	6,6,6	0.66	0
2	PO4	D	122	-	4,4,4	0.98	0	6,6,6	0.60	0
2	PO4	B	122	-	4,4,4	1.51	0	6,6,6	1.13	0

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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	120/126 (95%)	0.43	7 (5%) 29 25	33, 57, 90, 189	0
1	B	120/126 (95%)	0.69	14 (11%) 9 7	34, 66, 108, 137	0
1	C	111/126 (88%)	1.32	26 (23%) 2 1	45, 83, 207, 246	0
1	D	120/126 (95%)	1.63	40 (33%) 1 1	40, 90, 167, 263	0
All	All	471/504 (93%)	1.01	87 (18%) 3 3	33, 72, 160, 263	0

All (87) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	10	ALA	6.5
1	D	51	SER	5.9
1	C	52	ILE	5.7
1	D	8	ILE	5.0
1	C	111	SER	4.6
1	D	52	ILE	4.6
1	A	120	PHE	4.6
1	B	120	PHE	4.3
1	A	116	ILE	4.3
1	A	121	GLY	4.2
1	C	58	LYS	4.2
1	D	19	VAL	4.1
1	C	55	ASP	4.1
1	D	3	ALA	3.8
1	D	30	ILE	3.8
1	D	58	LYS	3.8
1	D	60	GLY	3.8
1	D	9	ASP	3.8
1	C	102	ILE	3.7
1	C	96	GLY	3.6
1	D	84	GLY	3.6

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Mol	Chain	Res	Type	RSRZ
1	D	120	PHE	3.6
1	D	32	LEU	3.5
1	C	107	VAL	3.5
1	C	84	GLY	3.5
1	D	54	ARG	3.4
1	D	2	GLN	3.3
1	B	1	SER	3.3
1	D	1	SER	3.2
1	C	60	GLY	3.2
1	D	87	GLU	3.2
1	C	56	GLY	3.2
1	C	81	GLY	3.1
1	B	119	THR	3.1
1	D	57	THR	3.1
1	C	68	PHE	3.0
1	D	4	VAL	3.0
1	C	63	THR	2.9
1	D	53	ASP	2.9
1	C	61	TYR	2.9
1	D	11	LYS	2.9
1	D	55	ASP	2.9
1	C	54	ARG	2.8
1	B	24	GLY	2.8
1	D	17	PHE	2.8
1	A	117	ALA	2.7
1	D	24	GLY	2.7
1	C	101	SER	2.7
1	D	56	GLY	2.7
1	C	80	SER	2.6
1	A	119	THR	2.6
1	C	79	ALA	2.6
1	C	103	ASN	2.6
1	C	106	ALA	2.6
1	C	62	ASP	2.5
1	B	13	VAL	2.5
1	B	33	ARG	2.5
1	D	72	LEU	2.5
1	A	57	THR	2.5
1	D	22	TYR	2.5
1	B	3	ALA	2.5
1	D	20	PHE	2.5
1	D	117	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	57	THR	2.5
1	A	93	PHE	2.4
1	D	13	VAL	2.4
1	D	66	ILE	2.4
1	D	68	PHE	2.4
1	B	26	LYS	2.4
1	C	83	ALA	2.3
1	B	57	THR	2.3
1	C	100	VAL	2.3
1	B	55	ASP	2.3
1	B	2	GLN	2.3
1	C	108	GLU	2.3
1	D	71	PRO	2.3
1	D	15	GLY	2.2
1	B	25	LYS	2.2
1	D	50	THR	2.2
1	D	59	SER	2.2
1	B	14	ASP	2.2
1	D	119	THR	2.2
1	B	90	LEU	2.2
1	D	7	ALA	2.1
1	D	31	LEU	2.1
1	D	91	GLU	2.0
1	C	17	PHE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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	PO4	C	122	5/5	0.53	0.17	89,89,90,91	0
2	PO4	D	122	5/5	0.91	0.18	53,55,56,57	0
2	PO4	A	122	5/5	0.97	0.09	31,31,34,35	0
2	PO4	B	122	5/5	0.98	0.11	29,29,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.