



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:06 PM UTC

PDB ID : 3K17 / pdb_00003k17
Title : Crystal structure of a Lin0012 protein from *Listeria innocua*
Authors : Palani, K.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center
for Structural Genomics (NYSGXRC)
Deposited on : 2009-09-25
Resolution : 2.10 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

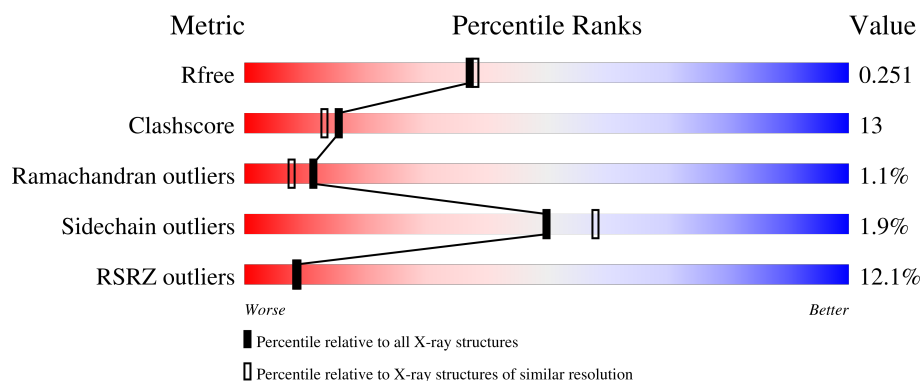
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	365	2% 77% 19% ..
1	B	365	12% 67% 28% ..
1	C	365	7% 70% 25% ..
1	D	365	26% 67% 26% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11491 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 Lin0012 protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	355	Total	C	N	O	S	Se	0	0	0
			2774	1776	459	528	3	8			
1	B	355	Total	C	N	O	S	Se	0	0	0
			2774	1776	459	528	3	8			
1	C	355	Total	C	N	O	S	Se	0	0	0
			2774	1776	459	528	3	8			
1	D	355	Total	C	N	O	S	Se	0	0	0
			2774	1776	459	528	3	8			

There are 40 discrepancies between the modelled and reference sequences:

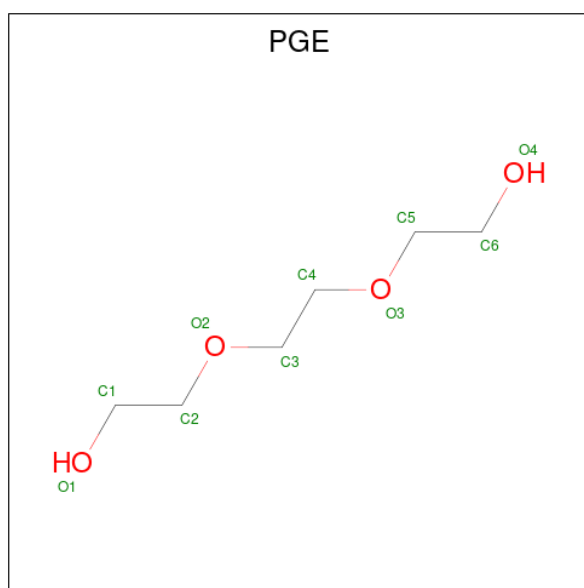
Chain	Residue	Modelled	Actual	Comment	Reference
A	3	MSE	-	expression tag	UNP Q92FU1
A	4	SER	-	expression tag	UNP Q92FU1
A	5	LEU	-	expression tag	UNP Q92FU1
A	361	GLY	-	expression tag	UNP Q92FU1
A	362	HIS	-	expression tag	UNP Q92FU1
A	363	HIS	-	expression tag	UNP Q92FU1
A	364	HIS	-	expression tag	UNP Q92FU1
A	365	HIS	-	expression tag	UNP Q92FU1
A	366	HIS	-	expression tag	UNP Q92FU1
A	367	HIS	-	expression tag	UNP Q92FU1
B	3	MSE	-	expression tag	UNP Q92FU1
B	4	SER	-	expression tag	UNP Q92FU1
B	5	LEU	-	expression tag	UNP Q92FU1
B	361	GLY	-	expression tag	UNP Q92FU1
B	362	HIS	-	expression tag	UNP Q92FU1
B	363	HIS	-	expression tag	UNP Q92FU1
B	364	HIS	-	expression tag	UNP Q92FU1
B	365	HIS	-	expression tag	UNP Q92FU1
B	366	HIS	-	expression tag	UNP Q92FU1
B	367	HIS	-	expression tag	UNP Q92FU1
C	3	MSE	-	expression tag	UNP Q92FU1

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Chain	Residue	Modelled	Actual	Comment	Reference
C	4	SER	-	expression tag	UNP Q92FU1
C	5	LEU	-	expression tag	UNP Q92FU1
C	361	GLY	-	expression tag	UNP Q92FU1
C	362	HIS	-	expression tag	UNP Q92FU1
C	363	HIS	-	expression tag	UNP Q92FU1
C	364	HIS	-	expression tag	UNP Q92FU1
C	365	HIS	-	expression tag	UNP Q92FU1
C	366	HIS	-	expression tag	UNP Q92FU1
C	367	HIS	-	expression tag	UNP Q92FU1
D	3	MSE	-	expression tag	UNP Q92FU1
D	4	SER	-	expression tag	UNP Q92FU1
D	5	LEU	-	expression tag	UNP Q92FU1
D	361	GLY	-	expression tag	UNP Q92FU1
D	362	HIS	-	expression tag	UNP Q92FU1
D	363	HIS	-	expression tag	UNP Q92FU1
D	364	HIS	-	expression tag	UNP Q92FU1
D	365	HIS	-	expression tag	UNP Q92FU1
D	366	HIS	-	expression tag	UNP Q92FU1
D	367	HIS	-	expression tag	UNP Q92FU1

- Molecule 2 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			10	6	4		

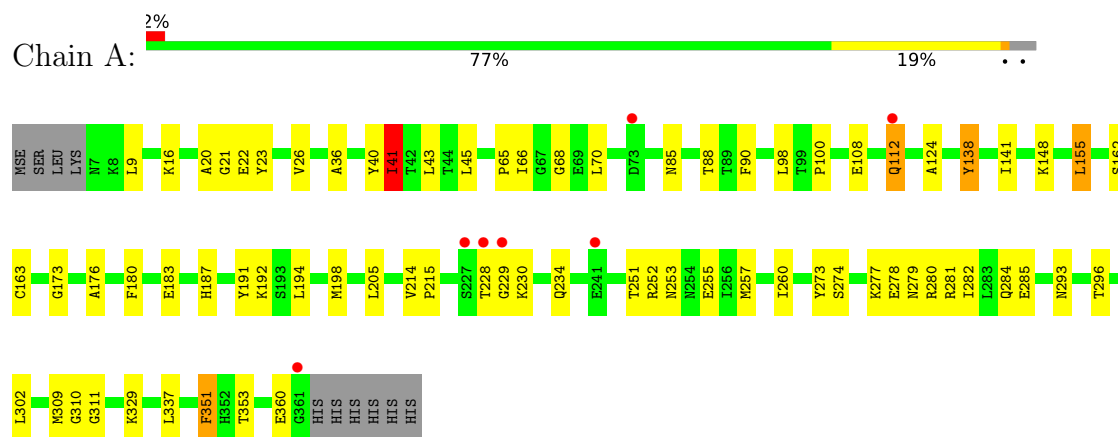
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	128	Total 128	O 128	0	0
3	B	69	Total 69	O 69	0	0
3	C	124	Total 124	O 124	0	0
3	D	64	Total 64	O 64	0	0

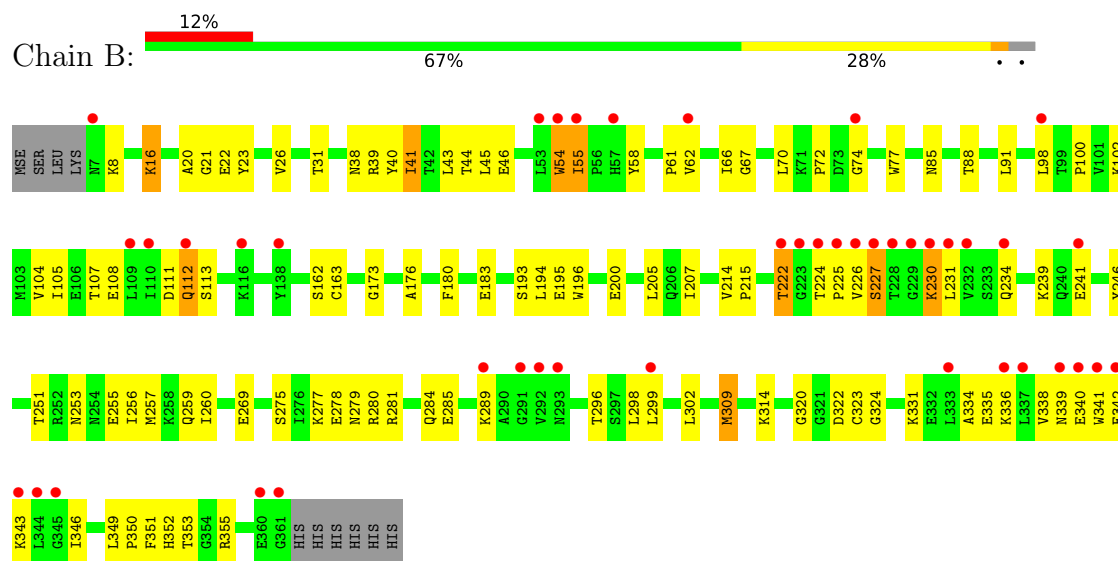
3 Residue-property plots [i](#)

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: Lin0012 protein

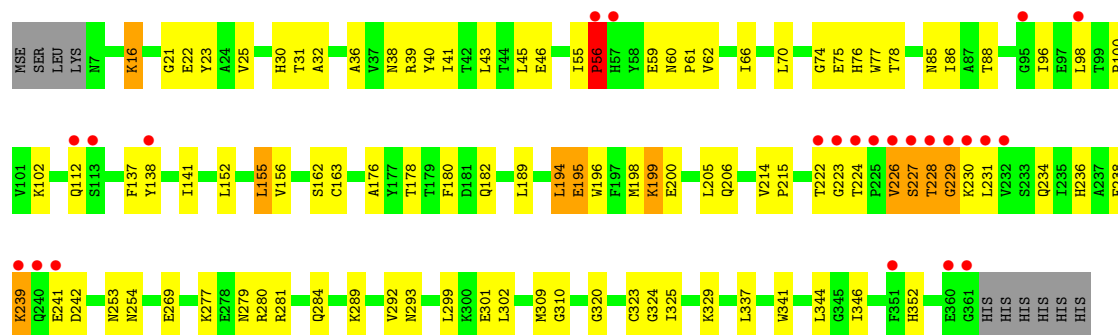


• Molecule 1: Lin0012 protein

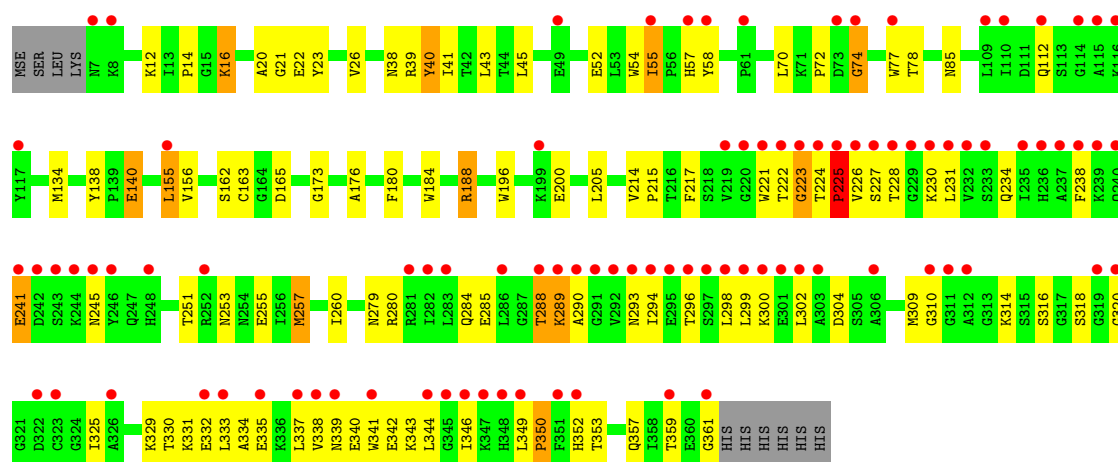


• Molecule 1: Lin0012 protein





● Molecule 1: Lin0012 protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	105.97Å 122.36Å 138.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.06 – 2.10 43.06 – 2.10	Depositor EDS
% Data completeness (in resolution range)	96.3 (43.06-2.10) 96.3 (43.06-2.10)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.00 (at 2.10Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.223 , 0.252 0.222 , 0.251	Depositor DCC
R_{free} test set	5099 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.285	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11491	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.73% 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: PGE

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	0.40	0/2829	0.88	8/3818 (0.2%)
1	B	0.35	0/2829	0.86	10/3818 (0.3%)
1	C	0.40	0/2829	0.91	12/3818 (0.3%)
1	D	0.34	0/2829	0.85	6/3818 (0.2%)
All	All	0.37	0/11316	0.88	36/15272 (0.2%)

There are no bond length outliers.

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	41	ILE	N-CA-C	-8.62	96.11	108.17
1	A	41	ILE	N-CA-C	-8.07	96.69	108.71
1	A	40	TYR	N-CA-C	7.40	121.55	109.85
1	D	40	TYR	N-CA-C	7.37	121.50	109.85
1	C	78	THR	N-CA-C	7.36	119.94	111.11
1	C	40	TYR	N-CA-C	7.22	120.86	109.81
1	C	324	GLY	N-CA-C	-6.91	100.11	111.38
1	B	40	TYR	N-CA-C	6.61	119.92	109.81
1	C	176	ALA	N-CA-C	-6.55	97.84	108.52
1	D	41	ILE	N-CA-C	-6.52	98.35	107.80
1	C	55	ILE	CA-C-N	6.12	127.50	119.84
1	C	55	ILE	C-N-CA	6.12	127.50	119.84
1	C	56	PRO	N-CA-C	6.05	124.93	112.47
1	A	351	PHE	N-CA-C	5.98	119.28	109.72
1	C	241	GLU	N-CA-C	5.88	117.49	111.14
1	B	108	GLU	N-CA-C	-5.71	105.62	112.59
1	B	55	ILE	CA-C-N	5.68	125.39	119.82
1	B	55	ILE	C-N-CA	5.68	125.39	119.82
1	C	323	CYS	N-CA-C	5.61	118.82	110.52
1	B	41	ILE	N-CA-C	-5.61	99.57	107.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	176	ALA	N-CA-C	-5.42	99.69	108.52
1	A	108	GLU	N-CA-C	-5.40	104.97	113.02
1	B	176	ALA	N-CA-C	-5.40	98.44	107.99
1	B	322	ASP	CB-CA-C	-5.33	109.44	115.79
1	C	36	ALA	N-CA-C	-5.31	102.66	110.52
1	C	39	ARG	N-CA-C	-5.31	99.78	108.76
1	B	323	CYS	N-CA-C	5.30	118.37	110.52
1	A	176	ALA	N-CA-C	-5.16	98.85	107.99
1	B	324	GLY	N-CA-C	-5.14	103.26	111.18
1	A	36	ALA	N-CA-C	-5.13	103.63	110.55
1	D	78	THR	N-CA-C	5.11	117.25	111.02
1	A	138	TYR	CA-C-N	5.06	124.84	119.28
1	A	138	TYR	C-N-CA	5.06	124.84	119.28
1	D	55	ILE	CA-C-N	5.05	126.16	119.84
1	D	55	ILE	C-N-CA	5.05	126.16	119.84
1	B	351	PHE	N-CA-C	5.01	117.66	109.59

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	2774	0	2762	53	0
1	B	2774	0	2762	88	0
1	C	2774	0	2762	74	0
1	D	2774	0	2762	83	0
2	A	10	0	14	4	0
3	A	128	0	0	0	0
3	B	69	0	0	1	0
3	C	124	0	0	2	0
3	D	64	0	0	0	0
All	All	11491	0	11062	296	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 13.

All (296) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:226:VAL:HG22	1:C:227:SER:H	1.20	1.01
1:D:245:ASN:HD22	1:D:289:LYS:HD2	1.22	1.01
1:D:257:MSE:HE2	1:D:260:ILE:HG21	1.41	0.98
1:C:21:GLY:HA2	1:C:253:ASN:HD21	1.37	0.90
1:A:70:LEU:H	1:A:85:ASN:HD21	1.18	0.88
1:C:70:LEU:H	1:C:85:ASN:HD21	1.23	0.84
1:B:70:LEU:H	1:B:85:ASN:HD21	1.25	0.82
1:C:21:GLY:H	1:C:279:ASN:HD21	1.26	0.81
1:D:245:ASN:ND2	1:D:289:LYS:HD2	1.95	0.81
1:D:70:LEU:H	1:D:85:ASN:HD21	1.28	0.81
1:B:331:LYS:O	1:B:335:GLU:HG3	1.83	0.78
1:A:9:LEU:HD13	1:A:360:GLU:HG2	1.66	0.77
1:A:20:ALA:HB1	1:A:257:MSE:HE1	1.66	0.76
1:A:41:ILE:HD12	1:A:124:ALA:HB3	1.67	0.75
1:D:39:ARG:HH21	1:D:349:LEU:HB3	1.49	0.75
1:D:296:THR:H	1:D:299:LEU:HD12	1.50	0.75
1:C:226:VAL:HG22	1:C:227:SER:N	2.00	0.75
1:D:184:TRP:O	1:D:188:ARG:HD2	1.86	0.75
1:B:21:GLY:HA2	1:B:253:ASN:HD21	1.51	0.74
1:A:21:GLY:HA2	1:A:253:ASN:HD21	1.52	0.74
1:A:21:GLY:H	1:A:279:ASN:HD21	1.36	0.73
1:D:257:MSE:HE2	1:D:260:ILE:CG2	2.18	0.73
1:D:21:GLY:HA2	1:D:253:ASN:HD21	1.54	0.72
1:D:257:MSE:HA	1:D:257:MSE:HE3	1.69	0.72
1:C:289:LYS:HE3	3:C:427:HOH:O	1.88	0.72
1:C:277:LYS:O	1:C:281:ARG:HG2	1.91	0.71
1:C:309:MSE:HE3	1:C:337:LEU:HA	1.72	0.70
1:A:277:LYS:O	1:A:281:ARG:HG2	1.92	0.70
1:B:21:GLY:H	1:B:279:ASN:HD21	1.38	0.70
1:A:310:GLY:O	1:A:329:LYS:HE3	1.92	0.69
1:D:257:MSE:HA	1:D:257:MSE:CE	2.23	0.68
1:A:194:LEU:H	2:A:1:PGE:H22	1.59	0.67
1:C:178:THR:HB	1:C:206:GLN:HB2	1.77	0.67
1:A:194:LEU:N	2:A:1:PGE:H22	2.10	0.67
1:C:30:HIS:CD2	1:C:254:ASN:HD21	2.13	0.67
1:D:279:ASN:ND2	1:D:314:LYS:HE3	2.08	0.67
1:C:155:LEU:HD12	1:C:155:LEU:C	2.21	0.66
1:A:230:LYS:HD2	1:B:241:GLU:OE2	1.96	0.66
1:A:281:ARG:O	1:A:285:GLU:HG2	1.96	0.66
1:B:8:LYS:HB3	1:B:46:GLU:HB3	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:LYS:O	1:A:234:GLN:HG3	1.95	0.65
1:D:52:GLU:CD	1:D:54:TRP:HE1	2.05	0.65
1:B:16:LYS:HD2	1:B:16:LYS:C	2.22	0.64
1:C:222:THR:O	1:C:224:THR:N	2.31	0.63
1:D:332:GLU:CD	1:D:332:GLU:H	2.05	0.63
1:C:196:TRP:CH2	1:C:200:GLU:HG2	2.34	0.63
1:A:183:GLU:HG2	1:A:187:HIS:CE1	2.34	0.63
1:B:340:GLU:HA	1:B:343:LYS:HE2	1.82	0.62
1:A:41:ILE:HD12	1:A:124:ALA:CB	2.29	0.62
1:D:329:LYS:HB2	1:D:333:LEU:HD12	1.81	0.62
1:B:309:MSE:HE3	1:B:309:MSE:HA	1.82	0.62
1:D:196:TRP:CH2	1:D:200:GLU:HG2	2.36	0.61
1:B:173:GLY:HA2	1:B:353:THR:CG2	2.30	0.61
1:A:173:GLY:HA2	1:A:353:THR:CG2	2.30	0.60
1:D:180:PHE:HB3	1:D:205:LEU:HB2	1.84	0.60
1:D:334:ALA:O	1:D:338:VAL:HG23	2.00	0.60
1:D:23:TYR:O	1:D:26:VAL:HG22	2.02	0.59
1:A:88:THR:HG23	1:A:98:LEU:CD1	2.32	0.59
1:D:16:LYS:HE2	1:D:16:LYS:C	2.28	0.59
1:B:277:LYS:O	1:B:281:ARG:HG2	2.01	0.58
1:B:302:LEU:C	1:B:302:LEU:HD23	2.28	0.58
1:C:195:GLU:O	1:C:199:LYS:HD2	2.02	0.58
1:B:38:ASN:HB3	1:B:352:HIS:HB2	1.85	0.58
1:B:336:LYS:O	1:B:339:ASN:HB3	2.03	0.58
1:B:111:ASP:C	1:B:113:SER:H	2.10	0.58
1:B:180:PHE:HB3	1:B:205:LEU:HB2	1.86	0.58
1:A:280:ARG:O	1:A:284:GLN:HG3	2.03	0.58
1:B:54:TRP:CZ3	1:B:102:LYS:HE3	2.39	0.58
1:D:285:GLU:O	1:D:289:LYS:HG2	2.04	0.58
1:C:66:ILE:HD13	1:C:100:PRO:HD3	1.86	0.57
1:B:23:TYR:O	1:B:26:VAL:HG22	2.04	0.57
1:B:222:THR:CG2	1:B:299:LEU:HD23	2.33	0.57
1:B:260:ILE:HG12	1:B:275:SER:HB3	1.86	0.57
1:C:155:LEU:HD12	1:C:156:VAL:N	2.19	0.57
1:D:221:TRP:CE2	1:D:223:GLY:HA2	2.38	0.57
1:A:112:GLN:CD	1:A:112:GLN:H	2.11	0.57
1:B:285:GLU:HG3	1:B:289:LYS:HE3	1.86	0.57
1:D:309:MSE:HE3	1:D:337:LEU:N	2.20	0.57
1:B:162:SER:O	1:B:163:CYS:HB2	2.05	0.57
1:B:39:ARG:HH12	1:B:349:LEU:HD22	1.69	0.57
1:B:222:THR:HG21	1:B:299:LEU:HD23	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:20:ALA:HB1	1:D:257:MSE:HE1	1.88	0.56
1:B:230:LYS:HB2	1:B:230:LYS:NZ	2.20	0.56
1:B:214:VAL:HB	1:B:215:PRO:HD3	1.85	0.56
1:C:226:VAL:CG2	1:C:227:SER:H	1.98	0.56
1:A:22:GLU:O	1:A:23:TYR:HB2	2.05	0.56
1:D:359:THR:HG22	1:D:361:GLY:H	1.69	0.56
1:A:162:SER:O	1:A:163:CYS:HB2	2.06	0.56
1:D:162:SER:O	1:D:163:CYS:HB2	2.06	0.56
1:D:280:ARG:O	1:D:284:GLN:HG3	2.06	0.56
1:C:231:LEU:HD23	1:C:292:VAL:HG21	1.88	0.55
1:D:74:GLY:HA3	1:D:77:TRP:HD1	1.70	0.55
1:D:55:ILE:HD11	1:D:58:TYR:CD2	2.41	0.55
1:A:45:LEU:C	1:A:45:LEU:HD23	2.32	0.55
1:A:138:TYR:CD2	1:A:141:ILE:HB	2.42	0.55
1:C:88:THR:HG23	1:C:98:LEU:CD1	2.36	0.55
1:C:45:LEU:C	1:C:45:LEU:HD23	2.31	0.54
1:B:20:ALA:HB1	1:B:279:ASN:HD22	1.72	0.54
1:D:38:ASN:HB3	1:D:352:HIS:HB2	1.90	0.54
1:D:288:THR:C	1:D:290:ALA:H	2.16	0.54
1:B:31:THR:O	1:B:257:MSE:HG2	2.06	0.54
1:B:45:LEU:C	1:B:45:LEU:HD23	2.32	0.54
1:C:194:LEU:O	1:C:198:MSE:HG2	2.08	0.54
1:C:228:THR:HG23	1:C:229:GLY:N	2.23	0.54
1:C:152:LEU:O	1:C:156:VAL:HG13	2.08	0.54
1:C:86:ILE:HG23	1:C:194:LEU:HD21	1.90	0.54
1:B:39:ARG:HG2	1:B:39:ARG:HH11	1.72	0.53
1:C:180:PHE:CE1	1:C:182:GLN:HG2	2.42	0.53
1:B:302:LEU:HB2	1:B:346:ILE:HD13	1.91	0.53
1:D:302:LEU:C	1:D:302:LEU:HD23	2.34	0.53
1:B:43:LEU:HD23	1:B:43:LEU:C	2.34	0.53
1:B:70:LEU:HG	1:B:72:PRO:HD3	1.91	0.53
1:B:43:LEU:HD23	1:B:44:THR:N	2.24	0.52
1:C:222:THR:O	1:C:224:THR:HG22	2.10	0.52
1:D:329:LYS:HB2	1:D:333:LEU:CD1	2.39	0.52
1:D:330:THR:OG1	1:D:333:LEU:HG	2.09	0.52
1:D:310:GLY:O	1:D:329:LYS:HE2	2.10	0.52
1:D:155:LEU:HD12	1:D:156:VAL:N	2.24	0.52
1:B:54:TRP:CD2	1:B:61:PRO:HB3	2.45	0.52
1:B:70:LEU:N	1:B:85:ASN:HD21	2.02	0.52
1:C:74:GLY:HA3	1:C:76:HIS:CE1	2.45	0.52
1:A:23:TYR:O	1:A:26:VAL:HG22	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:214:VAL:HB	1:C:215:PRO:HD3	1.92	0.51
1:D:43:LEU:C	1:D:43:LEU:HD23	2.35	0.51
1:D:155:LEU:HD12	1:D:155:LEU:C	2.34	0.51
1:D:226:VAL:HG22	1:D:320:GLY:O	2.10	0.51
1:D:45:LEU:C	1:D:45:LEU:HD23	2.36	0.51
1:D:224:THR:O	1:D:225:PRO:C	2.52	0.51
1:B:45:LEU:HD23	1:B:46:GLU:N	2.26	0.51
1:B:54:TRP:CE3	1:B:102:LYS:HE3	2.46	0.51
1:B:66:ILE:HD12	1:B:66:ILE:H	1.75	0.51
1:C:162:SER:O	1:C:163:CYS:HB2	2.10	0.51
1:D:12:LYS:HD3	1:D:357:GLN:NE2	2.27	0.50
1:D:316:SER:HB2	1:D:325:ILE:HG23	1.93	0.50
1:D:341:TRP:HB3	1:D:346:ILE:HB	1.92	0.50
1:D:55:ILE:HD11	1:D:58:TYR:HD2	1.76	0.50
1:A:214:VAL:N	1:A:215:PRO:CD	2.74	0.50
1:B:54:TRP:CZ2	1:B:61:PRO:HG3	2.47	0.50
1:A:20:ALA:HB1	1:A:279:ASN:HD22	1.76	0.50
1:D:173:GLY:HA2	1:D:353:THR:CG2	2.41	0.50
1:B:355:ARG:HG2	1:B:355:ARG:HH11	1.76	0.50
1:A:148:LYS:HB3	1:A:198:MSE:HE1	1.93	0.50
1:C:301:GLU:HB3	1:C:344:LEU:HD13	1.93	0.50
1:C:222:THR:HG21	1:C:299:LEU:HG	1.94	0.50
1:C:228:THR:O	1:C:230:LYS:N	2.45	0.50
1:C:226:VAL:HG13	1:C:227:SER:N	2.27	0.50
1:B:193:SER:OG	1:B:195:GLU:HG2	2.11	0.50
1:B:226:VAL:O	1:B:227:SER:HB3	2.11	0.50
1:B:239:LYS:HG3	1:B:246:TYR:CD2	2.46	0.50
1:B:88:THR:HG23	1:B:98:LEU:CD1	2.42	0.49
1:A:173:GLY:HA2	1:A:353:THR:HG21	1.93	0.49
1:A:252:ARG:NH2	1:A:278:GLU:OE1	2.46	0.49
1:D:214:VAL:HB	1:D:215:PRO:HD3	1.95	0.49
1:A:112:GLN:OE1	1:A:112:GLN:N	2.46	0.49
1:D:231:LEU:HA	1:D:234:GLN:OE1	2.13	0.49
1:C:62:VAL:HG13	1:C:77:TRP:CZ2	2.48	0.49
1:D:39:ARG:HH21	1:D:349:LEU:CB	2.23	0.49
1:B:280:ARG:O	1:B:284:GLN:HG3	2.12	0.49
1:D:251:THR:O	1:D:255:GLU:HG3	2.13	0.49
1:D:222:THR:HB	1:D:298:LEU:HD23	1.95	0.49
1:C:325:ILE:C	1:C:325:ILE:HD12	2.39	0.48
1:B:341:TRP:HB3	1:B:346:ILE:HB	1.96	0.48
1:C:239:LYS:HB2	1:C:239:LYS:NZ	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:THR:HG23	1:A:98:LEU:HD11	1.96	0.48
1:D:222:THR:HG21	1:D:299:LEU:HD23	1.96	0.48
1:C:310:GLY:O	1:C:329:LYS:HE2	2.13	0.48
1:B:256:ILE:O	1:B:259:GLN:HB2	2.14	0.48
1:D:74:GLY:HA3	1:D:77:TRP:CD1	2.49	0.48
1:A:194:LEU:H	2:A:1:PGE:C2	2.27	0.48
1:B:39:ARG:NH1	1:B:349:LEU:HD22	2.29	0.48
1:B:226:VAL:HG11	1:B:320:GLY:O	2.14	0.48
1:A:274:SER:O	1:A:278:GLU:HG3	2.14	0.47
1:D:300:LYS:HG2	1:D:304:ASP:OD2	2.13	0.47
1:C:215:PRO:HG3	1:C:269:GLU:HG2	1.96	0.47
1:B:230:LYS:HG2	1:B:234:GLN:OE1	2.14	0.47
1:B:251:THR:O	1:B:255:GLU:HG3	2.14	0.47
1:C:228:THR:HG23	1:C:229:GLY:H	1.78	0.47
1:A:155:LEU:C	1:A:155:LEU:HD12	2.39	0.47
1:A:277:LYS:NZ	1:A:311:GLY:O	2.47	0.47
1:C:341:TRP:HB3	1:C:346:ILE:HB	1.96	0.47
1:A:43:LEU:C	1:A:43:LEU:HD23	2.40	0.47
1:C:226:VAL:HG22	1:C:228:THR:H	1.79	0.47
1:A:112:GLN:CD	1:A:112:GLN:N	2.72	0.46
1:A:273:TYR:O	1:A:277:LYS:HG3	2.15	0.46
1:C:236:HIS:O	1:C:239:LYS:HG2	2.16	0.46
1:B:222:THR:OG1	1:B:298:LEU:HG	2.16	0.46
1:C:112:GLN:H	1:C:112:GLN:CD	2.23	0.46
1:C:309:MSE:HE3	1:C:337:LEU:CA	2.42	0.46
1:D:338:VAL:O	1:D:342:GLU:HG3	2.15	0.46
1:A:66:ILE:HD13	1:A:100:PRO:HD3	1.97	0.46
1:D:16:LYS:HD3	1:D:165:ASP:CB	2.46	0.46
1:C:43:LEU:C	1:C:43:LEU:HD23	2.41	0.46
1:B:91:LEU:CD1	1:B:98:LEU:HD21	2.46	0.46
1:B:230:LYS:O	1:B:234:GLN:HG3	2.14	0.46
1:C:38:ASN:ND2	1:C:352:HIS:HB2	2.31	0.46
1:D:331:LYS:NZ	1:D:331:LYS:HB3	2.30	0.46
1:C:292:VAL:HG22	1:C:293:ASN:H	1.80	0.46
1:B:260:ILE:CG1	1:B:275:SER:HB3	2.46	0.46
1:C:195:GLU:HG3	1:C:196:TRP:N	2.31	0.46
1:D:39:ARG:HG2	1:D:39:ARG:HH11	1.81	0.46
1:D:293:ASN:ND2	1:D:296:THR:HG22	2.31	0.46
1:D:332:GLU:HA	1:D:335:GLU:HG3	1.98	0.46
1:B:338:VAL:O	1:B:342:GLU:HG3	2.16	0.45
1:C:280:ARG:O	1:C:284:GLN:HG3	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:PHE:HB3	1:A:205:LEU:HB2	1.98	0.45
1:A:191:TYR:CE1	1:A:192:LYS:HE2	2.51	0.45
1:B:355:ARG:HG2	1:B:355:ARG:NH1	2.31	0.45
1:B:224:THR:O	1:B:225:PRO:C	2.59	0.45
1:C:292:VAL:HG22	1:C:293:ASN:N	2.31	0.45
1:B:55:ILE:HD11	1:B:58:TYR:CD2	2.52	0.45
1:A:65:PRO:HG2	1:A:68:GLY:HA3	1.98	0.45
1:D:294:ILE:HG23	1:D:318:SER:O	2.16	0.45
1:B:67:GLY:HA2	1:B:98:LEU:HD12	1.99	0.45
1:C:155:LEU:C	1:C:155:LEU:CD1	2.90	0.45
1:D:57:HIS:ND1	1:D:57:HIS:O	2.50	0.45
1:D:349:LEU:O	1:D:350:PRO:C	2.60	0.45
1:C:180:PHE:HB3	1:C:205:LEU:HB2	1.99	0.45
1:D:12:LYS:HD3	1:D:357:GLN:HE21	1.81	0.45
1:B:334:ALA:O	1:B:338:VAL:HG23	2.16	0.44
1:C:269:GLU:HG3	3:C:443:HOH:O	2.17	0.44
1:D:22:GLU:O	1:D:23:TYR:HB2	2.17	0.44
1:A:20:ALA:HB1	1:A:279:ASN:ND2	2.32	0.44
1:B:260:ILE:HG12	1:B:275:SER:CB	2.48	0.44
1:B:349:LEU:O	1:B:350:PRO:C	2.61	0.44
1:C:30:HIS:CD2	1:C:254:ASN:ND2	2.84	0.44
1:A:302:LEU:C	1:A:302:LEU:HD23	2.41	0.44
1:B:224:THR:O	1:B:224:THR:HG23	2.17	0.44
1:D:217:PHE:CE1	1:D:325:ILE:HB	2.53	0.44
1:B:256:ILE:CD1	1:B:278:GLU:HB3	2.48	0.44
1:D:70:LEU:CD1	1:D:72:PRO:HD3	2.47	0.44
1:D:140:GLU:H	1:D:140:GLU:CD	2.25	0.44
1:A:252:ARG:HG2	1:A:282:ILE:HG21	2.00	0.44
1:C:222:THR:HG21	1:C:299:LEU:CD2	2.47	0.44
1:D:339:ASN:O	1:D:343:LYS:HG3	2.18	0.43
1:D:16:LYS:HD3	1:D:165:ASP:HB3	2.00	0.43
1:B:173:GLY:HA2	1:B:353:THR:HG21	1.99	0.43
1:B:111:ASP:C	1:B:113:SER:N	2.75	0.43
1:C:21:GLY:HA2	1:C:253:ASN:ND2	2.18	0.43
1:D:230:LYS:O	1:D:234:GLN:HG3	2.18	0.43
1:B:296:THR:H	1:B:299:LEU:HD12	1.83	0.43
1:C:189:LEU:HD23	1:C:189:LEU:HA	1.90	0.43
1:C:238:PHE:CE1	1:C:242:ASP:HB3	2.53	0.43
1:B:339:ASN:O	1:B:343:LYS:HG3	2.18	0.43
1:C:230:LYS:O	1:C:234:GLN:HG3	2.18	0.43
2:A:1:PGE:H6	1:B:194:LEU:HB2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:207:ILE:HB	1:C:59:GLU:HB2	2.00	0.43
1:C:22:GLU:O	1:C:23:TYR:HB2	2.18	0.43
1:D:112:GLN:OE1	1:D:112:GLN:N	2.52	0.43
1:C:96:ILE:HG21	1:C:137:PHE:HB3	2.01	0.43
1:D:14:PRO:HB3	1:D:40:TYR:CE1	2.54	0.43
1:D:332:GLU:O	1:D:335:GLU:HB2	2.18	0.43
1:A:351:PHE:CD2	1:A:351:PHE:C	2.97	0.43
1:C:31:THR:HG22	1:C:32:ALA:N	2.34	0.43
1:A:21:GLY:N	1:A:279:ASN:HD21	2.12	0.42
1:C:226:VAL:HG22	1:C:228:THR:N	2.33	0.42
1:D:227:SER:O	1:D:228:THR:C	2.62	0.42
1:B:22:GLU:CD	1:B:314:LYS:HZ1	2.26	0.42
1:C:60:ASN:ND2	1:C:61:PRO:HD2	2.33	0.42
1:B:20:ALA:HB1	1:B:279:ASN:ND2	2.33	0.42
1:C:112:GLN:CD	1:C:112:GLN:N	2.77	0.42
1:B:88:THR:HG23	1:B:98:LEU:HD11	2.02	0.42
1:D:302:LEU:HB2	1:D:346:ILE:HD13	2.01	0.42
1:A:251:THR:O	1:A:255:GLU:HG3	2.19	0.42
1:A:309:MSE:HE3	1:A:337:LEU:HA	2.02	0.42
1:C:228:THR:O	1:C:229:GLY:C	2.62	0.42
1:D:293:ASN:HD21	1:D:296:THR:HG22	1.84	0.42
1:C:138:TYR:CD2	1:C:141:ILE:HB	2.54	0.42
1:D:238:PHE:O	1:D:241:GLU:HB3	2.19	0.42
1:D:341:TRP:O	1:D:344:LEU:N	2.52	0.42
1:B:222:THR:HG23	1:B:299:LEU:HD23	2.01	0.42
1:C:46:GLU:HG2	1:C:102:LYS:HB3	2.01	0.42
1:C:25:VAL:HB	1:C:32:ALA:HB2	2.02	0.42
1:B:41:ILE:HD13	1:B:107:THR:HB	2.02	0.41
1:A:257:MSE:O	1:A:260:ILE:HG22	2.20	0.41
1:A:228:THR:O	1:A:229:GLY:C	2.64	0.41
1:A:293:ASN:ND2	1:A:296:THR:HG22	2.35	0.41
1:C:75:GLU:O	1:C:75:GLU:HG2	2.20	0.41
1:D:134:MSE:O	1:D:138:TYR:C	2.63	0.41
1:B:43:LEU:HA	1:B:104:VAL:O	2.20	0.41
1:D:257:MSE:HE3	1:D:257:MSE:CA	2.46	0.41
1:B:55:ILE:HG22	1:B:105:ILE:HD12	2.02	0.41
1:B:74:GLY:HA3	1:B:77:TRP:HD1	1.85	0.41
1:B:214:VAL:HB	1:B:269:GLU:OE2	2.21	0.41
1:C:16:LYS:C	1:C:16:LYS:HD2	2.46	0.41
1:D:279:ASN:HD22	1:D:314:LYS:HE3	1.83	0.41
1:D:296:THR:HG23	1:D:320:GLY:HA2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:ILE:HG12	1:B:100:PRO:HD3	2.03	0.41
1:A:90:PHE:CE1	1:A:198:MSE:HG3	2.56	0.40
1:B:112:GLN:H	1:B:112:GLN:CD	2.29	0.40
1:B:196:TRP:CH2	1:B:200:GLU:HG2	2.57	0.40
1:C:302:LEU:HD23	1:C:302:LEU:C	2.46	0.40
1:B:231:LEU:N	1:B:231:LEU:HD12	2.36	0.40
1:C:231:LEU:HD11	1:C:320:GLY:HA3	2.04	0.40
1:B:62:VAL:HG13	1:B:77:TRP:CZ2	2.56	0.40
1:B:66:ILE:HD12	1:B:66:ILE:N	2.37	0.40
1:B:183:GLU:HG3	3:B:373:HOH:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	353/365 (97%)	342 (97%)	11 (3%)	0	100	100
1	B	353/365 (97%)	335 (95%)	16 (4%)	2 (1%)	21	18
1	C	353/365 (97%)	336 (95%)	11 (3%)	6 (2%)	7	3
1	D	353/365 (97%)	322 (91%)	24 (7%)	7 (2%)	6	2
All	All	1412/1460 (97%)	1335 (94%)	62 (4%)	15 (1%)	11	8

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	227	SER
1	C	223	GLY
1	C	226	VAL
1	C	228	THR
1	C	229	GLY

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Mol	Chain	Res	Type
1	D	223	GLY
1	B	112	GLN
1	C	227	SER
1	D	225	PRO
1	D	289	LYS
1	D	74	GLY
1	D	241	GLU
1	D	288	THR
1	C	56	PRO
1	D	350	PRO

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	298/299 (100%)	294 (99%)	4 (1%)	61	69
1	B	298/299 (100%)	293 (98%)	5 (2%)	53	62
1	C	298/299 (100%)	291 (98%)	7 (2%)	44	51
1	D	298/299 (100%)	291 (98%)	7 (2%)	44	51
All	All	1192/1196 (100%)	1169 (98%)	23 (2%)	50	58

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LYS
1	A	41	ILE
1	A	112	GLN
1	A	155	LEU
1	B	16	LYS
1	B	54	TRP
1	B	222	THR
1	B	230	LYS
1	B	309	MSE
1	C	16	LYS
1	C	56	PRO

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Mol	Chain	Res	Type
1	C	155	LEU
1	C	194	LEU
1	C	195	GLU
1	C	199	LYS
1	C	239	LYS
1	D	16	LYS
1	D	140	GLU
1	D	155	LEU
1	D	188	ARG
1	D	225	PRO
1	D	257	MSE
1	D	340	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	85	ASN
1	A	240	GLN
1	A	245	ASN
1	A	253	ASN
1	A	259	GLN
1	A	279	ASN
1	A	352	HIS
1	A	357	GLN
1	B	60	ASN
1	B	76	HIS
1	B	85	ASN
1	B	240	GLN
1	B	245	ASN
1	B	253	ASN
1	B	259	GLN
1	B	279	ASN
1	B	293	ASN
1	B	357	GLN
1	C	60	ASN
1	C	85	ASN
1	C	182	GLN
1	C	253	ASN
1	C	254	ASN
1	C	279	ASN
1	C	284	GLN

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Mol	Chain	Res	Type
1	C	339	ASN
1	C	348	HIS
1	D	85	ASN
1	D	131	ASN
1	D	245	ASN
1	D	248	HIS
1	D	253	ASN
1	D	259	GLN
1	D	293	ASN
1	D	339	ASN
1	D	348	HIS
1	D	357	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	PGE	A	1	-	9,9,9	0.93	0	8,8,8	2.57	3 (37%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PGE	A	1	-	-	3/7/7/7	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	PGE	O3-C4-C3	-5.00	87.57	110.35
2	A	1	PGE	C3-O2-C2	-4.28	94.52	113.26
2	A	1	PGE	O3-C5-C6	-2.14	100.67	110.11

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1	PGE	O2-C3-C4-O3
2	A	1	PGE	O1-C1-C2-O2
2	A	1	PGE	O3-C5-C6-O4

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1	PGE	4	0

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	347/365 (95%)	0.07	7 (2%) 65 67	12, 24, 45, 65	0
1	B	347/365 (95%)	0.80	43 (12%) 8 8	16, 37, 63, 81	0
1	C	347/365 (95%)	0.32	24 (6%) 23 24	13, 24, 51, 75	0
1	D	347/365 (95%)	1.23	94 (27%) 1 1	17, 40, 73, 90	0
All	All	1388/1460 (95%)	0.60	168 (12%) 8 9	12, 30, 66, 90	0

All (168) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	223	GLY	13.6
1	C	226	VAL	10.9
1	D	226	VAL	10.6
1	C	224	THR	10.3
1	C	227	SER	9.0
1	B	226	VAL	8.1
1	B	222	THR	7.3
1	D	228	THR	7.2
1	D	229	GLY	7.2
1	C	228	THR	7.1
1	C	229	GLY	5.6
1	D	224	THR	5.3
1	B	227	SER	5.3
1	D	231	LEU	5.2
1	B	225	PRO	5.1
1	D	227	SER	5.0
1	D	232	VAL	5.0
1	C	225	PRO	4.9
1	D	225	PRO	4.9
1	D	223	GLY	4.6
1	D	110	ILE	4.6

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Mol	Chain	Res	Type	RSRZ
1	B	224	THR	4.5
1	B	228	THR	4.5
1	B	229	GLY	4.5
1	D	361	GLY	4.4
1	D	289	LYS	4.3
1	D	245	ASN	4.1
1	B	54	TRP	4.0
1	D	221	TRP	3.9
1	D	294	ILE	3.9
1	D	235	ILE	3.9
1	B	223	GLY	3.8
1	D	345	GLY	3.8
1	D	155	LEU	3.7
1	D	282	ILE	3.7
1	D	298	LEU	3.7
1	B	361	GLY	3.7
1	B	339	ASN	3.6
1	C	351	PHE	3.6
1	D	243	SER	3.6
1	D	237	ALA	3.6
1	D	299	LEU	3.5
1	D	306	ALA	3.5
1	D	252	ARG	3.5
1	A	361	GLY	3.5
1	D	297	SER	3.5
1	D	311	GLY	3.4
1	D	338	VAL	3.3
1	C	231	LEU	3.3
1	D	57	HIS	3.3
1	A	241	GLU	3.3
1	D	238	PHE	3.2
1	D	288	THR	3.2
1	D	292	VAL	3.2
1	D	352	HIS	3.2
1	C	113	SER	3.2
1	C	239	LYS	3.1
1	D	283	LEU	3.1
1	D	346	ILE	3.1
1	B	337	LEU	3.0
1	C	138	TYR	3.0
1	D	117	TYR	3.0
1	D	310	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	222	THR	2.9
1	B	7	ASN	2.9
1	D	341	TRP	2.9
1	C	230	LYS	2.9
1	C	56	PRO	2.9
1	D	339	ASN	2.9
1	D	281	ARG	2.8
1	D	302	LEU	2.8
1	A	227	SER	2.8
1	D	359	THR	2.8
1	D	293	ASN	2.8
1	D	301	GLU	2.8
1	D	248	HIS	2.7
1	D	322	ASP	2.7
1	B	292	VAL	2.7
1	B	344	LEU	2.7
1	D	286	LEU	2.7
1	B	293	ASN	2.6
1	D	73	ASP	2.6
1	B	110	ILE	2.6
1	B	57	HIS	2.6
1	D	333	LEU	2.6
1	D	337	LEU	2.6
1	A	112	GLN	2.5
1	B	55	ILE	2.5
1	C	361	GLY	2.5
1	B	98	LEU	2.5
1	B	231	LEU	2.5
1	D	349	LEU	2.5
1	D	115	ALA	2.5
1	D	312	ALA	2.5
1	B	343	LYS	2.5
1	B	112	GLN	2.4
1	B	234	GLN	2.4
1	D	58	TYR	2.4
1	D	114	GLY	2.4
1	D	319	GLY	2.4
1	D	55	ILE	2.4
1	D	303	ALA	2.4
1	D	320	GLY	2.4
1	B	336	LYS	2.4
1	D	199	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	344	LEU	2.4
1	D	241	GLU	2.4
1	B	291	GLY	2.4
1	D	220	GLY	2.4
1	D	219	VAL	2.4
1	A	228	THR	2.4
1	D	246	TYR	2.3
1	D	326	ALA	2.3
1	D	240	GLN	2.3
1	B	232	VAL	2.3
1	D	112	GLN	2.3
1	B	53	LEU	2.3
1	B	109	LEU	2.3
1	B	299	LEU	2.3
1	D	74	GLY	2.3
1	D	109	LEU	2.3
1	D	242	ASP	2.3
1	B	360	GLU	2.3
1	C	232	VAL	2.3
1	C	241	GLU	2.3
1	A	73	ASP	2.3
1	B	74	GLY	2.3
1	D	116	LYS	2.3
1	D	291	GLY	2.3
1	D	296	THR	2.2
1	D	347	LYS	2.2
1	B	340	GLU	2.2
1	D	335	GLU	2.2
1	B	341	TRP	2.2
1	C	222	THR	2.2
1	B	138	TYR	2.2
1	D	49	GLU	2.2
1	C	240	GLN	2.2
1	D	290	ALA	2.2
1	D	61	PRO	2.1
1	B	62	VAL	2.1
1	C	112	GLN	2.1
1	C	57	HIS	2.1
1	D	236	HIS	2.1
1	D	230	LYS	2.1
1	D	332	GLU	2.1
1	D	77	TRP	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	351	PHE	2.1
1	A	229	GLY	2.1
1	C	98	LEU	2.1
1	D	244	LYS	2.1
1	D	300	LYS	2.1
1	D	348	HIS	2.1
1	C	95	GLY	2.1
1	B	230	LYS	2.1
1	B	333	LEU	2.1
1	C	360	GLU	2.1
1	D	295	GLU	2.1
1	B	116	LYS	2.1
1	B	289	LYS	2.1
1	D	323	CYS	2.0
1	B	241	GLU	2.0
1	B	342	GLU	2.0
1	D	233	SER	2.0
1	D	7	ASN	2.0
1	D	8	LYS	2.0
1	D	239	LYS	2.0
1	B	345	GLY	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	PGE	A	1	10/10	0.65	0.26	35,46,49,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.