



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 03:56 AM UTC

PDB ID : 2Z3R / pdb_00002z3r
Title : Crystal structure of the IL-15/IL-15Ra complex
Authors : Chirifu, M.; Yamagata, Y.; Davis, S.J.; Ikemizu, S.
Deposited on : 2007-06-05
Resolution : 2.00 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

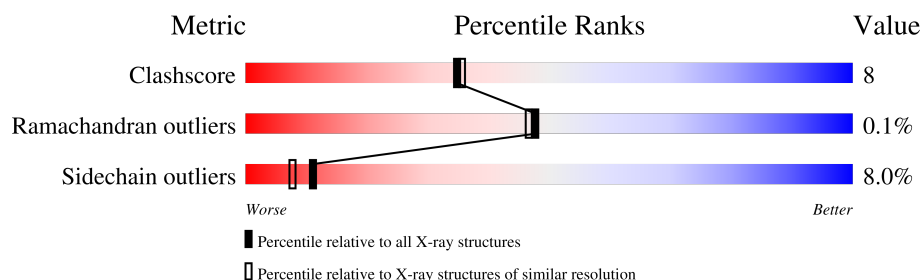
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

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



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

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

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	119	76% 15% 5%
1	C	119	76% 16% 5%
1	E	119	85% 14%
1	G	119	76% 20% ..
1	I	119	76% 21% ..
1	K	119	80% 18%
1	M	119	79% 18% ..
1	O	119	76% 18% ..

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Mol	Chain	Length	Quality of chain
2	B	107	
2	D	107	
2	F	107	
2	H	107	
2	J	107	
2	L	107	
2	N	107	
2	P	107	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12517 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 Interleukin-15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	113	Total	C	N	O	S	0	0	0
			883	556	140	179	8			
1	C	116	Total	C	N	O	S	0	0	0
			904	568	143	185	8			
1	E	119	Total	C	N	O	S	0	0	0
			924	578	148	190	8			
1	G	119	Total	C	N	O	S	0	0	0
			924	578	148	190	8			
1	I	116	Total	C	N	O	S	0	0	0
			904	568	143	185	8			
1	K	116	Total	C	N	O	S	0	0	0
			904	568	143	185	8			
1	M	116	Total	C	N	O	S	0	0	0
			904	568	143	185	8			
1	O	118	Total	C	N	O	S	0	0	0
			917	575	147	187	8			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	ALA	-	expression tag	UNP P40933
A	-3	MET	-	expression tag	UNP P40933
A	-2	ALA	-	expression tag	UNP P40933
A	-1	ILE	-	expression tag	UNP P40933
A	0	SER	-	expression tag	UNP P40933
C	-4	ALA	-	expression tag	UNP P40933
C	-3	MET	-	expression tag	UNP P40933
C	-2	ALA	-	expression tag	UNP P40933
C	-1	ILE	-	expression tag	UNP P40933
C	0	SER	-	expression tag	UNP P40933
E	-4	ALA	-	expression tag	UNP P40933
E	-3	MET	-	expression tag	UNP P40933
E	-2	ALA	-	expression tag	UNP P40933

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-1	ILE	-	expression tag	UNP P40933
E	0	SER	-	expression tag	UNP P40933
G	-4	ALA	-	expression tag	UNP P40933
G	-3	MET	-	expression tag	UNP P40933
G	-2	ALA	-	expression tag	UNP P40933
G	-1	ILE	-	expression tag	UNP P40933
G	0	SER	-	expression tag	UNP P40933
I	-4	ALA	-	expression tag	UNP P40933
I	-3	MET	-	expression tag	UNP P40933
I	-2	ALA	-	expression tag	UNP P40933
I	-1	ILE	-	expression tag	UNP P40933
I	0	SER	-	expression tag	UNP P40933
K	-4	ALA	-	expression tag	UNP P40933
K	-3	MET	-	expression tag	UNP P40933
K	-2	ALA	-	expression tag	UNP P40933
K	-1	ILE	-	expression tag	UNP P40933
K	0	SER	-	expression tag	UNP P40933
M	-4	ALA	-	expression tag	UNP P40933
M	-3	MET	-	expression tag	UNP P40933
M	-2	ALA	-	expression tag	UNP P40933
M	-1	ILE	-	expression tag	UNP P40933
M	0	SER	-	expression tag	UNP P40933
O	-4	ALA	-	expression tag	UNP P40933
O	-3	MET	-	expression tag	UNP P40933
O	-2	ALA	-	expression tag	UNP P40933
O	-1	ILE	-	expression tag	UNP P40933
O	0	SER	-	expression tag	UNP P40933

- Molecule 2 is a protein called Interleukin-15 receptor alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	74	Total	C	N	O	S	0	0	0
			584	367	106	106	5			
2	D	76	Total	C	N	O	S	0	0	0
			594	375	105	108	6			
2	F	73	Total	C	N	O	S	0	0	0
			573	361	102	105	5			
2	H	77	Total	C	N	O	S	0	0	0
			599	378	106	109	6			
2	J	78	Total	C	N	O	S	0	0	0
			610	384	110	110	6			
2	L	75	Total	C	N	O	S	0	0	0
			592	373	107	107	5			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	N	76	Total	C	N	O	S	0	0	0
			594	375	105	108	6			
2	P	72	Total	C	N	O	S	0	0	0
			564	356	100	103	5			

There are 40 discrepancies between the modelled and reference sequences:

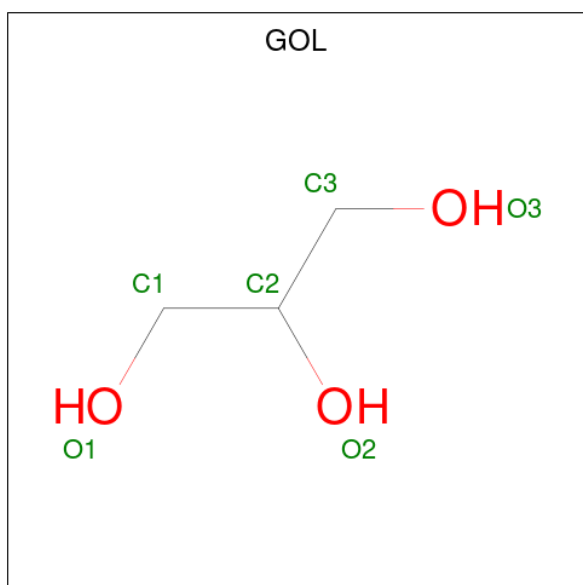
Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	ALA	-	expression tag	UNP Q13261
B	-3	MET	-	expression tag	UNP Q13261
B	-2	ALA	-	expression tag	UNP Q13261
B	-1	ILE	-	expression tag	UNP Q13261
B	0	SER	-	expression tag	UNP Q13261
D	-4	ALA	-	expression tag	UNP Q13261
D	-3	MET	-	expression tag	UNP Q13261
D	-2	ALA	-	expression tag	UNP Q13261
D	-1	ILE	-	expression tag	UNP Q13261
D	0	SER	-	expression tag	UNP Q13261
F	-4	ALA	-	expression tag	UNP Q13261
F	-3	MET	-	expression tag	UNP Q13261
F	-2	ALA	-	expression tag	UNP Q13261
F	-1	ILE	-	expression tag	UNP Q13261
F	0	SER	-	expression tag	UNP Q13261
H	-4	ALA	-	expression tag	UNP Q13261
H	-3	MET	-	expression tag	UNP Q13261
H	-2	ALA	-	expression tag	UNP Q13261
H	-1	ILE	-	expression tag	UNP Q13261
H	0	SER	-	expression tag	UNP Q13261
J	-4	ALA	-	expression tag	UNP Q13261
J	-3	MET	-	expression tag	UNP Q13261
J	-2	ALA	-	expression tag	UNP Q13261
J	-1	ILE	-	expression tag	UNP Q13261
J	0	SER	-	expression tag	UNP Q13261
L	-4	ALA	-	expression tag	UNP Q13261
L	-3	MET	-	expression tag	UNP Q13261
L	-2	ALA	-	expression tag	UNP Q13261
L	-1	ILE	-	expression tag	UNP Q13261
L	0	SER	-	expression tag	UNP Q13261
N	-4	ALA	-	expression tag	UNP Q13261
N	-3	MET	-	expression tag	UNP Q13261
N	-2	ALA	-	expression tag	UNP Q13261
N	-1	ILE	-	expression tag	UNP Q13261

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Chain	Residue	Modelled	Actual	Comment	Reference
N	0	SER	-	expression tag	UNP Q13261
P	-4	ALA	-	expression tag	UNP Q13261
P	-3	MET	-	expression tag	UNP Q13261
P	-2	ALA	-	expression tag	UNP Q13261
P	-1	ILE	-	expression tag	UNP Q13261
P	0	SER	-	expression tag	UNP Q13261

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	O	1	Total C O 6 3 3	0	0
3	O	1	Total C O 6 3 3	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	17	Total O 17 17	0	0
4	B	32	Total O 32 32	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	35	Total 35	O 35	0	0
4	D	26	Total 26	O 26	0	0
4	E	32	Total 32	O 32	0	0
4	F	43	Total 43	O 43	0	0
4	G	41	Total 41	O 41	0	0
4	H	43	Total 43	O 43	0	0
4	I	35	Total 35	O 35	0	0
4	J	37	Total 37	O 37	0	0
4	K	34	Total 34	O 34	0	0
4	L	45	Total 45	O 45	0	0
4	M	22	Total 22	O 22	0	0
4	N	21	Total 21	O 21	0	0
4	O	20	Total 20	O 20	0	0
4	P	36	Total 36	O 36	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. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

• Molecule 1: Interleukin-15

Chain A: 



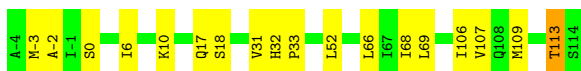
• Molecule 1: Interleukin-15

Chain C: 



• Molecule 1: Interleukin-15

Chain E: 



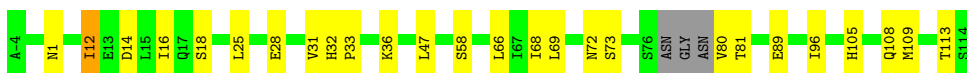
• Molecule 1: Interleukin-15

Chain G: 



• Molecule 1: Interleukin-15

Chain I: 



• Molecule 1: Interleukin-15

Chain K: 



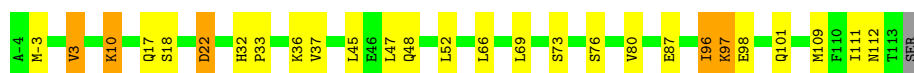
- Molecule 1: Interleukin-15

Chain M: 79% 18% . .



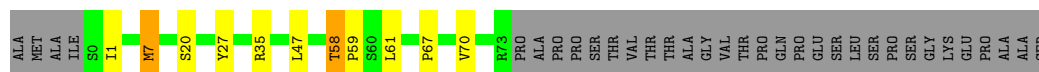
- Molecule 1: Interleukin-15

Chain O: 76% 18% . .



- Molecule 2: Interleukin-15 receptor alpha chain

Chain B: 59% 8% . 31%



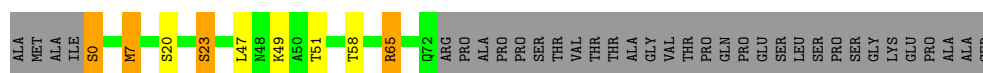
- Molecule 2: Interleukin-15 receptor alpha chain

Chain D: 55% 13% . 29%



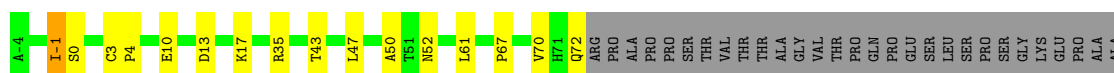
- Molecule 2: Interleukin-15 receptor alpha chain

Chain F: 60% 5% . 32%



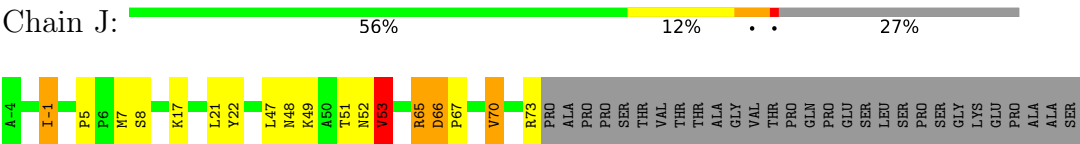
- Molecule 2: Interleukin-15 receptor alpha chain

Chain H: 57% 14% . 28%

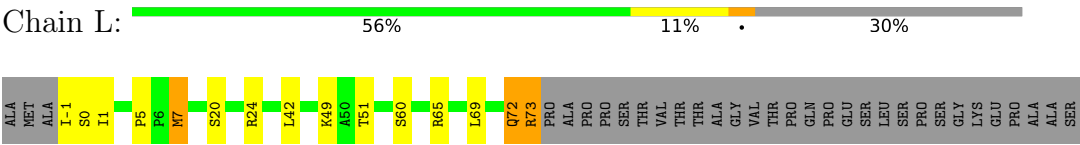


SER

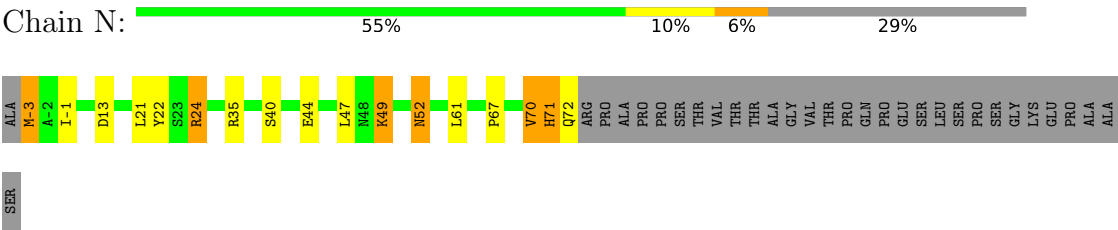
- Molecule 2: Interleukin-15 receptor alpha chain



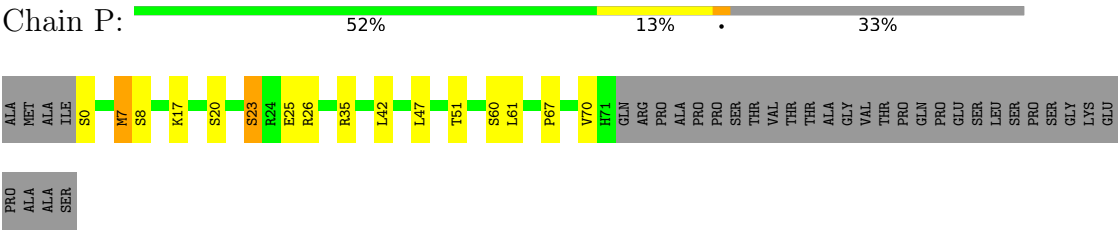
● Molecule 2: Interleukin-15 receptor alpha chain



● Molecule 2: Interleukin-15 receptor alpha chain



● Molecule 2: Interleukin-15 receptor alpha chain



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.80Å 127.03Å 191.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.93 – 2.00	Depositor
% Data completeness (in resolution range)	100.0 (30.93-2.00)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.205 , 0.250	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	12517	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.97	0/893	1.07	0/1205
1	C	1.11	0/914	1.11	0/1233
1	E	1.01	0/935	1.06	0/1263
1	G	1.06	0/935	1.18	3/1263 (0.2%)
1	I	1.19	2/914 (0.2%)	1.13	0/1233
1	K	1.13	0/914	1.13	1/1233 (0.1%)
1	M	1.01	0/914	1.05	0/1233
1	O	0.97	0/928	1.13	3/1255 (0.2%)
2	B	1.04	0/599	1.08	2/814 (0.2%)
2	D	1.10	1/609 (0.2%)	1.05	2/828 (0.2%)
2	F	1.14	0/588	1.07	1/800 (0.1%)
2	H	1.19	1/614 (0.2%)	1.12	0/835
2	J	1.15	1/625 (0.2%)	1.30	4/849 (0.5%)
2	L	1.13	1/607 (0.2%)	1.12	3/825 (0.4%)
2	N	1.05	0/609	1.05	1/828 (0.1%)
2	P	1.07	0/579	1.04	1/788 (0.1%)
All	All	1.08	6/12177 (0.0%)	1.11	21/16485 (0.1%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	ILE	C-O	-7.21	1.16	1.24
2	H	43	THR	CA-C	6.93	1.61	1.52
1	I	12	ILE	CA-CB	-6.14	1.47	1.54
2	J	-1	ILE	CA-CB	5.50	1.61	1.54
1	I	68	ILE	C-O	-5.22	1.18	1.24
2	L	42	LEU	C-O	-5.17	1.17	1.23

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	66	ASP	CA-C-N	9.76	130.44	119.32
2	J	66	ASP	C-N-CA	9.76	130.44	119.32
2	F	7	MET	CB-CG-SD	6.74	132.92	112.70
2	N	71	HIS	N-CA-C	6.64	120.94	108.65
2	D	71	HIS	N-CA-C	-6.47	105.51	113.41
2	L	5	PRO	CA-C-N	-5.95	114.48	120.31
2	L	5	PRO	C-N-CA	-5.95	114.48	120.31
2	L	60	SER	N-CA-C	-5.92	106.06	113.28
1	O	96	ILE	CB-CA-C	-5.78	104.47	112.04
1	K	43	PHE	N-CA-C	5.51	117.72	111.11
1	G	80	VAL	CA-C-N	5.41	131.43	121.70
1	G	80	VAL	C-N-CA	5.41	131.43	121.70
2	J	7	MET	CB-CG-SD	5.33	128.71	112.70
2	B	58	THR	CA-C-N	-5.24	115.18	120.31
2	B	58	THR	C-N-CA	-5.24	115.18	120.31
1	O	3	VAL	N-CA-C	-5.22	105.45	110.72
2	D	53	VAL	CB-CA-C	-5.16	105.04	111.08
2	P	42	LEU	N-CA-C	5.12	117.73	110.10
2	J	53	VAL	CB-CA-C	-5.08	104.61	110.91
1	G	50	ILE	N-CA-C	5.02	115.24	110.42
1	O	22	ASP	N-CA-CB	-5.02	103.69	110.67

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	883	0	876	11	0
1	C	904	0	897	22	0
1	E	924	0	913	10	0
1	G	924	0	913	26	0
1	I	904	0	897	20	0
1	K	904	0	897	21	0
1	M	904	0	897	13	0
1	O	917	0	908	17	0
2	B	584	0	583	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	594	0	595	7	0
2	F	573	0	570	6	0
2	H	599	0	600	8	0
2	J	610	0	613	11	0
2	L	592	0	594	5	0
2	N	594	0	595	9	0
2	P	564	0	562	10	0
3	A	12	0	16	1	0
3	O	12	0	16	0	0
4	A	17	0	0	0	0
4	B	32	0	0	0	0
4	C	35	0	0	2	0
4	D	26	0	0	0	0
4	E	32	0	0	0	0
4	F	43	0	0	2	0
4	G	41	0	0	1	0
4	H	43	0	0	2	0
4	I	35	0	0	0	0
4	J	37	0	0	2	0
4	K	34	0	0	2	0
4	L	45	0	0	2	0
4	M	22	0	0	0	0
4	N	21	0	0	0	0
4	O	20	0	0	2	0
4	P	36	0	0	2	0
All	All	12517	0	11942	180	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 8.

All (180) 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:M:36:LYS:HE3	1:M:113:THR:HG21	1.21	1.18
1:C:-1:ILE:HD11	1:C:114:SER:OG	1.48	1.12
1:G:80:VAL:HG13	1:G:81:THR:HA	1.04	1.03
2:H:17:LYS:HD2	4:H:144:HOH:O	1.59	1.01
1:M:36:LYS:HE3	1:M:113:THR:CG2	1.92	0.98
1:G:80:VAL:CG1	1:G:81:THR:HA	1.95	0.96
2:D:67:PRO:HA	2:D:70:VAL:HG22	1.50	0.93
1:I:1:ASN:HD21	1:I:72:ASN:HD22	1.17	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:96:ILE:HD13	4:G:120:HOH:O	1.79	0.82
1:E:-3:MET:HE3	1:E:-2:ALA:H	1.43	0.82
1:G:80:VAL:HG13	1:G:81:THR:CA	1.99	0.81
1:M:36:LYS:CE	1:M:113:THR:HG21	2.07	0.80
2:F:7:MET:HG2	4:F:137:HOH:O	1.81	0.80
1:G:101:GLN:HE22	1:K:-3:MET:HE1	1.48	0.80
1:C:75:SER:HB3	4:C:117:HOH:O	1.86	0.75
1:G:78:GLY:HA2	1:G:79:ASN:CB	2.17	0.74
2:H:13:ASP:HB2	4:H:142:HOH:O	1.88	0.73
1:G:78:GLY:HA2	1:G:79:ASN:HB2	1.70	0.73
1:C:-1:ILE:CD1	1:C:114:SER:OG	2.35	0.72
1:I:1:ASN:ND2	1:I:72:ASN:HD22	1.88	0.71
1:C:32:HIS:CD2	1:C:34:SER:H	2.10	0.70
1:G:105:HIS:HA	1:G:108:GLN:HE21	1.57	0.69
1:I:25:LEU:HD12	1:I:96:ILE:HD12	1.73	0.69
1:G:77:ASN:N	1:G:78:GLY:HA3	2.08	0.69
2:L:69:LEU:O	2:L:72:GLN:HG3	1.92	0.68
2:P:17:LYS:HE2	2:P:25:GLU:OE1	1.92	0.68
1:M:18:SER:HB2	1:O:18:SER:HB3	1.75	0.68
1:C:75:SER:CB	4:C:117:HOH:O	2.42	0.68
1:G:23:ALA:O	1:G:96:ILE:CD1	2.41	0.68
2:B:35:ARG:HG3	2:B:61:LEU:CD1	2.24	0.68
1:I:25:LEU:CD1	1:I:96:ILE:HD12	2.24	0.67
1:C:-1:ILE:HD11	1:C:114:SER:HG	1.59	0.67
1:I:1:ASN:HB2	1:I:73:SER:OG	1.94	0.66
1:O:66:LEU:C	1:O:66:LEU:HD23	2.22	0.65
1:E:32:HIS:ND1	1:E:33:PRO:HD2	2.13	0.64
1:O:22:ASP:HB3	2:P:26:ARG:HH11	1.62	0.64
1:K:31:VAL:CG1	1:K:109:MET:HE1	2.27	0.64
1:C:14:ASP:O	1:C:17:GLN:HB2	1.99	0.63
1:C:-3:MET:HE2	4:J:138:HOH:O	1.98	0.62
3:A:1004:GOL:O2	2:D:-3:MET:HE3	1.99	0.62
2:P:35:ARG:HG3	2:P:61:LEU:CD1	2.31	0.61
2:J:65:ARG:HG2	2:J:65:ARG:HH11	1.66	0.61
1:G:-4:ALA:HA	1:K:104:VAL:HG21	1.82	0.61
1:O:32:HIS:ND1	1:O:33:PRO:HD2	2.16	0.61
1:C:28:GLU:CG	1:C:31:VAL:HG22	2.31	0.61
1:E:68:ILE:HG13	2:H:-1:ILE:HD11	1.84	0.60
1:G:76:SER:C	1:G:78:GLY:HA3	2.27	0.59
1:I:89:GLU:HB2	2:J:67:PRO:HG3	1.86	0.58
1:G:2:TRP:HB2	1:G:111:ILE:HG13	1.83	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:33:PRO:O	1:I:36:LYS:HG2	2.04	0.58
1:M:6:ILE:HG13	1:M:107:VAL:HG11	1.86	0.58
1:O:87:GLU:HG3	2:P:67:PRO:HB3	1.86	0.58
1:G:23:ALA:O	1:G:96:ILE:HD13	2.04	0.58
1:K:36:LYS:CB	1:K:109:MET:HE3	2.34	0.57
2:L:20:SER:HB3	4:L:136:HOH:O	2.05	0.57
1:E:31:VAL:HG11	1:E:109:MET:HE1	1.87	0.57
1:K:-1:ILE:HD11	1:K:114:SER:HB2	1.86	0.56
1:O:96:ILE:HG12	4:O:1016:HOH:O	2.05	0.56
1:E:109:MET:O	1:E:113:THR:HB	2.05	0.56
2:B:58:THR:HG21	2:D:50:ALA:HB2	1.88	0.56
1:C:31:VAL:O	1:C:109:MET:HE1	2.05	0.56
2:P:7:MET:HE3	2:P:8:SER:O	2.07	0.55
2:B:7:MET:HE2	2:B:27:TYR:OH	2.06	0.55
1:C:75:SER:O	1:C:76:SER:O	2.25	0.55
2:F:0:SER:N	1:G:64:GLU:OE2	2.40	0.54
1:A:97:LYS:O	1:A:101:GLN:HG3	2.07	0.54
2:F:20:SER:O	2:F:23:SER:HB2	2.08	0.53
1:G:78:GLY:CA	1:G:79:ASN:CB	2.87	0.53
1:K:32:HIS:CD2	1:K:34:SER:H	2.26	0.53
2:J:65:ARG:HH11	2:J:65:ARG:CG	2.20	0.53
1:O:111:ILE:O	1:O:112:ASN:CB	2.57	0.53
1:C:11:LYS:HE2	1:C:62:THR:OG1	2.08	0.53
1:E:31:VAL:CG1	1:E:109:MET:HE1	2.37	0.53
1:I:18:SER:HB2	1:K:18:SER:HB2	1.91	0.53
1:K:-1:ILE:HD11	1:K:114:SER:CB	2.38	0.53
1:K:36:LYS:HB3	1:K:109:MET:HE3	1.90	0.53
2:N:-3:MET:HG2	2:N:52:ASN:HD21	1.73	0.53
1:I:12:ILE:O	1:I:16:ILE:HG23	2.08	0.53
2:D:66:ASP:O	2:D:70:VAL:HG13	2.09	0.53
1:C:28:GLU:HG3	1:C:31:VAL:HG22	1.92	0.52
1:M:14:ASP:O	1:M:17:GLN:HB2	2.09	0.52
1:O:45:LEU:O	1:O:48:GLN:HB2	2.08	0.52
1:G:6:ILE:O	1:G:10:LYS:HG2	2.10	0.52
1:C:32:HIS:HD2	1:C:34:SER:H	1.54	0.52
2:D:65:ARG:HB3	2:D:70:VAL:HG11	1.92	0.52
1:G:16:ILE:CD1	1:G:96:ILE:HG21	2.39	0.52
1:K:31:VAL:HG11	1:K:109:MET:HE1	1.91	0.52
1:K:32:HIS:CD2	1:K:33:PRO:HD2	2.45	0.51
2:L:7:MET:HG2	4:L:140:HOH:O	2.09	0.51
1:G:89:GLU:HB2	2:H:67:PRO:HG3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:97:LYS:CD	4:K:126:HOH:O	2.58	0.51
2:F:7:MET:HG2	4:F:108:HOH:O	2.10	0.51
2:J:21:LEU:O	2:J:22:TYR:HB2	2.11	0.51
1:O:10:LYS:HG2	4:O:1019:HOH:O	2.11	0.51
2:P:35:ARG:HG3	2:P:61:LEU:HD12	1.93	0.50
1:A:9:LEU:O	1:A:13:GLU:HG3	2.11	0.50
2:N:35:ARG:HG3	2:N:61:LEU:CD1	2.42	0.50
1:M:36:LYS:HB2	1:M:109:MET:HE3	1.92	0.50
2:N:49:LYS:HG2	2:P:60:SER:HB3	1.93	0.50
1:G:101:GLN:NE2	1:K:-3:MET:HE1	2.21	0.49
1:O:97:LYS:HD2	1:O:101:GLN:NE2	2.28	0.49
1:G:16:ILE:HD13	1:G:96:ILE:HG21	1.93	0.49
1:K:32:HIS:HD2	1:K:33:PRO:HD2	1.77	0.49
1:G:28:GLU:HG3	1:G:31:VAL:HG22	1.94	0.48
2:P:7:MET:HE1	4:P:120:HOH:O	2.14	0.48
2:F:65:ARG:HD3	2:F:65:ARG:HA	1.65	0.48
1:I:47:LEU:HD21	1:I:66:LEU:HD13	1.95	0.48
1:G:-4:ALA:HA	1:K:104:VAL:CG2	2.43	0.48
1:A:1:ASN:HB2	1:A:73:SER:OG	2.14	0.48
2:H:35:ARG:HG3	2:H:61:LEU:CD1	2.43	0.48
1:C:16:ILE:CD1	1:C:96:ILE:HG21	2.43	0.48
1:M:-1:ILE:O	1:M:3:VAL:HG23	2.14	0.48
2:L:72:GLN:O	2:L:73:ARG:HB3	2.13	0.47
1:I:32:HIS:ND1	1:I:33:PRO:HD2	2.28	0.47
1:O:73:SER:HA	1:O:76:SER:HB3	1.97	0.47
1:I:18:SER:CB	1:K:18:SER:HB2	2.44	0.47
2:J:48:ASN:C	2:J:48:ASN:HD22	2.22	0.47
1:A:10:LYS:O	1:A:10:LYS:HD3	2.15	0.47
2:B:35:ARG:HG3	2:B:61:LEU:HD13	1.95	0.47
1:A:66:LEU:C	1:A:66:LEU:HD23	2.41	0.46
1:E:18:SER:HB2	1:G:18:SER:HB2	1.98	0.46
1:M:89:GLU:HB2	2:N:67:PRO:HG3	1.98	0.46
1:A:33:PRO:O	1:A:36:LYS:HG2	2.16	0.46
1:I:25:LEU:CD1	1:I:96:ILE:CD1	2.92	0.46
1:O:36:LYS:HD2	1:O:109:MET:HE2	1.98	0.46
2:B:27:TYR:CD2	2:B:59:PRO:HG3	2.51	0.46
2:J:51:THR:HB	2:J:53:VAL:HG23	1.98	0.46
2:J:65:ARG:HB3	2:J:70:VAL:HG21	1.97	0.46
1:I:105:HIS:O	1:I:108:GLN:HB2	2.16	0.45
2:L:65:ARG:HA	2:L:65:ARG:HD3	1.67	0.45
2:P:20:SER:O	2:P:23:SER:HB3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:36:LYS:HB2	1:K:109:MET:HE3	1.98	0.45
2:N:70:VAL:CG1	2:N:70:VAL:O	2.64	0.45
2:J:66:ASP:HA	2:J:67:PRO:HD2	1.60	0.45
1:O:47:LEU:HD21	1:O:66:LEU:HD13	1.99	0.45
1:A:87:GLU:HG3	2:B:67:PRO:HB3	1.99	0.45
1:C:100:LEU:HB3	1:O:-3:MET:HE1	1.99	0.44
2:N:24:ARG:NH1	2:N:44:GLU:OE2	2.51	0.44
2:F:58:THR:HG21	2:H:50:ALA:HB2	2.00	0.44
2:H:-1:ILE:HG12	2:H:0:SER:N	2.33	0.44
1:K:75:SER:O	1:K:76:SER:C	2.61	0.44
1:O:37:VAL:HG21	1:O:80:VAL:HG11	2.00	0.44
1:A:47:LEU:HD21	1:A:66:LEU:HD13	1.98	0.44
1:M:87:GLU:HB2	1:M:90:GLU:HG3	1.99	0.44
1:I:16:ILE:CD1	1:I:96:ILE:HG21	2.48	0.43
2:N:40:SER:O	2:N:61:LEU:HB2	2.17	0.43
2:P:7:MET:HG2	4:P:122:HOH:O	2.17	0.43
1:O:97:LYS:HG3	1:O:98:GLU:N	2.30	0.43
1:E:66:LEU:C	1:E:66:LEU:HD23	2.43	0.43
1:M:11:LYS:HD3	1:M:62:THR:OG1	2.19	0.43
1:C:69:LEU:HD13	2:J:5:PRO:CD	2.49	0.42
2:H:3:CYS:HB3	2:H:4:PRO:HD2	2.00	0.42
1:I:28:GLU:CG	1:I:31:VAL:HG22	2.49	0.42
2:N:21:LEU:O	2:N:22:TYR:HB2	2.20	0.42
1:I:28:GLU:HG3	1:I:31:VAL:HG22	2.02	0.42
1:K:-1:ILE:CD1	1:K:114:SER:HB2	2.49	0.42
1:A:-3:MET:HE2	1:A:-3:MET:HB2	1.82	0.42
2:B:35:ARG:HG3	2:B:61:LEU:HD12	1.99	0.42
1:K:97:LYS:HD2	4:K:126:HOH:O	2.17	0.42
1:M:36:LYS:HB2	1:M:109:MET:CE	2.50	0.42
1:C:16:ILE:HA	1:C:19:MET:HG3	2.00	0.42
1:A:52:LEU:HD21	2:D:47:LEU:HD12	2.02	0.41
1:E:6:ILE:HG13	1:E:107:VAL:HG11	2.02	0.41
1:I:109:MET:HE2	1:I:109:MET:HB3	1.83	0.41
1:K:11:LYS:HE2	1:K:62:THR:OG1	2.20	0.41
1:M:12:ILE:O	1:M:16:ILE:HG23	2.21	0.41
2:D:6:PRO:HG3	2:D:16:VAL:HG21	2.02	0.41
1:C:28:GLU:HG3	1:C:31:VAL:CG2	2.50	0.41
2:N:70:VAL:O	2:N:70:VAL:HG12	2.20	0.41
1:E:106:ILE:HG13	1:E:109:MET:HE2	2.03	0.41
1:I:33:PRO:HA	1:I:109:MET:HE1	2.02	0.41
1:C:32:HIS:CD2	1:C:34:SER:OG	2.73	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:28:GLU:HG2	1:C:31:VAL:HG22	2.02	0.41
1:I:1:ASN:HD21	1:I:72:ASN:ND2	1.99	0.41
2:J:17:LYS:HD2	4:J:129:HOH:O	2.21	0.41
2:J:66:ASP:OD2	2:J:66:ASP:C	2.64	0.41
1:G:14:ASP:O	1:G:17:GLN:HB2	2.21	0.41
1:A:18:SER:HB2	1:C:18:SER:HB2	2.02	0.40
1:G:45:LEU:HD23	1:G:45:LEU:HA	1.91	0.40
1:O:3:VAL:CG2	1:O:111:ILE:HD11	2.52	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	109/119 (92%)	108 (99%)	1 (1%)	0	100	100
1	C	112/119 (94%)	109 (97%)	3 (3%)	0	100	100
1	E	117/119 (98%)	113 (97%)	4 (3%)	0	100	100
1	G	117/119 (98%)	110 (94%)	6 (5%)	1 (1%)	14	9
1	I	112/119 (94%)	108 (96%)	4 (4%)	0	100	100
1	K	112/119 (94%)	109 (97%)	3 (3%)	0	100	100
1	M	112/119 (94%)	109 (97%)	2 (2%)	1 (1%)	14	9
1	O	116/119 (98%)	114 (98%)	2 (2%)	0	100	100
2	B	72/107 (67%)	66 (92%)	6 (8%)	0	100	100
2	D	74/107 (69%)	73 (99%)	1 (1%)	0	100	100
2	F	71/107 (66%)	69 (97%)	2 (3%)	0	100	100
2	H	75/107 (70%)	75 (100%)	0	0	100	100
2	J	76/107 (71%)	74 (97%)	2 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	L	73/107 (68%)	70 (96%)	3 (4%)	0	100	100
2	N	74/107 (69%)	72 (97%)	2 (3%)	0	100	100
2	P	70/107 (65%)	69 (99%)	1 (1%)	0	100	100
All	All	1492/1808 (82%)	1448 (97%)	42 (3%)	2 (0%)	48	46

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	79	ASN
1	M	17	GLN

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	105/110 (96%)	94 (90%)	11 (10%)	6	4
1	C	108/110 (98%)	97 (90%)	11 (10%)	7	4
1	E	110/110 (100%)	104 (94%)	6 (6%)	19	17
1	G	110/110 (100%)	105 (96%)	5 (4%)	24	23
1	I	108/110 (98%)	102 (94%)	6 (6%)	19	16
1	K	108/110 (98%)	104 (96%)	4 (4%)	30	30
1	M	108/110 (98%)	103 (95%)	5 (5%)	24	22
1	O	109/110 (99%)	104 (95%)	5 (5%)	24	22
2	B	67/92 (73%)	62 (92%)	5 (8%)	12	9
2	D	68/92 (74%)	60 (88%)	8 (12%)	5	3
2	F	66/92 (72%)	60 (91%)	6 (9%)	9	6
2	H	68/92 (74%)	62 (91%)	6 (9%)	9	6
2	J	69/92 (75%)	60 (87%)	9 (13%)	4	2
2	L	68/92 (74%)	59 (87%)	9 (13%)	4	2
2	N	68/92 (74%)	58 (85%)	10 (15%)	3	2

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	P	65/92 (71%)	59 (91%)	6 (9%)	8 5
All	All	1405/1616 (87%)	1293 (92%)	112 (8%)	11 8

All (112) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-3	MET
1	A	10	LYS
1	A	18	SER
1	A	20	HIS
1	A	21	ILE
1	A	52	LEU
1	A	69	LEU
1	A	81	THR
1	A	92	GLU
1	A	96	ILE
1	A	111	ILE
2	B	1	ILE
2	B	7	MET
2	B	20	SER
2	B	47	LEU
2	B	70	VAL
1	C	-3	MET
1	C	-1	ILE
1	C	17	GLN
1	C	19	MET
1	C	20	HIS
1	C	69	LEU
1	C	80	VAL
1	C	81	THR
1	C	86	LYS
1	C	97	LYS
1	C	109	MET
2	D	10	GLU
2	D	13	ASP
2	D	47	LEU
2	D	49	LYS
2	D	52	ASN
2	D	53	VAL
2	D	62	LYS
2	D	71	HIS
1	E	0	SER

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Mol	Chain	Res	Type
1	E	10	LYS
1	E	17	GLN
1	E	52	LEU
1	E	69	LEU
1	E	113	THR
2	F	0	SER
2	F	23	SER
2	F	47	LEU
2	F	49	LYS
2	F	51	THR
2	F	65	ARG
1	G	17	GLN
1	G	20	HIS
1	G	79	ASN
1	G	83	SER
1	G	96	ILE
2	H	-1	ILE
2	H	10	GLU
2	H	47	LEU
2	H	52	ASN
2	H	70	VAL
2	H	72	GLN
1	I	14	ASP
1	I	58	SER
1	I	69	LEU
1	I	80	VAL
1	I	81	THR
1	I	113	THR
2	J	-1	ILE
2	J	8	SER
2	J	47	LEU
2	J	49	LYS
2	J	52	ASN
2	J	53	VAL
2	J	65	ARG
2	J	70	VAL
2	J	73	ARG
1	K	17	GLN
1	K	52	LEU
1	K	69	LEU
1	K	87	GLU
2	L	-1	ILE

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Mol	Chain	Res	Type
2	L	0	SER
2	L	1	ILE
2	L	7	MET
2	L	24	ARG
2	L	49	LYS
2	L	51	THR
2	L	72	GLN
2	L	73	ARG
1	M	31	VAL
1	M	58	SER
1	M	69	LEU
1	M	75	SER
1	M	80	VAL
2	N	-3	MET
2	N	-1	ILE
2	N	13	ASP
2	N	24	ARG
2	N	47	LEU
2	N	49	LYS
2	N	52	ASN
2	N	70	VAL
2	N	71	HIS
2	N	72	GLN
1	O	10	LYS
1	O	17	GLN
1	O	52	LEU
1	O	69	LEU
1	O	97	LYS
2	P	0	SER
2	P	7	MET
2	P	23	SER
2	P	47	LEU
2	P	51	THR
2	P	70	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	HIS
1	A	101	GLN
1	A	105	HIS
1	A	108	GLN

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Mol	Chain	Res	Type
2	B	30	ASN
2	B	71	HIS
1	C	32	HIS
1	C	48	GLN
1	C	65	ASN
2	D	11	HIS
2	D	30	ASN
1	E	20	HIS
1	E	48	GLN
1	E	108	GLN
2	F	30	ASN
1	G	20	HIS
1	G	48	GLN
1	G	101	GLN
1	G	108	GLN
2	H	30	ASN
2	H	55	HIS
1	I	48	GLN
1	I	65	ASN
1	I	72	ASN
1	I	101	GLN
2	J	30	ASN
2	J	48	ASN
2	J	55	HIS
1	K	20	HIS
1	K	32	HIS
1	K	112	ASN
2	L	72	GLN
1	M	48	GLN
1	M	108	GLN
2	N	52	ASN
2	N	55	HIS
1	O	1	ASN
1	O	17	GLN
1	O	20	HIS
1	O	79	ASN
1	O	101	GLN
1	O	108	GLN
2	P	30	ASN
2	P	55	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 ⓘ

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	$\# Z > 2$
3	GOL	O	1001	-	5,5,5	0.54	0	5,5,5	1.02	0
3	GOL	O	1002	-	5,5,5	0.39	0	5,5,5	0.31	0
3	GOL	A	1003	-	5,5,5	0.38	0	5,5,5	0.41	0
3	GOL	A	1004	-	5,5,5	0.42	0	5,5,5	0.54	0

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
3	GOL	O	1001	-	-	0/4/4/4	-
3	GOL	O	1002	-	-	3/4/4/4	-
3	GOL	A	1003	-	-	2/4/4/4	-
3	GOL	A	1004	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1004	GOL	C1-C2-C3-O3
3	A	1003	GOL	O1-C1-C2-C3
3	A	1004	GOL	O1-C1-C2-C3
3	O	1002	GOL	C1-C2-C3-O3
3	A	1003	GOL	O1-C1-C2-O2
3	A	1004	GOL	O1-C1-C2-O2
3	A	1004	GOL	O2-C2-C3-O3
3	O	1002	GOL	O2-C2-C3-O3
3	O	1002	GOL	O1-C1-C2-C3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1004	GOL	1	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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.