



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:03 PM UTC

PDB ID : 4RSU / pdb_00004rsu
Title : Crystal structure of the light and hvem complex
Authors : Liu, W.; Ramagoal, U.A.; Himmel, D.; Bonanno, J.B.; Nathenson, S.G.; Almo, S.C.; Atoms-to-Animals: The Immune Function Network (IFN); New York Structural Genomics Research Consortium (NYSGRG)
Deposited on : 2014-11-11
Resolution : 2.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

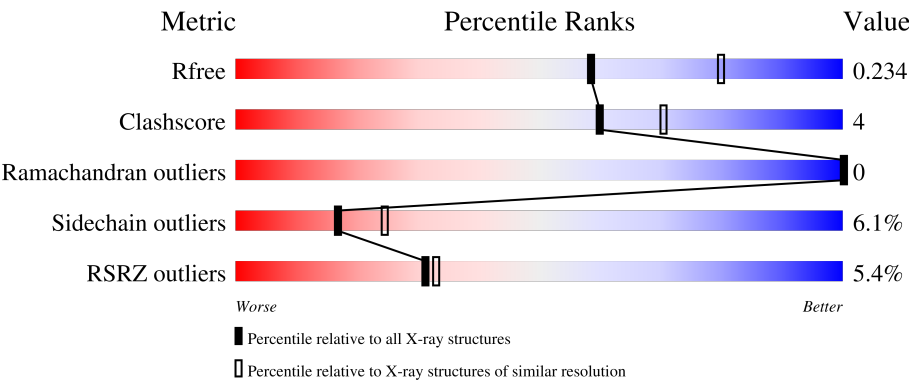
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	165	% 79% 10% • 10%
1	B	165	4% 76% 9% • 14%
1	C	165	4% 70% 13% • 16%
1	G	165	% 81% 8% • 9%
1	H	165	4% 77% 13% • 9%

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Mol	Chain	Length	Quality of chain
1	I	165	7% 76% 8% 5% 11%
2	D	134	3% 71% 7% 22%
2	E	134	7% 60% 14% 25%
2	F	134	7% 64% 7% 27%
2	J	134	3% 65% 12% 21%
2	K	134	8% 61% 17% 21%
2	L	134	8% 59% 18% 22%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11553 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Tumor necrosis factor ligand superfamily member 14, soluble form.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	149	Total	C	N	O	S	0	0	0
			1140	727	199	211	3			
1	B	142	Total	C	N	O	S	0	0	0
			1093	697	189	204	3			
1	C	138	Total	C	N	O	S	0	0	0
			1069	684	185	197	3			
1	G	150	Total	C	N	O	S	0	0	0
			1149	732	200	214	3			
1	H	150	Total	C	N	O	S	0	0	0
			1149	732	200	214	3			
1	I	147	Total	C	N	O	S	0	0	0
			1125	716	197	209	3			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	76	HIS	-	expression tag	UNP O43557
A	77	HIS	-	expression tag	UNP O43557
A	78	HIS	-	expression tag	UNP O43557
A	79	HIS	-	expression tag	UNP O43557
A	80	HIS	-	expression tag	UNP O43557
A	81	HIS	-	expression tag	UNP O43557
A	82	GLY	-	expression tag	UNP O43557
A	214	GLU	LYS	conflict	UNP O43557
B	76	HIS	-	expression tag	UNP O43557
B	77	HIS	-	expression tag	UNP O43557
B	78	HIS	-	expression tag	UNP O43557
B	79	HIS	-	expression tag	UNP O43557
B	80	HIS	-	expression tag	UNP O43557
B	81	HIS	-	expression tag	UNP O43557
B	82	GLY	-	expression tag	UNP O43557
B	214	GLU	LYS	conflict	UNP O43557

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Chain	Residue	Modelled	Actual	Comment	Reference
C	76	HIS	-	expression tag	UNP O43557
C	77	HIS	-	expression tag	UNP O43557
C	78	HIS	-	expression tag	UNP O43557
C	79	HIS	-	expression tag	UNP O43557
C	80	HIS	-	expression tag	UNP O43557
C	81	HIS	-	expression tag	UNP O43557
C	82	GLY	-	expression tag	UNP O43557
C	214	GLU	LYS	conflict	UNP O43557
G	76	HIS	-	expression tag	UNP O43557
G	77	HIS	-	expression tag	UNP O43557
G	78	HIS	-	expression tag	UNP O43557
G	79	HIS	-	expression tag	UNP O43557
G	80	HIS	-	expression tag	UNP O43557
G	81	HIS	-	expression tag	UNP O43557
G	82	GLY	-	expression tag	UNP O43557
G	214	GLU	LYS	conflict	UNP O43557
H	76	HIS	-	expression tag	UNP O43557
H	77	HIS	-	expression tag	UNP O43557
H	78	HIS	-	expression tag	UNP O43557
H	79	HIS	-	expression tag	UNP O43557
H	80	HIS	-	expression tag	UNP O43557
H	81	HIS	-	expression tag	UNP O43557
H	82	GLY	-	expression tag	UNP O43557
H	214	GLU	LYS	conflict	UNP O43557
I	76	HIS	-	expression tag	UNP O43557
I	77	HIS	-	expression tag	UNP O43557
I	78	HIS	-	expression tag	UNP O43557
I	79	HIS	-	expression tag	UNP O43557
I	80	HIS	-	expression tag	UNP O43557
I	81	HIS	-	expression tag	UNP O43557
I	82	GLY	-	expression tag	UNP O43557
I	214	GLU	LYS	conflict	UNP O43557

- Molecule 2 is a protein called Tumor necrosis factor receptor superfamily member 14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	105	Total	C	N	O	S	0	0	0
			761	458	136	149	18			
2	E	101	Total	C	N	O	S	0	0	0
			732	443	131	140	18			
2	F	98	Total	C	N	O	S	0	0	0
			705	425	126	136	18			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	J	106	Total	C	N	O	S	0	0	0
			766	461	137	150	18			
2	K	106	Total	C	N	O	S	0	0	0
			768	463	137	150	18			
2	L	104	Total	C	N	O	S	0	0	0
			741	445	133	145	18			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	37	ARG	-	expression tag	UNP Q92956
D	38	SER	-	expression tag	UNP Q92956
D	163	THR	-	expression tag	UNP Q92956
D	164	GLY	-	expression tag	UNP Q92956
D	165	HIS	-	expression tag	UNP Q92956
D	166	HIS	-	expression tag	UNP Q92956
D	167	HIS	-	expression tag	UNP Q92956
D	168	HIS	-	expression tag	UNP Q92956
D	169	HIS	-	expression tag	UNP Q92956
D	170	HIS	-	expression tag	UNP Q92956
E	37	ARG	-	expression tag	UNP Q92956
E	38	SER	-	expression tag	UNP Q92956
E	163	THR	-	expression tag	UNP Q92956
E	164	GLY	-	expression tag	UNP Q92956
E	165	HIS	-	expression tag	UNP Q92956
E	166	HIS	-	expression tag	UNP Q92956
E	167	HIS	-	expression tag	UNP Q92956
E	168	HIS	-	expression tag	UNP Q92956
E	169	HIS	-	expression tag	UNP Q92956
E	170	HIS	-	expression tag	UNP Q92956
F	37	ARG	-	expression tag	UNP Q92956
F	38	SER	-	expression tag	UNP Q92956
F	163	THR	-	expression tag	UNP Q92956
F	164	GLY	-	expression tag	UNP Q92956
F	165	HIS	-	expression tag	UNP Q92956
F	166	HIS	-	expression tag	UNP Q92956
F	167	HIS	-	expression tag	UNP Q92956
F	168	HIS	-	expression tag	UNP Q92956
F	169	HIS	-	expression tag	UNP Q92956
F	170	HIS	-	expression tag	UNP Q92956
J	37	ARG	-	expression tag	UNP Q92956
J	38	SER	-	expression tag	UNP Q92956

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Chain	Residue	Modelled	Actual	Comment	Reference
J	163	THR	-	expression tag	UNP Q92956
J	164	GLY	-	expression tag	UNP Q92956
J	165	HIS	-	expression tag	UNP Q92956
J	166	HIS	-	expression tag	UNP Q92956
J	167	HIS	-	expression tag	UNP Q92956
J	168	HIS	-	expression tag	UNP Q92956
J	169	HIS	-	expression tag	UNP Q92956
J	170	HIS	-	expression tag	UNP Q92956
K	37	ARG	-	expression tag	UNP Q92956
K	38	SER	-	expression tag	UNP Q92956
K	163	THR	-	expression tag	UNP Q92956
K	164	GLY	-	expression tag	UNP Q92956
K	165	HIS	-	expression tag	UNP Q92956
K	166	HIS	-	expression tag	UNP Q92956
K	167	HIS	-	expression tag	UNP Q92956
K	168	HIS	-	expression tag	UNP Q92956
K	169	HIS	-	expression tag	UNP Q92956
K	170	HIS	-	expression tag	UNP Q92956
L	37	ARG	-	expression tag	UNP Q92956
L	38	SER	-	expression tag	UNP Q92956
L	163	THR	-	expression tag	UNP Q92956
L	164	GLY	-	expression tag	UNP Q92956
L	165	HIS	-	expression tag	UNP Q92956
L	166	HIS	-	expression tag	UNP Q92956
L	167	HIS	-	expression tag	UNP Q92956
L	168	HIS	-	expression tag	UNP Q92956
L	169	HIS	-	expression tag	UNP Q92956
L	170	HIS	-	expression tag	UNP Q92956

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

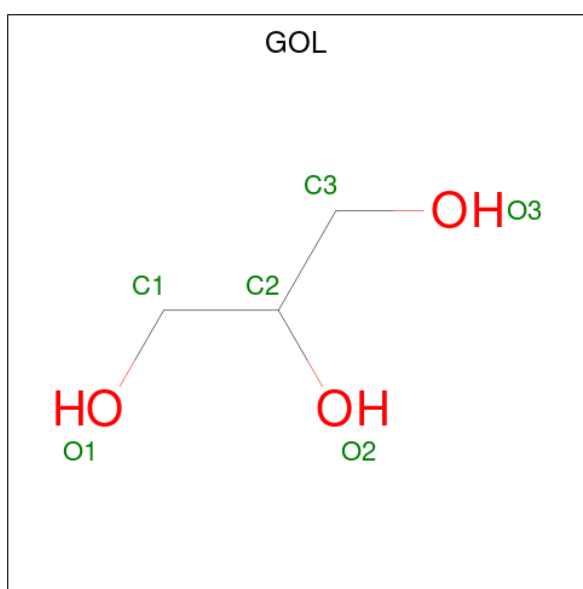
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Cl 2 2	0	0
3	B	2	Total Cl 2 2	0	0
3	C	2	Total Cl 2 2	0	0
3	E	1	Total Cl 1 1	0	0
3	G	3	Total Cl 3 3	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	1	Total	Cl	0	0
			1	1		
3	I	1	Total	Cl	0	0
			1	1		
3	K	1	Total	Cl	0	0
			1	1		
3	L	1	Total	Cl	0	0
			1	1		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



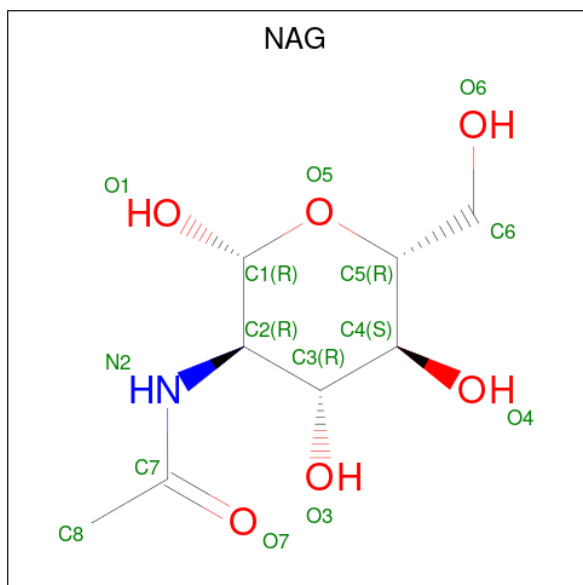
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		
4	F	1	Total	C	O	0	0
			6	3	3		
4	G	1	Total	C	O	0	0
			6	3	3		
4	H	1	Total	C	O	0	0
			6	3	3		
4	I	1	Total	C	O	0	0
			6	3	3		

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

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	D	1	Total	C	N	O	0	0
			14	8	1	5		
5	E	1	Total	C	N	O	0	0
			14	8	1	5		
5	J	1	Total	C	N	O	0	0
			14	8	1	5		
5	K	1	Total	C	N	O	0	0
			14	8	1	5		
5	L	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	39	Total	O	0	0
			39	39		
6	B	16	Total	O	0	0
			16	16		
6	C	12	Total	O	0	0
			12	12		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	17	Total 17	O 17	0	0
6	E	4	Total 4	O 4	0	0
6	F	16	Total 16	O 16	0	0
6	G	47	Total 47	O 47	0	0
6	H	19	Total 19	O 19	0	0
6	I	21	Total 21	O 21	0	0
6	J	14	Total 14	O 14	0	0
6	K	4	Total 4	O 4	0	0
6	L	14	Total 14	O 14	0	0

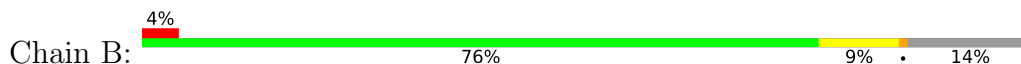
3 Residue-property plots [i](#)

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.

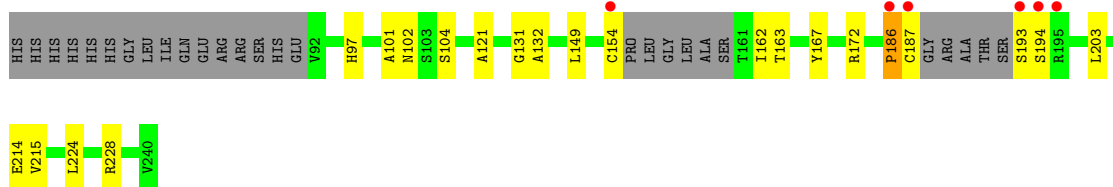
- Molecule 1: Tumor necrosis factor ligand superfamily member 14, soluble form



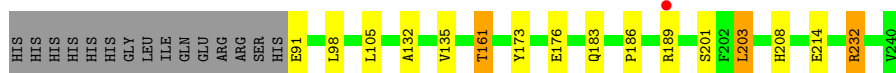
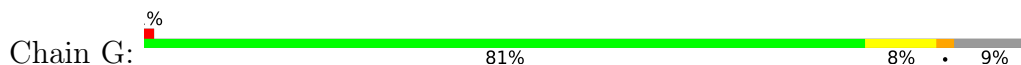
- Molecule 1: Tumor necrosis factor ligand superfamily member 14, soluble form



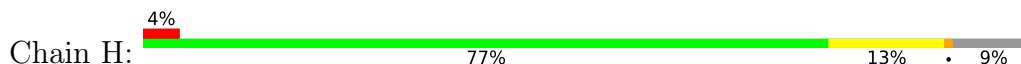
- Molecule 1: Tumor necrosis factor ligand superfamily member 14, soluble form

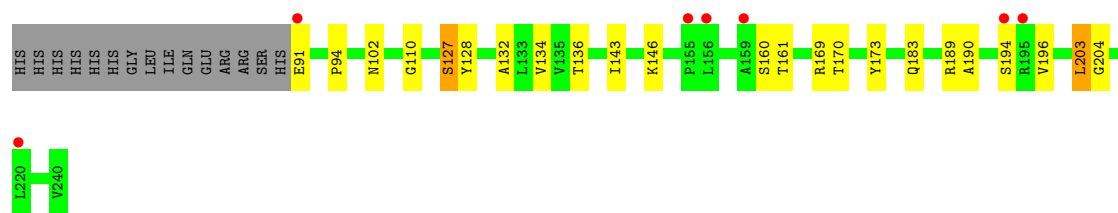


- Molecule 1: Tumor necrosis factor ligand superfamily member 14, soluble form

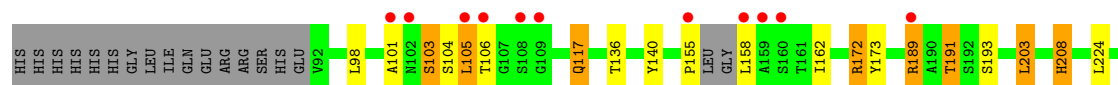
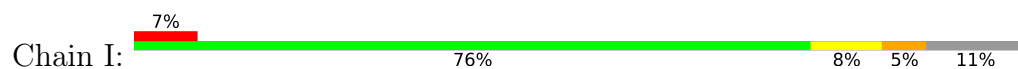


- Molecule 1: Tumor necrosis factor ligand superfamily member 14, soluble form

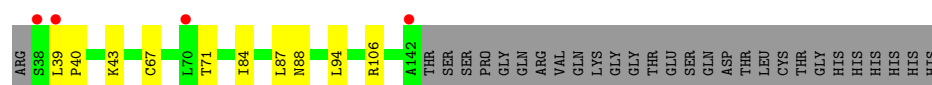




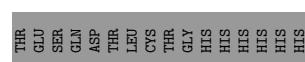
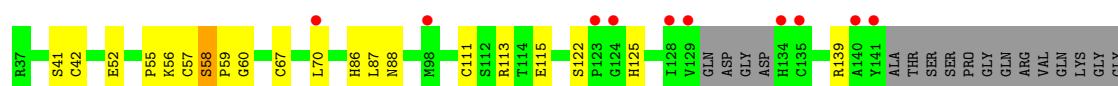
- Molecule 1: Tumor necrosis factor ligand superfamily member 14, soluble form



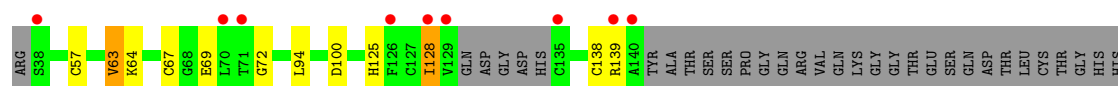
- Molecule 2: Tumor necrosis factor receptor superfamily member 14



- Molecule 2: Tumor necrosis factor receptor superfamily member 14

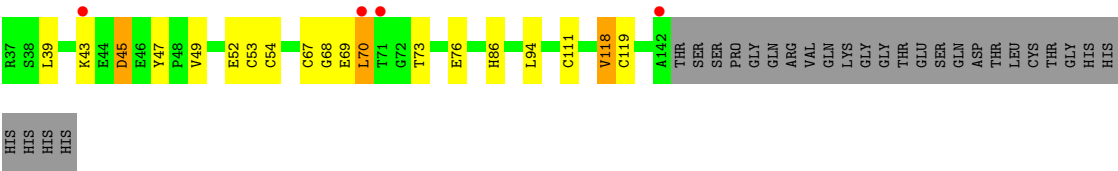


- Molecule 2: Tumor necrosis factor receptor superfamily member 14



- Molecule 2: Tumor necrosis factor receptor superfamily member 14

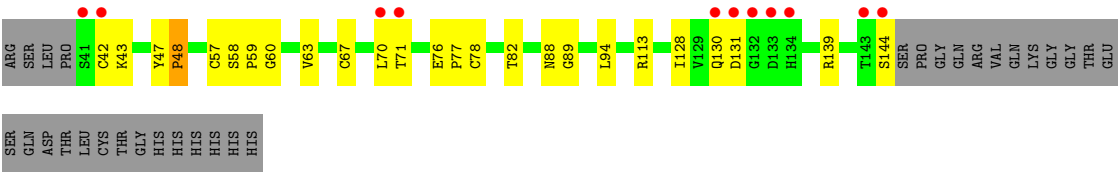




● Molecule 2: Tumor necrosis factor receptor superfamily member 14



● Molecule 2: Tumor necrosis factor receptor superfamily member 14



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	111.65Å 113.60Å 163.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.92 – 2.30 48.92 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.92-2.30) 99.8 (48.92-2.30)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.64 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.188 , 0.231 0.194 , 0.234	Depositor DCC
R_{free} test set	4631 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	35.0	Xtriage
Anisotropy	0.022	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 26.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.012 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11553	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.91% 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: GOL, NAG, CL

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	1.39	7/1167 (0.6%)	1.11	2/1585 (0.1%)
1	B	1.28	0/1117	1.09	0/1514
1	C	1.32	1/1093 (0.1%)	1.13	1/1481 (0.1%)
1	G	1.47	4/1176 (0.3%)	1.18	3/1597 (0.2%)
1	H	1.28	1/1176 (0.1%)	1.16	7/1597 (0.4%)
1	I	1.36	3/1151 (0.3%)	1.15	2/1562 (0.1%)
2	D	1.32	1/777 (0.1%)	1.12	1/1052 (0.1%)
2	E	1.14	1/747 (0.1%)	1.14	5/1010 (0.5%)
2	F	1.28	2/718 (0.3%)	1.17	1/970 (0.1%)
2	J	1.38	4/782 (0.5%)	1.30	6/1059 (0.6%)
2	K	1.24	1/784 (0.1%)	1.29	8/1061 (0.8%)
2	L	1.36	3/755 (0.4%)	1.15	0/1022
All	All	1.33	28/11443 (0.2%)	1.16	36/15510 (0.2%)

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
1	C	0	1
1	I	0	1
All	All	0	3

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	180	LEU	C-O	-7.42	1.15	1.23
1	I	117	GLN	C-O	-6.77	1.15	1.24
1	G	135	VAL	CA-C	6.74	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	88	ASN	CA-C	-6.28	1.44	1.52
2	J	86	HIS	CA-C	-6.18	1.44	1.52
2	J	119	CYS	C-O	-6.03	1.16	1.23
1	I	203	LEU	C-O	-5.91	1.16	1.23
1	H	94	PRO	CA-C	5.84	1.59	1.52
2	K	42	CYS	N-CA	5.69	1.53	1.46
1	A	128	TYR	C-O	-5.66	1.17	1.23
1	C	172	ARG	C-O	-5.62	1.16	1.24
1	I	208	HIS	C-O	-5.61	1.17	1.24
2	L	42	CYS	N-CA	5.59	1.52	1.46
1	A	152	VAL	C-O	-5.58	1.17	1.24
1	G	176	GLU	C-O	-5.46	1.17	1.23
1	A	219	VAL	C-O	-5.45	1.18	1.24
1	A	218	ARG	C-O	-5.37	1.17	1.23
1	A	105	LEU	C-O	-5.37	1.16	1.23
1	G	201	SER	C-O	-5.36	1.18	1.24
1	A	222	GLU	N-CA	-5.35	1.39	1.46
2	L	48	PRO	C-O	-5.35	1.18	1.23
2	F	100	ASP	N-CA	-5.29	1.39	1.46
2	L	88	ASN	N-CA	-5.24	1.39	1.46
2	J	111	CYS	C-O	-5.22	1.17	1.23
2	D	94	LEU	CA-C	-5.12	1.46	1.52
2	F	63	VAL	C-O	-5.09	1.19	1.24
1	G	161	THR	C-O	-5.04	1.17	1.23
2	J	53	CYS	CA-C	5.01	1.59	1.52

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	110	GLY	CA-C-N	-7.59	112.86	120.98
1	H	110	GLY	C-N-CA	-7.59	112.86	120.98
2	E	42	CYS	N-CA-C	-7.35	100.81	110.43
2	K	76	GLU	CA-C-N	-6.96	112.82	119.85
2	K	76	GLU	C-N-CA	-6.96	112.82	119.85
1	H	173	TYR	CA-C-N	-6.92	112.52	120.89
1	H	173	TYR	C-N-CA	-6.92	112.52	120.89
2	J	68	GLY	N-CA-C	-6.88	104.53	112.79
2	K	39	LEU	CA-C-N	-6.67	113.81	120.21
2	K	39	LEU	C-N-CA	-6.67	113.81	120.21
2	K	54	CYS	CA-C-N	-6.53	113.23	119.76
2	K	54	CYS	C-N-CA	-6.53	113.23	119.76
1	I	173	TYR	CA-C-N	-6.32	113.44	120.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	173	TYR	C-N-CA	-6.32	113.44	120.45
2	J	76	GLU	CA-C-N	-6.28	114.15	120.31
2	J	76	GLU	C-N-CA	-6.28	114.15	120.31
2	J	39	LEU	CA-C-N	-6.25	113.53	119.85
2	J	39	LEU	C-N-CA	-6.25	113.53	119.85
2	J	67	CYS	CB-CA-C	-5.82	99.34	109.64
2	F	67	CYS	N-CA-CB	5.82	118.58	110.26
1	A	232	ARG	NE-CZ-NH1	-5.81	115.69	121.50
1	A	195	ARG	N-CA-C	5.61	116.78	108.60
2	D	87	LEU	N-CA-C	-5.56	102.40	110.24
1	G	173	TYR	CA-C-N	-5.54	114.06	120.04
1	G	173	TYR	C-N-CA	-5.54	114.06	120.04
1	G	232	ARG	NE-CZ-NH1	-5.32	116.18	121.50
1	H	203	LEU	CB-CA-C	-5.32	98.37	110.07
1	C	131	GLY	N-CA-C	-5.21	108.49	114.69
2	E	55	PRO	CB-CA-C	-5.10	104.89	111.46
2	K	78	CYS	CA-C-N	-5.08	115.15	120.38
2	K	78	CYS	C-N-CA	-5.08	115.15	120.38
1	H	136	THR	CB-CA-C	-5.05	101.80	110.24
2	E	58	SER	CA-C-N	-5.03	114.60	119.78
2	E	58	SER	C-N-CA	-5.03	114.60	119.78
2	E	57	CYS	N-CA-CB	-5.03	102.72	110.56
1	H	127	SER	CB-CA-C	-5.00	101.12	111.22

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	191	THR	Peptide
1	C	186	PRO	Peptide
1	I	103	SER	Peptide

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	1140	0	1129	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1093	0	1071	8	0
1	C	1069	0	1052	11	0
1	G	1149	0	1135	6	0
1	H	1149	0	1135	13	0
1	I	1125	0	1105	22	0
2	D	761	0	702	2	0
2	E	732	0	679	6	0
2	F	705	0	662	9	0
2	J	766	0	704	6	0
2	K	768	0	708	6	0
2	L	741	0	680	9	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	E	1	0	0	0	0
3	G	3	0	0	0	0
3	H	1	0	0	0	0
3	I	1	0	0	0	0
3	K	1	0	0	0	0
3	L	1	0	0	0	0
4	B	6	0	8	0	0
4	C	6	0	8	0	0
4	D	6	0	8	0	0
4	F	6	0	8	2	0
4	G	6	0	8	0	0
4	H	6	0	8	0	0
4	I	12	0	16	0	0
5	D	14	0	13	0	0
5	E	14	0	13	0	0
5	J	14	0	13	0	0
5	K	14	0	13	0	0
5	L	14	0	13	0	0
6	A	39	0	0	0	0
6	B	16	0	0	0	0
6	C	12	0	0	0	0
6	D	17	0	0	0	0
6	E	4	0	0	1	0
6	F	16	0	0	3	0
6	G	47	0	0	2	0
6	H	19	0	0	0	0
6	I	21	0	0	0	0
6	J	14	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	K	4	0	0	0	0
6	L	14	0	0	0	0
All	All	11553	0	10891	96	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 4.

All (96) 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:I:189:ARG:HH21	1:I:189:ARG:HB3	1.29	0.97
1:I:101:ALA:N	1:I:104:SER:OG	2.04	0.91
1:I:105:LEU:HD13	1:I:227:LEU:HD11	1.58	0.85
1:H:189:ARG:HD3	1:I:189:ARG:HG3	1.60	0.83
1:I:117:GLN:O	2:K:92:LYS:NZ	2.12	0.82
2:K:59:PRO:HD3	2:K:87:LEU:HD23	1.62	0.81
1:C:149:LEU:HD13	1:C:162:ILE:HD13	1.69	0.74
2:F:128:ILE:HD12	2:F:139:ARG:CB	2.19	0.72
2:F:128:ILE:HD12	2:F:139:ARG:HB3	1.72	0.69
2:F:128:ILE:CD1	2:F:139:ARG:HB3	2.24	0.68
2:L:128:ILE:HD13	2:L:139:ARG:HG3	1.76	0.67
1:H:161:THR:HG21	1:H:183:GLN:HE21	1.59	0.67
1:H:161:THR:HG21	1:H:183:GLN:NE2	2.11	0.65
2:F:72:GLY:O	6:F:314:HOH:O	2.14	0.65
1:H:190:ALA:O	1:I:191:THR:HG21	1.96	0.65
1:I:105:LEU:CD1	1:I:227:LEU:HD11	2.27	0.65
2:F:128:ILE:CD1	2:F:139:ARG:CB	2.75	0.64
1:B:186:PRO:HB2	1:B:187:CYS:SG	2.38	0.63
1:I:189:ARG:HH21	1:I:189:ARG:CB	2.08	0.62
1:G:208:HIS:HD2	6:G:445:HOH:O	1.81	0.61
1:B:162:ILE:HD11	1:B:186:PRO:HB3	1.82	0.61
2:J:47:TYR:OH	6:J:313:HOH:O	2.15	0.60
1:I:189:ARG:HB3	1:I:189:ARG:NH2	2.08	0.60
2:E:60:GLY:HA2	2:E:113:ARG:O	2.03	0.58
4:F:201:GOL:H12	6:F:303:HOH:O	2.03	0.57
1:B:161:THR:HG21	1:B:183:GLN:HE21	1.67	0.57
2:F:125:HIS:HB3	2:F:138:CYS:HB3	1.87	0.57
1:H:161:THR:CG2	1:H:183:GLN:HE21	2.17	0.57
2:J:43:LYS:HE2	2:J:45:ASP:OD1	2.06	0.55
2:L:76:GLU:HG3	2:L:77:PRO:HD2	1.89	0.55
1:B:192:SER:HB3	1:B:195:ARG:HB2	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:155:PRO:O	1:I:158:LEU:N	2.41	0.54
1:H:127:SER:HB2	1:H:134:VAL:HB	1.90	0.54
1:G:132:ALA:HB1	1:G:214:GLU:HB3	1.91	0.53
1:I:101:ALA:O	1:I:104:SER:N	2.34	0.53
1:I:104:SER:C	1:I:105:LEU:HD12	2.34	0.52
1:G:186:PRO:O	1:I:189:ARG:NH1	2.42	0.52
2:E:125:HIS:HA	2:E:139:ARG:O	2.10	0.52
1:A:108:SER:HB2	2:J:118:VAL:HG11	1.92	0.52
1:C:228:ARG:HH11	1:C:228:ARG:HG3	1.75	0.52
1:H:91:GLU:HA	1:H:91:GLU:OE2	2.10	0.51
2:F:128:ILE:CD1	2:F:139:ARG:HB2	2.39	0.51
1:G:161:THR:HG23	1:G:183:GLN:HG2	1.92	0.51
2:D:39:LEU:HG	2:D:40:PRO:HD2	1.92	0.51
1:I:101:ALA:O	1:I:104:SER:HB2	2.10	0.51
2:K:133:ASP:HB3	2:K:134:HIS:CD2	2.46	0.51
2:L:57:CYS:SG	2:L:63:VAL:HG22	2.50	0.51
1:A:127:SER:HB2	1:A:134:VAL:HB	1.93	0.51
2:L:58:SER:HB3	2:L:59:PRO:HD2	1.95	0.49
2:J:70:LEU:H	2:J:70:LEU:HD22	1.78	0.48
2:F:57:CYS:SG	2:F:63:VAL:HG22	2.52	0.48
1:I:103:SER:O	1:I:105:LEU:N	2.46	0.48
1:C:132:ALA:HB1	1:C:214:GLU:HB3	1.96	0.47
1:C:228:ARG:HG3	1:C:228:ARG:NH1	2.28	0.47
1:B:161:THR:HG21	1:B:183:GLN:NE2	2.28	0.47
1:H:189:ARG:HD3	1:I:189:ARG:CG	2.38	0.47
1:I:189:ARG:CB	1:I:189:ARG:NH2	2.73	0.47
1:H:196:VAL:O	1:H:196:VAL:HG13	2.14	0.46
1:C:101:ALA:N	1:C:104:SER:OG	2.39	0.45
1:I:101:ALA:H	1:I:104:SER:CB	2.24	0.45
1:I:172:ARG:HG2	2:L:89:GLY:O	2.17	0.45
2:J:49:VAL:HG21	2:J:54:CYS:SG	2.56	0.45
1:G:203:LEU:HD23	1:H:146:LYS:HE2	1.99	0.45
1:A:173:TYR:CD1	1:A:174:PRO:HD2	2.52	0.44
1:B:143:ILE:HG22	1:B:237:ALA:HB2	1.99	0.44
2:F:128:ILE:HD12	2:F:139:ARG:HB2	1.96	0.44
1:C:162:ILE:CD1	1:C:186:PRO:HB3	2.47	0.44
1:H:128:TYR:HA	1:H:132:ALA:O	2.18	0.44
1:C:167:TYR:O	1:C:215:VAL:HA	2.18	0.43
2:L:94:LEU:HD12	2:L:94:LEU:N	2.33	0.43
1:C:162:ILE:HD11	1:C:224:LEU:HD13	2.00	0.43
1:C:194:SER:O	2:E:139:ARG:NH1	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:161:THR:HG23	1:G:183:GLN:CG	2.48	0.43
2:K:133:ASP:CB	2:K:134:HIS:CD2	3.01	0.43
1:B:168:LYS:HB2	1:B:179:LEU:HD11	2.01	0.43
2:D:84:ILE:HG21	2:D:88:ASN:ND2	2.33	0.43
2:L:78:CYS:HB3	2:L:82:THR:OG1	2.19	0.43
1:C:186:PRO:HG2	1:C:187:CYS:N	2.34	0.43
1:H:170:THR:O	1:H:170:THR:HG23	2.19	0.43
1:A:189:ARG:HG3	6:G:407:HOH:O	2.18	0.42
1:I:140:TYR:CE1	1:I:208:HIS:HB2	2.54	0.42
1:H:143:ILE:O	1:H:204:GLY:HA2	2.20	0.42
1:I:162:ILE:HD11	1:I:224:LEU:HD13	2.00	0.42
2:K:60:GLY:HA2	2:K:113:ARG:O	2.20	0.42
1:B:164:HIS:CE1	1:B:203:LEU:HD12	2.54	0.42
2:E:56:LYS:NZ	6:E:404:HOH:O	2.52	0.42
1:I:101:ALA:HB3	1:I:104:SER:OG	2.19	0.41
2:L:47:TYR:HA	2:L:48:PRO:HD3	1.81	0.41
2:J:70:LEU:HD22	2:J:70:LEU:N	2.35	0.41
2:K:104:GLY:O	2:K:138:CYS:HB2	2.19	0.41
4:F:201:GOL:H31	6:F:302:HOH:O	2.20	0.41
1:A:128:TYR:HA	1:A:132:ALA:O	2.21	0.41
2:E:86:HIS:O	2:E:87:LEU:C	2.64	0.41
2:L:60:GLY:HA2	2:L:113:ARG:O	2.21	0.40
1:C:97:HIS:O	1:C:121:ALA:HA	2.22	0.40
2:E:58:SER:O	2:E:59:PRO:C	2.63	0.40

There are no symmetry-related clashes.

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	147/165 (89%)	144 (98%)	3 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	136/165 (82%)	132 (97%)	4 (3%)	0	100	100
1	C	132/165 (80%)	126 (96%)	6 (4%)	0	100	100
1	G	148/165 (90%)	145 (98%)	3 (2%)	0	100	100
1	H	148/165 (90%)	143 (97%)	5 (3%)	0	100	100
1	I	143/165 (87%)	139 (97%)	4 (3%)	0	100	100
2	D	103/134 (77%)	102 (99%)	1 (1%)	0	100	100
2	E	97/134 (72%)	96 (99%)	1 (1%)	0	100	100
2	F	94/134 (70%)	93 (99%)	1 (1%)	0	100	100
2	J	104/134 (78%)	102 (98%)	2 (2%)	0	100	100
2	K	104/134 (78%)	101 (97%)	3 (3%)	0	100	100
2	L	102/134 (76%)	99 (97%)	3 (3%)	0	100	100
All	All	1458/1794 (81%)	1422 (98%)	36 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	120/135 (89%)	115 (96%)	5 (4%)	26	40
1	B	116/135 (86%)	113 (97%)	3 (3%)	40	59
1	C	113/135 (84%)	108 (96%)	5 (4%)	25	38
1	G	121/135 (90%)	115 (95%)	6 (5%)	22	33
1	H	121/135 (90%)	116 (96%)	5 (4%)	27	41
1	I	118/135 (87%)	108 (92%)	10 (8%)	10	13
2	D	86/111 (78%)	82 (95%)	4 (5%)	23	35
2	E	83/111 (75%)	76 (92%)	7 (8%)	10	14
2	F	81/111 (73%)	77 (95%)	4 (5%)	22	34
2	J	86/111 (78%)	79 (92%)	7 (8%)	11	14

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	K	86/111 (78%)	75 (87%)	11 (13%)	4	4
2	L	82/111 (74%)	75 (92%)	7 (8%)	10	13
All	All	1213/1476 (82%)	1139 (94%)	74 (6%)	17	24

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	193	SER
1	A	194	SER
1	A	195	ARG
1	A	203	LEU
1	A	232	ARG
1	B	105	LEU
1	B	160	SER
1	B	195	ARG
1	C	102	ASN
1	C	154	CYS
1	C	163	THR
1	C	193	SER
1	C	203	LEU
2	D	43	LYS
2	D	67	CYS
2	D	71	THR
2	D	106	ARG
2	E	41	SER
2	E	52	GLU
2	E	67	CYS
2	E	70	LEU
2	E	111	CYS
2	E	115	GLU
2	E	122	SER
2	F	64	LYS
2	F	69	GLU
2	F	94	LEU
2	F	128	ILE
1	G	91	GLU
1	G	98	LEU
1	G	105	LEU
1	G	189	ARG
1	G	203	LEU
1	G	232	ARG

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Mol	Chain	Res	Type
1	H	102	ASN
1	H	160	SER
1	H	169	ARG
1	H	194	SER
1	H	203	LEU
1	I	98	LEU
1	I	105	LEU
1	I	106	THR
1	I	136	THR
1	I	172	ARG
1	I	189	ARG
1	I	191	THR
1	I	193	SER
1	I	203	LEU
1	I	228	ARG
2	J	45	ASP
2	J	52	GLU
2	J	69	GLU
2	J	70	LEU
2	J	73	THR
2	J	94	LEU
2	J	118	VAL
2	K	38	SER
2	K	39	LEU
2	K	44	GLU
2	K	51	SER
2	K	62	ARG
2	K	70	LEU
2	K	94	LEU
2	K	98	MET
2	K	106	ARG
2	K	129	VAL
2	K	131	ASP
2	L	43	LYS
2	L	67	CYS
2	L	70	LEU
2	L	71	THR
2	L	130	GLN
2	L	131	ASP
2	L	144	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	102	ASN
1	B	129	HIS
1	B	183	GLN
2	E	134	HIS
2	F	125	HIS
1	G	184	GLN
1	G	208	HIS
1	I	102	ASN
1	I	129	HIS
2	J	86	HIS
2	K	134	HIS

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](#)

Of 27 ligands modelled in this entry, 14 are monoatomic - leaving 13 for Mogul analysis.

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	$\# Z > 2$
4	GOL	H	302	-	5,5,5	0.94	0	5,5,5	0.48	0
5	NAG	J	201	2	14,14,15	0.73	0	17,19,21	2.38	5 (29%)
4	GOL	I	302	-	5,5,5	0.54	0	5,5,5	1.23	1 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GOL	D	201	-	5,5,5	0.42	0	5,5,5	0.86	0
4	GOL	B	303	-	5,5,5	0.49	0	5,5,5	0.42	0
5	NAG	L	202	2	14,14,15	1.28	2 (14%)	17,19,21	1.77	4 (23%)
4	GOL	C	302	-	5,5,5	0.45	0	5,5,5	0.76	0
5	NAG	D	202	2	14,14,15	1.32	1 (7%)	17,19,21	2.62	6 (35%)
5	NAG	K	202	2	14,14,15	0.80	1 (7%)	17,19,21	2.10	6 (35%)
4	GOL	F	201	-	5,5,5	0.27	0	5,5,5	1.48	2 (40%)
4	GOL	I	303	-	5,5,5	0.51	0	5,5,5	1.00	0
4	GOL	G	303	-	5,5,5	0.31	0	5,5,5	0.88	0
5	NAG	E	202	2	14,14,15	0.97	1 (7%)	17,19,21	1.43	2 (11%)

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
4	GOL	H	302	-	-	4/4/4/4	-
5	NAG	J	201	2	-	0/6/23/26	0/1/1/1
4	GOL	I	302	-	-	4/4/4/4	-
4	GOL	D	201	-	-	0/4/4/4	-
4	GOL	B	303	-	-	2/4/4/4	-
5	NAG	L	202	2	-	2/6/23/26	0/1/1/1
4	GOL	C	302	-	-	4/4/4/4	-
5	NAG	D	202	2	-	0/6/23/26	0/1/1/1
5	NAG	K	202	2	-	2/6/23/26	0/1/1/1
4	GOL	F	201	-	-	2/4/4/4	-
4	GOL	I	303	-	-	2/4/4/4	-
4	GOL	G	303	-	-	4/4/4/4	-
5	NAG	E	202	2	-	0/6/23/26	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	202	NAG	C1-C2	4.18	1.58	1.52
5	L	202	NAG	C1-C2	3.13	1.56	1.52
5	L	202	NAG	O5-C1	-3.06	1.38	1.43
5	E	202	NAG	C1-C2	2.92	1.56	1.52
5	K	202	NAG	C1-C2	2.23	1.55	1.52

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	202	NAG	C1-O5-C5	8.48	123.55	112.19
5	J	201	NAG	C1-O5-C5	7.70	122.51	112.19
5	K	202	NAG	C1-O5-C5	4.33	117.99	112.19
5	L	202	NAG	O5-C1-C2	-4.28	104.67	111.29
5	K	202	NAG	O5-C5-C6	3.83	115.11	107.66
5	K	202	NAG	C2-N2-C7	3.60	127.73	122.90
5	E	202	NAG	C1-O5-C5	3.42	116.78	112.19
5	L	202	NAG	C3-C4-C5	-3.13	104.55	110.23
5	L	202	NAG	C1-O5-C5	3.01	116.22	112.19
5	D	202	NAG	C8-C7-N2	2.92	120.96	116.12
5	J	201	NAG	O5-C5-C4	2.91	117.91	110.83
5	J	201	NAG	C3-C4-C5	-2.77	105.21	110.23
5	L	202	NAG	O5-C5-C6	2.66	112.85	107.66
5	K	202	NAG	C1-C2-N2	2.60	114.54	110.43
5	D	202	NAG	O5-C1-C2	-2.52	107.39	111.29
5	E	202	NAG	O4-C4-C5	2.52	115.53	109.32
5	D	202	NAG	O7-C7-C8	-2.52	117.57	122.05
4	I	302	GOL	O2-C2-C3	2.48	119.44	109.18
5	J	201	NAG	C1-C2-N2	-2.47	106.54	110.43
5	K	202	NAG	O7-C7-C8	-2.43	117.72	122.05
5	K	202	NAG	C6-C5-C4	-2.28	107.41	113.02
5	J	201	NAG	O4-C4-C5	2.26	114.89	109.32
4	F	201	GOL	O2-C2-C3	2.15	118.09	109.18
4	F	201	GOL	C3-C2-C1	-2.09	104.12	111.80
5	D	202	NAG	O5-C5-C6	2.03	111.62	107.66
5	D	202	NAG	C3-C4-C5	-2.03	106.55	110.23

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	303	GOL	O1-C1-C2-O2
4	B	303	GOL	O1-C1-C2-C3
4	C	302	GOL	C1-C2-C3-O3
4	G	303	GOL	O1-C1-C2-O2
4	G	303	GOL	O1-C1-C2-C3
4	G	303	GOL	C1-C2-C3-O3
4	H	302	GOL	O1-C1-C2-C3
4	I	302	GOL	C1-C2-C3-O3
5	L	202	NAG	C4-C5-C6-O6
5	K	202	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
5	K	202	NAG	C4-C5-C6-O6
5	L	202	NAG	O5-C5-C6-O6
4	I	302	GOL	O1-C1-C2-O2
4	C	302	GOL	O1-C1-C2-C3
4	F	201	GOL	C1-C2-C3-O3
4	H	302	GOL	C1-C2-C3-O3
4	I	302	GOL	O1-C1-C2-C3
4	I	303	GOL	C1-C2-C3-O3
4	C	302	GOL	O1-C1-C2-O2
4	C	302	GOL	O2-C2-C3-O3
4	G	303	GOL	O2-C2-C3-O3
4	H	302	GOL	O1-C1-C2-O2
4	I	302	GOL	O2-C2-C3-O3
4	F	201	GOL	O2-C2-C3-O3
4	H	302	GOL	O2-C2-C3-O3
4	I	303	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	F	201	GOL	2	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	149/165 (90%)	-0.46	1 (0%) 84 85	20, 27, 44, 82	0
1	B	142/165 (86%)	-0.14	6 (4%) 40 42	20, 31, 57, 91	0
1	C	138/165 (83%)	-0.05	6 (4%) 40 41	20, 31, 62, 85	0
1	G	150/165 (90%)	-0.51	1 (0%) 84 85	18, 25, 45, 73	0
1	H	150/165 (90%)	-0.11	7 (4%) 36 38	20, 32, 58, 73	0
1	I	147/165 (89%)	0.04	11 (7%) 20 22	20, 30, 70, 89	0
2	D	105/134 (78%)	-0.08	4 (3%) 44 46	23, 33, 61, 90	0
2	E	101/134 (75%)	0.80	10 (9%) 13 14	33, 52, 91, 108	0
2	F	98/134 (73%)	0.37	9 (9%) 14 16	25, 38, 81, 100	0
2	J	106/134 (79%)	0.27	4 (3%) 44 46	25, 36, 63, 91	0
2	K	106/134 (79%)	0.87	11 (10%) 11 13	32, 52, 88, 113	0
2	L	104/134 (77%)	0.24	11 (10%) 11 12	23, 36, 76, 137	0
All	All	1496/1794 (83%)	0.05	81 (5%) 31 33	18, 33, 73, 137	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	129	VAL	6.2
2	L	133	ASP	5.9
2	K	142	ALA	5.5
1	I	106	THR	5.3
1	I	159	ALA	5.1
2	K	129	VAL	4.9
2	K	134	HIS	4.7
1	B	154	CYS	4.6
2	J	70	LEU	4.4
1	I	155	PRO	4.3
2	J	142	ALA	4.2

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Mol	Chain	Res	Type	RSRZ
2	E	141	TYR	4.2
2	F	140	ALA	3.9
1	B	191	THR	3.9
2	E	134	HIS	3.9
1	I	158	LEU	3.8
2	L	131	ASP	3.7
2	L	132	GLY	3.6
2	E	129	VAL	3.6
1	C	154	CYS	3.5
1	C	186	PRO	3.4
2	K	70	LEU	3.4
1	G	189	ARG	3.4
2	L	41	SER	3.4
2	L	70	LEU	3.2
1	I	109	GLY	3.2
1	C	187	CYS	3.2
1	C	195	ARG	3.2
2	F	38	SER	3.1
2	L	130	GLN	3.1
2	D	70	LEU	3.1
1	C	193	SER	3.1
1	I	108	SER	3.1
2	E	140	ALA	3.1
2	L	134	HIS	3.0
1	I	105	LEU	3.0
1	I	160	SER	3.0
2	F	135	CYS	2.9
1	B	160	SER	2.9
2	F	128	ILE	2.9
1	H	220	LEU	2.8
2	K	132	GLY	2.8
1	B	91	GLU	2.8
2	E	123	PRO	2.8
2	F	126	PHE	2.8
1	A	195	ARG	2.8
2	L	144	SER	2.7
2	K	71	THR	2.7
1	H	155	PRO	2.7
1	I	101	ALA	2.7
2	K	141	TYR	2.6
1	C	194	SER	2.6
2	J	43	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
2	E	70	LEU	2.6
1	H	195	ARG	2.6
2	K	123	PRO	2.6
2	E	98	MET	2.5
2	F	71	THR	2.5
2	L	143	THR	2.5
1	H	159	ALA	2.5
2	D	142	ALA	2.4
2	E	128	ILE	2.4
1	H	156	LEU	2.4
2	F	70	LEU	2.4
2	F	139	ARG	2.4
1	I	102	ASN	2.3
2	E	124	GLY	2.3
2	L	71	THR	2.3
1	H	194	SER	2.3
2	K	140	ALA	2.3
2	J	71	THR	2.2
1	B	192	SER	2.2
2	D	38	SER	2.2
2	K	75	CYS	2.1
1	I	189	ARG	2.1
2	E	135	CYS	2.1
2	K	68	GLY	2.0
1	H	91	GLU	2.0
1	B	187	CYS	2.0
2	L	42	CYS	2.0
2	D	39	LEU	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(Å ²)	Q<0.9
5	NAG	K	202	14/15	0.65	0.17	71,82,93,93	0
5	NAG	D	202	14/15	0.69	0.19	73,87,99,112	0
5	NAG	J	201	14/15	0.74	0.14	54,70,85,88	0
5	NAG	L	202	14/15	0.75	0.15	64,78,86,87	0
5	NAG	E	202	14/15	0.76	0.14	74,84,93,94	0
4	GOL	H	302	6/6	0.85	0.20	36,37,38,40	0
4	GOL	B	303	6/6	0.91	0.13	44,48,49,53	0
4	GOL	I	302	6/6	0.92	0.10	33,33,36,37	0
4	GOL	C	302	6/6	0.94	0.10	33,35,37,38	0
4	GOL	F	201	6/6	0.94	0.10	38,40,42,48	0
3	CL	E	201	1/1	0.94	0.12	45,45,45,45	0
4	GOL	G	303	6/6	0.96	0.08	40,44,45,46	0
3	CL	G	304	1/1	0.96	0.09	43,43,43,43	0
4	GOL	D	201	6/6	0.97	0.08	31,33,34,36	0
3	CL	B	302	1/1	0.97	0.06	48,48,48,48	0
4	GOL	I	303	6/6	0.97	0.08	32,35,37,37	0
3	CL	C	301	1/1	0.97	0.11	40,40,40,40	0
3	CL	L	201	1/1	0.98	0.05	46,46,46,46	0
3	CL	C	303	1/1	0.98	0.06	42,42,42,42	0
3	CL	A	301	1/1	0.98	0.04	29,29,29,29	0
3	CL	G	302	1/1	0.98	0.06	35,35,35,35	0
3	CL	A	302	1/1	0.98	0.04	37,37,37,37	0
3	CL	H	301	1/1	0.98	0.05	33,33,33,33	0
3	CL	K	201	1/1	0.98	0.07	46,46,46,46	0
3	CL	I	301	1/1	0.99	0.03	37,37,37,37	0
3	CL	B	301	1/1	0.99	0.03	31,31,31,31	0
3	CL	G	301	1/1	1.00	0.02	25,25,25,25	0

6.5 Other polymers ⓘ

There are no such residues in this entry.