



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:55 PM UTC

PDB ID : 7KGU / pdb_00007kgu
Title : Structure of 2Q1-Fab, an antibody selective for IDH2R140Q-HLA-B*07:02
Authors : Miller, M.S.; Aytenfisu, T.Y.; Wright, K.M.; Gabelli, S.B.
Deposited on : 2020-10-18
Resolution : 2.40 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

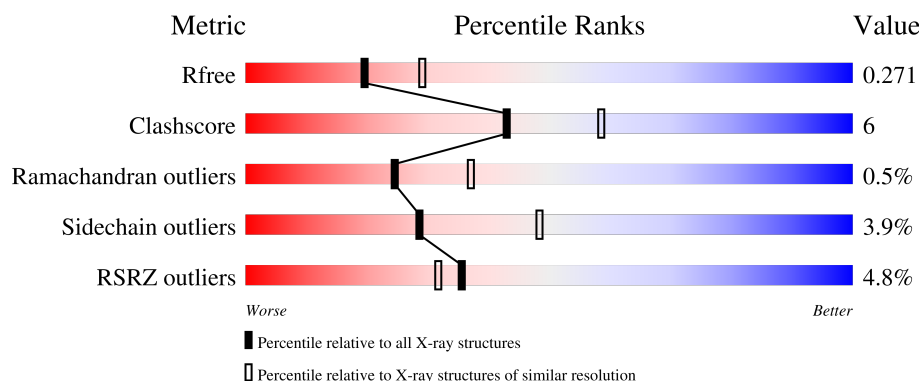
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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

Mol	Chain	Length	Quality of chain
1	A	214	9% 85% 13% ..
1	C	214	9% 89% 9% ..
1	E	214	9% 86% 10% ..
1	L	214	2% 82% 13% ..
2	B	218	6% 81% 14% ..

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Mol	Chain	Length	Quality of chain
2	D	218	
2	F	218	
2	H	218	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	BTB	A	302	-	-	X	-
5	SO4	D	304	-	-	X	-

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 13756 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.

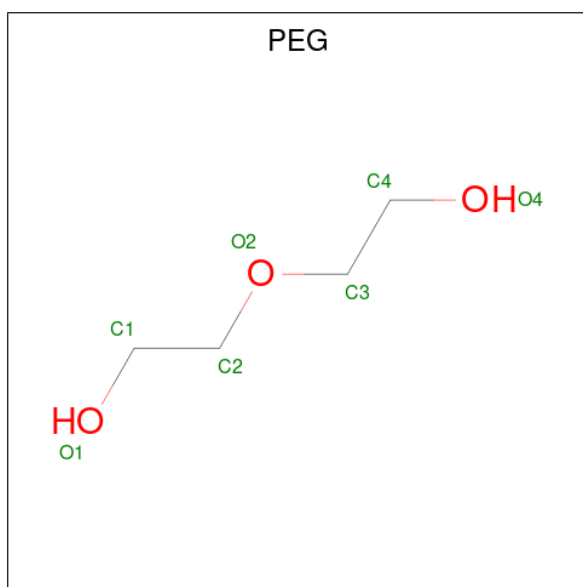
- Molecule 1 is a protein called Light Chain, Fab fragment, IGG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	212	Total	C	N	O	S	0	0	0
			1633	1025	273	330	5			
1	C	212	Total	C	N	O	S	0	0	0
			1633	1025	273	330	5			
1	E	210	Total	C	N	O	S	0	0	0
			1625	1021	271	328	5			
1	L	211	Total	C	N	O	S	0	0	0
			1629	1023	272	329	5			

- Molecule 2 is a protein called Heavy Chain, Fab, IGG, Fab.

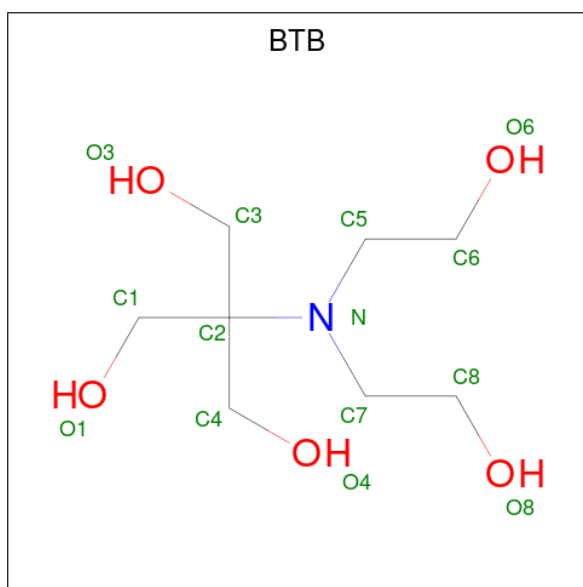
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	212	Total	C	N	O	S	0	0	0
			1600	1018	265	310	7			
2	D	212	Total	C	N	O	S	0	0	0
			1597	1016	265	309	7			
2	F	210	Total	C	N	O	S	0	0	0
			1592	1014	263	308	7			
2	H	211	Total	C	N	O	S	0	0	0
			1591	1013	264	307	7			

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



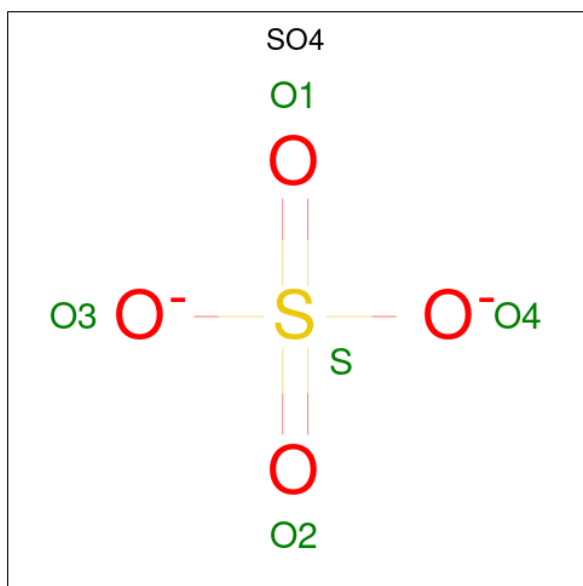
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		
3	C	1	Total	C	O	0	0
			7	4	3		
3	D	1	Total	C	O	0	0
			7	4	3		
3	D	1	Total	C	O	0	0
			7	4	3		
3	F	1	Total	C	O	0	0
			7	4	3		
3	H	1	Total	C	O	0	0
			7	4	3		
3	L	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is 2-[BIS-(2-HYDROXY-ETHYL)-AMINO]-2-HYDROXYMETHYL-PROPAN E-1,3-DIOL (CCD ID: BTB) (formula: C₈H₁₉NO₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		

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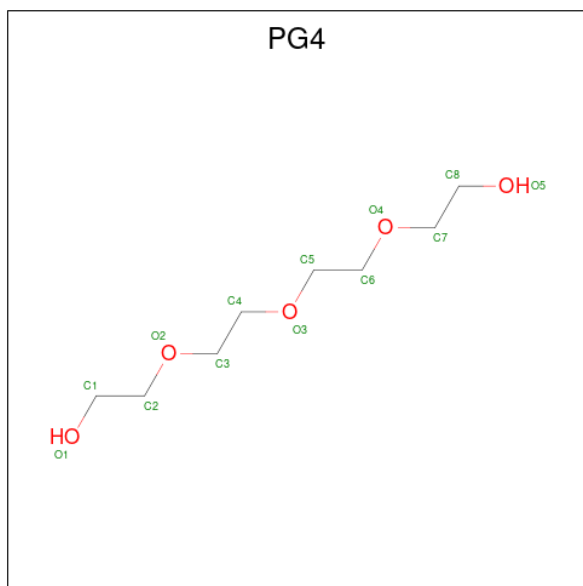
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	C	1	Total	O	S	0	0
			5	4	1		
5	C	1	Total	O	S	0	0
			5	4	1		
5	C	1	Total	O	S	0	0
			5	4	1		
5	D	1	Total	O	S	0	0
			5	4	1		
5	D	1	Total	O	S	0	0
			5	4	1		
5	D	1	Total	O	S	0	0
			5	4	1		
5	D	1	Total	O	S	0	0
			5	4	1		
5	D	1	Total	O	S	0	0
			5	4	1		
5	D	1	Total	O	S	0	0
			5	4	1		
5	F	1	Total	O	S	0	0
			5	4	1		
5	F	1	Total	O	S	0	0
			5	4	1		
5	F	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	F	1	Total	O	S	0	0
			5	4	1		
5	F	1	Total	O	S	0	0
			5	4	1		
5	F	1	Total	O	S	0	0
			5	4	1		
5	H	1	Total	O	S	0	0
			5	4	1		
5	H	1	Total	O	S	0	0
			5	4	1		
5	H	1	Total	O	S	0	0
			5	4	1		
5	H	1	Total	O	S	0	0
			5	4	1		
5	L	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			13	8	5		
6	B	1	Total	C	O	0	0
			13	8	5		

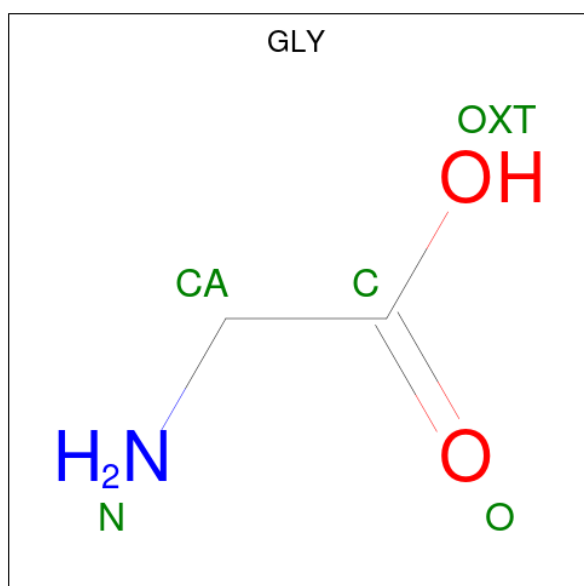
- Molecule 7 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	F	1	Total Cl 1 1	0	0
7	L	1	Total Cl 1 1	0	0

- Molecule 8 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	H	1	Total K 1 1	0	0

- Molecule 9 is GLYCINE (CCD ID: GLY) (formula: C₂H₅NO₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	L	1	Total C N O 4 2 1 1	0	0

- Molecule 10 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	A	67	Total O 67 67	0	0
10	B	84	Total O 84 84	0	0
10	C	65	Total O 65 65	0	0

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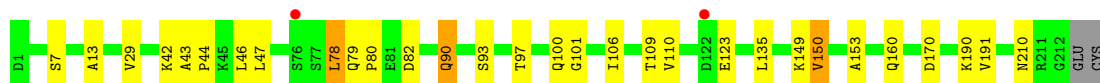
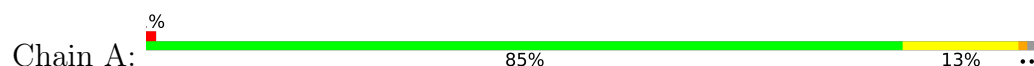
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	D	100	Total 100	O 100	0	0
10	E	46	Total 46	O 46	0	0
10	F	82	Total 82	O 82	0	0
10	H	91	Total 91	O 91	0	0
10	L	60	Total 60	O 60	0	0

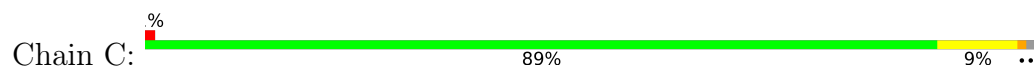
3 Residue-property plots [i](#)

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.

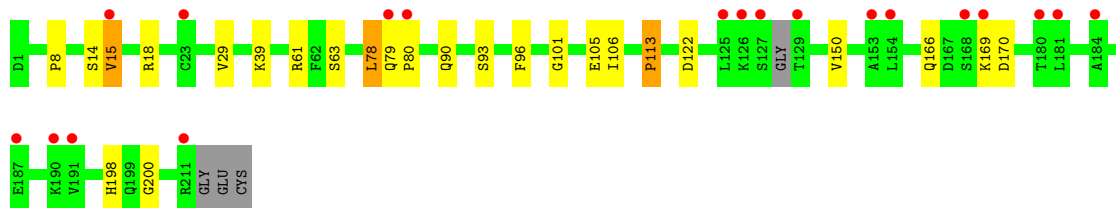
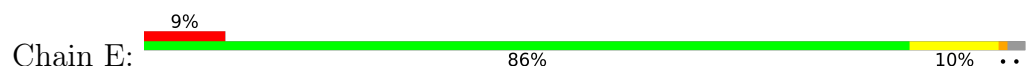
- Molecule 1: Light Chain, Fab fragment, IGG



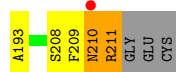
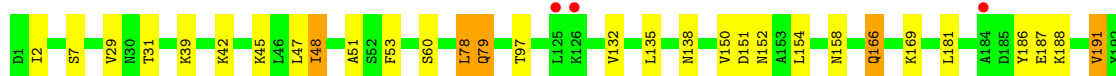
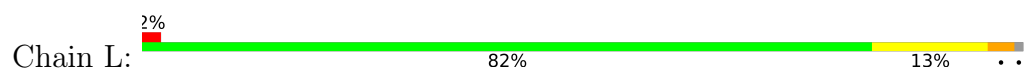
- Molecule 1: Light Chain, Fab fragment, IGG



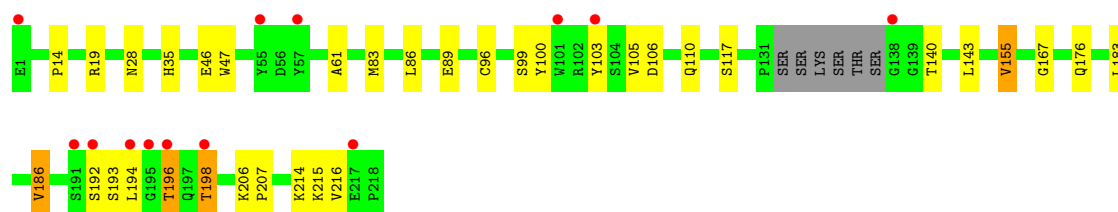
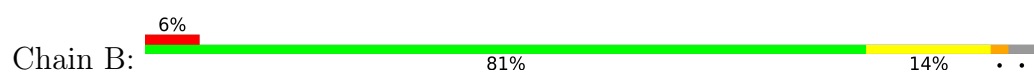
- Molecule 1: Light Chain, Fab fragment, IGG



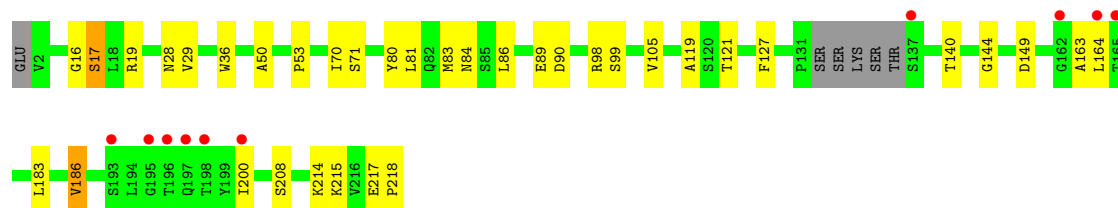
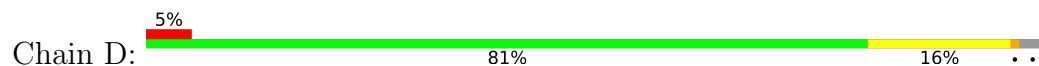
- Molecule 1: Light Chain, Fab fragment, IGG



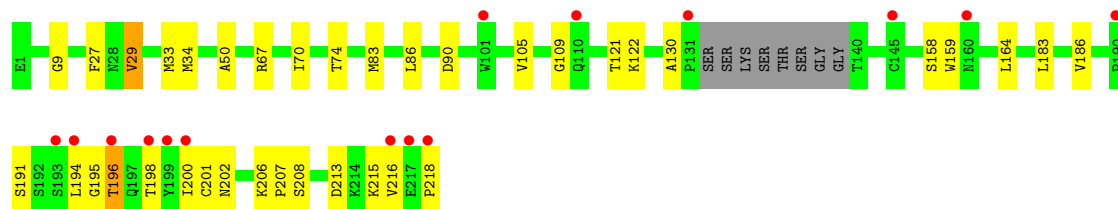
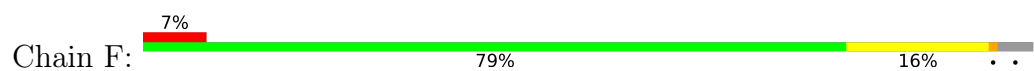
- Molecule 2: Heavy Chain, Fab, IGG, Fab



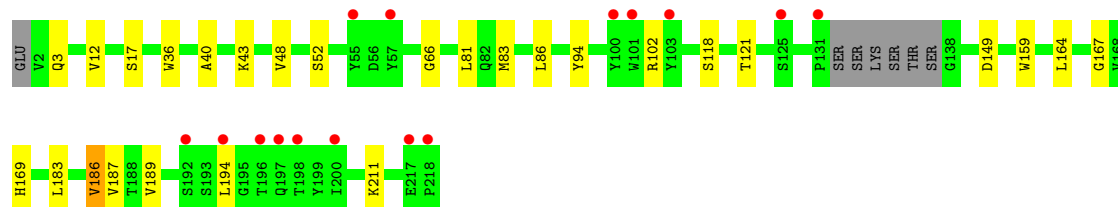
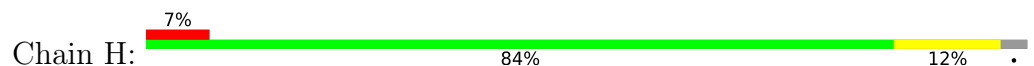
- Molecule 2: Heavy Chain, Fab, IGG, Fab



- Molecule 2: Heavy Chain, Fab, IGG, Fab



- Molecule 2: Heavy Chain, Fab, IGG, Fab



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	38.46Å 263.95Å 91.24Å 90.00° 99.53° 90.00°	Depositor
Resolution (Å)	45.53 – 2.40 45.53 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.8 (45.53-2.40) 98.8 (45.53-2.40)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0257	Depositor
R, R_{free}	0.208 , 0.269 0.213 , 0.271	Depositor DCC
R_{free} test set	3437 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	33.9	Xtriage
Anisotropy	0.351	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 41.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.039 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13756	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.94% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BTB, SO4, K, PG4, CL, PEG

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	1.11	0/1669	1.37	3/2267 (0.1%)
1	C	1.17	2/1669 (0.1%)	1.44	4/2267 (0.2%)
1	E	1.13	2/1660 (0.1%)	1.46	7/2254 (0.3%)
1	L	1.13	0/1665	1.46	2/2262 (0.1%)
2	B	1.17	1/1642 (0.1%)	1.36	2/2241 (0.1%)
2	D	1.21	6/1639 (0.4%)	1.40	6/2237 (0.3%)
2	F	1.19	1/1634 (0.1%)	1.43	8/2231 (0.4%)
2	H	1.21	5/1633 (0.3%)	1.40	4/2229 (0.2%)
All	All	1.16	17/13211 (0.1%)	1.42	36/17988 (0.2%)

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	130	ALA	C-N	9.54	1.49	1.34
2	D	71	SER	C-O	8.10	1.32	1.23
2	B	155	VAL	C-O	6.90	1.30	1.24
2	H	66	GLY	C-O	6.61	1.30	1.23
2	D	36	TRP	C-O	6.15	1.31	1.24
1	C	30	ASN	C-O	5.95	1.30	1.24
2	H	3	GLN	C-O	5.89	1.31	1.23
2	D	16	GLY	C-O	5.84	1.29	1.23
2	H	36	TRP	C-O	5.77	1.30	1.24
2	D	81	LEU	C-O	5.76	1.30	1.24
1	E	169	LYS	C-O	5.54	1.28	1.23
2	H	81	LEU	C-O	5.43	1.30	1.23
2	D	144	GLY	C-O	5.38	1.29	1.23
2	D	119	ALA	C-O	5.25	1.30	1.23
2	H	118	SER	C-O	5.16	1.30	1.24
1	C	80	PRO	C-O	-5.01	1.18	1.24
1	E	113	PRO	C-O	-5.01	1.18	1.23

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	27	PHE	CA-C-O	-6.51	114.30	121.33
1	C	187	GLU	CB-CG-CD	6.16	123.07	112.60
2	H	149	ASP	CA-CB-CG	6.14	118.74	112.60
1	E	170	ASP	CA-CB-CG	6.12	118.72	112.60
2	D	90	ASP	CA-CB-CG	5.98	118.58	112.60
2	F	67	ARG	CB-CA-C	-5.78	101.02	110.85
1	E	8	PRO	CA-C-N	5.65	128.17	120.54
1	E	8	PRO	C-N-CA	5.65	128.17	120.54
2	F	109	GLY	CA-C-N	5.49	128.19	120.28
2	F	109	GLY	C-N-CA	5.49	128.19	120.28
1	E	63	SER	CA-C-N	5.47	125.83	121.61
1	E	63	SER	C-N-CA	5.47	125.83	121.61
2	F	122	LYS	CB-CA-C	5.42	118.15	110.24
1	L	166	GLN	N-CA-CB	5.42	117.97	110.17
2	H	12	VAL	CA-C-N	5.39	127.88	120.39
2	H	12	VAL	C-N-CA	5.39	127.88	120.39
2	D	90	ASP	N-CA-C	-5.38	106.71	113.28
2	H	48	VAL	N-CA-C	-5.22	107.07	111.56
2	B	61	ALA	CA-C-O	-5.20	115.89	121.56
1	C	206	THR	CA-CB-OG1	-5.19	101.81	109.60
2	D	186	VAL	N-CA-CB	-5.18	102.95	111.45
1	A	170	ASP	CA-CB-CG	5.18	117.78	112.60
1	E	101	GLY	CA-C-O	-5.17	118.66	122.23
2	B	89	GLU	CB-CG-CD	5.17	121.38	112.60
2	F	213	ASP	CB-CA-C	5.16	119.15	109.71
1	A	44	PRO	CA-C-O	-5.15	115.58	121.56
2	D	163	ALA	CA-C-O	-5.15	115.39	120.90
1	C	91	VAL	CA-C-N	5.13	127.42	120.44
1	C	91	VAL	C-N-CA	5.13	127.42	120.44
1	L	31	THR	CA-CB-OG1	-5.12	101.92	109.60
2	D	164	LEU	CA-C-N	5.11	130.26	122.24
2	D	164	LEU	C-N-CA	5.11	130.26	122.24
2	F	90	ASP	N-CA-C	-5.09	107.07	113.28
1	E	18	ARG	CA-C-O	-5.07	115.48	121.16
2	F	74	THR	CA-C-O	-5.05	115.17	120.63
1	A	101	GLY	CA-C-O	-5.03	118.76	122.23

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	1633	0	1589	23	1
1	C	1633	0	1589	17	0
1	E	1625	0	1582	17	0
1	L	1629	0	1586	27	1
2	B	1600	0	1557	26	0
2	D	1597	0	1553	15	0
2	F	1592	0	1550	24	0
2	H	1591	0	1548	20	0
3	A	7	0	10	0	0
3	C	7	0	10	3	0
3	D	14	0	20	0	0
3	F	7	0	10	0	0
3	H	7	0	10	0	0
3	L	7	0	10	3	0
4	A	14	0	19	11	0
5	A	20	0	0	0	0
5	B	30	0	0	2	0
5	C	15	0	0	0	0
5	D	40	0	0	2	0
5	F	30	0	0	1	0
5	H	25	0	0	1	0
5	L	5	0	0	0	0
6	B	26	0	36	5	0
7	F	1	0	0	0	0
7	L	1	0	0	1	0
8	H	1	0	0	0	0
9	L	4	0	2	0	0
10	A	67	0	0	1	0
10	B	84	0	0	1	0
10	C	65	0	0	5	0
10	D	100	0	0	1	0
10	E	46	0	0	1	0
10	F	82	0	0	0	0
10	H	91	0	0	1	0
10	L	60	0	0	2	0
All	All	13756	0	12681	160	1

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 6.

All (160) 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)
2:F:194:LEU:HD11	2:F:218:PRO:CG	1.80	1.09
2:F:194:LEU:CD1	2:F:218:PRO:HG3	1.84	1.08
2:F:194:LEU:HD21	2:F:218:PRO:HG2	1.42	0.97
1:L:151:ASP:OD1	1:L:191:VAL:HG23	1.64	0.97
1:C:9:SER:H	3:C:301:PEG:H21	1.33	0.92
1:A:123:GLU:HB3	10:A:455:HOH:O	1.70	0.90
2:B:83:MET:HE2	2:B:86:LEU:HD21	1.63	0.80
2:F:194:LEU:HD21	2:F:218:PRO:CG	2.12	0.80
2:F:194:LEU:HD11	2:F:218:PRO:HG3	0.90	0.79
1:L:211:ARG:NH1	1:L:211:ARG:HG3	1.97	0.79
1:A:43:ALA:HB3	4:A:302:BTB:H61	1.65	0.79
2:F:83:MET:HE2	2:F:86:LEU:HD21	1.64	0.78
1:C:8:PRO:HA	3:C:301:PEG:H31	1.67	0.77
1:L:211:ARG:HG3	1:L:211:ARG:HH11	1.51	0.76
1:L:211:ARG:CD	1:L:211:ARG:H	1.97	0.75
1:A:42:LYS:HA	4:A:302:BTB:H82	1.69	0.73
1:E:80:PRO:HA	1:E:106:ILE:HD13	1.70	0.73
1:C:178:THR:HG22	10:C:435:HOH:O	1.88	0.72
1:L:211:ARG:H	1:L:211:ARG:HD2	1.54	0.72
2:D:28:ASN:N	5:D:304:SO4:O2	2.23	0.72
2:F:194:LEU:CD2	2:F:218:PRO:HG2	2.20	0.72
2:H:194:LEU:HD13	2:H:194:LEU:O	1.93	0.69
1:E:79:GLN:HB3	1:E:80:PRO:HD2	1.74	0.69
1:C:198:HIS:CD2	1:C:200:GLY:H	2.11	0.69
4:A:302:BTB:H71	2:B:110:GLN:HG2	1.75	0.68
1:L:45:LYS:HE3	3:L:302:PEG:H42	1.74	0.68
1:L:187:GLU:HA	1:L:187:GLU:OE1	1.94	0.67
2:H:40:ALA:HB3	2:H:43:LYS:HD2	1.77	0.67
1:A:190:LYS:HD2	1:A:210:ASN:HB3	1.77	0.66
1:L:187:GLU:HA	1:L:211:ARG:HH22	1.61	0.65
2:H:83:MET:HB3	2:H:86:LEU:HD21	1.79	0.64
1:A:78:LEU:O	1:A:82:ASP:OD2	2.17	0.64
2:D:83:MET:HE2	2:D:86:LEU:HD21	1.81	0.62
2:B:183:LEU:HD12	2:B:183:LEU:C	2.24	0.62
1:L:210:ASN:ND2	1:L:210:ASN:H	1.96	0.61
2:H:211:LYS:NZ	5:H:406:SO4:O4	2.31	0.60
1:E:90:GLN:HE21	1:E:93:SER:H	1.50	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:198:HIS:CD2	1:E:200:GLY:H	2.19	0.59
1:E:96:PHE:HE2	2:F:33:MET:HE1	1.68	0.59
1:E:15:VAL:HG22	1:E:15:VAL:O	2.04	0.57
2:F:183:LEU:HD12	2:F:183:LEU:C	2.29	0.57
1:L:211:ARG:HD2	1:L:211:ARG:N	2.17	0.57
1:C:103:LYS:CG	10:C:424:HOH:O	2.52	0.57
2:D:29:VAL:HG22	2:D:53:PRO:HG2	1.87	0.56
1:E:96:PHE:CE2	2:F:33:MET:HE1	2.40	0.56
2:H:183:LEU:C	2:H:183:LEU:HD12	2.31	0.56
2:B:35:HIS:HD2	2:B:47:TRP:HE1	1.54	0.56
6:B:502:PG4:H41	2:H:17:SER:N	2.22	0.55
1:C:103:LYS:HD2	10:C:425:HOH:O	2.05	0.55
1:E:106:ILE:H	1:E:166:GLN:HE22	1.53	0.55
2:H:186:VAL:HG11	1:L:135:LEU:HD22	1.88	0.55
1:A:135:LEU:CD2	2:B:186:VAL:HG11	2.36	0.55
2:F:33:MET:HE2	2:F:50:ALA:CB	2.37	0.55
2:B:83:MET:HE2	2:B:86:LEU:CD2	2.35	0.54
1:E:106:ILE:N	1:E:166:GLN:HE22	2.06	0.53
2:D:183:LEU:C	2:D:183:LEU:HD12	2.33	0.53
1:E:15:VAL:O	1:E:15:VAL:HG13	2.08	0.53
2:H:159:TRP:HB3	2:H:164:LEU:HD21	1.91	0.52
1:C:124:GLN:HE22	1:C:131:SER:HB2	1.74	0.52
1:A:135:LEU:HD22	2:B:186:VAL:HG11	1.90	0.52
1:C:15:VAL:HG21	1:C:80:PRO:HG3	1.91	0.52
2:F:195:GLY:O	2:F:196:THR:HG23	2.10	0.52
1:L:45:LYS:CE	3:L:302:PEG:H42	2.39	0.52
2:H:189:VAL:HG12	10:H:521:HOH:O	2.10	0.52
6:B:502:PG4:H41	2:H:17:SER:H	1.74	0.52
1:C:103:LYS:HG2	10:C:424:HOH:O	2.09	0.52
2:D:83:MET:HE2	2:D:86:LEU:CD2	2.40	0.52
1:A:109:THR:HG22	1:A:110:VAL:O	2.10	0.51
1:A:42:LYS:HB3	4:A:302:BTB:H42	1.93	0.51
2:F:206:LYS:N	2:F:207:PRO:CD	2.74	0.51
6:B:502:PG4:C4	2:H:17:SER:HB2	2.41	0.51
6:B:502:PG4:H41	2:H:17:SER:HB2	1.93	0.51
1:C:9:SER:N	3:C:301:PEG:H21	2.14	0.51
1:L:152:ASN:HB2	10:L:421:HOH:O	2.11	0.51
1:L:150:VAL:HA	1:L:191:VAL:O	2.10	0.51
1:A:150:VAL:HA	1:A:191:VAL:O	2.10	0.50
2:F:9:GLY:N	5:F:307:SO4:O3	2.45	0.50
1:E:106:ILE:H	1:E:166:GLN:NE2	2.09	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:211:ARG:HH11	1:L:211:ARG:CG	2.21	0.50
2:B:206:LYS:N	2:B:207:PRO:CD	2.75	0.49
4:A:302:BTB:H52	2:B:110:GLN:HG2	1.94	0.49
1:C:206:THR:O	1:C:207:LYS:HE2	2.11	0.49
2:D:17:SER:OG	2:D:84:ASN:OD1	2.30	0.49
2:D:99:SER:HB3	2:D:105:VAL:HG12	1.94	0.49
2:B:193:SER:HA	2:B:196:THR:HB	1.95	0.49
1:A:80:PRO:HA	1:A:106:ILE:HG12	1.93	0.49
4:A:302:BTB:H71	2:B:110:GLN:HE21	1.78	0.49
1:C:103:LYS:HG3	10:C:424:HOH:O	2.13	0.48
1:E:90:GLN:NE2	1:E:93:SER:H	2.11	0.48
1:A:46:LEU:HD22	2:B:106:ASP:HA	1.95	0.48
1:C:124:GLN:HA	2:D:127:PHE:CE1	2.49	0.48
4:A:302:BTB:C7	2:B:110:GLN:HE21	2.27	0.47
1:A:190:LYS:O	1:A:210:ASN:HA	2.15	0.47
4:A:302:BTB:H52	2:B:110:GLN:HA	1.95	0.47
2:F:159:TRP:CZ3	2:F:201:CYS:HB3	2.49	0.47
2:B:35:HIS:HE1	2:B:99:SER:HB3	1.79	0.47
2:F:159:TRP:CH2	2:F:201:CYS:HB3	2.49	0.47
1:L:169:LYS:NZ	10:L:401:HOH:O	2.48	0.47
1:L:209:PHE:CD1	1:L:209:PHE:C	2.92	0.47
2:D:19:ARG:HD3	2:D:80:TYR:HB3	1.97	0.47
1:L:42:LYS:HD3	3:L:302:PEG:H32	1.97	0.46
1:L:210:ASN:ND2	1:L:210:ASN:N	2.59	0.46
2:B:28:ASN:N	5:D:304:SO4:O1	2.41	0.46
1:A:90:GLN:HE21	1:A:93:SER:H	1.64	0.46
2:H:40:ALA:CB	2:H:43:LYS:HD2	2.45	0.46
2:H:169:HIS:HE1	1:L:138:ASN:HD21	1.63	0.46
2:B:35:HIS:O	2:B:96:CYS:HA	2.17	0.45
2:D:83:MET:HB3	2:D:86:LEU:HD21	2.00	0.44
1:E:113:PRO:HD3	1:E:198:HIS:CD2	2.53	0.44
1:A:106:ILE:N	1:A:106:ILE:HD12	2.32	0.44
1:A:90:GLN:OE1	1:A:97:THR:OG1	2.36	0.44
4:A:302:BTB:C5	2:B:110:GLN:HG2	2.47	0.43
2:F:33:MET:HE2	2:F:50:ALA:HB1	2.00	0.43
2:H:194:LEU:HD13	2:H:194:LEU:C	2.43	0.43
1:A:160:GLN:HE22	2:B:176:GLN:HA	1.83	0.43
2:B:167:GLY:HA2	5:B:506:SO4:O2	2.18	0.43
2:H:83:MET:CE	2:H:94:TYR:CZ	3.01	0.43
1:A:13:ALA:HB3	1:A:78:LEU:HD22	2.00	0.43
1:L:193:ALA:HB2	1:L:208:SER:HB3	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:19:ARG:NH1	10:D:407:HOH:O	2.51	0.43
2:F:121:THR:HG21	2:F:208:SER:HA	2.00	0.43
2:B:100:TYR:O	2:B:103:TYR:HB2	2.18	0.43
1:E:105:GLU:HB3	10:E:334:HOH:O	2.19	0.42
2:F:29:VAL:HG23	2:F:34:MET:SD	2.60	0.42
2:F:158:SER:OG	2:F:202:ASN:HB2	2.19	0.42
6:B:502:PG4:H41	2:H:17:SER:CB	2.48	0.42
2:H:83:MET:HE1	2:H:94:TYR:CZ	2.55	0.42
2:D:217:GLU:HB2	2:D:218:PRO:HD2	2.01	0.42
2:F:194:LEU:CD1	2:F:218:PRO:CG	2.66	0.42
2:D:121:THR:HG22	2:D:208:SER:HB3	2.01	0.42
2:F:83:MET:HE2	2:F:86:LEU:CD2	2.42	0.42
1:C:198:HIS:HD2	1:C:200:GLY:H	1.61	0.42
1:L:2:ILE:O	1:L:97:THR:OG1	2.31	0.42
1:A:90:GLN:NE2	1:A:93:SER:O	2.53	0.41
2:H:167:GLY:O	2:H:187:VAL:HA	2.20	0.41
2:D:200:ILE:HA	2:D:214:LYS:O	2.19	0.41
1:E:78:LEU:HD11	1:E:106:ILE:HG12	2.01	0.41
1:A:149:LYS:HA	1:A:153:ALA:O	2.20	0.41
2:B:14:PRO:HD3	2:B:117:SER:C	2.44	0.41
2:F:50:ALA:C	2:F:70:ILE:HD13	2.46	0.41
1:A:42:LYS:HD3	4:A:302:BTB:O1	2.21	0.41
1:C:90:GLN:NE2	1:C:93:SER:H	2.18	0.41
2:D:50:ALA:C	2:D:70:ILE:HD13	2.46	0.41
2:B:198:THR:HB	2:B:215:LYS:NZ	2.36	0.41
1:C:15:VAL:CG2	1:C:80:PRO:HG3	2.49	0.41
1:C:124:GLN:HE22	1:C:131:SER:CB	2.34	0.41
1:E:29:VAL:HG11	1:E:90:GLN:HG3	2.03	0.41
2:B:46:GLU:HA	5:B:507:SO4:O3	2.21	0.41
1:E:150:VAL:O	1:E:150:VAL:HG23	2.20	0.41
4:A:302:BTB:H11	4:A:302:BTB:H72	1.89	0.40
1:L:51:ALA:O	7:L:304:CL:CL	2.76	0.40
1:A:78:LEU:O	1:A:82:ASP:CG	2.65	0.40
2:F:159:TRP:HB3	2:F:164:LEU:HD21	2.02	0.40
2:H:186:VAL:HG11	1:L:135:LEU:CD2	2.49	0.40
1:L:188:LYS:HB3	1:L:188:LYS:HE2	1.95	0.40
2:B:155:VAL:CG1	2:B:183:LEU:HD21	2.51	0.40
1:L:48:ILE:HA	1:L:53:PHE:O	2.20	0.40
1:L:186:TYR:O	1:L:211:ARG:NH2	2.55	0.40
1:A:29:VAL:HG11	1:A:90:GLN:HG3	2.04	0.40
2:B:28:ASN:ND2	10:B:606:HOH:O	2.52	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:THR:CG2	1:L:79:GLN:NE2[1_556]	1.68	0.52

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	210/214 (98%)	199 (95%)	9 (4%)	2 (1%)	12	20
1	C	210/214 (98%)	200 (95%)	9 (4%)	1 (0%)	24	37
1	E	206/214 (96%)	197 (96%)	8 (4%)	1 (0%)	24	37
1	L	209/214 (98%)	197 (94%)	10 (5%)	2 (1%)	12	20
2	B	208/218 (95%)	200 (96%)	7 (3%)	1 (0%)	24	37
2	D	208/218 (95%)	204 (98%)	3 (1%)	1 (0%)	24	37
2	F	206/218 (94%)	198 (96%)	7 (3%)	1 (0%)	24	37
2	H	207/218 (95%)	201 (97%)	6 (3%)	0	100	100
All	All	1664/1728 (96%)	1596 (96%)	59 (4%)	9 (0%)	24	37

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	15	VAL
2	B	196	THR
1	C	122	ASP
1	A	78	LEU
2	D	149	ASP
2	F	196	THR
1	A	79	GLN
1	L	78	LEU
1	L	79	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	187/189 (99%)	182 (97%)	5 (3%)	39	62
1	C	187/189 (99%)	183 (98%)	4 (2%)	47	69
1	E	187/189 (99%)	182 (97%)	5 (3%)	39	62
1	L	187/189 (99%)	172 (92%)	15 (8%)	11	19
2	B	176/182 (97%)	166 (94%)	10 (6%)	18	33
2	D	176/182 (97%)	170 (97%)	6 (3%)	32	54
2	F	176/182 (97%)	168 (96%)	8 (4%)	24	42
2	H	175/182 (96%)	171 (98%)	4 (2%)	44	66
All	All	1451/1484 (98%)	1394 (96%)	57 (4%)	28	48

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	SER
1	A	47	LEU
1	A	90	GLN
1	A	100	GLN
1	A	150	VAL
2	B	19	ARG
2	B	105	VAL
2	B	140	THR
2	B	143	LEU
2	B	186	VAL
2	B	192	SER
2	B	194	LEU
2	B	198	THR
2	B	214	LYS
2	B	216	VAL
1	C	23	CYS
1	C	56	SER
1	C	63	SER
1	C	150	VAL

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Mol	Chain	Res	Type
2	D	17	SER
2	D	89	GLU
2	D	98	ARG
2	D	140	THR
2	D	186	VAL
2	D	215	LYS
1	E	14	SER
1	E	39	LYS
1	E	61	ARG
1	E	78	LEU
1	E	122	ASP
2	F	29	VAL
2	F	105	VAL
2	F	186	VAL
2	F	191	SER
2	F	198	THR
2	F	200	ILE
2	F	215	LYS
2	F	216	VAL
2	H	52	SER
2	H	102	ARG
2	H	121	THR
2	H	186	VAL
1	L	7	SER
1	L	29	VAL
1	L	39	LYS
1	L	47	LEU
1	L	48	ILE
1	L	60	SER
1	L	78	LEU
1	L	132	VAL
1	L	154	LEU
1	L	158	ASN
1	L	166	GLN
1	L	181	LEU
1	L	191	VAL
1	L	210	ASN
1	L	211	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	137	ASN
1	A	138	ASN
1	A	152	ASN
1	A	160	GLN
1	A	210	ASN
2	B	13	GLN
2	B	35	HIS
2	B	110	GLN
2	B	197	GLN
1	C	90	GLN
1	C	124	GLN
1	C	138	ASN
1	C	160	GLN
1	C	198	HIS
2	D	3	GLN
2	D	169	HIS
1	E	38	GLN
1	E	79	GLN
1	E	198	HIS
2	F	3	GLN
2	F	39	GLN
2	F	197	GLN
2	H	3	GLN
2	H	110	GLN
1	L	37	GLN
1	L	79	GLN
1	L	124	GLN
1	L	137	ASN
1	L	138	ASN
1	L	158	ASN
1	L	198	HIS
1	L	210	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 47 ligands modelled in this entry, 3 are monoatomic - leaving 44 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	$\# Z > 2$
3	PEG	F	301	-	6,6,6	0.20	0	5,5,5	0.14	0
3	PEG	D	301	-	6,6,6	0.17	0	5,5,5	0.29	0
5	SO4	D	307	-	4,4,4	0.31	0	6,6,6	0.10	0
5	SO4	F	305	-	4,4,4	0.35	0	6,6,6	0.09	0
5	SO4	F	302	-	4,4,4	0.45	0	6,6,6	0.13	0
5	SO4	F	307	-	4,4,4	0.32	0	6,6,6	0.09	0
9	GLY	L	301	-	3,3,4	0.70	0	1,2,4	0.72	0
5	SO4	C	303	-	4,4,4	0.36	0	6,6,6	0.10	0
4	BTB	A	302	-	13,13,13	1.64	3 (23%)	7,16,16	1.23	1 (14%)
5	SO4	D	305	-	4,4,4	0.35	0	6,6,6	0.07	0
5	SO4	H	404	-	4,4,4	0.30	0	6,6,6	0.09	0
3	PEG	A	301	-	6,6,6	0.26	0	5,5,5	0.17	0
5	SO4	B	508	-	4,4,4	0.31	0	6,6,6	0.08	0
5	SO4	C	302	-	4,4,4	0.32	0	6,6,6	0.09	0
5	SO4	D	304	-	4,4,4	0.35	0	6,6,6	0.08	0
5	SO4	A	304	-	4,4,4	0.30	0	6,6,6	0.16	0
3	PEG	D	302	-	6,6,6	0.21	0	5,5,5	0.17	0
6	PG4	B	502	-	12,12,12	0.32	0	11,11,11	0.13	0
5	SO4	F	304	-	4,4,4	0.31	0	6,6,6	0.07	0
5	SO4	D	309	-	4,4,4	0.30	0	6,6,6	0.13	0
5	SO4	C	304	-	4,4,4	0.32	0	6,6,6	0.14	0
5	SO4	H	402	-	4,4,4	0.32	0	6,6,6	0.06	0
6	PG4	B	501	-	12,12,12	0.23	0	11,11,11	0.20	0
5	SO4	D	303	-	4,4,4	0.28	0	6,6,6	0.21	0
5	SO4	B	507	-	4,4,4	0.36	0	6,6,6	0.07	0
3	PEG	C	301	-	6,6,6	0.21	0	5,5,5	0.31	0
5	SO4	A	303	-	4,4,4	0.37	0	6,6,6	0.13	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	SO4	B	505	-	4,4,4	0.28	0	6,6,6	0.05	0
5	SO4	F	306	-	4,4,4	0.28	0	6,6,6	0.06	0
5	SO4	D	306	-	4,4,4	0.25	0	6,6,6	0.09	0
5	SO4	A	306	-	4,4,4	0.33	0	6,6,6	0.07	0
5	SO4	H	405	-	4,4,4	0.43	0	6,6,6	0.14	0
5	SO4	D	310	-	4,4,4	0.33	0	6,6,6	0.06	0
5	SO4	A	305	-	4,4,4	0.35	0	6,6,6	0.10	0
3	PEG	L	302	-	6,6,6	0.16	0	5,5,5	0.17	0
5	SO4	L	303	-	4,4,4	0.35	0	6,6,6	0.08	0
5	SO4	B	504	-	4,4,4	0.31	0	6,6,6	0.11	0
5	SO4	D	308	-	4,4,4	0.32	0	6,6,6	0.09	0
5	SO4	H	403	-	4,4,4	0.27	0	6,6,6	0.09	0
5	SO4	B	503	-	4,4,4	0.36	0	6,6,6	0.19	0
5	SO4	B	506	-	4,4,4	0.32	0	6,6,6	0.14	0
5	SO4	H	406	-	4,4,4	0.31	0	6,6,6	0.07	0
5	SO4	F	303	-	4,4,4	0.33	0	6,6,6	0.05	0
3	PEG	H	401	-	6,6,6	0.17	0	5,5,5	0.13	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
6	PG4	B	501	-	-	5/10/10/10	-
3	PEG	D	302	-	-	0/4/4/4	-
3	PEG	F	301	-	-	0/4/4/4	-
3	PEG	D	301	-	-	1/4/4/4	-
6	PG4	B	502	-	-	6/10/10/10	-
3	PEG	C	301	-	-	2/4/4/4	-
9	GLY	L	301	-	-	0/0/1/2	-
4	BTB	A	302	-	-	10/21/21/21	-
3	PEG	A	301	-	-	2/4/4/4	-
3	PEG	L	302	-	-	0/4/4/4	-
3	PEG	H	401	-	-	0/4/4/4	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	302	BTB	C2-N	4.22	1.56	1.48
4	A	302	BTB	C5-N	2.50	1.51	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	302	BTB	C7-N	2.26	1.51	1.48

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	302	BTB	O4-C4-C2	2.25	116.70	111.40

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	302	BTB	C1-C2-C4-O4
4	A	302	BTB	C3-C2-C4-O4
4	A	302	BTB	N-C2-C4-O4
4	A	302	BTB	C1-C2-N-C7
4	A	302	BTB	C3-C2-N-C7
4	A	302	BTB	C4-C2-N-C7
4	A	302	BTB	C6-C5-N-C2
6	B	501	PG4	O2-C3-C4-O3
3	A	301	PEG	O2-C3-C4-O4
3	C	301	PEG	O2-C3-C4-O4
6	B	501	PG4	O1-C1-C2-O2
6	B	502	PG4	O4-C7-C8-O5
6	B	502	PG4	O1-C1-C2-O2
3	A	301	PEG	O1-C1-C2-O2
3	C	301	PEG	C4-C3-O2-C2
6	B	502	PG4	C1-C2-O2-C3
6	B	502	PG4	O3-C5-C6-O4
4	A	302	BTB	C4-C2-N-C5
6	B	501	PG4	O4-C7-C8-O5
6	B	502	PG4	C6-C5-O3-C4
4	A	302	BTB	C4-C2-C3-O3
6	B	501	PG4	O3-C5-C6-O4
6	B	502	PG4	C3-C4-O3-C5
6	B	501	PG4	C3-C4-O3-C5
4	A	302	BTB	C3-C2-N-C5
3	D	301	PEG	O1-C1-C2-O2

There are no ring outliers.

9 monomers are involved in 28 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	F	307	SO4	1	0
4	A	302	BTB	11	0
5	D	304	SO4	2	0
6	B	502	PG4	5	0
5	B	507	SO4	1	0
3	C	301	PEG	3	0
3	L	302	PEG	3	0
5	B	506	SO4	1	0
5	H	406	SO4	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 ⓘ

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	212/214 (99%)	0.04	2 (0%) 81 78	23, 41, 65, 80	0
1	C	212/214 (99%)	0.13	3 (1%) 73 69	20, 41, 73, 88	0
1	E	210/214 (98%)	0.68	19 (9%) 15 12	29, 57, 95, 119	0
1	L	211/214 (98%)	0.11	4 (1%) 66 62	22, 39, 81, 96	0
2	B	212/218 (97%)	0.01	13 (6%) 27 23	16, 30, 72, 92	0
2	D	212/218 (97%)	0.01	10 (4%) 36 32	17, 29, 72, 93	0
2	F	210/218 (96%)	0.14	15 (7%) 22 18	17, 33, 87, 105	0
2	H	211/218 (96%)	0.11	15 (7%) 22 18	16, 31, 77, 94	0
All	All	1690/1728 (97%)	0.15	81 (4%) 35 32	16, 38, 79, 119	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	154	LEU	5.1
2	D	164	LEU	4.9
2	H	57	TYR	4.7
1	E	187	GLU	4.5
2	B	138	GLY	4.3
2	B	194	LEU	4.1
2	F	198	THR	4.0
2	H	101	TRP	3.9
1	E	125	LEU	3.8
2	F	196	THR	3.7
2	D	196	THR	3.6
1	L	184	ALA	3.6
1	E	181	LEU	3.6
2	B	196	THR	3.5
2	H	100	TYR	3.5
2	B	192	SER	3.4

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Mol	Chain	Res	Type	RSRZ
2	F	218	PRO	3.3
2	F	200	ILE	3.3
2	F	199	TYR	3.2
2	F	216	VAL	3.2
1	E	153	ALA	3.1
1	E	15	VAL	3.0
2	B	55	TYR	2.9
2	H	192	SER	2.9
2	B	103	TYR	2.7
2	F	194	LEU	2.7
2	D	198	THR	2.7
1	E	126	LYS	2.7
2	H	55	TYR	2.6
2	F	101	TRP	2.6
2	D	193	SER	2.6
2	B	57	TYR	2.6
2	H	196	THR	2.6
2	F	193	SER	2.6
2	F	131	PRO	2.5
1	L	125	LEU	2.5
1	E	191	VAL	2.5
1	E	168	SER	2.5
1	E	127	SER	2.4
2	H	218	PRO	2.4
2	H	103	TYR	2.4
2	H	194	LEU	2.4
2	B	101	TRP	2.4
1	E	184	ALA	2.4
2	B	191	SER	2.4
1	E	129	THR	2.4
1	E	80	PRO	2.4
2	H	217	GLU	2.4
1	C	212	GLY	2.3
2	B	195	GLY	2.3
2	D	200	ILE	2.3
2	H	200	ILE	2.3
1	L	126	LYS	2.3
2	B	217	GLU	2.3
2	F	145	CYS	2.3
2	B	198	THR	2.3
1	E	79	GLN	2.3
2	F	190	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	128	GLY	2.2
1	C	125	LEU	2.2
1	E	169	LYS	2.2
2	D	137	SER	2.2
2	F	217	GLU	2.2
2	D	162	GLY	2.1
2	H	198	THR	2.1
2	H	125	SER	2.1
1	E	190	LYS	2.1
2	F	110	GLN	2.1
2	H	197	GLN	2.1
2	D	195	GLY	2.1
2	H	131	PRO	2.1
1	L	210	ASN	2.1
1	E	211	ARG	2.0
2	F	160	ASN	2.0
2	B	1	GLU	2.0
1	E	180	THR	2.0
2	D	165	THR	2.0
1	A	122	ASP	2.0
1	E	23	CYS	2.0
2	D	197	GLN	2.0
1	A	76	SER	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.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	SO4	F	306	5/5	0.54	0.17	76,80,83,92	0
5	SO4	B	508	5/5	0.70	0.12	79,81,90,91	0
3	PEG	A	301	7/7	0.71	0.17	67,73,76,76	0
5	SO4	D	309	5/5	0.72	0.16	74,77,84,88	0
5	SO4	A	306	5/5	0.72	0.12	87,90,93,96	0
5	SO4	D	307	5/5	0.75	0.12	72,84,86,87	0
9	GLY	L	301	4/5	0.75	0.17	73,75,75,76	0
5	SO4	H	406	5/5	0.76	0.16	64,69,71,81	0
4	BTB	A	302	14/14	0.76	0.19	68,72,85,85	0
5	SO4	D	310	5/5	0.77	0.13	90,91,94,95	0
5	SO4	B	506	5/5	0.78	0.15	65,71,78,82	0
6	PG4	B	502	13/13	0.78	0.20	44,68,78,78	0
5	SO4	A	305	5/5	0.78	0.10	87,91,94,95	0
3	PEG	L	302	7/7	0.79	0.19	62,65,70,71	0
6	PG4	B	501	13/13	0.80	0.19	51,58,71,71	0
3	PEG	D	302	7/7	0.80	0.19	49,53,58,62	0
3	PEG	C	301	7/7	0.80	0.15	53,58,60,60	0
3	PEG	D	301	7/7	0.82	0.18	49,56,66,73	0
5	SO4	C	302	5/5	0.82	0.10	85,86,92,93	0
3	PEG	F	301	7/7	0.83	0.20	62,65,69,69	0
5	SO4	F	307	5/5	0.83	0.12	79,83,87,88	0
5	SO4	A	304	5/5	0.83	0.13	68,73,76,77	0
5	SO4	H	404	5/5	0.84	0.15	71,80,83,84	0
3	PEG	H	401	7/7	0.85	0.14	49,51,52,57	0
5	SO4	B	507	5/5	0.85	0.19	83,85,94,106	0
5	SO4	H	402	5/5	0.86	0.18	67,73,80,81	0
5	SO4	D	306	5/5	0.87	0.23	62,62,63,65	0
5	SO4	H	403	5/5	0.89	0.14	70,70,77,77	0
5	SO4	C	304	5/5	0.89	0.14	74,79,80,84	0
5	SO4	F	303	5/5	0.90	0.17	63,65,70,74	0
5	SO4	F	304	5/5	0.91	0.10	55,60,65,69	0
5	SO4	D	305	5/5	0.91	0.09	54,56,58,61	0
5	SO4	A	303	5/5	0.92	0.13	51,53,55,58	0
5	SO4	C	303	5/5	0.92	0.24	51,56,62,65	0
5	SO4	B	504	5/5	0.92	0.13	53,57,62,66	0
5	SO4	D	304	5/5	0.92	0.27	37,50,57,59	0
5	SO4	B	505	5/5	0.92	0.17	56,62,63,65	0
7	CL	L	304	1/1	0.93	0.20	62,62,62,62	0
5	SO4	F	305	5/5	0.93	0.12	74,76,77,82	0
5	SO4	D	308	5/5	0.94	0.12	64,66,68,74	0
5	SO4	L	303	5/5	0.96	0.07	49,49,53,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	SO4	B	503	5/5	0.97	0.06	31,33,36,37	0
5	SO4	F	302	5/5	0.97	0.07	33,34,34,38	0
8	K	H	407	1/1	0.97	0.15	52,52,52,52	0
5	SO4	H	405	5/5	0.97	0.10	39,39,42,46	0
5	SO4	D	303	5/5	0.98	0.06	29,31,34,35	0
7	CL	F	308	1/1	0.99	0.05	21,21,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.