



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 05:47 pm BST

PDB ID : 5MV4
Title : ACC1 Fab fragment in complex with citrullinated CII616-639 epitope of collagen type II (ptm23)
Authors : Dobritsch, D.; Holmdahl, R.; Ge, C.
Deposited on : 2017-01-15
Resolution : 2.90 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

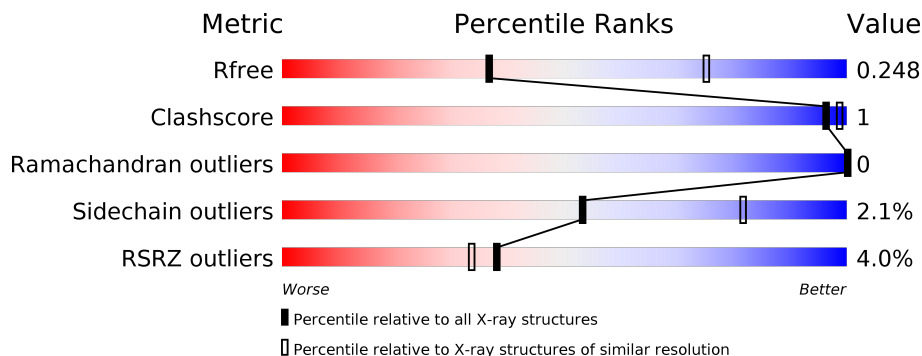
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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.



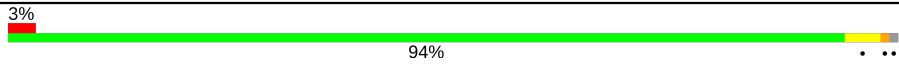
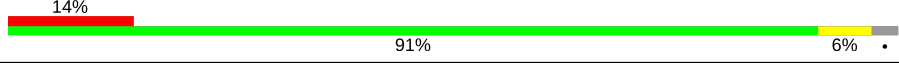
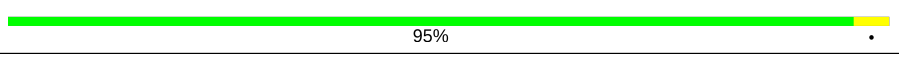
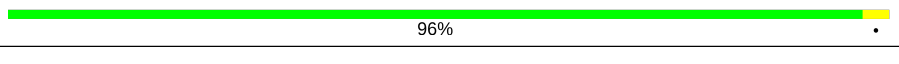
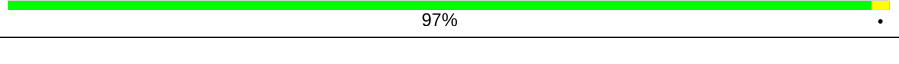
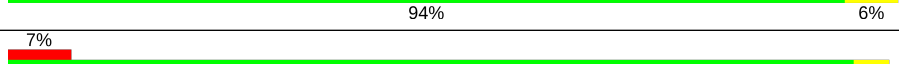
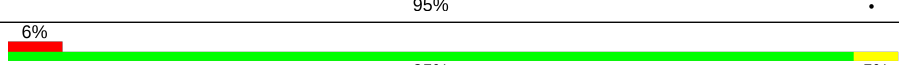
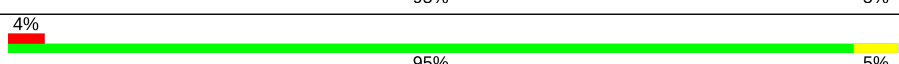
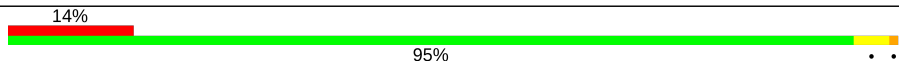






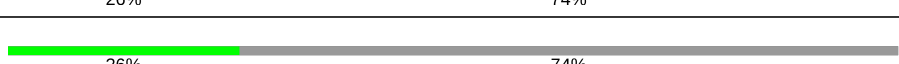

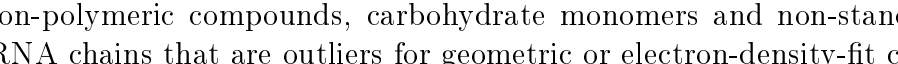
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 on 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	218	97%
1	C	218	95% 5%
1	F	218	92% 6%
1	I	218	97%
1	L	218	95% 7%
1	O	218	95% 2%

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Mol	Chain	Length	Quality of chain
1	R	218	
1	U	218	
2	B	218	
2	D	218	
2	G	218	
2	J	218	
2	M	218	
2	P	218	
2	S	218	
2	V	218	
3	E	46	
3	H	46	
3	K	46	
3	N	46	
3	Q	46	
3	T	46	
3	W	46	
3	X	46	

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
3	HYP	T	21	-	-	-	X
3	HYP	W	21	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 27296 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 ACC1 antibody Fab fragment, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	218	Total 1640	C 1034	N 273	O 325	S 8	0	0	0
1	C	218	Total 1640	C 1034	N 273	O 325	S 8	0	0	0
1	F	213	Total 1612	C 1019	N 268	O 317	S 8	0	0	0
1	I	218	Total 1640	C 1034	N 273	O 325	S 8	0	0	0
1	L	216	Total 1632	C 1030	N 271	O 323	S 8	0	0	0
1	O	216	Total 1629	C 1028	N 271	O 322	S 8	0	0	0
1	R	215	Total 1625	C 1026	N 270	O 321	S 8	0	0	0
1	U	212	Total 1608	C 1017	N 267	O 317	S 7	0	0	0

- Molecule 2 is a protein called ACC1 antibody Fab fragment, light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	217	Total 1660	C 1035	N 281	O 338	S 6	0	0	0
2	D	217	Total 1660	C 1035	N 281	O 338	S 6	0	0	0
2	G	218	Total 1666	C 1038	N 282	O 339	S 7	0	0	0
2	J	218	Total 1666	C 1038	N 282	O 339	S 7	0	0	0
2	M	217	Total 1660	C 1035	N 281	O 338	S 6	0	0	0
2	P	218	Total 1666	C 1038	N 282	O 339	S 7	0	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	S	218	Total	C	N	O	S	0	0	0
			1666	1038	282	339	7			
2	V	218	Total	C	N	O	S	0	0	0
			1666	1038	282	339	7			

- Molecule 3 is a protein called synthetic peptide containing the citrullinated collagen type II epitope CII616-639, Collagen alpha-1(II) chain, synthetic peptide containing the citrullinated collagen type II epitope CII616-639.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	X	12	Total	C	N	O	0	0	0
			82	48	15	19			
3	E	12	Total	C	N	O	0	0	0
			82	48	15	19			
3	H	13	Total	C	N	O	0	0	0
			87	51	16	20			
3	K	13	Total	C	N	O	0	0	0
			87	51	16	20			
3	N	12	Total	C	N	O	0	0	0
			82	48	15	19			
3	Q	12	Total	C	N	O	0	0	0
			82	48	15	19			
3	T	12	Total	C	N	O	0	0	0
			82	48	15	19			
3	W	12	Total	C	N	O	0	0	0
			82	48	15	19			

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	X	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	H	1	Total	O	S	0	0
			5	4	1		
4	I	1	Total	O	S	0	0
			5	4	1		
4	L	1	Total	O	S	0	0
			5	4	1		
4	Q	1	Total	O	S	0	0
			5	4	1		
4	T	1	Total	O	S	0	0
			5	4	1		
4	W	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	23	Total	O	0	0
			23	23		
5	B	34	Total	O	0	0
			34	34		
5	C	15	Total	O	0	0
			15	15		
5	D	35	Total	O	0	0
			35	35		

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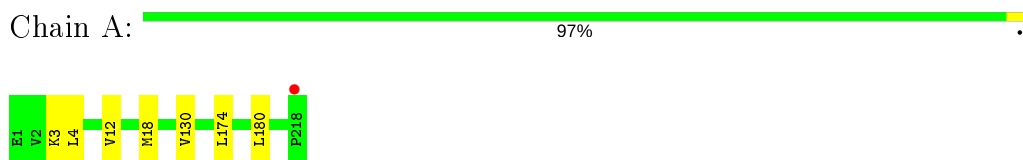
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	F	18	Total O 18 18	0	0
5	G	22	Total O 22 22	0	0
5	H	3	Total O 3 3	0	0
5	I	21	Total O 21 21	0	0
5	J	15	Total O 15 15	0	0
5	K	4	Total O 4 4	0	0
5	L	13	Total O 13 13	0	0
5	M	4	Total O 4 4	0	0
5	O	11	Total O 11 11	0	0
5	P	8	Total O 8 8	0	0
5	Q	1	Total O 1 1	0	0
5	R	8	Total O 8 8	0	0
5	S	10	Total O 10 10	0	0
5	U	3	Total O 3 3	0	0
5	V	6	Total O 6 6	0	0

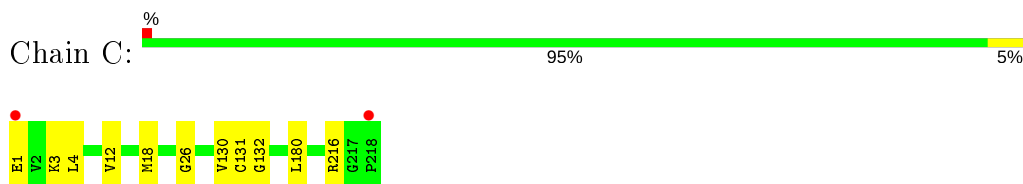
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

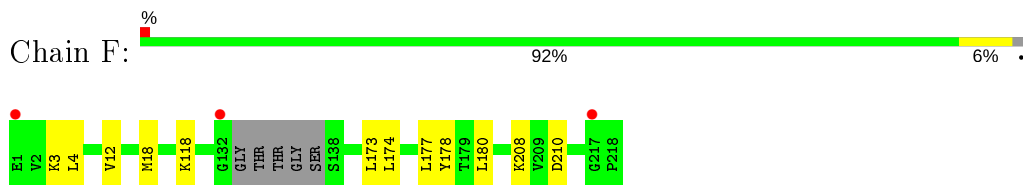
- Molecule 1: ACC1 antibody Fab fragment, heavy chain



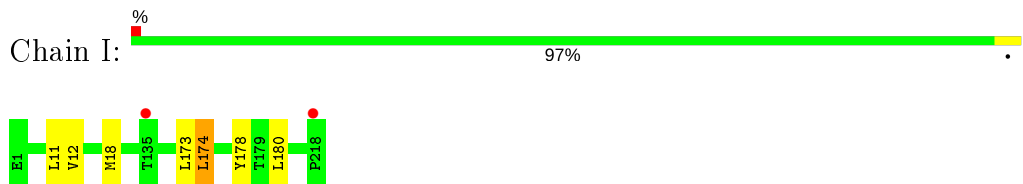
- Molecule 1: ACC1 antibody Fab fragment, heavy chain



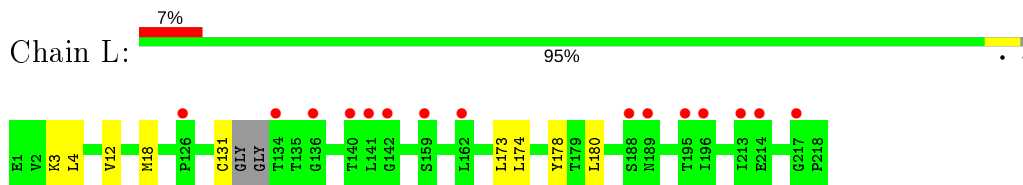
- Molecule 1: ACC1 antibody Fab fragment, heavy chain



- Molecule 1: ACC1 antibody Fab fragment, heavy chain



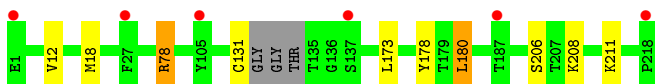
- Molecule 1: ACC1 antibody Fab fragment, heavy chain



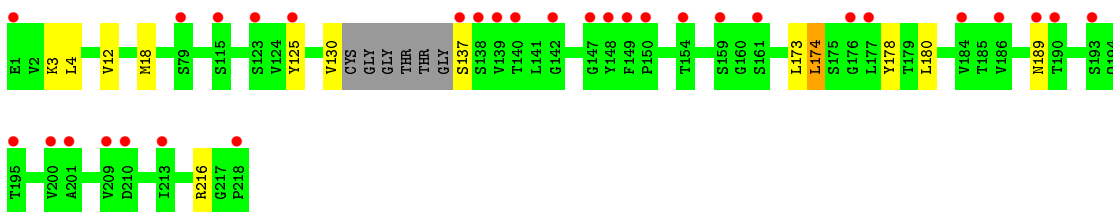
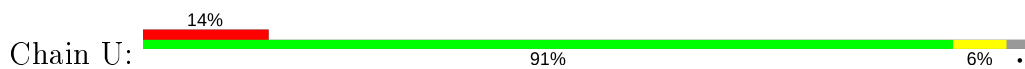
- Molecule 1: ACC1 antibody Fab fragment, heavy chain



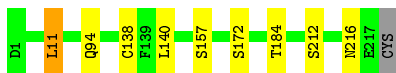
- Molecule 1: ACC1 antibody Fab fragment, heavy chain



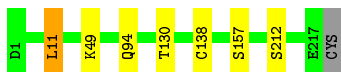
- Molecule 1: ACC1 antibody Fab fragment, heavy chain



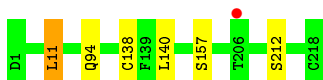
- Molecule 2: ACC1 antibody Fab fragment, light chain



- Molecule 2: ACC1 antibody Fab fragment, light chain



- Molecule 2: ACC1 antibody Fab fragment, light chain



- Molecule 2: ACC1 antibody Fab fragment, light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	119.24Å 91.60Å 182.36Å 90.00° 91.52° 90.00°	Depositor
Resolution (Å)	32.80 – 2.90 32.80 – 2.90	Depositor EDS
% Data completeness (in resolution range)	96.8 (32.80-2.90) 96.9 (32.80-2.90)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 2.90Å)	Xtrriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.232 , 0.247 0.235 , 0.248	Depositor DCC
R_{free} test set	4191 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	38.7	Xtrriage
Anisotropy	0.627	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 36.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	0.036 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	27296	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 42.04 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.1622e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: HYP, SO4

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	0.50	0/1681	0.71	1/2295 (0.0%)
1	C	0.51	0/1681	0.72	1/2295 (0.0%)
1	F	0.53	0/1652	0.72	2/2254 (0.1%)
1	I	0.54	0/1681	0.74	2/2295 (0.1%)
1	L	0.53	0/1672	0.72	1/2282 (0.0%)
1	O	0.50	0/1669	0.70	0/2277
1	R	0.50	0/1665	0.76	4/2272 (0.2%)
1	U	0.47	0/1648	0.70	2/2249 (0.1%)
2	B	0.53	0/1697	0.71	1/2302 (0.0%)
2	D	0.51	0/1697	0.70	1/2302 (0.0%)
2	G	0.50	0/1703	0.70	1/2310 (0.0%)
2	J	0.53	0/1703	0.69	0/2310
2	M	0.55	0/1697	0.72	1/2302 (0.0%)
2	P	0.51	0/1703	0.68	1/2310 (0.0%)
2	S	0.48	0/1703	0.68	1/2310 (0.0%)
2	V	0.47	0/1703	0.70	3/2310 (0.1%)
3	E	0.85	0/66	1.17	0/86
3	H	0.77	0/71	1.01	0/93
3	K	0.77	0/71	1.11	0/93
3	N	0.81	0/66	1.22	1/86 (1.2%)
3	Q	0.80	0/66	1.13	1/86 (1.2%)
3	T	0.67	0/66	1.03	1/86 (1.2%)
3	W	0.72	0/66	1.06	0/86
3	X	0.75	0/66	1.04	0/86
All	All	0.52	0/27493	0.72	25/37377 (0.1%)

There are no bond length outliers.

All (25) bond angle outliers are listed below:

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	R	180	LEU	CB-CG-CD2	6.91	122.74	111.00
2	V	192	ARG	NE-CZ-NH1	6.78	123.69	120.30
2	V	218	CYS	CA-CB-SG	6.21	125.17	114.00
2	D	11	LEU	CA-CB-CG	6.20	129.57	115.30
2	S	11	LEU	CA-CB-CG	6.17	129.50	115.30
1	R	78	ARG	NE-CZ-NH2	6.16	123.38	120.30
2	G	11	LEU	CA-CB-CG	6.05	129.21	115.30
2	B	11	LEU	CA-CB-CG	5.94	128.96	115.30
1	R	131	CYS	CA-CB-SG	5.94	124.69	114.00
2	V	11	LEU	CA-CB-CG	5.94	128.96	115.30
3	N	24	ARG	NE-CZ-NH1	5.92	123.26	120.30
2	P	11	LEU	CA-CB-CG	5.84	128.72	115.30
1	F	177	LEU	CB-CG-CD1	5.80	120.86	111.00
1	U	216	ARG	NE-CZ-NH2	5.79	123.19	120.30
1	I	11	LEU	CB-CG-CD2	-5.54	101.57	111.00
3	Q	24	ARG	NE-CZ-NH1	5.53	123.07	120.30
1	R	180	LEU	CA-CB-CG	5.51	127.97	115.30
2	M	192	ARG	NE-CZ-NH1	5.41	123.00	120.30
1	C	216	ARG	NE-CZ-NH2	5.34	122.97	120.30
1	F	174	LEU	CB-CG-CD2	5.26	119.95	111.00
1	A	174	LEU	CA-CB-CG	5.24	127.35	115.30
1	L	174	LEU	CA-CB-CG	5.21	127.28	115.30
3	T	24	ARG	NE-CZ-NH1	5.17	122.89	120.30
1	U	174	LEU	CA-CB-CG	5.14	127.13	115.30
1	I	174	LEU	CA-CB-CG	5.13	127.09	115.30

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	1640	0	1601	2	0
1	C	1640	0	1601	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	1612	0	1574	5	0
1	I	1640	0	1600	3	0
1	L	1632	0	1594	3	0
1	O	1629	0	1589	5	0
1	R	1625	0	1586	8	0
1	U	1608	0	1572	12	0
2	B	1660	0	1593	2	0
2	D	1660	0	1593	1	0
2	G	1666	0	1597	1	0
2	J	1666	0	1597	6	0
2	M	1660	0	1593	6	0
2	P	1666	0	1597	4	0
2	S	1666	0	1597	6	0
2	V	1666	0	1598	5	0
3	E	82	0	73	0	0
3	H	87	0	77	0	0
3	K	87	0	77	0	0
3	N	82	0	72	0	0
3	Q	82	0	72	0	0
3	T	82	0	72	0	0
3	W	82	0	72	0	0
3	X	82	0	72	0	0
4	E	5	0	0	0	0
4	H	5	0	0	0	0
4	I	5	0	0	0	0
4	L	5	0	0	0	0
4	Q	5	0	0	0	0
4	T	5	0	0	0	0
4	W	5	0	0	0	0
4	X	5	0	0	0	0
5	A	23	0	0	0	0
5	B	34	0	0	0	0
5	C	15	0	0	0	0
5	D	35	0	0	0	0
5	F	18	0	0	0	0
5	G	22	0	0	0	0
5	H	3	0	0	0	0
5	I	21	0	0	0	0
5	J	15	0	0	0	0
5	K	4	0	0	0	0
5	L	13	0	0	0	0
5	M	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	O	11	0	0	0	0
5	P	8	0	0	0	0
5	Q	1	0	0	0	0
5	R	8	0	0	0	0
5	S	10	0	0	0	0
5	U	3	0	0	0	0
5	V	6	0	0	0	0
All	All	27296	0	26069	55	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 1.

All (55) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:3:LYS:HG3	2:M:206:THR:HB	1.67	0.76
1:I:174:LEU:HD13	2:J:164:LEU:HD11	1.67	0.76
2:J:206:THR:HG23	1:R:208:LYS:HB2	1.74	0.70
2:P:194:ASN:OD1	2:P:215:ARG:HG2	1.95	0.66
1:R:211:LYS:NZ	2:S:127:GLU:OE2	2.25	0.66
2:P:193:HIS:O	2:P:215:ARG:HD2	1.98	0.63
1:U:130:VAL:C	2:V:218:CYS:SG	2.78	0.61
1:C:3:LYS:CG	2:M:206:THR:HB	2.32	0.60
1:C:3:LYS:HG3	2:M:206:THR:CB	2.34	0.57
2:D:130:THR:O	1:O:27:PHE:HA	2.05	0.57
1:I:12:VAL:HB	1:I:18:MET:CE	2.37	0.55
1:F:12:VAL:HB	1:F:18:MET:CE	2.37	0.55
1:U:174:LEU:HD13	2:V:164:LEU:HD11	1.89	0.54
1:R:12:VAL:HB	1:R:18:MET:CE	2.38	0.54
1:U:12:VAL:HB	1:U:18:MET:CE	2.39	0.53
2:S:58:GLN:OE1	1:U:189:ASN:HB2	2.08	0.53
1:O:12:VAL:HB	1:O:18:MET:CE	2.39	0.52
1:F:208:LYS:NZ	1:F:210:ASP:OD1	2.43	0.52
1:C:12:VAL:HB	1:C:18:MET:CE	2.39	0.52
1:A:12:VAL:HB	1:A:18:MET:CE	2.39	0.51
1:L:12:VAL:HB	1:L:18:MET:CE	2.40	0.51
2:J:11:LEU:HD12	2:J:13:VAL:HG23	1.92	0.51
1:C:130:VAL:HG12	1:C:132:GLY:H	1.78	0.48
1:F:173:LEU:HD13	1:F:178:TYR:CZ	2.51	0.46
2:J:217:GLU:CG	2:J:218:CYS:N	2.78	0.46
1:U:125:TYR:HB3	2:V:125:SER:OG	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:173:LEU:HD13	1:L:178:TYR:CZ	2.51	0.45
2:M:11:LEU:HD12	2:M:13:VAL:HG23	1.98	0.45
1:O:130:VAL:HG21	2:P:213:PHE:HB3	1.98	0.45
1:R:173:LEU:HD13	1:R:178:TYR:CZ	2.52	0.44
1:U:173:LEU:HD13	1:U:178:TYR:CZ	2.53	0.44
1:U:125:TYR:CE1	2:V:128:GLN:HA	2.53	0.44
1:C:1:GLU:H2	1:R:78:ARG:HG3	1.83	0.44
1:I:173:LEU:HD13	1:I:178:TYR:CZ	2.54	0.43
1:O:3:LYS:C	1:O:4:LEU:HD12	2.40	0.42
2:S:61:GLY:CA	1:U:137:SER:HB3	2.49	0.42
1:L:3:LYS:C	1:L:4:LEU:HD12	2.40	0.42
1:C:3:LYS:C	1:C:4:LEU:HD12	2.40	0.42
1:O:131:CYS:HB3	2:P:218:CYS:HA	2.02	0.42
1:C:3:LYS:HB2	2:M:206:THR:CG2	2.49	0.42
2:B:216:ASN:HB2	1:F:118:LYS:HB2	2.03	0.41
1:C:26:GLY:HA2	1:R:78:ARG:HG2	2.00	0.41
2:S:61:GLY:HA2	1:U:137:SER:HB3	2.02	0.41
2:J:140:LEU:N	2:J:140:LEU:HD12	2.36	0.41
1:U:3:LYS:C	1:U:4:LEU:HD12	2.40	0.41
1:F:3:LYS:C	1:F:4:LEU:HD12	2.41	0.41
2:S:61:GLY:C	1:U:137:SER:HB3	2.41	0.41
1:A:3:LYS:C	1:A:4:LEU:HD12	2.41	0.41
2:M:11:LEU:HD12	2:M:13:VAL:CG2	2.51	0.40
1:U:130:VAL:HG22	2:V:123:PRO:HD3	2.03	0.40
2:B:140:LEU:N	2:B:140:LEU:HD12	2.36	0.40
2:J:206:THR:HG21	1:R:206:SER:O	2.21	0.40
2:G:140:LEU:N	2:G:140:LEU:HD12	2.37	0.40
1:C:1:GLU:N	1:R:78:ARG:HG3	2.36	0.40
2:S:140:LEU:HD12	2:S:140:LEU:N	2.36	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	216/218 (99%)	210 (97%)	6 (3%)	0	100	100
1	C	216/218 (99%)	210 (97%)	6 (3%)	0	100	100
1	F	209/218 (96%)	205 (98%)	4 (2%)	0	100	100
1	I	216/218 (99%)	211 (98%)	5 (2%)	0	100	100
1	L	212/218 (97%)	207 (98%)	5 (2%)	0	100	100
1	O	212/218 (97%)	207 (98%)	5 (2%)	0	100	100
1	R	211/218 (97%)	207 (98%)	4 (2%)	0	100	100
1	U	208/218 (95%)	203 (98%)	5 (2%)	0	100	100
2	B	215/218 (99%)	210 (98%)	5 (2%)	0	100	100
2	D	215/218 (99%)	211 (98%)	4 (2%)	0	100	100
2	G	216/218 (99%)	212 (98%)	4 (2%)	0	100	100
2	J	216/218 (99%)	212 (98%)	4 (2%)	0	100	100
2	M	215/218 (99%)	211 (98%)	4 (2%)	0	100	100
2	P	216/218 (99%)	212 (98%)	4 (2%)	0	100	100
2	S	216/218 (99%)	212 (98%)	4 (2%)	0	100	100
2	V	216/218 (99%)	211 (98%)	5 (2%)	0	100	100
3	E	9/46 (20%)	8 (89%)	1 (11%)	0	100	100
3	H	10/46 (22%)	9 (90%)	1 (10%)	0	100	100
3	K	10/46 (22%)	9 (90%)	1 (10%)	0	100	100
3	N	9/46 (20%)	8 (89%)	1 (11%)	0	100	100
3	Q	9/46 (20%)	8 (89%)	1 (11%)	0	100	100
3	T	9/46 (20%)	8 (89%)	1 (11%)	0	100	100
3	W	9/46 (20%)	8 (89%)	1 (11%)	0	100	100
3	X	9/46 (20%)	8 (89%)	1 (11%)	0	100	100
All	All	3499/3856 (91%)	3417 (98%)	82 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	185/185 (100%)	183 (99%)	2 (1%)	73	92
1	C	185/185 (100%)	183 (99%)	2 (1%)	73	92
1	F	182/185 (98%)	181 (100%)	1 (0%)	88	96
1	I	185/185 (100%)	184 (100%)	1 (0%)	88	96
1	L	185/185 (100%)	183 (99%)	2 (1%)	73	92
1	O	184/185 (100%)	181 (98%)	3 (2%)	62	86
1	R	184/185 (100%)	183 (100%)	1 (0%)	88	96
1	U	182/185 (98%)	181 (100%)	1 (0%)	88	96
2	B	187/188 (100%)	180 (96%)	7 (4%)	34	68
2	D	187/188 (100%)	181 (97%)	6 (3%)	39	73
2	G	188/188 (100%)	183 (97%)	5 (3%)	44	77
2	J	188/188 (100%)	182 (97%)	6 (3%)	39	73
2	M	187/188 (100%)	182 (97%)	5 (3%)	44	77
2	P	188/188 (100%)	182 (97%)	6 (3%)	39	73
2	S	188/188 (100%)	181 (96%)	7 (4%)	34	68
2	V	188/188 (100%)	181 (96%)	7 (4%)	34	68
3	E	6/15 (40%)	6 (100%)	0	100	100
3	H	6/15 (40%)	6 (100%)	0	100	100
3	K	6/15 (40%)	6 (100%)	0	100	100
3	N	6/15 (40%)	6 (100%)	0	100	100
3	Q	6/15 (40%)	6 (100%)	0	100	100
3	T	6/15 (40%)	6 (100%)	0	100	100
3	W	6/15 (40%)	6 (100%)	0	100	100
3	X	6/15 (40%)	6 (100%)	0	100	100
All	All	3021/3104 (97%)	2959 (98%)	62 (2%)	53	81

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	130	VAL
1	A	180	LEU
2	B	11	LEU
2	B	94	GLN

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Mol	Chain	Res	Type
2	B	138	CYS
2	B	157	SER
2	B	172	SER
2	B	184	THR
2	B	212	SER
1	C	131	CYS
1	C	180	LEU
2	D	11	LEU
2	D	49	LYS
2	D	94	GLN
2	D	138	CYS
2	D	157	SER
2	D	212	SER
1	F	180	LEU
2	G	11	LEU
2	G	94	GLN
2	G	138	CYS
2	G	157	SER
2	G	212	SER
1	I	180	LEU
2	J	49	LYS
2	J	94	GLN
2	J	138	CYS
2	J	147	ASP
2	J	157	SER
2	J	212	SER
1	L	131	CYS
1	L	180	LEU
2	M	49	LYS
2	M	94	GLN
2	M	138	CYS
2	M	157	SER
2	M	212	SER
1	O	130	VAL
1	O	131	CYS
1	O	180	LEU
2	P	11	LEU
2	P	49	LYS
2	P	94	GLN
2	P	138	CYS
2	P	157	SER
2	P	212	SER

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Mol	Chain	Res	Type
1	R	180	LEU
2	S	11	LEU
2	S	49	LYS
2	S	94	GLN
2	S	126	SER
2	S	138	CYS
2	S	157	SER
2	S	212	SER
1	U	180	LEU
2	V	11	LEU
2	V	94	GLN
2	V	138	CYS
2	V	157	SER
2	V	192	ARG
2	V	212	SER
2	V	218	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	194	GLN
2	B	161	ASN
2	D	161	ASN
2	G	161	ASN
2	J	161	ASN
2	M	149	ASN
2	M	161	ASN
2	P	161	ASN
2	S	161	ASN
2	V	161	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

16 non-standard protein/DNA/RNA residues 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	HYP	H	30	3	6,8,9	0.56	0	5,10,12	2.00	2 (40%)
3	HYP	N	30	3	6,8,9	0.59	0	5,10,12	2.05	2 (40%)
3	HYP	X	21	3	6,8,9	0.78	0	5,10,12	2.61	2 (40%)
3	HYP	E	30	3	6,8,9	1.14	0	5,10,12	1.54	1 (20%)
3	HYP	T	30	3	6,8,9	0.57	0	5,10,12	2.14	2 (40%)
3	HYP	K	30	3	6,8,9	0.76	0	5,10,12	1.91	2 (40%)
3	HYP	T	21	3	6,8,9	0.71	0	5,10,12	2.56	1 (20%)
3	HYP	W	21	3	6,8,9	0.79	0	5,10,12	2.85	2 (40%)
3	HYP	X	30	3	6,8,9	0.85	0	5,10,12	1.99	2 (40%)
3	HYP	Q	21	3	6,8,9	0.74	0	5,10,12	2.69	1 (20%)
3	HYP	Q	30	3	6,8,9	0.58	0	5,10,12	2.04	2 (40%)
3	HYP	N	21	3	6,8,9	0.56	0	5,10,12	2.70	1 (20%)
3	HYP	W	30	3	6,8,9	0.46	0	5,10,12	2.04	2 (40%)
3	HYP	H	21	3	6,8,9	0.62	0	5,10,12	2.68	1 (20%)
3	HYP	K	21	3	6,8,9	0.62	0	5,10,12	2.63	3 (60%)
3	HYP	E	21	3	6,8,9	0.64	0	5,10,12	2.58	2 (40%)

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	HYP	H	30	3	-	0/0/11/13	0/1/1/1
3	HYP	N	30	3	-	0/0/11/13	0/1/1/1
3	HYP	X	21	3	-	0/0/11/13	0/1/1/1
3	HYP	E	30	3	-	0/0/11/13	0/1/1/1
3	HYP	T	30	3	-	0/0/11/13	0/1/1/1
3	HYP	K	30	3	-	0/0/11/13	0/1/1/1
3	HYP	T	21	3	-	0/0/11/13	0/1/1/1
3	HYP	W	21	3	-	0/0/11/13	0/1/1/1
3	HYP	X	30	3	-	0/0/11/13	0/1/1/1
3	HYP	Q	21	3	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	HYP	Q	30	3	-	0/0/11/13	0/1/1/1
3	HYP	N	21	3	-	0/0/11/13	0/1/1/1
3	HYP	W	30	3	-	0/0/11/13	0/1/1/1
3	HYP	H	21	3	-	0/0/11/13	0/1/1/1
3	HYP	K	21	3	-	0/0/11/13	0/1/1/1
3	HYP	E	21	3	-	0/0/11/13	0/1/1/1

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	W	21	HYP	CB-CG-CD	5.83	110.42	103.27
3	Q	21	HYP	CB-CG-CD	5.71	110.27	103.27
3	N	21	HYP	CB-CG-CD	5.66	110.21	103.27
3	H	21	HYP	CB-CG-CD	5.49	110.00	103.27
3	T	21	HYP	CB-CG-CD	5.16	109.60	103.27
3	X	21	HYP	CB-CG-CD	5.07	109.48	103.27
3	K	21	HYP	CB-CG-CD	4.99	109.38	103.27
3	E	21	HYP	CB-CG-CD	4.89	109.27	103.27
3	T	30	HYP	CB-CG-CD	3.90	108.05	103.27
3	N	30	HYP	CB-CG-CD	3.80	107.93	103.27
3	H	30	HYP	CB-CG-CD	3.78	107.91	103.27
3	Q	30	HYP	CB-CG-CD	3.76	107.87	103.27
3	W	30	HYP	CB-CG-CD	3.75	107.87	103.27
3	X	30	HYP	CB-CG-CD	3.61	107.69	103.27
3	K	30	HYP	CB-CG-CD	3.46	107.50	103.27
3	E	30	HYP	CB-CG-CD	2.58	106.43	103.27
3	T	30	HYP	O-C-CA	-2.51	118.20	124.78
3	Q	30	HYP	O-C-CA	-2.43	118.40	124.78
3	X	30	HYP	O-C-CA	-2.40	118.49	124.78
3	K	30	HYP	O-C-CA	-2.40	118.49	124.78
3	N	30	HYP	O-C-CA	-2.39	118.51	124.78
3	W	30	HYP	O-C-CA	-2.34	118.63	124.78
3	W	21	HYP	O-C-CA	-2.15	119.13	124.78
3	E	21	HYP	O-C-CA	-2.14	119.17	124.78
3	K	21	HYP	O-C-CA	-2.12	119.22	124.78
3	H	30	HYP	O-C-CA	-2.12	119.22	124.78
3	X	21	HYP	O-C-CA	-2.06	119.38	124.78
3	K	21	HYP	CG-CB-CA	2.00	106.49	103.96

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

8 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	$\# Z > 2$
4	SO4	Q	101	-	4,4,4	0.33	0	6,6,6	0.27	0
4	SO4	X	101	-	4,4,4	0.32	0	6,6,6	0.38	0
4	SO4	T	101	-	4,4,4	0.38	0	6,6,6	0.31	0
4	SO4	E	101	-	4,4,4	0.33	0	6,6,6	0.13	0
4	SO4	H	101	-	4,4,4	0.35	0	6,6,6	0.28	0
4	SO4	L	301	-	4,4,4	0.38	0	6,6,6	0.17	0
4	SO4	I	301	-	4,4,4	0.39	0	6,6,6	0.16	0
4	SO4	W	101	-	4,4,4	0.34	0	6,6,6	0.21	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data i

6.1 Protein, DNA and RNA chains i

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	218/218 (100%)	-0.06	1 (0%) 91 91	24, 42, 69, 79	0
1	C	218/218 (100%)	-0.12	2 (0%) 84 84	30, 43, 67, 80	0
1	F	213/218 (97%)	0.00	3 (1%) 75 75	27, 45, 67, 93	0
1	I	218/218 (100%)	-0.13	2 (0%) 84 84	27, 40, 59, 79	0
1	L	216/218 (99%)	0.35	15 (6%) 16 13	27, 49, 97, 116	0
1	O	216/218 (99%)	0.12	4 (1%) 66 65	28, 47, 85, 108	0
1	R	215/218 (98%)	0.07	6 (2%) 53 49	34, 49, 73, 95	0
1	U	212/218 (97%)	0.79	31 (14%) 2 1	41, 76, 110, 127	0
2	B	217/218 (99%)	-0.23	0 100 100	22, 39, 52, 57	0
2	D	217/218 (99%)	-0.27	0 100 100	24, 38, 51, 59	0
2	G	218/218 (100%)	0.15	1 (0%) 91 91	25, 50, 86, 101	0
2	J	218/218 (100%)	-0.02	0 100 100	28, 42, 58, 73	0
2	M	217/218 (99%)	0.49	16 (7%) 14 11	27, 59, 106, 113	0
2	P	218/218 (100%)	0.42	14 (6%) 19 15	29, 63, 110, 119	0
2	S	218/218 (100%)	0.29	9 (4%) 37 32	32, 56, 87, 97	0
2	V	218/218 (100%)	0.81	31 (14%) 2 2	32, 74, 120, 144	0
3	E	10/46 (21%)	0.27	0 100 100	39, 45, 56, 58	0
3	H	11/46 (23%)	0.27	1 (9%) 9 6	48, 52, 55, 55	0
3	K	11/46 (23%)	0.51	1 (9%) 9 6	43, 49, 56, 62	0
3	N	10/46 (21%)	0.56	2 (20%) 1 0	48, 54, 68, 75	0
3	Q	10/46 (21%)	0.65	1 (10%) 7 5	73, 78, 96, 106	0
3	T	10/46 (21%)	0.64	1 (10%) 7 5	55, 58, 66, 69	0
3	W	10/46 (21%)	0.76	2 (20%) 1 0	66, 70, 80, 80	0
3	X	10/46 (21%)	0.48	0 100 100	37, 42, 52, 61	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
All	All	3549/3856 (92%)	0.17	143 (4%) 38 33	22, 47, 98, 144	0

All (143) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	U	139	VAL	6.1
2	V	159	ARG	5.2
1	U	190	THR	5.1
2	P	210	VAL	4.9
2	V	218	CYS	4.5
1	L	196	ILE	4.3
1	L	141	LEU	4.3
1	U	140	THR	4.2
1	U	193	SER	4.2
1	U	218	PRO	4.2
2	V	216	ASN	4.1
2	M	196	TYR	3.8
1	C	218	PRO	3.8
1	U	189	ASN	3.8
2	S	218	CYS	3.7
1	L	134	THR	3.7
1	L	195	THR	3.7
2	V	188	ASP	3.7
1	U	1	GLU	3.6
2	M	214	ASN	3.5
1	L	188	SER	3.5
1	U	209	VAL	3.4
2	V	138	CYS	3.3
2	V	154	ILE	3.3
3	N	32	PRO	3.2
2	M	32	TYR	3.1
1	U	201	ALA	3.1
3	W	32	PRO	3.1
1	U	154	THR	3.0
2	V	196	TYR	3.0
1	I	218	PRO	3.0
1	A	218	PRO	3.0
1	L	142	GLY	2.9
2	S	158	GLU	2.9
1	L	140	THR	2.9
1	U	138	SER	2.9
2	M	157	SER	2.9

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Mol	Chain	Res	Type	RSRZ
2	V	158	GLU	2.9
2	P	154	ILE	2.9
1	C	1	GLU	2.9
1	U	200	VAL	2.9
1	L	214	GLU	2.9
2	V	160	GLN	2.8
1	L	162	LEU	2.8
1	U	159	SER	2.8
2	V	125	SER	2.8
1	U	150	PRO	2.8
2	V	198	CYS	2.8
2	M	206	THR	2.8
2	M	185	LEU	2.8
3	H	33	ALA	2.8
2	P	192	ARG	2.8
3	Q	32	PRO	2.8
1	U	213	ILE	2.7
1	U	161	SER	2.7
2	V	131	SER	2.7
1	U	148	TYR	2.7
1	L	189	ASN	2.7
3	T	32	PRO	2.7
2	M	207	SER	2.7
2	V	185	LEU	2.7
1	F	1	GLU	2.7
2	P	208	PRO	2.7
1	R	1	GLU	2.6
2	S	197	THR	2.6
1	U	176	GLY	2.6
2	P	162	GLY	2.6
2	M	134	ALA	2.5
1	U	210	ASP	2.5
1	L	126	PRO	2.5
2	S	217	GLU	2.5
2	V	186	THR	2.5
1	U	137	SER	2.5
2	P	207	SER	2.5
1	O	218	PRO	2.5
2	V	153	LYS	2.5
1	U	149	PHE	2.5
2	V	202	HIS	2.5
1	F	132	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
2	V	217	GLU	2.5
2	P	216	ASN	2.5
2	P	32	TYR	2.4
2	M	126	SER	2.4
2	V	192	ARG	2.4
1	O	189	ASN	2.4
2	V	130	THR	2.4
2	V	152	TRP	2.4
2	V	197	THR	2.4
2	V	213	PHE	2.4
1	U	79	SER	2.4
1	F	217	GLY	2.4
1	L	159	SER	2.4
1	R	27	PHE	2.4
2	M	188	ASP	2.3
1	O	195	THR	2.3
1	R	137	SER	2.3
2	V	183	LEU	2.3
3	N	31	GLY	2.3
2	V	137	VAL	2.3
2	V	32	TYR	2.3
1	U	123	SER	2.2
2	G	206	THR	2.2
2	V	126	SER	2.2
2	P	209	ILE	2.2
1	U	186	VAL	2.2
1	R	218	PRO	2.2
2	P	152	TRP	2.2
2	V	214	ASN	2.2
1	U	195	THR	2.2
3	K	33	ALA	2.2
1	U	115	SER	2.2
2	P	159	ARG	2.2
2	M	192	ARG	2.2
1	U	142	GLY	2.2
2	P	218	CYS	2.2
2	V	60	SER	2.1
2	V	120	SER	2.1
2	M	210	VAL	2.1
2	M	34	ILE	2.1
2	M	217	GLU	2.1
1	O	193	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	L	217	GLY	2.1
1	R	187	THR	2.1
2	M	31	ASN	2.1
2	V	31	ASN	2.1
2	S	216	ASN	2.1
2	S	130	THR	2.0
2	S	134	ALA	2.0
2	S	162	GLY	2.0
3	W	22	GLY	2.0
2	P	194	ASN	2.0
1	L	213	ILE	2.0
2	P	24	ARG	2.0
1	L	136	GLY	2.0
1	U	147	GLY	2.0
2	S	173	LYS	2.0
2	V	195	SER	2.0
1	R	105	TYR	2.0
1	I	135	THR	2.0
1	U	184	VAL	2.0
1	U	177	LEU	2.0
2	M	135	SER	2.0
1	U	125	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
3	HYP	X	21	8/9	0.73	0.30	48,49,50,50	0
3	HYP	W	21	8/9	0.73	0.42	77,78,78,79	0
3	HYP	Q	21	8/9	0.73	0.30	83,85,92,92	0
3	HYP	T	21	8/9	0.75	0.40	65,67,67,67	0
3	HYP	K	21	8/9	0.77	0.31	50,52,59,61	0
3	HYP	N	21	8/9	0.81	0.28	57,61,65,65	0
3	HYP	H	21	8/9	0.85	0.30	50,52,56,58	0
3	HYP	W	30	8/9	0.86	0.26	70,72,72,72	0
3	HYP	E	21	8/9	0.87	0.20	50,52,52,53	0
3	HYP	Q	30	8/9	0.90	0.26	85,89,92,93	0
3	HYP	X	30	8/9	0.91	0.18	42,44,47,48	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	HYP	T	30	8/9	0.91	0.23	59,61,61,61	0
3	HYP	N	30	8/9	0.93	0.18	58,61,62,63	0
3	HYP	E	30	8/9	0.94	0.17	45,46,48,49	0
3	HYP	K	30	8/9	0.94	0.18	51,52,60,61	0
3	HYP	H	30	8/9	0.95	0.17	52,53,53,53	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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(\AA^2)	Q<0.9
4	SO4	W	101	5/5	0.91	0.17	83,84,88,89	0
4	SO4	T	101	5/5	0.92	0.20	64,68,69,70	0
4	SO4	L	301	5/5	0.95	0.19	62,65,69,70	0
4	SO4	Q	101	5/5	0.95	0.18	65,65,68,69	0
4	SO4	E	101	5/5	0.96	0.12	61,63,65,65	0
4	SO4	H	101	5/5	0.96	0.17	43,43,46,49	0
4	SO4	I	301	5/5	0.98	0.14	47,49,50,53	0
4	SO4	X	101	5/5	0.98	0.11	40,41,41,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.