



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 06:58 PM GMT

PDB ID : 4N3E
Title : Crystal structure of Hyp-1, a St John's wort PR-10 protein, in complex with 8-anilino-1-naphthalene sulfonate (ANS)
Authors : Sliwiak, J.; Dauter, Z.; Mccoy, A.J.; Read, R.J.; Jaskolski, M.
Deposited on : 2013-10-07
Resolution : 2.43 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

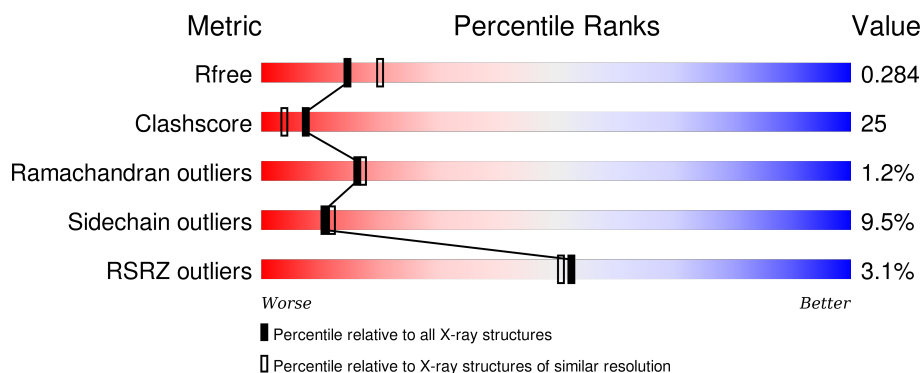
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.43 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}	91344	1003 (2.46-2.42)
Clashscore	102246	1071 (2.46-2.42)
Ramachandran outliers	100387	1065 (2.46-2.42)
Sidechain outliers	100360	1065 (2.46-2.42)
RSRZ outliers	91569	1005 (2.46-2.42)

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. 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	<div> <div>5%</div> <div>38% 50% 8% .</div> </div>
1	B	165	<div> <div>5%</div> <div>47% 42% 8% .</div> </div>
1	C	165	<div> <div>4%</div> <div>54% 35% 8% .</div> </div>
1	D	165	<div> <div>5%</div> <div>61% 28% 6% . .</div> </div>
1	E	165	<div> <div>5%</div> <div>59% 31% 6% .</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	165	
1	G	165	
1	H	165	
1	I	165	
1	J	165	
1	K	165	
1	L	165	
1	M	165	
1	N	165	
1	O	165	
1	P	165	
1	Q	165	
1	R	165	
1	S	165	
1	T	165	
1	U	165	
1	V	165	
1	W	165	
1	X	165	
1	Y	165	
1	Z	165	
1	a	165	
1	b	165	

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
2	2AN	A	201	-	-	X	-
2	2AN	A	203	-	-	X	-
2	2AN	A	204	-	-	X	-
2	2AN	B	201	-	-	X	-
2	2AN	B	202	-	-	X	-
2	2AN	B	203	-	-	X	-
2	2AN	C	202	-	-	X	-
2	2AN	C	203	-	-	X	-
2	2AN	C	204	-	-	X	-
2	2AN	D	202	-	-	X	-
2	2AN	D	203	-	-	X	-
2	2AN	E	202	-	-	X	-
2	2AN	E	203	-	-	X	X
2	2AN	E	204	-	-	X	-
2	2AN	E	205	-	-	X	-
2	2AN	F	201	-	-	X	-
2	2AN	G	202	-	-	X	-
2	2AN	G	204	-	-	X	X
2	2AN	G	205	-	-	X	-
2	2AN	I	201	-	-	X	-
2	2AN	I	203	-	-	X	X
2	2AN	I	205	-	-	X	-
2	2AN	J	201	-	-	-	X
2	2AN	J	202	-	-	X	-
2	2AN	K	202	-	-	X	-
2	2AN	K	203	-	-	X	-
2	2AN	L	202	-	-	X	-
2	2AN	L	204	-	-	X	-
2	2AN	M	202	-	-	X	-
2	2AN	M	203	-	-	X	-
2	2AN	N	202	-	-	X	-
2	2AN	N	203	-	-	X	-
2	2AN	Q	201	-	-	X	-
2	2AN	Q	203	-	-	X	-
2	2AN	R	201	-	-	X	-
2	2AN	R	203	-	-	-	X
2	2AN	S	201	-	-	X	-
2	2AN	U	201	-	-	X	-
2	2AN	U	202	-	-	-	X
2	2AN	W	201	-	-	X	-
2	2AN	W	202	-	-	X	-
2	2AN	W	203	-	-	-	X
2	2AN	W	204	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	2AN	X	201	-	-	-	X
2	2AN	Y	202	-	-	X	-
2	2AN	Y	204	-	-	X	-
2	2AN	Z	202	-	-	X	X
2	2AN	Z	203	-	-	X	-
2	2AN	b	203	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 37252 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 Phenolic oxidative coupling protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	159	Total	C	N	O	S	0	1	0
			1269	820	204	241	4			
1	B	159	Total	C	N	O	S	1	0	0
			1256	813	203	236	4			
1	C	159	Total	C	N	O	S	0	0	0
			1256	813	203	236	4			
1	D	159	Total	C	N	O	S	0	0	0
			1257	814	203	236	4			
1	E	159	Total	C	N	O	S	1	0	0
			1260	815	203	238	4			
1	F	159	Total	C	N	O	S	0	0	0
			1257	814	202	237	4			
1	G	159	Total	C	N	O	S	0	0	0
			1259	815	202	238	4			
1	H	159	Total	C	N	O	S	0	0	0
			1260	815	203	238	4			
1	I	159	Total	C	N	O	S	0	0	0
			1255	813	202	236	4			
1	J	159	Total	C	N	O	S	0	0	0
			1255	812	203	236	4			
1	K	159	Total	C	N	O	S	0	0	0
			1260	815	203	238	4			
1	L	159	Total	C	N	O	S	1	0	0
			1260	815	203	238	4			
1	M	159	Total	C	N	O	S	0	0	0
			1256	813	203	236	4			
1	N	159	Total	C	N	O	S	0	1	0
			1269	820	204	241	4			
1	O	159	Total	C	N	O	S	0	0	0
			1258	814	203	237	4			
1	P	159	Total	C	N	O	S	0	0	0
			1260	815	203	238	4			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	159	Total	C	N	O	S	0	0	0
			1257	812	203	238	4			
1	R	159	Total	C	N	O	S	1	0	0
			1260	815	203	238	4			
1	S	159	Total	C	N	O	S	0	0	0
			1257	812	203	238	4			
1	T	159	Total	C	N	O	S	0	6	0
			1307	844	210	249	4			
1	U	159	Total	C	N	O	S	1	0	0
			1260	815	203	238	4			
1	V	159	Total	C	N	O	S	0	0	0
			1257	814	203	236	4			
1	W	159	Total	C	N	O	S	0	0	0
			1258	814	203	237	4			
1	X	159	Total	C	N	O	S	0	0	0
			1260	815	203	238	4			
1	Y	159	Total	C	N	O	S	1	0	0
			1260	815	203	238	4			
1	Z	159	Total	C	N	O	S	0	0	0
			1256	813	203	236	4			
1	a	159	Total	C	N	O	S	0	1	0
			1270	821	206	239	4			
1	b	159	Total	C	N	O	S	1	0	0
			1254	813	202	235	4			

There are 252 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
A	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
A	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
A	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
A	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
A	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
A	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
A	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
A	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
B	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
B	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
B	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
B	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
B	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
B	0	THR	-	EXPRESSION TAG	UNP Q8H1L1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
B	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
B	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
C	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
C	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
C	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
C	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
C	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
C	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
C	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
C	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
C	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
D	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
D	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
D	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
D	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
D	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
D	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
D	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
D	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
D	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
E	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
E	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
E	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
E	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
E	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
E	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
E	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
E	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
E	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
F	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
F	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
F	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
F	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
F	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
F	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
F	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
F	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
F	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
G	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
G	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
G	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
G	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
G	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
G	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
G	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
G	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
H	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
H	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
H	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
H	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
H	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
H	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
H	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
H	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
H	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
I	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
I	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
I	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
I	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
I	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
I	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
I	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
I	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
I	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
J	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
J	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
J	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
J	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
J	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
J	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
J	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
J	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
J	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
K	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
K	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
K	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
K	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
K	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
K	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
K	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
K	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
K	119	THR	SER	SEE REMARK 999	UNP Q8H1L1

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Chain	Residue	Modelled	Actual	Comment	Reference
L	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
L	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
L	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
L	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
L	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
L	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
L	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
L	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
L	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
M	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
M	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
M	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
M	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
M	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
M	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
M	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
M	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
M	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
N	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
N	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
N	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
N	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
N	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
N	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
N	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
N	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
N	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
O	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
O	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
O	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
O	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
O	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
O	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
O	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
O	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
O	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
P	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
P	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
P	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
P	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
P	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
P	0	THR	-	EXPRESSION TAG	UNP Q8H1L1

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Chain	Residue	Modelled	Actual	Comment	Reference
P	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
P	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
P	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
Q	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
Q	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
Q	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
Q	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
Q	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
Q	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
Q	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
Q	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
Q	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
R	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
R	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
R	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
R	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
R	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
R	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
R	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
R	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
R	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
S	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
S	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
S	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
S	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
S	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
S	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
S	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
S	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
S	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
T	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
T	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
T	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
T	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
T	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
T	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
T	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
T	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
T	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
U	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
U	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
U	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1

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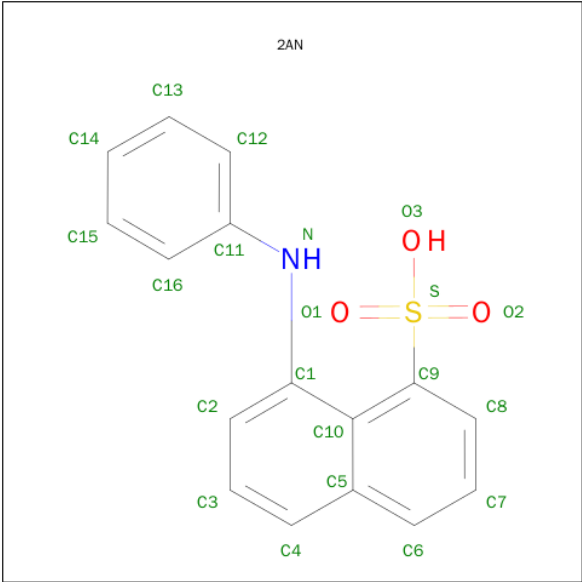
Chain	Residue	Modelled	Actual	Comment	Reference
U	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
U	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
U	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
U	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
U	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
U	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
V	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
V	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
V	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
V	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
V	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
V	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
V	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
V	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
V	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
W	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
W	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
W	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
W	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
W	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
W	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
W	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
W	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
W	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
X	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
X	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
X	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
X	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
X	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
X	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
X	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
X	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
X	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
Y	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
Y	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
Y	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
Y	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
Y	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
Y	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
Y	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
Y	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
Y	119	THR	SER	SEE REMARK 999	UNP Q8H1L1

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Chain	Residue	Modelled	Actual	Comment	Reference
Z	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
Z	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
Z	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
Z	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
Z	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
Z	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
Z	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
Z	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
Z	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
a	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
a	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
a	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
a	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
a	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
a	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
a	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
a	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
a	119	THR	SER	SEE REMARK 999	UNP Q8H1L1
b	-5	GLY	-	EXPRESSION TAG	UNP Q8H1L1
b	-4	ILE	-	EXPRESSION TAG	UNP Q8H1L1
b	-3	ASP	-	EXPRESSION TAG	UNP Q8H1L1
b	-2	PRO	-	EXPRESSION TAG	UNP Q8H1L1
b	-1	PHE	-	EXPRESSION TAG	UNP Q8H1L1
b	0	THR	-	EXPRESSION TAG	UNP Q8H1L1
b	14	ILE	THR	SEE REMARK 999	UNP Q8H1L1
b	86	LEU	ILE	SEE REMARK 999	UNP Q8H1L1
b	119	THR	SER	SEE REMARK 999	UNP Q8H1L1

- Molecule 2 is 8-ANILINO-1-NAPHTHALENE SULFONATE (three-letter code: 2AN) (formula: C₁₆H₁₃NO₃S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	A	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	A	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	A	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	B	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	B	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	B	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	B	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	C	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	C	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	C	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	C	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	D	1	Total	C	N	O	S	0	0
			21	16	1	3	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	D	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	D	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	E	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	E	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	E	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	E	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	E	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	F	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	G	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	G	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	G	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	G	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	G	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	H	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	H	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	H	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	I	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	I	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	I	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	I	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	I	1	Total 21	C 16	N 1	O 3	S 1	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	J	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	J	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	K	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	K	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	K	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	L	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	L	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	L	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	L	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	L	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	M	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	M	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	M	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	N	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	N	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	N	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	N	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	O	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	O	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	P	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	P	1	Total 21	C 16	N 1	O 3	S 1	0	0

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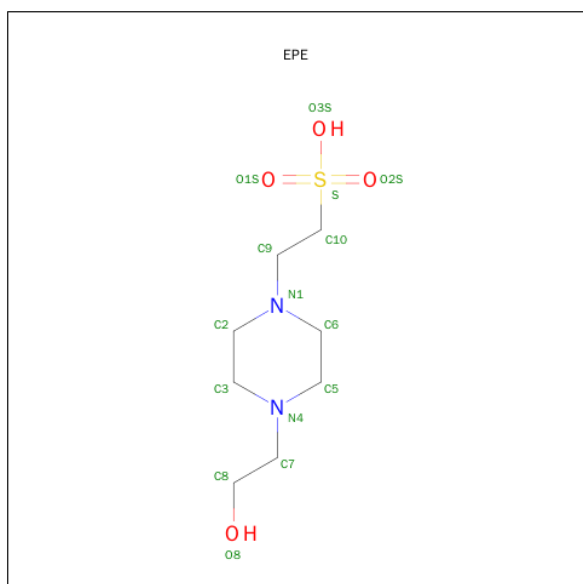
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2	Q	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	R	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	R	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	R	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	R	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	S	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	U	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	U	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	U	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	U	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	U	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	V	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	V	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	W	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	W	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	W	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	W	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	X	1	Total 21	C 16	N 1	O 3	S 1	0	0
2	X	1	Total 21	C 16	N 1	O 3	S 1	0	0

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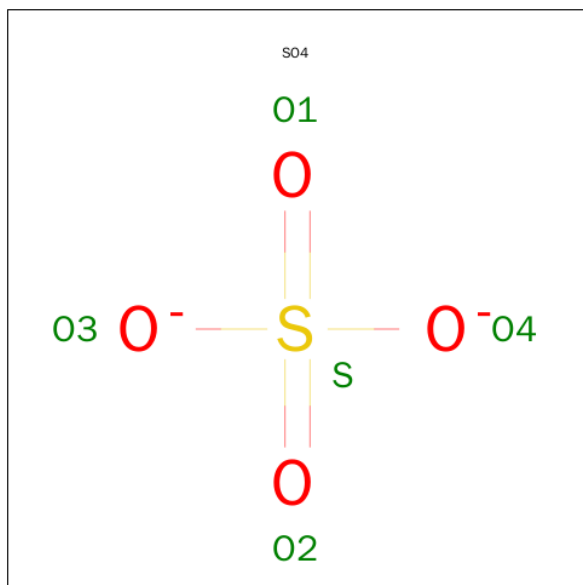
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	X	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	Y	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	Y	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	Y	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	Y	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	Z	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	Z	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	Z	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	a	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	b	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	b	1	Total	C	N	O	S	0	0
			21	16	1	3	1		
2	b	1	Total	C	N	O	S	0	0
			21	16	1	3	1		

- Molecule 3 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	C	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
3	Z	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	V	1	Total	O	S	0	0
			5	4	1		
4	X	1	Total	O	S	0	0
			5	4	1		
4	b	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	5	Total	O	0	0
			5	5		
5	C	2	Total	O	0	0
			2	2		
5	D	1	Total	O	0	0
			1	1		
5	E	2	Total	O	0	0
			2	2		
5	F	2	Total	O	0	0
			2	2		

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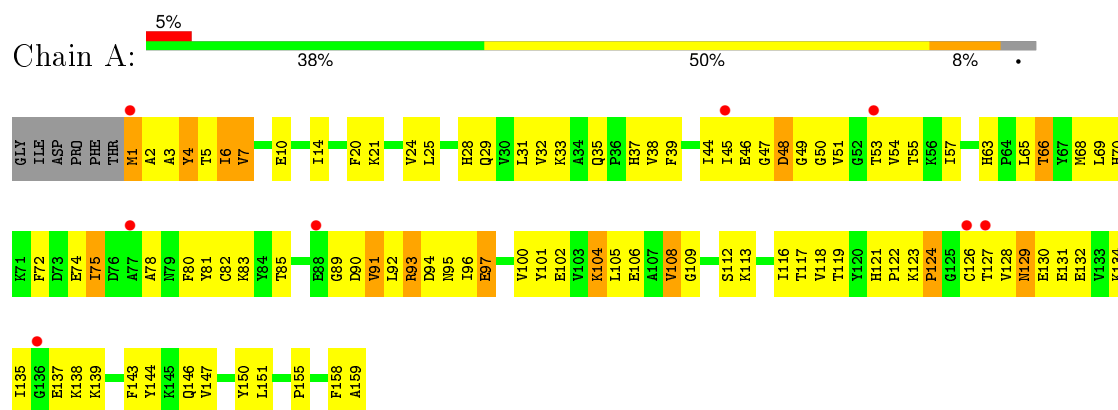
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	G	1	Total 1	O 1	0	0
5	H	4	Total 4	O 4	0	0
5	I	2	Total 2	O 2	0	0
5	L	2	Total 2	O 2	0	0
5	R	2	Total 2	O 2	0	0
5	T	1	Total 1	O 1	0	0
5	U	1	Total 1	O 1	0	0
5	W	2	Total 2	O 2	0	0
5	Y	2	Total 2	O 2	0	0
5	Z	3	Total 3	O 3	0	0
5	a	2	Total 2	O 2	0	0
5	b	1	Total 1	O 1	0	0

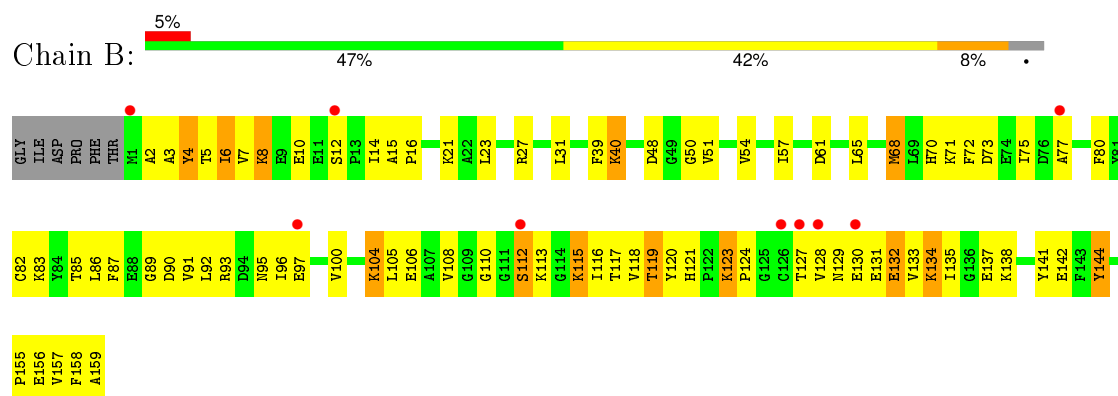
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

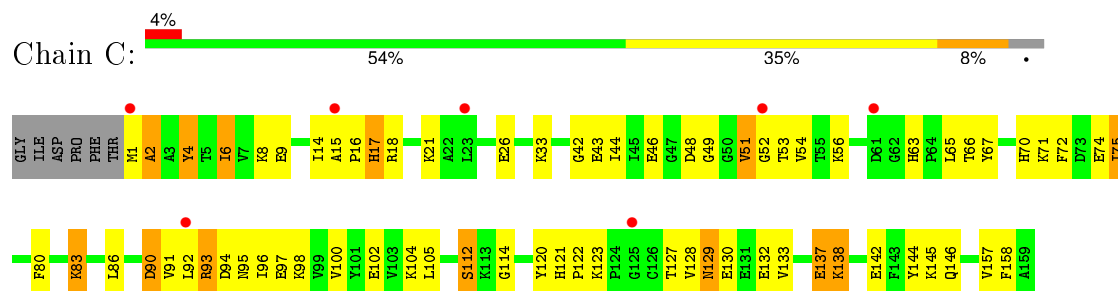
- Molecule 1: Phenolic oxidative coupling protein



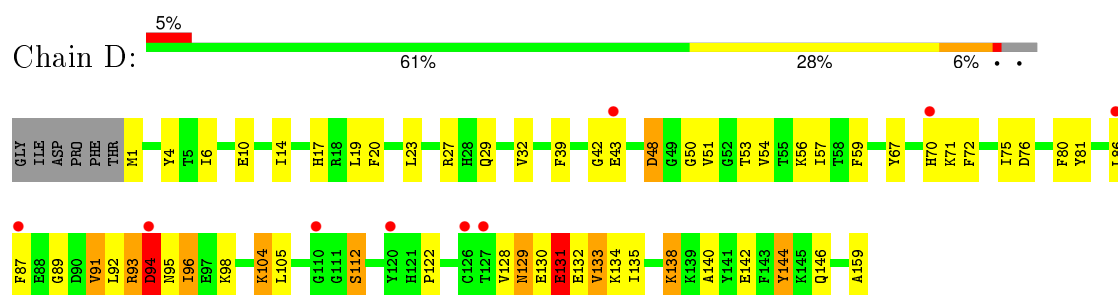
- Molecule 1: Phenolic oxidative coupling protein



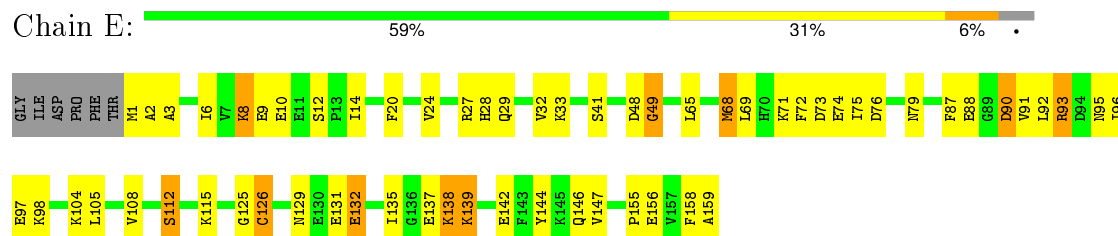
- Molecule 1: Phenolic oxidative coupling protein



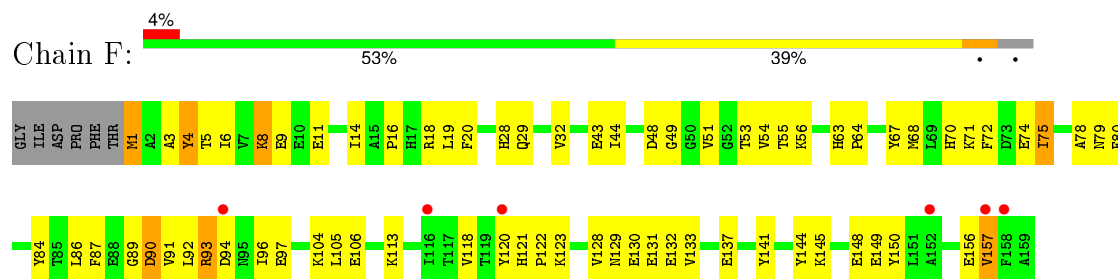
- Molecule 1: Phenolic oxidative coupling protein



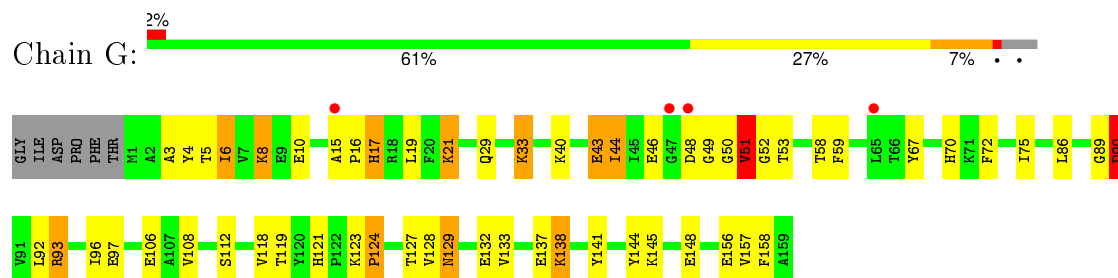
- Molecule 1: Phenolic oxidative coupling protein



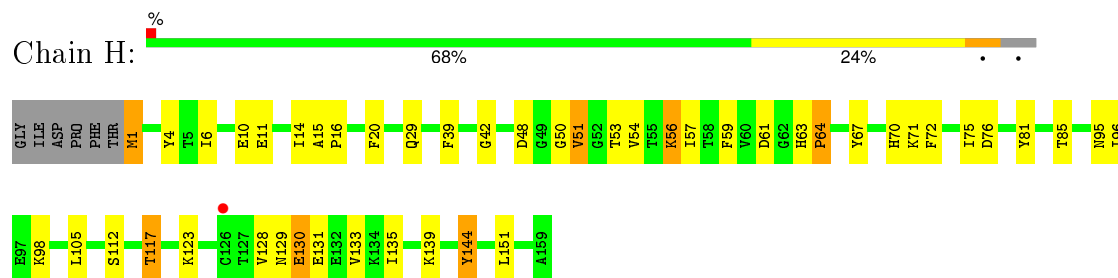
- Molecule 1: Phenolic oxidative coupling protein



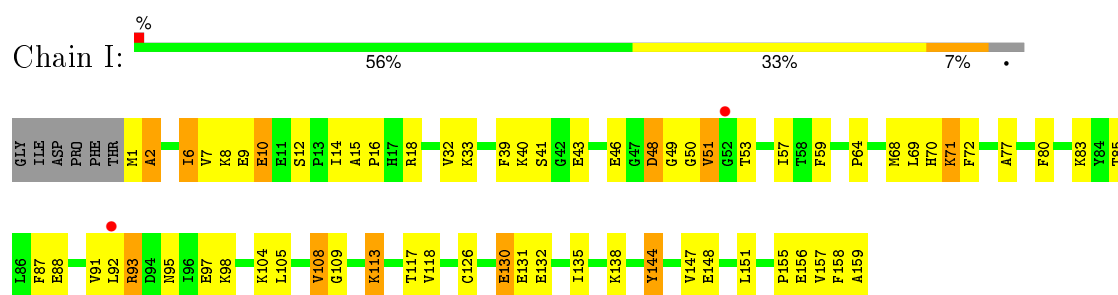
- Molecule 1: Phenolic oxidative coupling protein



- Molecule 1: Phenolic oxidative coupling protein



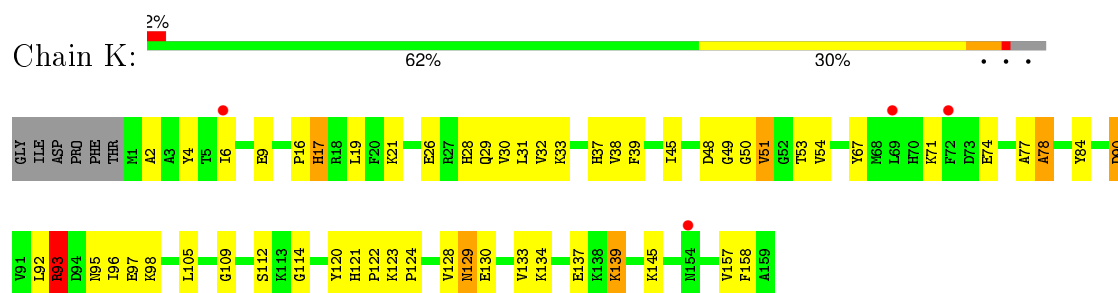
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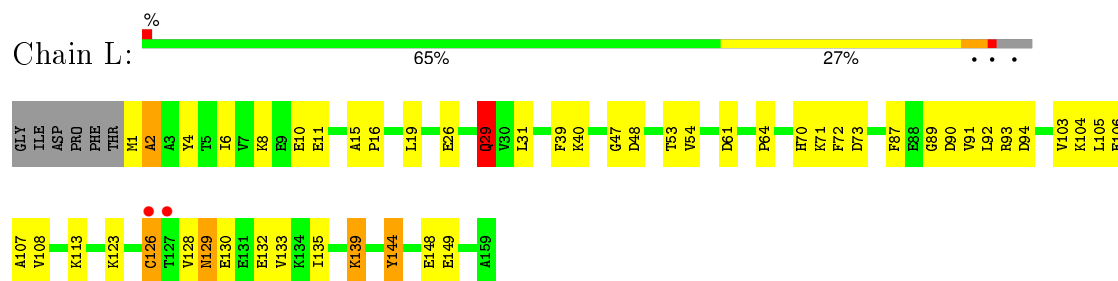
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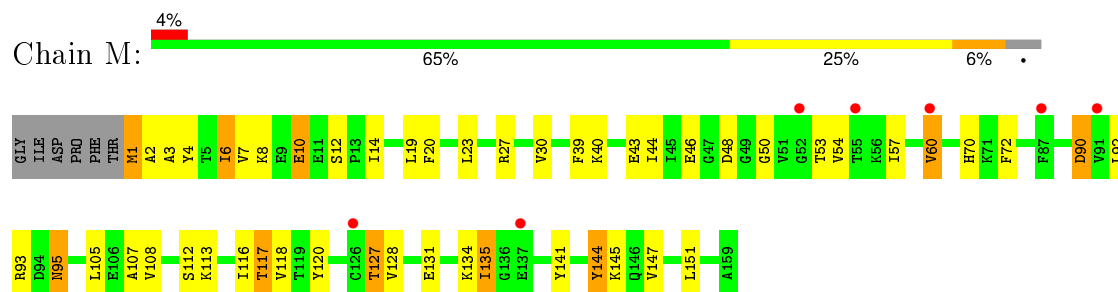
- Molecule 1: Phenolic oxidative coupling protein



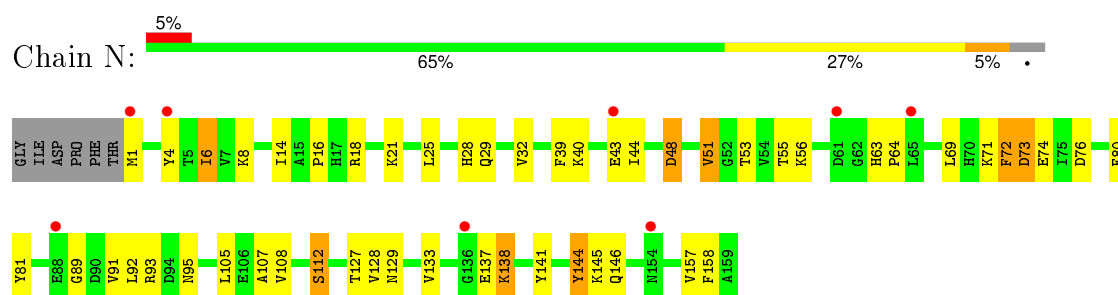
- Molecule 1: Phenolic oxidative coupling protein



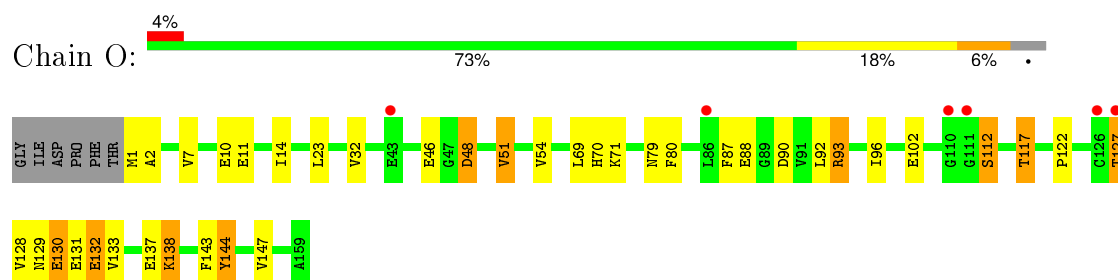
- Molecule 1: Phenolic oxidative coupling protein



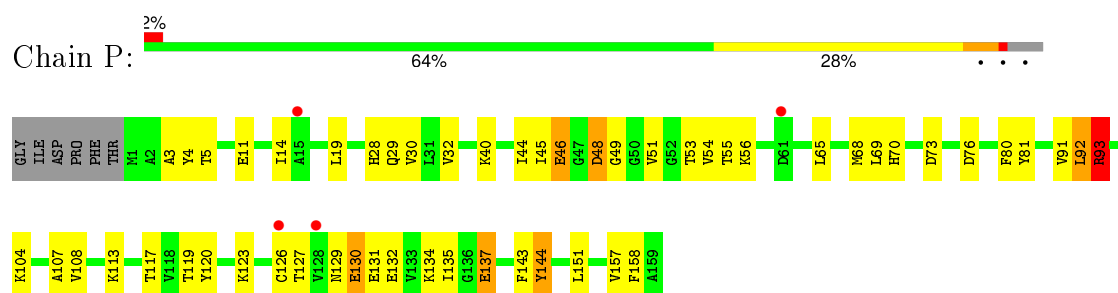
- Molecule 1: Phenolic oxidative coupling protein



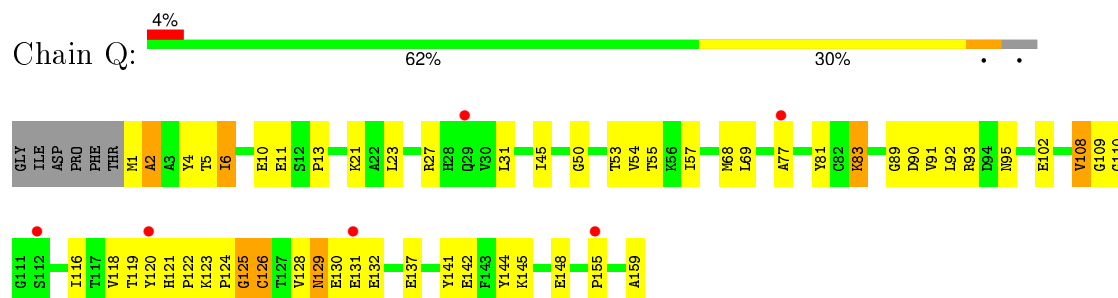
- Molecule 1: Phenolic oxidative coupling protein



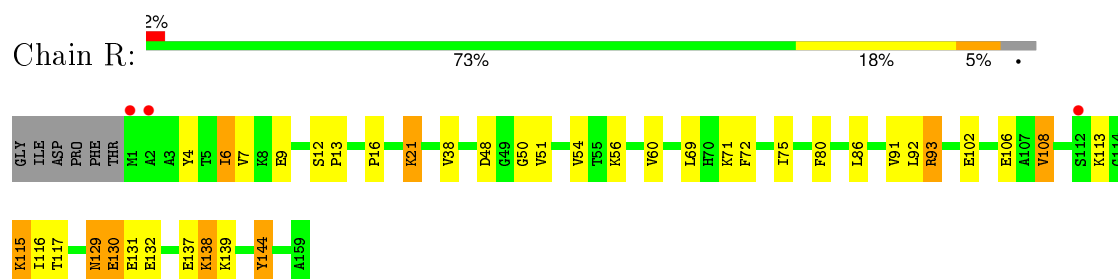
- Molecule 1: Phenolic oxidative coupling protein



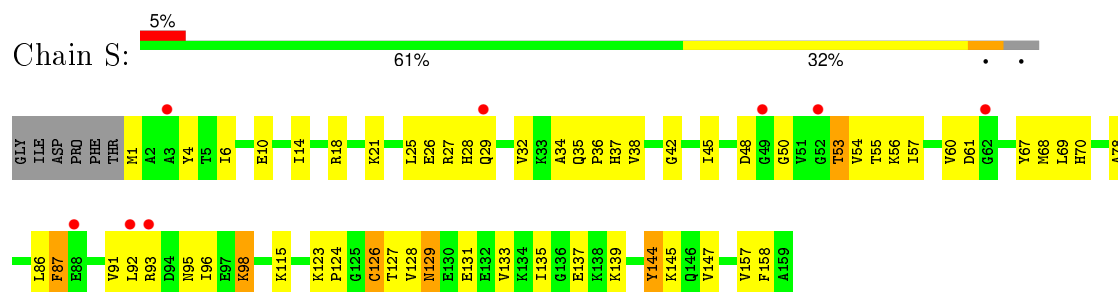
- Molecule 1: Phenolic oxidative coupling protein



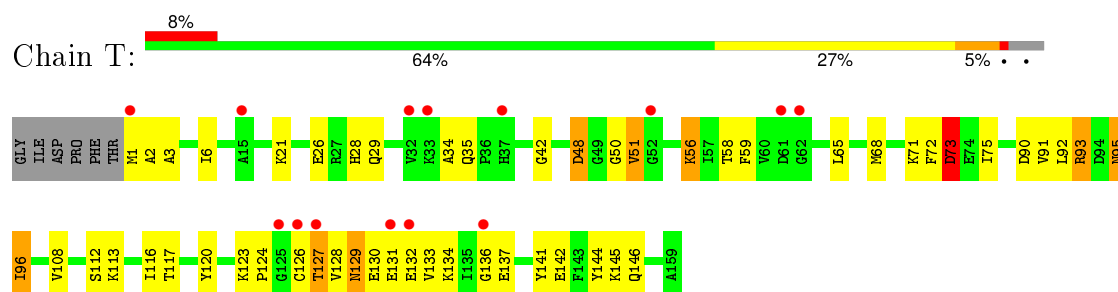
- Molecule 1: Phenolic oxidative coupling protein



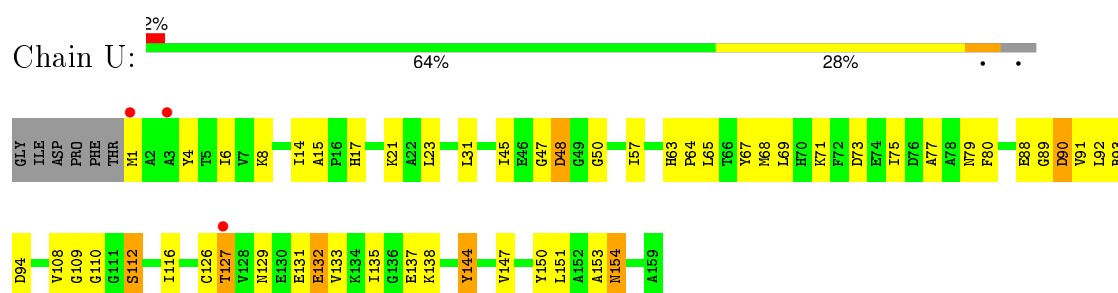
- Molecule 1: Phenolic oxidative coupling protein



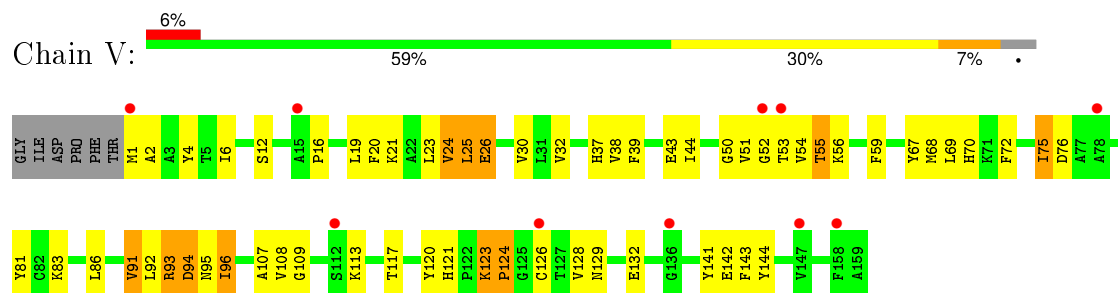
- Molecule 1: Phenolic oxidative coupling protein



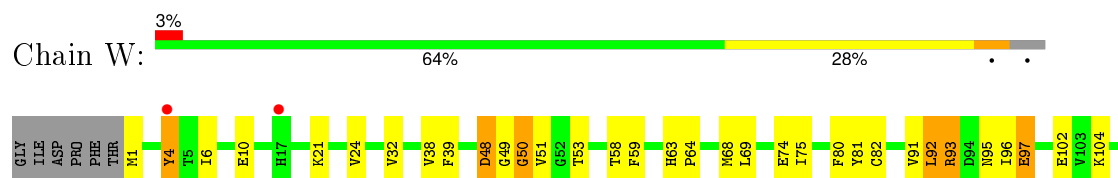
- Molecule 1: Phenolic oxidative coupling protein



- Molecule 1: Phenolic oxidative coupling protein

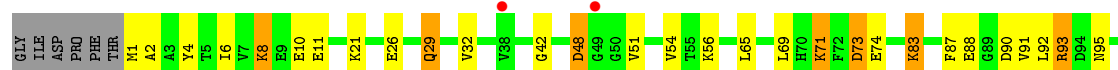


- Molecule 1: Phenolic oxidative coupling protein





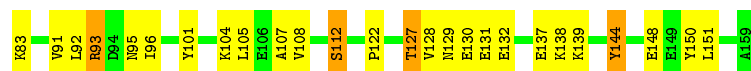
- Molecule 1: Phenolic oxidative coupling protein



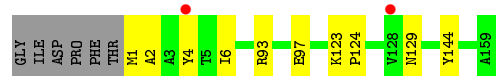
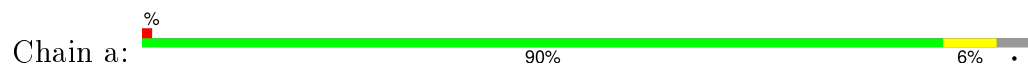
- Molecule 1: Phenolic oxidative coupling protein



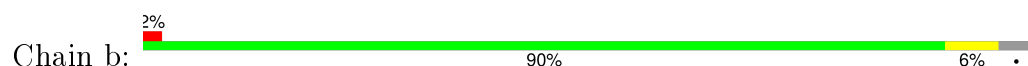
- Molecule 1: Phenolic oxidative coupling protein



- Molecule 1: Phenolic oxidative coupling protein



- Molecule 1: Phenolic oxidative coupling protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	146.29Å 146.29Å 298.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.87 – 2.43 48.87 – 2.43	Depositor EDS
% Data completeness (in resolution range)	99.7 (48.87-2.43) 99.7 (48.87-2.43)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.69 (at 2.42Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.223 , 0.278 0.229 , 0.284	Depositor DCC
R_{free} test set	3078 reflections (1.33%)	DCC
Wilson B-factor (Å ²)	49.9	Xtriage
Anisotropy	0.143	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 22.3	EDS
Estimated twinning fraction	0.251 for H, K, L 0.292 for -K, -H, -L 0.249 for -H, -K, L 0.209 for K, H, -L 0.489 for k,h,-l 0.489 for -k,-h,-l 0.489 for -h,-k,l	Xtriage
Reported twinning fraction	0.251 for H, K, L 0.292 for -K, -H, -L 0.249 for -H, -K, L 0.209 for K, H, -L	Depositor
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	2 of 235350 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	37252	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 67.36 % 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 5.2079e-06. The detected translational NCS is most likely*

¹ Intensities estimated from amplitudes.

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

also responsible for the elevated intensity ratio.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: 2AN, EPE, 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	0.74	0/1298	1.03	0/1753
1	B	0.81	0/1285	1.08	2/1736 (0.1%)
1	C	0.72	0/1285	1.04	1/1736 (0.1%)
1	D	0.69	0/1286	0.97	0/1737
1	E	0.93	0/1289	1.17	4/1741 (0.2%)
1	F	0.61	0/1286	0.93	2/1737 (0.1%)
1	G	0.97	1/1288 (0.1%)	1.17	4/1740 (0.2%)
1	H	0.77	0/1289	1.06	2/1741 (0.1%)
1	I	0.82	0/1284	1.09	2/1735 (0.1%)
1	J	0.93	0/1284	1.18	4/1734 (0.2%)
1	K	0.68	0/1289	1.01	0/1741
1	L	0.89	0/1289	1.17	2/1741 (0.1%)
1	M	0.75	0/1285	1.08	1/1736 (0.1%)
1	N	0.76	0/1298	1.05	0/1753
1	O	0.72	0/1287	1.04	1/1738 (0.1%)
1	P	0.70	0/1289	1.07	2/1741 (0.1%)
1	Q	0.70	0/1286	1.04	1/1737 (0.1%)
1	R	0.81	0/1289	1.11	0/1741
1	S	0.64	0/1286	0.99	0/1737
1	T	0.61	0/1337	0.96	2/1806 (0.1%)
1	U	0.91	0/1289	1.19	3/1741 (0.2%)
1	V	0.59	0/1286	0.94	0/1737
1	W	0.86	0/1287	1.14	2/1738 (0.1%)
1	X	0.78	0/1289	1.07	1/1741 (0.1%)
1	Y	0.81	0/1289	1.15	3/1741 (0.2%)
1	Z	0.82	0/1285	1.08	3/1736 (0.2%)
1	a	0.75	0/1300	1.06	0/1756
1	b	0.89	0/1283	1.18	1/1733 (0.1%)
All	All	0.78	1/36117 (0.0%)	1.08	43/48784 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	148	GLU	CG-CD	-5.36	1.44	1.51

The worst 5 of 43 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	U	109	GLY	N-CA-C	-5.90	98.36	113.10
1	E	6	ILE	CB-CA-C	-5.87	99.87	111.60
1	G	51	VAL	CB-CA-C	-5.78	100.41	111.40
1	I	6	ILE	CB-CA-C	-5.76	100.08	111.60
1	W	50	GLY	N-CA-C	-5.75	98.74	113.10

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1269	0	1260	133	0
1	B	1256	0	1251	97	0
1	C	1256	0	1251	74	0
1	D	1257	0	1253	58	0
1	E	1260	0	1255	54	0
1	F	1257	0	1251	66	0
1	G	1259	0	1250	59	0
1	H	1260	0	1255	40	0
1	I	1255	0	1246	71	0
1	J	1255	0	1250	58	0
1	K	1260	0	1255	57	0
1	L	1260	0	1255	57	0
1	M	1256	0	1251	52	0
1	N	1269	0	1260	50	0
1	O	1258	0	1250	30	0
1	P	1260	0	1255	39	0
1	Q	1257	0	1246	47	0
1	R	1260	0	1255	39	0
1	S	1257	0	1246	57	0
1	T	1307	0	1286	37	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	U	1260	0	1255	45	0
1	V	1257	0	1253	53	0
1	W	1258	0	1250	53	0
1	X	1260	0	1255	46	0
1	Y	1260	0	1255	47	0
1	Z	1256	0	1251	55	0
1	a	1270	0	1261	0	0
1	b	1254	0	1249	0	0
2	A	84	0	49	34	0
2	B	105	0	63	37	0
2	C	84	0	49	28	0
2	D	63	0	36	22	0
2	E	105	0	62	45	0
2	F	21	0	13	8	0
2	G	105	0	60	36	0
2	H	63	0	39	17	0
2	I	105	0	62	36	0
2	J	42	0	25	13	0
2	K	63	0	36	30	0
2	L	105	0	63	31	0
2	M	63	0	37	24	0
2	N	84	0	51	32	0
2	O	42	0	24	10	0
2	P	42	0	24	10	0
2	Q	63	0	37	21	0
2	R	84	0	49	25	0
2	S	21	0	13	7	0
2	U	105	0	63	25	0
2	V	42	0	24	11	0
2	W	84	0	49	34	0
2	X	63	0	39	14	0
2	Y	84	0	49	24	0
2	Z	63	0	37	25	0
2	a	21	0	12	0	0
2	b	63	0	38	0	0
3	C	15	0	18	0	0
3	Z	15	0	18	0	0
4	V	5	0	0	0	0
4	X	5	0	0	0	0
4	b	5	0	0	0	0
5	A	5	0	0	1	0
5	C	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	D	1	0	0	0	0
5	E	2	0	0	0	0
5	F	2	0	0	0	0
5	G	1	0	0	1	0
5	H	4	0	0	0	0
5	I	2	0	0	0	0
5	L	2	0	0	0	0
5	R	2	0	0	0	0
5	T	1	0	0	0	0
5	U	1	0	0	0	0
5	W	2	0	0	0	0
5	Y	2	0	0	0	0
5	Z	3	0	0	0	0
5	a	2	0	0	0	0
5	b	1	0	0	0	0
All	All	37252	0	36249	1870	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 25.

The worst 5 of 1870 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:L:128:VAL:CG1	1:L:129:ASN:HB3	1.48	1.40
1:A:123:LYS:HD2	1:A:124:PRO:CD	4.03	1.38
1:A:123:LYS:CD	1:A:124:PRO:HD2	3.44	1.37
1:A:37:HIS:CD2	1:A:38:VAL:HG23	1.62	1.34
1:A:96:ILE:CD1	1:A:128:VAL:HG22	1.68	1.22

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	158/165 (96%)	148 (94%)	7 (4%)	3 (2%)	10	8
1	B	157/165 (95%)	143 (91%)	12 (8%)	2 (1%)	15	15
1	C	157/165 (95%)	139 (88%)	16 (10%)	2 (1%)	15	15
1	D	157/165 (95%)	140 (89%)	14 (9%)	3 (2%)	10	8
1	E	157/165 (95%)	142 (90%)	12 (8%)	3 (2%)	10	8
1	F	157/165 (95%)	145 (92%)	11 (7%)	1 (1%)	30	36
1	G	157/165 (95%)	145 (92%)	9 (6%)	3 (2%)	10	8
1	H	157/165 (95%)	150 (96%)	7 (4%)	0	100	100
1	I	157/165 (95%)	145 (92%)	10 (6%)	2 (1%)	15	15
1	J	157/165 (95%)	143 (91%)	12 (8%)	2 (1%)	15	15
1	K	157/165 (95%)	143 (91%)	10 (6%)	4 (2%)	7	4
1	L	157/165 (95%)	149 (95%)	6 (4%)	2 (1%)	15	15
1	M	157/165 (95%)	147 (94%)	10 (6%)	0	100	100
1	N	158/165 (96%)	145 (92%)	13 (8%)	0	100	100
1	O	157/165 (95%)	147 (94%)	10 (6%)	0	100	100
1	P	157/165 (95%)	139 (88%)	14 (9%)	4 (2%)	7	4
1	Q	157/165 (95%)	142 (90%)	13 (8%)	2 (1%)	15	15
1	R	157/165 (95%)	146 (93%)	9 (6%)	2 (1%)	15	15
1	S	157/165 (95%)	141 (90%)	15 (10%)	1 (1%)	30	36
1	T	163/165 (99%)	151 (93%)	11 (7%)	1 (1%)	30	36
1	U	157/165 (95%)	145 (92%)	11 (7%)	1 (1%)	30	36
1	V	157/165 (95%)	139 (88%)	16 (10%)	2 (1%)	15	15
1	W	157/165 (95%)	151 (96%)	4 (2%)	2 (1%)	15	15
1	X	157/165 (95%)	148 (94%)	7 (4%)	2 (1%)	15	15
1	Y	157/165 (95%)	145 (92%)	11 (7%)	1 (1%)	30	36
1	Z	157/165 (95%)	142 (90%)	12 (8%)	3 (2%)	10	8
1	a	158/165 (96%)	143 (90%)	13 (8%)	2 (1%)	15	15
1	b	157/165 (95%)	141 (90%)	15 (10%)	1 (1%)	30	36
All	All	4405/4620 (95%)	4044 (92%)	310 (7%)	51 (1%)	16	17

5 of 51 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	94	ASP
1	D	131	GLU
1	E	126	CYS
1	I	2	ALA
1	J	3	ALA

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	135/139 (97%)	118 (87%)	17 (13%)	5	4
1	B	133/139 (96%)	118 (89%)	15 (11%)	7	7
1	C	133/139 (96%)	117 (88%)	16 (12%)	6	5
1	D	133/139 (96%)	118 (89%)	15 (11%)	7	7
1	E	134/139 (96%)	124 (92%)	10 (8%)	17	22
1	F	133/139 (96%)	121 (91%)	12 (9%)	12	13
1	G	133/139 (96%)	114 (86%)	19 (14%)	4	3
1	H	134/139 (96%)	126 (94%)	8 (6%)	24	33
1	I	132/139 (95%)	120 (91%)	12 (9%)	12	13
1	J	133/139 (96%)	121 (91%)	12 (9%)	12	13
1	K	134/139 (96%)	123 (92%)	11 (8%)	14	18
1	L	134/139 (96%)	121 (90%)	13 (10%)	10	11
1	M	133/139 (96%)	120 (90%)	13 (10%)	10	11
1	N	135/139 (97%)	122 (90%)	13 (10%)	10	11
1	O	133/139 (96%)	121 (91%)	12 (9%)	12	13
1	P	134/139 (96%)	123 (92%)	11 (8%)	14	18
1	Q	133/139 (96%)	124 (93%)	9 (7%)	20	26
1	R	134/139 (96%)	122 (91%)	12 (9%)	12	13
1	S	133/139 (96%)	123 (92%)	10 (8%)	17	22
1	T	139/139 (100%)	121 (87%)	18 (13%)	5	4

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	U	134/139 (96%)	119 (89%)	15 (11%)	7	7
1	V	133/139 (96%)	118 (89%)	15 (11%)	7	7
1	W	133/139 (96%)	125 (94%)	8 (6%)	24	33
1	X	134/139 (96%)	121 (90%)	13 (10%)	10	11
1	Y	134/139 (96%)	118 (88%)	16 (12%)	6	5
1	Z	133/139 (96%)	120 (90%)	13 (10%)	10	11
1	a	135/139 (97%)	127 (94%)	8 (6%)	24	34
1	b	132/139 (95%)	124 (94%)	8 (6%)	23	32
All	All	3743/3892 (96%)	3389 (90%)	354 (10%)	11	12

5 of 354 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	M	1	MET
1	P	29	GLN
1	Z	5	THR
1	M	43	GLU
1	N	72	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 44 such sidechains are listed below:

Mol	Chain	Res	Type
1	J	29	GLN
1	L	35	GLN
1	X	154	ASN
1	J	95	ASN
1	K	79	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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

94 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$
2	2AN	A	201	-	23,23,23	0.76	0	30,33,33	1.43	4 (13%)
2	2AN	A	202	-	23,23,23	0.90	2 (8%)	30,33,33	1.22	4 (13%)
2	2AN	A	203	-	23,23,23	0.75	0	30,33,33	1.07	3 (10%)
2	2AN	A	204	-	23,23,23	0.89	0	30,33,33	1.30	5 (16%)
2	2AN	B	201	-	23,23,23	0.59	0	30,33,33	1.08	2 (6%)
2	2AN	B	202	-	23,23,23	0.56	0	30,33,33	1.25	4 (13%)
2	2AN	B	203	-	23,23,23	0.57	0	30,33,33	1.11	4 (13%)
2	2AN	B	204	-	23,23,23	0.76	0	30,33,33	1.13	4 (13%)
2	2AN	B	205	-	23,23,23	0.55	0	30,33,33	1.13	4 (13%)
2	2AN	C	201	-	23,23,23	0.90	1 (4%)	30,33,33	1.20	4 (13%)
2	2AN	C	202	-	23,23,23	0.73	0	30,33,33	1.09	4 (13%)
2	2AN	C	203	-	23,23,23	0.64	0	30,33,33	1.25	3 (10%)
2	2AN	C	204	-	23,23,23	0.83	1 (4%)	30,33,33	1.22	5 (16%)
3	EPE	C	205	-	14,15,15	1.07	1 (7%)	18,20,20	3.98	13 (72%)
2	2AN	D	201	-	23,23,23	0.67	0	30,33,33	1.04	3 (10%)
2	2AN	D	202	-	23,23,23	0.64	0	30,33,33	1.07	4 (13%)
2	2AN	D	203	-	23,23,23	0.85	1 (4%)	30,33,33	1.09	3 (10%)
2	2AN	E	201	-	23,23,23	1.28	3 (13%)	30,33,33	1.56	7 (23%)
2	2AN	E	202	-	23,23,23	0.59	0	30,33,33	1.27	4 (13%)
2	2AN	E	203	-	23,23,23	0.74	1 (4%)	30,33,33	1.49	4 (13%)
2	2AN	E	204	-	23,23,23	0.83	1 (4%)	30,33,33	1.24	4 (13%)
2	2AN	E	205	-	23,23,23	0.78	0	30,33,33	1.16	5 (16%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	2AN	F	201	-	23,23,23	0.75	0	30,33,33	1.20	3 (10%)
2	2AN	G	201	-	23,23,23	1.55	3 (13%)	30,33,33	1.42	4 (13%)
2	2AN	G	202	-	23,23,23	0.98	1 (4%)	30,33,33	1.16	4 (13%)
2	2AN	G	203	-	23,23,23	0.72	0	30,33,33	1.07	2 (6%)
2	2AN	G	204	-	23,23,23	1.01	1 (4%)	30,33,33	1.25	4 (13%)
2	2AN	G	205	-	23,23,23	0.50	0	30,33,33	1.17	4 (13%)
2	2AN	H	201	-	23,23,23	0.80	0	30,33,33	1.25	5 (16%)
2	2AN	H	202	-	23,23,23	0.73	0	30,33,33	1.04	2 (6%)
2	2AN	H	203	-	23,23,23	0.85	1 (4%)	30,33,33	1.23	3 (10%)
2	2AN	I	201	-	23,23,23	0.81	1 (4%)	30,33,33	1.05	2 (6%)
2	2AN	I	202	-	23,23,23	1.08	1 (4%)	30,33,33	1.06	3 (10%)
2	2AN	I	203	-	23,23,23	0.61	0	30,33,33	1.22	4 (13%)
2	2AN	I	204	-	23,23,23	0.69	0	30,33,33	1.48	6 (20%)
2	2AN	I	205	-	23,23,23	0.54	0	30,33,33	1.14	4 (13%)
2	2AN	J	201	-	23,23,23	0.93	1 (4%)	30,33,33	1.00	4 (13%)
2	2AN	J	202	-	23,23,23	0.89	1 (4%)	30,33,33	1.17	4 (13%)
2	2AN	K	201	-	23,23,23	0.88	0	30,33,33	1.22	3 (10%)
2	2AN	K	202	-	23,23,23	0.47	0	30,33,33	1.18	3 (10%)
2	2AN	K	203	-	23,23,23	0.47	0	30,33,33	1.20	5 (16%)
2	2AN	L	201	-	23,23,23	0.80	1 (4%)	30,33,33	1.09	3 (10%)
2	2AN	L	202	-	23,23,23	0.80	0	30,33,33	1.03	2 (6%)
2	2AN	L	203	-	23,23,23	0.79	0	30,33,33	1.32	5 (16%)
2	2AN	L	204	-	23,23,23	0.64	0	30,33,33	1.25	5 (16%)
2	2AN	L	205	-	23,23,23	0.62	0	30,33,33	1.15	3 (10%)
2	2AN	M	201	-	23,23,23	0.82	1 (4%)	30,33,33	1.04	2 (6%)
2	2AN	M	202	-	23,23,23	0.82	1 (4%)	30,33,33	1.08	2 (6%)
2	2AN	M	203	-	23,23,23	0.52	0	30,33,33	1.21	4 (13%)
2	2AN	N	201	-	23,23,23	0.71	0	30,33,33	1.07	3 (10%)
2	2AN	N	202	-	23,23,23	0.78	0	30,33,33	1.15	3 (10%)
2	2AN	N	203	-	23,23,23	0.83	0	30,33,33	1.36	4 (13%)
2	2AN	N	204	-	23,23,23	1.09	2 (8%)	30,33,33	1.24	4 (13%)
2	2AN	O	201	-	23,23,23	0.84	1 (4%)	30,33,33	1.17	3 (10%)
2	2AN	O	202	-	23,23,23	0.85	1 (4%)	30,33,33	1.11	3 (10%)
2	2AN	P	201	-	23,23,23	0.74	0	30,33,33	1.13	4 (13%)
2	2AN	P	202	-	23,23,23	0.64	0	30,33,33	1.29	4 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	2AN	Q	201	-	23,23,23	0.66	0	30,33,33	1.20	4 (13%)
2	2AN	Q	202	-	23,23,23	0.58	0	30,33,33	1.16	4 (13%)
2	2AN	Q	203	-	23,23,23	0.61	0	30,33,33	1.11	4 (13%)
2	2AN	R	201	-	23,23,23	0.71	0	30,33,33	1.31	4 (13%)
2	2AN	R	202	-	23,23,23	1.07	1 (4%)	30,33,33	1.22	3 (10%)
2	2AN	R	203	-	23,23,23	0.67	0	30,33,33	1.29	4 (13%)
2	2AN	R	204	-	23,23,23	0.76	0	30,33,33	1.10	4 (13%)
2	2AN	S	201	-	23,23,23	0.83	1 (4%)	30,33,33	1.04	3 (10%)
2	2AN	U	201	-	23,23,23	0.70	1 (4%)	30,33,33	1.20	4 (13%)
2	2AN	U	202	-	23,23,23	0.94	1 (4%)	30,33,33	1.32	2 (6%)
2	2AN	U	203	-	23,23,23	0.72	0	30,33,33	1.19	5 (16%)
2	2AN	U	204	-	23,23,23	0.69	0	30,33,33	1.04	3 (10%)
2	2AN	U	205	-	23,23,23	0.52	0	30,33,33	1.06	2 (6%)
2	2AN	V	201	-	23,23,23	0.65	0	30,33,33	1.09	2 (6%)
2	2AN	V	202	-	23,23,23	0.63	0	30,33,33	1.32	5 (16%)
4	SO4	V	203	-	4,4,4	0.79	0	6,6,6	1.11	0
2	2AN	W	201	-	23,23,23	0.86	1 (4%)	30,33,33	1.31	5 (16%)
2	2AN	W	202	-	23,23,23	0.77	1 (4%)	30,33,33	1.02	3 (10%)
2	2AN	W	203	-	23,23,23	1.08	2 (8%)	30,33,33	1.16	4 (13%)
2	2AN	W	204	-	23,23,23	0.84	1 (4%)	30,33,33	1.36	4 (13%)
2	2AN	X	201	-	23,23,23	0.69	0	30,33,33	1.05	3 (10%)
2	2AN	X	202	-	23,23,23	0.52	0	30,33,33	1.18	4 (13%)
2	2AN	X	203	-	23,23,23	0.56	0	30,33,33	1.24	2 (6%)
4	SO4	X	204	-	4,4,4	0.59	0	6,6,6	0.44	0
2	2AN	Y	201	-	23,23,23	1.07	3 (13%)	30,33,33	1.16	2 (6%)
2	2AN	Y	202	-	23,23,23	0.70	0	30,33,33	1.02	2 (6%)
2	2AN	Y	203	-	23,23,23	0.53	0	30,33,33	1.08	2 (6%)
2	2AN	Y	204	-	23,23,23	0.57	0	30,33,33	1.11	3 (10%)
2	2AN	Z	201	-	23,23,23	0.77	0	30,33,33	1.21	4 (13%)
2	2AN	Z	202	-	23,23,23	0.83	0	30,33,33	1.28	6 (20%)
2	2AN	Z	203	-	23,23,23	0.59	0	30,33,33	1.13	3 (10%)
3	EPE	Z	204	-	14,15,15	0.84	0	18,20,20	4.03	11 (61%)
2	2AN	a	201	-	23,23,23	0.53	0	30,33,33	1.02	3 (10%)
2	2AN	b	201	-	23,23,23	0.58	0	30,33,33	1.18	3 (10%)
2	2AN	b	202	-	23,23,23	1.13	1 (4%)	30,33,33	1.17	2 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	2AN	b	203	-	23,23,23	0.65	0	30,33,33	1.06	3 (10%)
4	SO4	b	204	-	4,4,4	0.50	0	6,6,6	1.73	1 (16%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	2AN	A	201	-	-	0/10/10/10	0/3/3/3
2	2AN	A	202	-	-	0/10/10/10	0/3/3/3
2	2AN	A	203	-	-	0/10/10/10	0/3/3/3
2	2AN	A	204	-	-	0/10/10/10	0/3/3/3
2	2AN	B	201	-	-	0/10/10/10	0/3/3/3
2	2AN	B	202	-	-	0/10/10/10	0/3/3/3
2	2AN	B	203	-	-	0/10/10/10	0/3/3/3
2	2AN	B	204	-	-	0/10/10/10	0/3/3/3
2	2AN	B	205	-	-	0/10/10/10	0/3/3/3
2	2AN	C	201	-	-	0/10/10/10	0/3/3/3
2	2AN	C	202	-	-	0/10/10/10	0/3/3/3
2	2AN	C	203	-	-	0/10/10/10	0/3/3/3
2	2AN	C	204	-	-	0/10/10/10	0/3/3/3
3	EPE	C	205	-	-	0/9/19/19	0/1/1/1
2	2AN	D	201	-	-	0/10/10/10	0/3/3/3
2	2AN	D	202	-	-	0/10/10/10	0/3/3/3
2	2AN	D	203	-	-	0/10/10/10	0/3/3/3
2	2AN	E	201	-	-	0/10/10/10	0/3/3/3
2	2AN	E	202	-	-	0/10/10/10	0/3/3/3
2	2AN	E	203	-	-	0/10/10/10	0/3/3/3
2	2AN	E	204	-	-	0/10/10/10	0/3/3/3
2	2AN	E	205	-	-	0/10/10/10	0/3/3/3
2	2AN	F	201	-	-	0/10/10/10	0/3/3/3
2	2AN	G	201	-	-	0/10/10/10	0/3/3/3
2	2AN	G	202	-	-	0/10/10/10	0/3/3/3
2	2AN	G	203	-	-	0/10/10/10	0/3/3/3
2	2AN	G	204	-	-	0/10/10/10	0/3/3/3
2	2AN	G	205	-	-	0/10/10/10	0/3/3/3
2	2AN	H	201	-	-	0/10/10/10	0/3/3/3
2	2AN	H	202	-	-	0/10/10/10	0/3/3/3
2	2AN	H	203	-	-	0/10/10/10	0/3/3/3
2	2AN	I	201	-	-	0/10/10/10	0/3/3/3
2	2AN	I	202	-	-	0/10/10/10	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	2AN	I	203	-	-	0/10/10/10	0/3/3/3
2	2AN	I	204	-	-	0/10/10/10	0/3/3/3
2	2AN	I	205	-	-	0/10/10/10	0/3/3/3
2	2AN	J	201	-	-	0/10/10/10	0/3/3/3
2	2AN	J	202	-	-	0/10/10/10	0/3/3/3
2	2AN	K	201	-	-	0/10/10/10	0/3/3/3
2	2AN	K	202	-	-	0/10/10/10	0/3/3/3
2	2AN	K	203	-	-	0/10/10/10	0/3/3/3
2	2AN	L	201	-	-	0/10/10/10	0/3/3/3
2	2AN	L	202	-	-	0/10/10/10	0/3/3/3
2	2AN	L	203	-	-	0/10/10/10	0/3/3/3
2	2AN	L	204	-	-	0/10/10/10	0/3/3/3
2	2AN	L	205	-	-	0/10/10/10	0/3/3/3
2	2AN	M	201	-	-	0/10/10/10	0/3/3/3
2	2AN	M	202	-	-	0/10/10/10	0/3/3/3
2	2AN	M	203	-	-	0/10/10/10	0/3/3/3
2	2AN	N	201	-	-	0/10/10/10	0/3/3/3
2	2AN	N	202	-	-	0/10/10/10	0/3/3/3
2	2AN	N	203	-	-	0/10/10/10	0/3/3/3
2	2AN	N	204	-	-	0/10/10/10	0/3/3/3
2	2AN	O	201	-	-	0/10/10/10	0/3/3/3
2	2AN	O	202	-	-	0/10/10/10	0/3/3/3
2	2AN	P	201	-	-	0/10/10/10	0/3/3/3
2	2AN	P	202	-	-	0/10/10/10	0/3/3/3
2	2AN	Q	201	-	-	0/10/10/10	0/3/3/3
2	2AN	Q	202	-	-	0/10/10/10	0/3/3/3
2	2AN	Q	203	-	-	0/10/10/10	0/3/3/3
2	2AN	R	201	-	-	0/10/10/10	0/3/3/3
2	2AN	R	202	-	-	0/10/10/10	0/3/3/3
2	2AN	R	203	-	-	0/10/10/10	0/3/3/3
2	2AN	R	204	-	-	0/10/10/10	0/3/3/3
2	2AN	S	201	-	-	0/10/10/10	0/3/3/3
2	2AN	U	201	-	-	0/10/10/10	0/3/3/3
2	2AN	U	202	-	-	0/10/10/10	0/3/3/3
2	2AN	U	203	-	-	0/10/10/10	0/3/3/3
2	2AN	U	204	-	-	0/10/10/10	0/3/3/3
2	2AN	U	205	-	-	0/10/10/10	0/3/3/3
2	2AN	V	201	-	-	0/10/10/10	0/3/3/3
2	2AN	V	202	-	-	0/10/10/10	0/3/3/3
4	SO4	V	203	-	-	0/0/0/0	0/0/0/0
2	2AN	W	201	-	-	0/10/10/10	0/3/3/3
2	2AN	W	202	-	-	0/10/10/10	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	2AN	W	203	-	-	0/10/10/10	0/3/3/3
2	2AN	W	204	-	-	0/10/10/10	0/3/3/3
2	2AN	X	201	-	-	0/10/10/10	0/3/3/3
2	2AN	X	202	-	-	0/10/10/10	0/3/3/3
2	2AN	X	203	-	-	0/10/10/10	0/3/3/3
4	SO4	X	204	-	-	0/0/0/0	0/0/0/0
2	2AN	Y	201	-	-	0/10/10/10	0/3/3/3
2	2AN	Y	202	-	-	0/10/10/10	0/3/3/3
2	2AN	Y	203	-	-	0/10/10/10	0/3/3/3
2	2AN	Y	204	-	-	0/10/10/10	0/3/3/3
2	2AN	Z	201	-	-	0/10/10/10	0/3/3/3
2	2AN	Z	202	-	-	0/10/10/10	0/3/3/3
2	2AN	Z	203	-	-	0/10/10/10	0/3/3/3
3	EPE	Z	204	-	-	0/9/19/19	0/1/1/1
2	2AN	a	201	-	-	0/10/10/10	0/3/3/3
2	2AN	b	201	-	-	0/10/10/10	0/3/3/3
2	2AN	b	202	-	-	0/10/10/10	0/3/3/3
2	2AN	b	203	-	-	0/10/10/10	0/3/3/3
4	SO4	b	204	-	-	0/0/0/0	0/0/0/0

The worst 5 of 41 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	201	2AN	C9-S	-3.93	1.72	1.78
2	R	202	2AN	C9-S	-3.89	1.72	1.78
2	b	202	2AN	C9-S	-3.56	1.73	1.78
2	N	204	2AN	C16-C11	-3.30	1.33	1.39
2	G	204	2AN	C13-C12	-3.23	1.32	1.38

The worst 5 of 343 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Z	204	EPE	O2S-S-C10	-10.76	97.72	106.91
3	C	205	EPE	C2-C3-N4	-5.26	101.22	110.63
3	Z	204	EPE	O3S-S-O2S	-4.13	101.99	111.61
3	C	205	EPE	O3S-S-O1S	-3.82	102.71	111.61
4	b	204	SO4	O2-S-O1	-3.70	97.78	109.50

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

85 monomers are involved in 599 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	201	2AN	17	0
2	A	202	2AN	3	0
2	A	203	2AN	7	0
2	A	204	2AN	7	0
2	B	201	2AN	8	0
2	B	202	2AN	10	0
2	B	203	2AN	7	0
2	B	204	2AN	6	0
2	B	205	2AN	6	0
2	C	201	2AN	4	0
2	C	202	2AN	8	0
2	C	203	2AN	7	0
2	C	204	2AN	9	0
2	D	201	2AN	6	0
2	D	202	2AN	9	0
2	D	203	2AN	7	0
2	E	201	2AN	6	0
2	E	202	2AN	11	0
2	E	203	2AN	14	0
2	E	204	2AN	7	0
2	E	205	2AN	8	0
2	F	201	2AN	8	0
2	G	201	2AN	5	0
2	G	202	2AN	11	0
2	G	203	2AN	2	0
2	G	204	2AN	12	0
2	G	205	2AN	7	0
2	H	201	2AN	6	0
2	H	202	2AN	5	0
2	H	203	2AN	6	0
2	I	201	2AN	8	0
2	I	202	2AN	6	0
2	I	203	2AN	12	0
2	I	204	2AN	6	0
2	I	205	2AN	7	0
2	J	201	2AN	6	0
2	J	202	2AN	7	0
2	K	201	2AN	4	0
2	K	202	2AN	8	0
2	K	203	2AN	18	0
2	L	201	2AN	3	0
2	L	202	2AN	8	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	L	203	2AN	5	0
2	L	204	2AN	10	0
2	L	205	2AN	5	0
2	M	201	2AN	6	0
2	M	202	2AN	7	0
2	M	203	2AN	11	0
2	N	201	2AN	4	0
2	N	202	2AN	7	0
2	N	203	2AN	16	0
2	N	204	2AN	5	0
2	O	201	2AN	6	0
2	O	202	2AN	4	0
2	P	201	2AN	6	0
2	P	202	2AN	4	0
2	Q	201	2AN	9	0
2	Q	202	2AN	5	0
2	Q	203	2AN	7	0
2	R	201	2AN	13	0
2	R	202	2AN	4	0
2	R	203	2AN	6	0
2	R	204	2AN	2	0
2	S	201	2AN	7	0
2	U	201	2AN	10	0
2	U	202	2AN	2	0
2	U	203	2AN	6	0
2	U	204	2AN	4	0
2	U	205	2AN	3	0
2	V	201	2AN	6	0
2	V	202	2AN	5	0
2	W	201	2AN	8	0
2	W	202	2AN	12	0
2	W	203	2AN	6	0
2	W	204	2AN	8	0
2	X	201	2AN	4	0
2	X	202	2AN	5	0
2	X	203	2AN	5	0
2	Y	201	2AN	4	0
2	Y	202	2AN	7	0
2	Y	203	2AN	6	0
2	Y	204	2AN	7	0
2	Z	201	2AN	5	0
2	Z	202	2AN	12	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	Z	203	2AN	8	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	159/165 (96%)	0.19	8 (5%) 32 32	27, 48, 69, 87	1 (0%)
1	B	159/165 (96%)	0.17	9 (5%) 27 25	25, 46, 77, 108	1 (0%)
1	C	159/165 (96%)	0.25	7 (4%) 38 37	29, 50, 73, 84	0
1	D	159/165 (96%)	0.23	9 (5%) 27 25	29, 50, 83, 123	0
1	E	159/165 (96%)	-0.17	0 100 100	20, 35, 57, 76	2 (1%)
1	F	159/165 (96%)	0.32	6 (3%) 44 44	34, 56, 82, 98	2 (1%)
1	G	159/165 (96%)	0.02	4 (2%) 61 59	20, 39, 64, 83	3 (1%)
1	H	159/165 (96%)	0.04	1 (0%) 90 91	24, 45, 65, 77	2 (1%)
1	I	159/165 (96%)	-0.01	2 (1%) 79 80	21, 40, 66, 83	4 (2%)
1	J	159/165 (96%)	-0.01	0 100 100	17, 40, 63, 72	1 (0%)
1	K	159/165 (96%)	0.16	4 (2%) 61 59	34, 48, 71, 77	3 (1%)
1	L	159/165 (96%)	-0.08	2 (1%) 79 80	19, 37, 62, 83	3 (1%)
1	M	159/165 (96%)	0.13	7 (4%) 38 37	25, 46, 79, 91	1 (0%)
1	N	159/165 (96%)	0.23	8 (5%) 32 32	23, 47, 72, 96	3 (1%)
1	O	159/165 (96%)	0.10	6 (3%) 44 44	28, 47, 77, 92	2 (1%)
1	P	159/165 (96%)	0.13	4 (2%) 61 59	29, 46, 70, 83	2 (1%)
1	Q	159/165 (96%)	0.24	6 (3%) 44 44	26, 50, 75, 86	2 (1%)
1	R	159/165 (96%)	0.03	3 (1%) 70 69	23, 44, 66, 107	2 (1%)
1	S	159/165 (96%)	0.40	8 (5%) 32 32	31, 57, 86, 99	2 (1%)
1	T	159/165 (96%)	0.43	14 (8%) 12 10	36, 54, 88, 115	1 (0%)
1	U	159/165 (96%)	-0.12	3 (1%) 70 69	17, 37, 55, 103	3 (1%)
1	V	159/165 (96%)	0.53	10 (6%) 23 21	35, 61, 86, 103	0
1	W	159/165 (96%)	0.09	5 (3%) 52 51	23, 40, 67, 74	2 (1%)
1	X	159/165 (96%)	0.03	3 (1%) 70 69	25, 44, 67, 80	2 (1%)

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	Y	159/165 (96%)	-0.13	1 (0%)	90 91	20, 39, 58, 78	3 (1%)
1	Z	159/165 (96%)	0.02	2 (1%)	79 80	27, 44, 63, 77	2 (1%)
1	a	159/165 (96%)	-0.03	2 (1%)	79 80	28, 43, 65, 76	3 (1%)
1	b	159/165 (96%)	-0.11	3 (1%)	70 69	20, 38, 59, 69	2 (1%)
All	All	4452/4620 (96%)	0.11	137 (3%)	52 51	17, 45, 74, 123	54 (1%)

The worst 5 of 137 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	127	THR	7.0
1	T	126	CYS	6.8
1	A	127	THR	6.5
1	V	112	SER	6.5
1	S	62	GLY	6.5

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 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	2AN	R	203	21/21	0.86	0.30	7.50	21,26,39,41	21
2	2AN	Z	202	21/21	0.82	0.29	4.33	21,29,35,38	21
2	2AN	W	203	21/21	0.88	0.32	4.29	16,21,29,33	21
2	2AN	U	202	21/21	0.91	0.25	3.49	24,31,42,48	0
2	2AN	G	204	21/21	0.87	0.24	2.71	22,32,42,45	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	2AN	X	201	21/21	0.94	0.30	2.67	37,41,48,50	0
2	2AN	I	203	21/21	0.96	0.17	2.50	23,32,55,59	0
2	2AN	b	203	21/21	0.91	0.19	2.34	29,43,52,55	0
2	2AN	E	203	21/21	0.93	0.21	2.18	23,32,48,61	0
2	2AN	J	201	21/21	0.92	0.23	2.12	20,23,29,33	21
2	2AN	L	205	21/21	0.93	0.20	1.79	37,44,50,54	0
2	2AN	E	204	21/21	0.88	0.19	1.66	35,46,58,73	0
2	2AN	H	201	21/21	0.94	0.19	1.59	26,31,43,45	0
2	2AN	W	204	21/21	0.84	0.21	1.52	41,53,68,75	0
2	2AN	N	203	21/21	0.86	0.20	1.49	41,48,58,62	0
2	2AN	X	203	21/21	0.94	0.22	1.47	22,31,53,61	0
2	2AN	X	202	21/21	0.89	0.23	1.37	27,32,37,44	21
2	2AN	L	204	21/21	0.91	0.18	1.36	31,40,46,52	0
2	2AN	Z	201	21/21	0.95	0.20	1.35	26,36,46,48	0
2	2AN	R	201	21/21	0.87	0.22	1.19	35,45,68,70	0
2	2AN	S	201	21/21	0.92	0.22	1.18	27,33,38,52	0
2	2AN	G	203	21/21	0.94	0.22	1.17	25,34,56,60	0
2	2AN	b	201	21/21	0.90	0.20	1.16	34,44,49,52	0
2	2AN	J	202	21/21	0.95	0.19	1.11	22,27,39,44	0
2	2AN	E	205	21/21	0.95	0.20	1.03	31,42,54,63	0
2	2AN	R	204	21/21	0.93	0.21	0.98	30,39,58,61	0
2	2AN	E	201	21/21	0.94	0.18	0.97	15,21,28,38	0
2	2AN	Y	202	21/21	0.93	0.17	0.95	27,32,35,38	0
2	2AN	A	201	21/21	0.95	0.18	0.85	25,34,46,71	0
2	2AN	W	201	21/21	0.91	0.24	0.83	35,44,54,63	0
2	2AN	M	202	21/21	0.92	0.18	0.82	26,30,37,43	0
2	2AN	N	201	21/21	0.91	0.22	0.76	24,29,34,37	0
2	2AN	I	201	21/21	0.89	0.18	0.73	21,28,35,42	0
2	2AN	I	204	21/21	0.89	0.19	0.71	42,57,68,81	0
3	EPE	Z	204	15/15	0.94	0.19	0.70	36,44,54,54	0
2	2AN	E	202	21/21	0.89	0.16	0.68	23,31,39,53	0
3	EPE	C	205	15/15	0.93	0.18	0.67	29,34,52,57	0
2	2AN	C	201	21/21	0.91	0.20	0.62	28,39,47,61	0
2	2AN	C	203	21/21	0.92	0.19	0.59	41,51,70,81	0
2	2AN	B	205	21/21	0.93	0.16	0.56	39,54,64,68	0
2	2AN	C	204	21/21	0.94	0.17	0.54	33,39,48,56	0
2	2AN	Q	203	21/21	0.88	0.21	0.53	70,81,96,100	0
2	2AN	P	201	21/21	0.95	0.18	0.53	25,30,38,41	21
2	2AN	b	202	21/21	0.94	0.20	0.50	19,23,34,40	0
2	2AN	N	204	21/21	0.92	0.20	0.50	25,32,44,52	0
2	2AN	O	201	21/21	0.93	0.17	0.49	28,35,40,43	0
2	2AN	M	203	21/21	0.92	0.20	0.47	42,53,70,74	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	2AN	a	201	21/21	0.92	0.17	0.46	25,36,46,49	0
2	2AN	R	202	21/21	0.94	0.18	0.43	21,29,48,53	0
2	2AN	Y	201	21/21	0.95	0.17	0.43	22,24,30,36	0
2	2AN	D	202	21/21	0.95	0.16	0.33	32,39,44,48	0
2	2AN	H	203	21/21	0.96	0.16	0.28	22,28,34,38	0
2	2AN	B	204	21/21	0.93	0.16	0.25	29,40,59,67	0
2	2AN	U	204	21/21	0.95	0.15	0.22	24,26,34,45	0
2	2AN	D	201	21/21	0.96	0.18	0.22	37,45,53,56	0
2	2AN	A	203	21/21	0.93	0.18	0.13	31,43,46,54	0
2	2AN	O	202	21/21	0.96	0.18	0.10	23,28,33,37	0
2	2AN	I	205	21/21	0.90	0.18	-0.00	46,60,65,70	0
2	2AN	F	201	21/21	0.91	0.17	-0.01	31,37,47,56	0
2	2AN	H	202	21/21	0.96	0.17	-0.02	25,36,40,43	0
2	2AN	Z	203	21/21	0.94	0.17	-0.04	35,43,56,64	0
2	2AN	M	201	21/21	0.97	0.17	-0.05	30,35,38,45	0
2	2AN	L	203	21/21	0.97	0.14	-0.05	18,31,40,48	0
2	2AN	K	201	21/21	0.95	0.17	-0.06	28,33,36,42	0
2	2AN	Y	203	21/21	0.90	0.16	-0.07	32,38,47,48	0
2	2AN	G	201	21/21	0.96	0.16	-0.08	15,23,26,34	0
2	2AN	P	202	21/21	0.92	0.15	-0.12	40,51,66,71	0
2	2AN	Y	204	21/21	0.93	0.16	-0.13	30,47,51,57	0
2	2AN	C	202	21/21	0.98	0.15	-0.18	24,39,49,53	0
2	2AN	D	203	21/21	0.90	0.21	-0.19	40,48,58,62	0
2	2AN	L	201	21/21	0.96	0.15	-0.21	19,26,30,37	0
2	2AN	B	203	21/21	0.95	0.15	-0.21	38,44,50,56	0
2	2AN	K	203	21/21	0.92	0.16	-0.22	33,42,66,74	0
2	2AN	I	202	21/21	0.95	0.16	-0.23	19,23,30,38	0
2	2AN	L	202	21/21	0.95	0.14	-0.25	23,28,32,33	0
2	2AN	B	201	21/21	0.96	0.15	-0.29	27,34,44,47	0
2	2AN	A	204	21/21	0.96	0.15	-0.39	35,43,51,54	0
2	2AN	U	203	21/21	0.94	0.14	-0.40	24,38,46,50	0
2	2AN	G	202	21/21	0.97	0.15	-0.40	18,27,34,41	0
2	2AN	U	201	21/21	0.95	0.14	-0.40	28,34,43,47	0
2	2AN	A	202	21/21	0.95	0.17	-0.42	23,35,40,41	0
2	2AN	Q	201	21/21	0.96	0.15	-0.49	30,37,45,51	0
2	2AN	Q	202	21/21	0.94	0.15	-0.56	35,50,64,68	0
2	2AN	W	202	21/21	0.96	0.13	-0.59	24,30,35,40	0
2	2AN	V	202	21/21	0.94	0.17	-0.59	34,46,79,82	0
2	2AN	B	202	21/21	0.94	0.13	-0.60	28,34,53,63	0
2	2AN	N	202	21/21	0.97	0.13	-0.63	25,30,46,54	0
2	2AN	U	205	21/21	0.93	0.15	-0.67	37,47,60,78	0
2	2AN	K	202	21/21	0.93	0.14	-0.73	43,46,53,59	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	SO4	V	203	5/5	0.97	0.14	-0.84	35,48,55,57	0
2	2AN	V	201	21/21	0.94	0.13	-1.22	44,51,56,61	0
2	2AN	G	205	21/21	0.85	0.23	-	52,66,74,85	0
4	SO4	X	204	5/5	0.96	0.10	-	39,42,44,49	0
4	SO4	b	204	5/5	0.97	0.10	-	32,36,40,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.