



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:19 PM GMT

PDB ID : 4EFD
Title : Crystal Structure of an M17 aminopeptidase from Trypanosoma Brucei, Tb427tmp.02.4440
Authors : Wernimont, A.K.; Osman, K.T.; Loppnau, P.; Arrowsmith, C.H.; Edwards, A.M.; Bountra, C.; Hui, R.; Lin, Y.H.; Structural Genomics Consortium (SGC)
Deposited on : 2012-03-29
Resolution : 2.45 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
<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

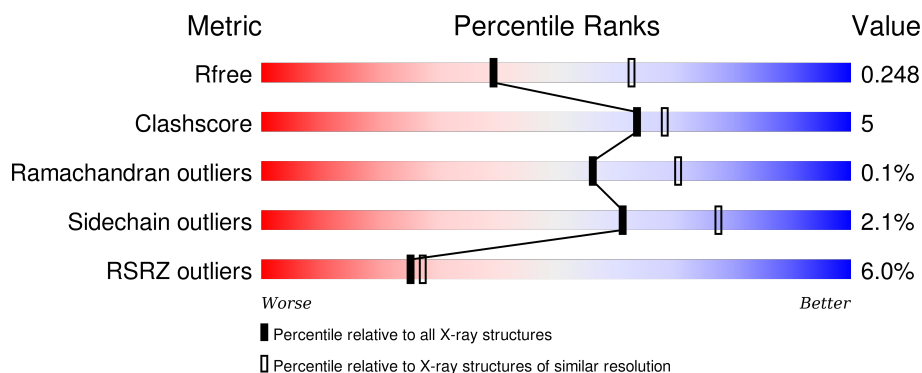
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	4776 (2.50-2.42)
Clashscore	102246	1030 (2.48-2.44)
Ramachandran outliers	100387	1024 (2.48-2.44)
Sidechain outliers	100360	1024 (2.48-2.44)
RSRZ outliers	91569	4787 (2.50-2.42)

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	522	<div> <div>4%</div> <div>88%</div> <div>11%</div> <div>..</div> </div>
1	B	522	<div> <div>7%</div> <div>88%</div> <div>10%</div> <div>..</div> </div>
1	C	522	<div> <div>8%</div> <div>86%</div> <div>12%</div> <div>..</div> </div>
1	D	522	<div> <div>6%</div> <div>89%</div> <div>9%</div> <div>..</div> </div>
1	E	522	<div> <div>5%</div> <div>89%</div> <div>9%</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	522	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	GOL	A	604	-	-	-	X
4	GOL	B	606	-	-	-	X
4	GOL	B	607	-	-	-	X
5	UNX	A	607	-	-	-	X
5	UNX	B	608	-	-	-	X
5	UNX	D	604	-	-	-	X
5	UNX	E	603	-	-	-	X
5	UNX	F	606	-	-	X	X

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 23792 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 Aminopeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	517	Total	C	N	O	S	0	4	0
			3836	2423	656	735	22			
1	B	515	Total	C	N	O	S	0	9	0
			3819	2411	648	739	21			
1	C	514	Total	C	N	O	S	0	9	0
			3808	2405	647	734	22			
1	D	516	Total	C	N	O	S	0	2	0
			3799	2401	643	733	22			
1	E	515	Total	C	N	O	S	0	3	0
			3809	2404	651	732	22			
1	F	516	Total	C	N	O	S	0	5	0
			3827	2419	647	739	22			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	2	Total Mn 2 2	0	0
3	A	1	Total Mn 1 1	0	0
3	D	1	Total Mn 1 1	0	0
3	C	2	Total Mn 2 2	0	0
3	F	2	Total Mn 2 2	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 5 is UNKNOWN ATOM OR ION (three-letter code: UNX) (formula: X).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	D	1	Total X 1 1	0	0
5	E	1	Total X 1 1	0	0
5	B	1	Total X 1 1	0	0
5	C	1	Total X 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total 1	X 1	0	0
5	F	1	Total 1	X 1	0	0

- Molecule 6 is SODIUM ION (three-letter code: NA) (formula: Na).

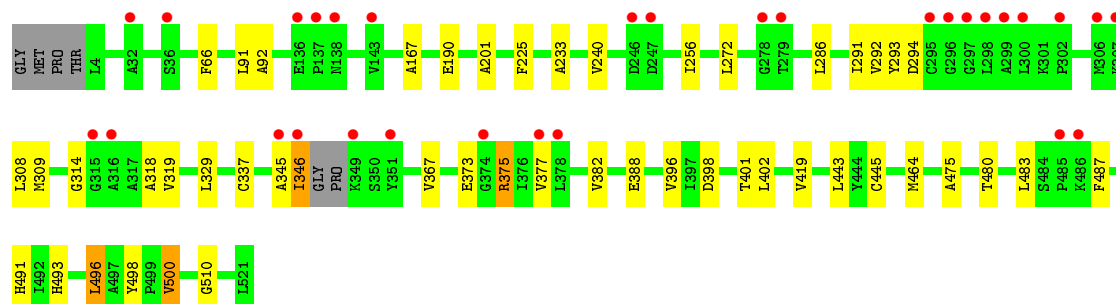
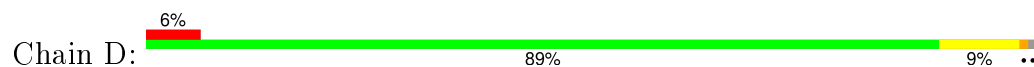
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	1	Total 1	Na 1	0	0
6	F	1	Total 1	Na 1	0	0

- Molecule 7 is water.

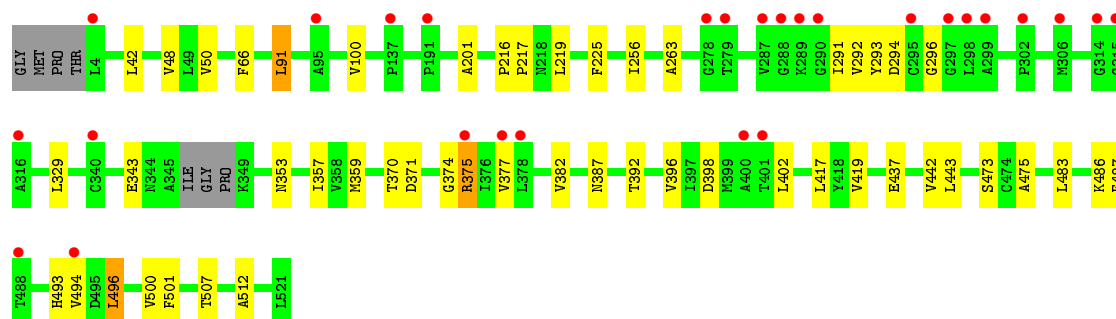
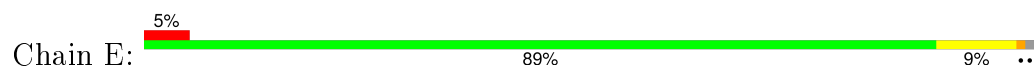
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	100	Total 100	O 100	0	0
7	B	131	Total 131	O 131	0	0
7	C	82	Total 82	O 82	0	0
7	D	97	Total 97	O 97	0	0
7	E	161	Total 162	O 162	0	1
7	F	198	Total 198	O 198	0	0



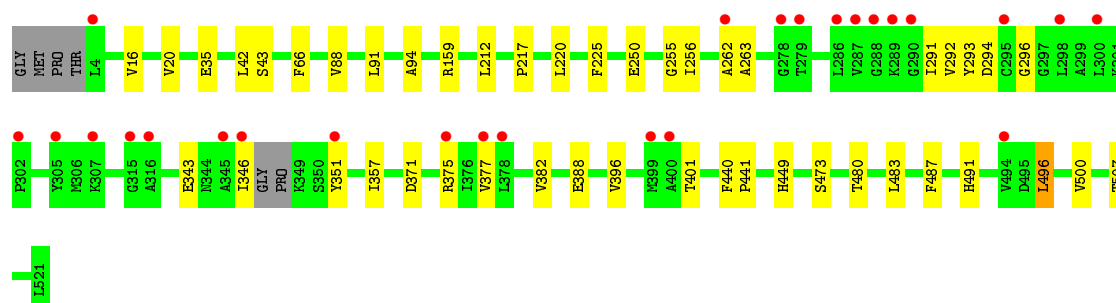
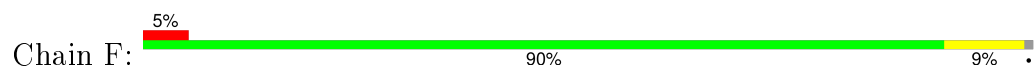
• Molecule 1: Aminopeptidase



• Molecule 1: Aminopeptidase



• Molecule 1: Aminopeptidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	161.92Å 161.74Å 176.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.45 29.87 – 2.45	Depositor EDS
% Data completeness (in resolution range)	99.3 (30.00-2.45) 99.3 (29.87-2.45)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.20 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.227 , 0.253 0.225 , 0.248	Depositor DCC
R_{free} test set	8505 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	46.7	Xtriage
Anisotropy	0.040	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.2	EDS
Estimated twinning fraction	0.015 for k,h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 168491 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	23792	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.03% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: UNX, GOL, MN, SO4, NA

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	0.28	0/3915	0.45	1/5337 (0.0%)
1	B	0.27	0/3917	0.44	1/5343 (0.0%)
1	C	0.27	0/3885	0.43	1/5302 (0.0%)
1	D	0.28	0/3874	0.45	0/5288
1	E	0.28	0/3887	0.45	0/5299
1	F	0.28	0/3905	0.45	0/5323
All	All	0.28	0/23383	0.45	3/31892 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	500	VAL	CG1-CB-CG2	5.52	119.73	110.90
1	C	500	VAL	CG1-CB-CG2	5.29	119.37	110.90
1	A	500	VAL	CG1-CB-CG2	5.01	118.91	110.90

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	3836	0	3746	37	0
1	B	3819	0	3716	41	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	3808	0	3669	52	0
1	D	3799	0	3677	40	0
1	E	3809	0	3713	39	0
1	F	3827	0	3731	32	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
2	C	10	0	0	0	0
2	D	10	0	0	0	0
2	E	10	0	0	0	0
2	F	10	0	0	0	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	1	0	0	0	0
3	F	2	0	0	0	0
4	A	18	0	24	1	0
4	B	18	0	24	0	0
4	C	12	0	16	1	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
5	E	1	0	0	0	0
5	F	1	0	0	2	0
6	C	1	0	0	0	0
6	F	1	0	0	0	0
7	A	100	0	0	0	0
7	B	131	0	0	0	0
7	C	82	0	0	0	0
7	D	97	0	0	0	0
7	E	162	0	0	1	0
7	F	198	0	0	0	0
All	All	23792	0	22316	236	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 5.

All (236) 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:372[B]:ALA:O	1:C:375[B]:ARG:HD3	1.35	1.22

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:375[B]:ARG:CG	1:F:375[B]:ARG:HH11	1.59	1.15
1:F:375[B]:ARG:HH11	1:F:375[B]:ARG:HG2	1.05	1.13
1:E:375[A]:ARG:HH11	1:E:375[A]:ARG:HG2	1.02	1.12
1:B:375[B]:ARG:CD	1:B:375[B]:ARG:H	1.58	1.10
1:C:375[B]:ARG:HD2	1:C:375[B]:ARG:N	1.63	1.08
1:B:375[B]:ARG:HH11	1:B:375[B]:ARG:CG	1.69	1.06
1:B:375[B]:ARG:HG3	1:B:375[B]:ARG:NH1	1.57	1.06
1:E:375[A]:ARG:CD	1:E:375[A]:ARG:H	1.71	1.04
1:D:496:LEU:O	1:D:500:VAL:HG22	1.58	1.04
1:C:496:LEU:O	1:C:500:VAL:HG22	1.58	1.03
1:E:375[A]:ARG:HD2	1:E:375[A]:ARG:H	1.24	1.03
1:D:345:ALA:HB1	1:D:346:ILE:HG23	1.39	1.03
1:B:496:LEU:O	1:B:500:VAL:HG22	1.59	1.02
1:B:375[B]:ARG:H	1:B:375[B]:ARG:HD3	1.26	1.00
1:C:375[B]:ARG:H	1:C:375[B]:ARG:CD	1.74	0.98
1:C:375[B]:ARG:HD2	1:C:375[B]:ARG:H	0.82	0.96
1:E:375[A]:ARG:CG	1:E:375[A]:ARG:HH11	1.77	0.96
1:F:375[B]:ARG:HH12	5:F:606:UNX:UNK	1.02	0.94
1:C:226:THR:HG22	1:C:270:VAL:HG21	1.49	0.94
1:F:375[B]:ARG:HG2	1:F:375[B]:ARG:NH1	1.75	0.93
1:E:375[A]:ARG:HD2	1:E:375[A]:ARG:N	1.84	0.93
1:B:375[B]:ARG:CD	1:B:375[B]:ARG:N	2.35	0.89
1:B:375[B]:ARG:HG3	1:B:375[B]:ARG:HH11	0.77	0.87
1:F:375[B]:ARG:NH1	5:F:606:UNX:UNK	1.70	0.87
1:C:372[B]:ALA:O	1:C:375[B]:ARG:CD	2.22	0.86
1:E:375[A]:ARG:HG2	1:E:375[A]:ARG:NH1	1.84	0.84
1:D:345:ALA:HB1	1:D:346:ILE:CG2	2.07	0.84
1:A:319:VAL:HG12	1:A:337:CYS:SG	2.22	0.80
1:E:357:ILE:HD11	1:E:359:MET:SD	2.21	0.80
1:D:319:VAL:HG12	1:D:337:CYS:SG	2.21	0.79
1:A:240:VAL:HG22	1:A:272:LEU:HD22	1.67	0.76
1:D:496:LEU:O	1:D:500:VAL:CG2	2.34	0.74
1:F:382:VAL:HG13	1:F:483:LEU:HD23	1.67	0.74
1:B:496:LEU:O	1:B:500:VAL:CG2	2.34	0.74
1:C:496:LEU:O	1:C:500:VAL:CG2	2.34	0.74
1:B:375[B]:ARG:N	1:B:375[B]:ARG:HD2	2.03	0.73
1:E:42:LEU:HD22	1:E:100:VAL:HG11	1.71	0.73
1:B:375[B]:ARG:H	1:B:375[B]:ARG:HD2	1.52	0.72
1:C:372[B]:ALA:O	1:C:374[B]:GLY:N	2.27	0.68
1:C:357:ILE:HD11	1:C:359:MET:SD	2.35	0.67
1:D:240:VAL:HG22	1:D:272:LEU:HD22	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:375[B]:ARG:CG	1:F:375[B]:ARG:NH1	2.32	0.66
1:F:382:VAL:CG1	1:F:483:LEU:HD23	2.26	0.65
1:F:375[B]:ARG:HH11	1:F:375[B]:ARG:HG3	1.57	0.64
1:B:382:VAL:HG13	1:B:483:LEU:HD23	1.79	0.64
1:B:357:ILE:HD11	1:B:359:MET:SD	2.37	0.64
1:D:314:GLY:HA3	1:D:500:VAL:HG11	1.80	0.64
1:C:88:VAL:HG11	1:C:91:LEU:HD12	1.80	0.63
1:E:382:VAL:HG13	1:E:483:LEU:HD23	1.80	0.63
1:B:240:VAL:HG22	1:B:272:LEU:HD22	1.80	0.63
1:C:35[B]:GLU:C	1:C:35[B]:GLU:OE1	2.36	0.62
1:A:289:LYS:HE3	1:A:374:GLY:HA3	1.81	0.62
1:A:286:LEU:HB3	1:A:319:VAL:HG13	1.82	0.61
1:A:496:LEU:O	1:A:500:VAL:HG12	2.01	0.60
1:E:375[A]:ARG:CG	1:E:375[A]:ARG:NH1	2.47	0.60
1:C:375[B]:ARG:CD	1:C:375[B]:ARG:N	2.45	0.58
1:E:496:LEU:O	1:E:500:VAL:HG12	2.04	0.57
1:F:496:LEU:O	1:F:500:VAL:HG12	2.04	0.57
1:B:382:VAL:CG1	1:B:483:LEU:HD23	2.34	0.57
1:A:382:VAL:CG1	1:A:483:LEU:HD23	2.35	0.57
1:A:113:ARG:NH1	1:D:419:VAL:O	2.39	0.56
1:C:38:GLN:O	1:C:41:VAL:HG22	2.06	0.56
1:F:396:VAL:HG23	1:F:487:PHE:CE1	2.41	0.56
1:C:314:GLY:HA3	1:C:500:VAL:HG11	1.87	0.55
1:D:345:ALA:CB	1:D:346:ILE:HG23	2.26	0.55
1:A:382:VAL:HG13	1:A:483:LEU:HD23	1.89	0.55
1:F:256:ILE:HA	1:F:377:VAL:HG12	1.89	0.55
1:C:374[B]:GLY:C	1:C:376[B]:ILE:N	2.60	0.55
1:D:382:VAL:CG1	1:D:483:LEU:HD23	2.37	0.55
1:B:42:LEU:HD22	1:B:100:VAL:HG11	1.87	0.54
1:E:396:VAL:HG23	1:E:487:PHE:CE1	2.42	0.54
1:E:375[A]:ARG:HD3	1:E:375[A]:ARG:H	1.65	0.54
1:C:382:VAL:HG13	1:C:483:LEU:HD23	1.89	0.54
1:C:256:ILE:HA	1:C:377:VAL:HG12	1.90	0.54
1:C:376[B]:ILE:HD11	1:C:462:ASN:HB2	1.91	0.53
1:D:66:PHE:HA	1:D:91:LEU:HD11	1.91	0.52
1:E:42:LEU:HD13	1:E:91:LEU:HD22	1.90	0.52
1:A:113:ARG:NH2	4:A:605:GOL:O1	2.42	0.52
1:D:256:ILE:HA	1:D:377:VAL:HG12	1.90	0.52
1:C:66:PHE:HA	1:C:91:LEU:HD11	1.92	0.52
1:F:292:VAL:HG12	1:F:293:TYR:CD2	2.46	0.51
1:D:480:THR:HG22	1:D:491:HIS:CD2	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:374:GLY:O	1:A:377:VAL:HG22	2.11	0.51
1:A:35[B]:GLU:HA	1:A:35[B]:GLU:OE1	2.10	0.51
1:E:292:VAL:HG12	1:E:293:TYR:CD2	2.46	0.51
1:A:42:LEU:HD13	1:A:91:LEU:HD22	1.91	0.51
1:D:401:THR:CG2	1:D:500:VAL:HG21	2.40	0.51
1:A:66:PHE:HA	1:A:91:LEU:HD11	1.93	0.51
1:E:382:VAL:CG1	1:E:483:LEU:HD23	2.41	0.50
1:D:292:VAL:HG12	1:D:293:TYR:CD2	2.46	0.50
1:B:292:VAL:HG12	1:B:293:TYR:CD2	2.46	0.50
1:B:66:PHE:HA	1:B:91:LEU:HD11	1.91	0.50
1:D:319:VAL:CG1	1:D:337:CYS:SG	2.97	0.50
1:F:66:PHE:HA	1:F:91:LEU:HD11	1.94	0.50
1:A:351:TYR:O	1:A:370:THR:HG21	2.12	0.50
1:A:292:VAL:HG12	1:A:293:TYR:CD2	2.46	0.50
1:B:373:GLU:HG2	1:B:376:ILE:HD12	1.94	0.49
1:E:392:THR:HG22	1:E:486:LYS:HE3	1.94	0.49
1:B:396:VAL:HG23	1:B:487:PHE:CE1	2.46	0.49
1:D:402:LEU:HD23	1:D:475:ALA:HB1	1.95	0.49
1:C:374[B]:GLY:C	1:C:376[B]:ILE:H	2.16	0.49
1:B:314:GLY:HA3	1:B:500:VAL:HG11	1.95	0.49
1:C:292:VAL:HG12	1:C:293:TYR:CD2	2.47	0.49
1:C:374[B]:GLY:O	1:C:376[B]:ILE:N	2.46	0.49
1:D:396:VAL:HG23	1:D:487:PHE:CE1	2.47	0.49
1:B:352:ARG:NH2	1:B:373:GLU:OE1	2.45	0.49
1:B:353[B]:ASN:HD22	1:D:309:MET:CE	2.26	0.49
1:B:401:THR:CG2	1:B:500:VAL:HG21	2.42	0.48
1:B:401:THR:HG22	1:B:500:VAL:HG21	1.95	0.48
1:C:401:THR:HG22	1:C:500:VAL:HG21	1.95	0.48
1:C:442:VAL:HG12	1:C:473:SER:OG	2.14	0.48
1:A:217:PRO:HB3	1:A:292:VAL:HG13	1.96	0.48
1:D:401:THR:HG22	1:D:500:VAL:HG21	1.96	0.48
1:D:286:LEU:HB3	1:D:319:VAL:HG13	1.96	0.48
1:D:375:ARG:HG3	1:D:375:ARG:HH11	1.79	0.48
1:C:442:VAL:HG11	1:C:472:VAL:HB	1.95	0.48
1:C:382:VAL:CG1	1:C:483:LEU:HD23	2.45	0.47
1:D:318:ALA:HB1	1:D:510:GLY:HA3	1.96	0.47
1:A:296:GLY:HA3	1:A:371:ASP:HB3	1.96	0.47
1:C:201:ALA:HB3	1:C:329:LEU:HD11	1.95	0.47
1:C:374[B]:GLY:O	1:C:377:VAL:HG22	2.14	0.47
1:D:382:VAL:HG13	1:D:483:LEU:HD23	1.96	0.47
1:E:387:ASN:HA	1:E:486:LYS:HE2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:437:GLU:CD	1:E:512:ALA:HB3	2.35	0.47
1:B:353[B]:ASN:HD22	1:D:309:MET:HE2	1.79	0.47
1:D:240:VAL:HG22	1:D:272:LEU:CD2	2.44	0.47
1:C:35[A]:GLU:HA	1:C:35[A]:GLU:OE2	2.15	0.46
1:A:399:MET:SD	1:A:514:LEU:HD11	2.55	0.46
1:F:42:LEU:HD23	1:F:91:LEU:HD22	1.97	0.46
1:F:262:ALA:HB2	1:F:351:TYR:HE2	1.80	0.46
1:B:364:THR:HG21	1:D:167:ALA:HB1	1.98	0.46
1:F:346:ILE:HD12	1:F:346:ILE:N	2.29	0.46
1:A:151:LEU:HD13	1:A:516:GLU:OE1	2.15	0.46
1:E:48:VAL:HG12	1:E:50:VAL:HG23	1.97	0.46
1:B:296:GLY:HA3	1:B:371:ASP:HB3	1.96	0.46
1:E:217:PRO:HB3	1:E:292:VAL:HG13	1.96	0.46
1:A:256:ILE:HA	1:A:377:VAL:HG12	1.97	0.46
1:A:318:ALA:HB1	1:A:510:GLY:HA3	1.96	0.46
1:C:351:TYR:O	1:C:370:THR:HG21	2.16	0.46
1:E:500:VAL:HG23	7:E:809:HOH:O	2.15	0.46
1:E:216:PRO:HD2	1:E:219:LEU:HD12	1.96	0.46
1:C:226:THR:HG21	1:C:268:ARG:HD3	1.98	0.46
1:B:281:VAL:HG12	1:B:332:PRO:O	2.16	0.45
1:B:201:ALA:HB3	1:B:329:LEU:HD11	1.98	0.45
1:B:374:GLY:O	1:B:377:VAL:HG22	2.15	0.45
1:C:401:THR:CG2	1:C:500:VAL:HG21	2.47	0.45
1:E:66:PHE:HA	1:E:91:LEU:HD11	1.99	0.45
1:C:91:LEU:HD23	1:C:92:ALA:N	2.31	0.45
1:A:465[B]:GLU:HA	1:A:465[B]:GLU:OE1	2.17	0.45
1:C:226:THR:HG21	1:C:268:ARG:CD	2.46	0.45
1:C:217:PRO:HB3	1:C:292:VAL:HG13	1.99	0.45
1:F:296:GLY:HA3	1:F:371:ASP:HB3	1.98	0.45
1:E:501:PHE:HA	1:E:507:THR:HG23	1.99	0.45
1:E:357:ILE:HD12	1:E:357:ILE:C	2.37	0.45
1:D:233:ALA:HB2	1:D:240:VAL:HG21	1.99	0.45
1:E:353:ASN:HD22	1:E:370:THR:HG23	1.82	0.44
1:E:291:ILE:HB	1:E:294:ASP:HB3	1.99	0.44
1:A:51:GLY:HA3	1:A:146:LEU:HD12	1.99	0.44
1:E:256:ILE:HA	1:E:377:VAL:HG12	2.00	0.44
1:D:91:LEU:HD23	1:D:92:ALA:N	2.32	0.44
1:A:442:VAL:HG13	1:A:473:SER:OG	2.17	0.44
1:A:216:PRO:HD2	1:A:219:LEU:HD12	1.99	0.44
1:D:402:LEU:HD22	1:D:493:HIS:CE1	2.52	0.44
1:C:291:ILE:HB	1:C:294:ASP:HB3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:151:LEU:HD13	1:C:516:GLU:OE1	2.17	0.44
1:B:256:ILE:HA	1:B:377:VAL:HG12	2.00	0.43
1:B:291:ILE:HB	1:B:294:ASP:HB3	2.00	0.43
1:A:240:VAL:HG22	1:A:272:LEU:CD2	2.41	0.43
1:E:402:LEU:HD23	1:E:475:ALA:HB1	2.00	0.43
1:B:440:PHE:CD1	1:E:443:LEU:HD23	2.53	0.43
1:E:201:ALA:HB3	1:E:329:LEU:HD11	1.99	0.43
1:D:201:ALA:HB3	1:D:329:LEU:HD11	1.99	0.43
1:F:480:THR:HG22	1:F:491:HIS:ND1	2.33	0.43
1:B:240:VAL:HG22	1:B:272:LEU:CD2	2.48	0.43
1:A:374:GLY:HA2	1:A:377:VAL:HG22	2.01	0.43
1:F:88:VAL:HG11	1:F:91:LEU:HD12	2.00	0.43
1:F:291:ILE:HB	1:F:294:ASP:HB3	2.00	0.43
1:E:263:ALA:CB	1:E:343:GLU:HB3	2.49	0.43
1:C:225:PHE:CE1	1:C:339:LEU:HD13	2.54	0.43
1:A:281:VAL:HG22	1:A:332:PRO:O	2.19	0.43
1:D:291:ILE:HB	1:D:294:ASP:HB3	2.00	0.42
1:F:16:VAL:O	1:F:20:VAL:HG23	2.19	0.42
1:C:296:GLY:HA3	1:C:371:ASP:HB3	2.00	0.42
1:E:442:VAL:HG13	1:E:473:SER:OG	2.18	0.42
1:F:255:GLY:HA2	1:F:357:ILE:HD12	2.01	0.42
1:F:43:SER:O	1:F:94:ALA:HB3	2.20	0.42
1:A:291:ILE:HB	1:A:294:ASP:HB3	2.00	0.42
1:A:396:VAL:HG23	1:A:487:PHE:CE1	2.54	0.42
1:B:357:ILE:C	1:B:357:ILE:HD12	2.40	0.42
1:D:375:ARG:HD2	1:D:402:LEU:HD11	2.02	0.42
1:C:240:VAL:HG23	1:C:272:LEU:HD23	2.01	0.42
1:C:396:VAL:HG23	1:C:487:PHE:CE1	2.54	0.42
1:F:159:ARG:HG3	1:F:507:THR:O	2.20	0.42
1:D:375:ARG:N	1:D:375:ARG:HD3	2.35	0.42
1:D:402:LEU:HD22	1:D:493:HIS:HE1	1.84	0.42
1:E:417:LEU:CD2	1:E:494:VAL:HG22	2.50	0.42
1:D:233:ALA:HB2	1:D:240:VAL:CG2	2.50	0.42
1:D:443:LEU:HD11	1:D:445:CYS:HB3	2.02	0.42
1:B:443:LEU:HD22	1:B:444:TYR:N	2.35	0.42
1:E:392:THR:HG22	1:E:486:LYS:CE	2.50	0.41
1:E:374:GLY:O	1:E:377:VAL:HG22	2.20	0.41
1:D:367:VAL:HG11	1:D:373:GLU:OE2	2.20	0.41
1:B:48:VAL:HG12	1:B:50:VAL:HG23	2.02	0.41
1:F:263:ALA:HB2	1:F:343:GLU:O	2.20	0.41
1:B:429:LEU:HD21	1:B:441:PRO:HD3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:35[B]:GLU:O	1:C:35[B]:GLU:OE1	2.38	0.41
1:B:398:ASP:OD2	1:B:493:HIS:ND1	2.50	0.41
1:C:88:VAL:CG1	1:C:91:LEU:HD12	2.47	0.41
1:F:217:PRO:HB3	1:F:292:VAL:HG13	2.02	0.41
1:A:42:LEU:HD22	1:A:100:VAL:HG11	2.02	0.41
1:F:401:THR:CG2	1:F:500:VAL:HG11	2.50	0.41
1:D:308:LEU:HD21	1:D:498:TYR:CZ	2.56	0.41
1:C:42:LEU:HD22	1:C:100:VAL:HG11	2.02	0.41
1:C:480:THR:HG22	1:C:491:HIS:CE1	2.56	0.41
1:E:296:GLY:HA3	1:E:371:ASP:HB3	2.02	0.41
1:B:442:VAL:HG21	1:B:472:VAL:HG11	2.01	0.41
1:A:335:LEU:HD11	1:A:518:PHE:CZ	2.56	0.41
1:A:480:THR:HG22	1:A:491:HIS:CD2	2.55	0.41
1:A:432:GLY:HA2	1:A:513:LEU:HD13	2.02	0.41
1:F:440:PHE:CD1	1:F:441:PRO:HD2	2.55	0.41
1:C:42:LEU:HD13	1:C:91:LEU:CD2	2.51	0.41
1:A:91:LEU:HD23	1:A:92:ALA:N	2.36	0.41
1:A:402:LEU:CD2	1:A:475:ALA:HB1	2.51	0.41
1:C:320:PHE:CE2	1:C:324:LEU:HD11	2.56	0.41
1:C:113[A]:ARG:NH2	4:C:607:GOL:O3	2.54	0.41
1:E:402:LEU:HD22	1:E:493:HIS:CE1	2.56	0.41
1:D:294:ASP:OD2	1:D:373:GLU:O	2.39	0.41
1:F:440:PHE:CG	1:F:441:PRO:HD2	2.56	0.40
1:F:449:HIS:CE1	1:F:473:SER:HB3	2.56	0.40
1:C:12:LEU:O	1:C:16:VAL:HG23	2.21	0.40
1:F:212:LEU:O	1:F:220:LEU:HD22	2.21	0.40
1:C:417:LEU:CD2	1:C:494:VAL:HG22	2.51	0.40
1:B:449:HIS:CE1	1:B:473:SER:HB3	2.56	0.40
1:C:480:THR:HG22	1:C:491:HIS:ND1	2.36	0.40
1:A:375[A]:ARG:HG2	1:A:375[A]:ARG:H	1.44	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	517/522 (99%)	505 (98%)	12 (2%)	0	100	100
1	B	520/522 (100%)	505 (97%)	15 (3%)	0	100	100
1	C	517/522 (99%)	498 (96%)	13 (2%)	6 (1%)	16	17
1	D	514/522 (98%)	501 (98%)	13 (2%)	0	100	100
1	E	514/522 (98%)	502 (98%)	12 (2%)	0	100	100
1	F	517/522 (99%)	498 (96%)	19 (4%)	0	100	100
All	All	3099/3132 (99%)	3009 (97%)	84 (3%)	6 (0%)	56	64

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	373[A]	GLU
1	C	373[B]	GLU
1	C	375[A]	ARG
1	C	375[B]	ARG
1	C	374[A]	GLY
1	C	374[B]	GLY

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	394/411 (96%)	386 (98%)	8 (2%)	63	78
1	B	393/411 (96%)	380 (97%)	13 (3%)	45	61
1	C	384/411 (93%)	375 (98%)	9 (2%)	58	74
1	D	386/411 (94%)	376 (97%)	10 (3%)	54	70
1	E	391/411 (95%)	384 (98%)	7 (2%)	66	81
1	F	393/411 (96%)	388 (99%)	5 (1%)	76	86
All	All	2341/2466 (95%)	2289 (98%)	52 (2%)	61	75

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62	GLN
1	A	225	PHE
1	A	244	CYS
1	A	375[A]	ARG
1	A	375[B]	ARG
1	A	388	GLU
1	A	442	VAL
1	A	496	LEU
1	B	35	GLU
1	B	91	LEU
1	B	225	PHE
1	B	250	GLU
1	B	312[A]	ASP
1	B	312[B]	ASP
1	B	388	GLU
1	B	398	ASP
1	B	419	VAL
1	B	442	VAL
1	B	467	ARG
1	B	496	LEU
1	B	500	VAL
1	C	250	GLU
1	C	282	LYS
1	C	353	ASN
1	C	388	GLU
1	C	398	ASP
1	C	419	VAL
1	C	465	GLU
1	C	496	LEU
1	C	500	VAL
1	D	190[A]	GLU
1	D	190[B]	GLU
1	D	225	PHE
1	D	346	ILE
1	D	375	ARG
1	D	388	GLU
1	D	398	ASP
1	D	464	MET
1	D	496	LEU
1	D	500	VAL
1	E	91	LEU
1	E	225	PHE

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Mol	Chain	Res	Type
1	E	375[A]	ARG
1	E	375[B]	ARG
1	E	398	ASP
1	E	419	VAL
1	E	496	LEU
1	F	35	GLU
1	F	225	PHE
1	F	250	GLU
1	F	388	GLU
1	F	496	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	GLN
1	A	148	ASN
1	A	311	HIS
1	A	334	GLN
1	A	353	ASN
1	A	491	HIS
1	B	334	GLN
1	B	504	ASN
1	C	62	GLN
1	C	148	ASN
1	C	230	GLN
1	C	334	GLN
1	D	230	GLN
1	D	334	GLN
1	D	491	HIS
1	E	334	GLN
1	E	353	ASN
1	E	449	HIS
1	F	207	GLN
1	F	334	GLN
1	F	504	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 36 ligands modelled in this entry, 6 are unknown and 10 are monoatomic - leaving 20 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	601	-	4,4,4	0.22	0	6,6,6	0.10	0
2	SO4	A	602	-	4,4,4	0.23	0	6,6,6	0.08	0
4	GOL	A	604	-	5,5,5	0.33	0	5,5,5	0.30	0
4	GOL	A	605	-	5,5,5	0.38	0	5,5,5	0.21	0
4	GOL	A	606	-	5,5,5	0.35	0	5,5,5	0.21	0
2	SO4	B	601	-	4,4,4	0.23	0	6,6,6	0.08	0
2	SO4	B	602	-	4,4,4	0.23	0	6,6,6	0.09	0
4	GOL	B	605	-	5,5,5	0.35	0	5,5,5	0.26	0
4	GOL	B	606	-	5,5,5	0.36	0	5,5,5	0.19	0
4	GOL	B	607	-	5,5,5	0.38	0	5,5,5	0.09	0
2	SO4	C	601	-	4,4,4	0.24	0	6,6,6	0.09	0
2	SO4	C	602	-	4,4,4	0.23	0	6,6,6	0.07	0
4	GOL	C	606	-	5,5,5	0.35	0	5,5,5	0.14	0
4	GOL	C	607	-	5,5,5	0.32	0	5,5,5	0.26	0
2	SO4	D	601	-	4,4,4	0.23	0	6,6,6	0.07	0
2	SO4	D	602	-	4,4,4	0.24	0	6,6,6	0.06	0
2	SO4	E	601	-	4,4,4	0.24	0	6,6,6	0.09	0
2	SO4	E	602	-	4,4,4	0.24	0	6,6,6	0.09	0
2	SO4	F	601	-	4,4,4	0.25	0	6,6,6	0.07	0
2	SO4	F	602	-	4,4,4	0.21	0	6,6,6	0.08	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SO4	A	601	-	-	0/0/0/0	0/0/0/0
2	SO4	A	602	-	-	0/0/0/0	0/0/0/0
4	GOL	A	604	-	-	0/4/4/4	0/0/0/0
4	GOL	A	605	-	-	0/4/4/4	0/0/0/0
4	GOL	A	606	-	-	0/4/4/4	0/0/0/0
2	SO4	B	601	-	-	0/0/0/0	0/0/0/0
2	SO4	B	602	-	-	0/0/0/0	0/0/0/0
4	GOL	B	605	-	-	0/4/4/4	0/0/0/0
4	GOL	B	606	-	-	0/4/4/4	0/0/0/0
4	GOL	B	607	-	-	0/4/4/4	0/0/0/0
2	SO4	C	601	-	-	0/0/0/0	0/0/0/0
2	SO4	C	602	-	-	0/0/0/0	0/0/0/0
4	GOL	C	606	-	-	0/4/4/4	0/0/0/0
4	GOL	C	607	-	-	0/4/4/4	0/0/0/0
2	SO4	D	601	-	-	0/0/0/0	0/0/0/0
2	SO4	D	602	-	-	0/0/0/0	0/0/0/0
2	SO4	E	601	-	-	0/0/0/0	0/0/0/0
2	SO4	E	602	-	-	0/0/0/0	0/0/0/0
2	SO4	F	601	-	-	0/0/0/0	0/0/0/0
2	SO4	F	602	-	-	0/0/0/0	0/0/0/0

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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	605	GOL	1	0
4	C	607	GOL	1	0

5.7 Other polymers ⓘ

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 ⓘ

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	517/522 (99%)	0.22	20 (3%)	43 47	35, 48, 59, 65	0
1	B	515/522 (98%)	0.27	39 (7%)	17 17	31, 43, 54, 65	0
1	C	514/522 (98%)	0.39	43 (8%)	14 14	34, 53, 62, 66	0
1	D	516/522 (98%)	0.22	30 (5%)	26 29	35, 47, 58, 65	0
1	E	515/522 (98%)	0.19	27 (5%)	31 34	30, 42, 53, 65	0
1	F	516/522 (98%)	0.07	26 (5%)	32 36	26, 37, 49, 65	0
All	All	3093/3132 (98%)	0.23	185 (5%)	25 27	26, 44, 58, 66	0

All (185) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	298	LEU	7.5
1	A	302	PRO	5.6
1	B	302	PRO	5.5
1	C	296	GLY	5.4
1	F	346	ILE	4.7
1	E	137	PRO	4.7
1	E	302	PRO	4.7
1	D	300	LEU	4.7
1	D	298	LEU	4.5
1	D	302	PRO	4.5
1	A	348	PRO	4.3
1	B	297	GLY	4.2
1	D	351	TYR	4.1
1	B	303	ALA	4.1
1	B	351	TYR	4.1
1	E	278	GLY	4.0
1	E	279	THR	4.0
1	C	295	CYS	4.0
1	D	307	LYS	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	315	GLY	3.8
1	F	288	GLY	3.7
1	E	288	GLY	3.6
1	C	316	ALA	3.6
1	B	312[A]	ASP	3.5
1	F	287	VAL	3.5
1	C	278	GLY	3.4
1	F	290	GLY	3.4
1	B	318	ALA	3.4
1	E	287	VAL	3.4
1	E	295	CYS	3.4
1	E	306	MET	3.3
1	D	295	CYS	3.3
1	E	375[A]	ARG	3.3
1	A	290	GLY	3.3
1	C	319	VAL	3.3
1	C	288	GLY	3.2
1	D	297	GLY	3.2
1	B	319	VAL	3.2
1	D	346	ILE	3.2
1	C	375[A]	ARG	3.2
1	B	137	PRO	3.2
1	C	488	THR	3.2
1	A	351	TYR	3.2
1	E	315	GLY	3.1
1	D	246	ASP	3.1
1	E	298	LEU	3.1
1	C	297	GLY	3.1
1	B	304	ASP	3.1
1	C	4	LEU	3.0
1	E	4	LEU	3.0
1	E	289	LYS	3.0
1	C	137	PRO	3.0
1	B	143	VAL	3.0
1	D	279	THR	3.0
1	C	305	TYR	2.9
1	E	290	GLY	2.9
1	B	299	ALA	2.9
1	F	345	ALA	2.9
1	D	377	VAL	2.9
1	F	278	GLY	2.9
1	A	4	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	F	298	LEU	2.9
1	F	302	PRO	2.9
1	B	316	ALA	2.9
1	D	374	GLY	2.9
1	E	299	ALA	2.9
1	C	304	ASP	2.9
1	C	279	THR	2.8
1	B	288	GLY	2.8
1	B	296	GLY	2.8
1	F	400	ALA	2.8
1	B	236	LEU	2.8
1	C	300	LEU	2.8
1	D	315	GLY	2.8
1	C	346	ILE	2.8
1	F	494	VAL	2.8
1	C	299	ALA	2.8
1	C	290	GLY	2.7
1	C	246	ASP	2.7
1	A	297	GLY	2.7
1	B	250	GLU	2.7
1	D	137	PRO	2.7
1	F	378	LEU	2.7
1	C	312[A]	ASP	2.7
1	F	289	LYS	2.7
1	A	295	CYS	2.7
1	D	306	MET	2.7
1	D	278	GLY	2.6
1	E	377	VAL	2.6
1	F	375[A]	ARG	2.6
1	C	244	CYS	2.6
1	C	307	LYS	2.6
1	B	279	THR	2.6
1	D	138	ASN	2.6
1	B	295	CYS	2.6
1	E	95	ALA	2.6
1	B	48	VAL	2.6
1	C	289	LYS	2.6
1	E	297	GLY	2.6
1	C	318	ALA	2.6
1	D	345	ALA	2.6
1	D	136	GLU	2.5
1	A	306	MET	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	350	SER	2.5
1	A	378	LEU	2.5
1	C	351	TYR	2.5
1	B	246	ASP	2.5
1	D	486	LYS	2.5
1	B	340	CYS	2.5
1	D	316	ALA	2.5
1	D	36	SER	2.5
1	C	302	PRO	2.5
1	C	378	LEU	2.5
1	F	300	LEU	2.5
1	E	314	GLY	2.4
1	E	494	VAL	2.4
1	A	288	GLY	2.4
1	F	279	THR	2.4
1	A	374	GLY	2.4
1	C	298	LEU	2.4
1	A	279	THR	2.4
1	D	349	LYS	2.4
1	D	296	GLY	2.4
1	F	286	LEU	2.4
1	B	289	LYS	2.4
1	C	309	MET	2.4
1	E	316	ALA	2.4
1	F	399	MET	2.4
1	E	340	CYS	2.3
1	B	399	MET	2.3
1	B	136	GLU	2.3
1	B	467	ARG	2.3
1	E	378	LEU	2.3
1	F	377	VAL	2.3
1	B	430	ARG	2.3
1	C	247	ASP	2.3
1	B	300	LEU	2.3
1	A	287	VAL	2.3
1	D	247	ASP	2.3
1	C	37	ASN	2.3
1	B	5	PRO	2.3
1	C	234	LYS	2.3
1	A	278	GLY	2.3
1	D	32	ALA	2.2
1	B	290	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	F	315	GLY	2.2
1	C	486	LYS	2.2
1	C	139	THR	2.2
1	E	488	THR	2.2
1	C	337	CYS	2.2
1	C	317	ALA	2.2
1	F	307	LYS	2.2
1	B	286	LEU	2.2
1	D	378	LEU	2.2
1	C	275	THR	2.2
1	A	301	LYS	2.2
1	F	305	TYR	2.1
1	F	351	TYR	2.1
1	B	145	VAL	2.1
1	C	90	GLY	2.1
1	E	400	ALA	2.1
1	F	316	ALA	2.1
1	C	276	PRO	2.1
1	D	485	PRO	2.1
1	F	295	CYS	2.1
1	A	494	VAL	2.1
1	B	494	VAL	2.1
1	D	299	ALA	2.1
1	C	286	LEU	2.1
1	B	278	GLY	2.1
1	C	287	VAL	2.1
1	D	143	VAL	2.1
1	B	378	LEU	2.1
1	E	191	PRO	2.1
1	A	316	ALA	2.1
1	B	154	ALA	2.0
1	A	298	LEU	2.0
1	E	401	THR	2.0
1	B	317	ALA	2.0
1	F	262	ALA	2.0
1	A	192	PRO	2.0
1	C	280	PRO	2.0
1	B	206	THR	2.0
1	F	4	LEU	2.0
1	A	349	LYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no 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
5	UNX	A	607	1/1	0.84	0.85	9.12	30,30,30,30	0
4	GOL	B	606	6/6	0.92	0.24	6.85	52,52,52,52	0
5	UNX	D	604	1/1	0.70	0.59	6.40	30,30,30,30	0
4	GOL	B	607	6/6	0.70	0.32	6.37	73,73,73,73	0
5	UNX	B	608	1/1	0.35	0.71	5.96	30,30,30,30	0
4	GOL	A	604	6/6	0.71	0.28	5.12	72,72,72,72	0
5	UNX	F	606	1/1	0.63	0.57	3.48	30,30,30,30	0
5	UNX	E	603	1/1	0.75	0.48	2.01	30,30,30,30	0
4	GOL	A	605	6/6	0.83	0.22	1.57	68,68,68,68	0
4	GOL	C	607	6/6	0.84	0.20	1.20	61,61,62,62	0
4	GOL	C	606	6/6	0.80	0.23	1.10	62,62,62,62	0
5	UNX	C	608	1/1	0.44	0.39	0.95	30,30,30,30	0
2	SO4	F	601	5/5	0.94	0.15	-0.71	64,64,64,64	0
2	SO4	A	601	5/5	0.96	0.14	-0.89	66,66,66,66	0
2	SO4	B	602	5/5	0.96	0.10	-1.29	74,74,74,74	0
2	SO4	D	602	5/5	0.95	0.11	-1.36	68,68,68,68	0
2	SO4	C	601	5/5	0.96	0.09	-1.48	70,70,70,70	0
2	SO4	E	601	5/5	0.98	0.09	-1.87	59,59,59,59	0
2	SO4	E	602	5/5	0.85	0.31	-	98,98,98,98	0
2	SO4	C	602	5/5	0.89	0.22	-	103,103,103,103	0
4	GOL	A	606	6/6	0.70	0.40	-	59,59,59,59	0
6	NA	C	603	1/1	0.88	0.17	-	67,67,67,67	0
2	SO4	D	601	5/5	0.91	0.15	-	104,104,104,104	0
3	MN	C	604	1/1	0.95	0.07	-	82,82,82,82	0
2	SO4	B	601	5/5	0.90	0.17	-	92,92,92,92	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	MN	D	603	1/1	0.91	0.04	-	94,94,94,94	0
3	MN	F	603	1/1	0.98	0.26	-	85,85,85,85	0
6	NA	F	605	1/1	0.84	0.14	-	47,47,47,47	0
3	MN	B	604	1/1	0.94	0.10	-	80,80,80,80	0
3	MN	C	605	1/1	0.95	0.16	-	100,100,100,100	0
2	SO4	F	602	5/5	0.92	0.23	-	76,76,76,76	0
4	GOL	B	605	6/6	0.81	0.19	-	82,82,82,82	0
3	MN	B	603	1/1	0.85	0.06	-	92,92,92,92	0
3	MN	F	604	1/1	0.86	0.09	-	99,99,99,99	0
2	SO4	A	602	5/5	0.93	0.23	-	93,93,94,94	0
3	MN	A	603	1/1	0.84	0.08	-	95,95,95,95	0

6.5 Other polymers [i](#)

There are no such residues in this entry.