Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 10:12 AM GMT

PDB ID	:	3L7J
Title	:	Structure of the Wall Teichoic Acid Polymerase TagF, H444N variant
Authors	:	Strynadka, N.C.J.; Lovering, A.L.
Deposited on	:	2009-12-28
Resolution	:	2.81 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

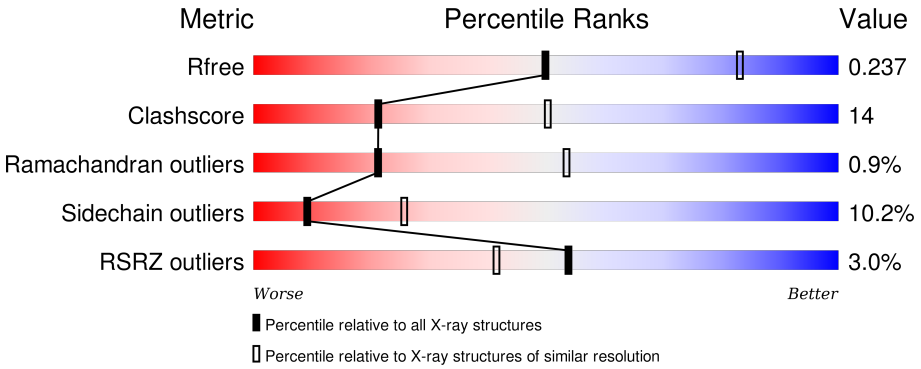
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.81 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	2676 (2.84-2.80)
Clashscore	102246	3124 (2.84-2.80)
Ramachandran outliers	100387	3072 (2.84-2.80)
Sidechain outliers	100360	3074 (2.84-2.80)
RSRZ outliers	91569	2690 (2.84-2.80)

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 <=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	729	39% 15% 44%
1	B	729	35% 18% 43%
1	C	729	38% 16% 44%
1	D	729	37% 16% 44%

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	A	732	-	-	-	X
2	SO4	B	730	-	-	X	-
3	CL	B	732	-	-	X	-
3	CL	B	736	-	-	X	-
3	CL	C	733	-	-	X	-
3	CL	D	731	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13970 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 Teichoic acid biosynthesis protein F.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	410	Total	C	N	O	S	0	0	0
			3462	2224	577	650	11			
1	B	412	Total	C	N	O	S	0	0	0
			3477	2233	581	652	11			
1	C	411	Total	C	N	O	S	0	0	0
			3467	2227	578	651	11			
1	D	411	Total	C	N	O	S	0	0	0
			3472	2230	580	651	11			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	444	ASN	HIS	ENGINEERED	UNP Q5HLM5
A	722	LEU	-	EXPRESSION TAG	UNP Q5HLM5
A	723	GLU	-	EXPRESSION TAG	UNP Q5HLM5
A	724	HIS	-	EXPRESSION TAG	UNP Q5HLM5
A	725	HIS	-	EXPRESSION TAG	UNP Q5HLM5
A	726	HIS	-	EXPRESSION TAG	UNP Q5HLM5
A	727	HIS	-	EXPRESSION TAG	UNP Q5HLM5
A	728	HIS	-	EXPRESSION TAG	UNP Q5HLM5
A	729	HIS	-	EXPRESSION TAG	UNP Q5HLM5
B	444	ASN	HIS	ENGINEERED	UNP Q5HLM5
B	722	LEU	-	EXPRESSION TAG	UNP Q5HLM5
B	723	GLU	-	EXPRESSION TAG	UNP Q5HLM5
B	724	HIS	-	EXPRESSION TAG	UNP Q5HLM5
B	725	HIS	-	EXPRESSION TAG	UNP Q5HLM5
B	726	HIS	-	EXPRESSION TAG	UNP Q5HLM5
B	727	HIS	-	EXPRESSION TAG	UNP Q5HLM5
B	728	HIS	-	EXPRESSION TAG	UNP Q5HLM5
B	729	HIS	-	EXPRESSION TAG	UNP Q5HLM5
C	444	ASN	HIS	ENGINEERED	UNP Q5HLM5
C	722	LEU	-	EXPRESSION TAG	UNP Q5HLM5
C	723	GLU	-	EXPRESSION TAG	UNP Q5HLM5

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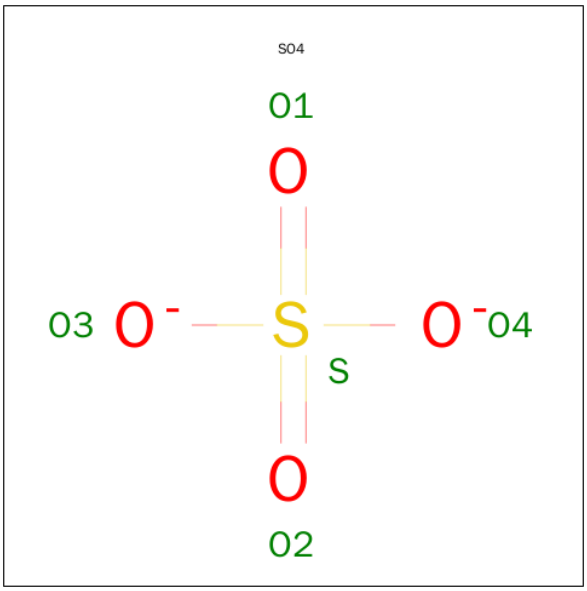
Chain	Residue	Modelled	Actual	Comment	Reference
C	724	HIS	-	EXPRESSION TAG	UNP Q5HLM5
C	725	HIS	-	EXPRESSION TAG	UNP Q5HLM5
C	726	HIS	-	EXPRESSION TAG	UNP Q5HLM5
C	727	HIS	-	EXPRESSION TAG	UNP Q5HLM5
C	728	HIS	-	EXPRESSION TAG	UNP Q5HLM5
C	729	HIS	-	EXPRESSION TAG	UNP Q5HLM5
D	444	ASN	HIS	ENGINEERED	UNP Q5HLM5
D	722	LEU	-	EXPRESSION TAG	UNP Q5HLM5
D	723	GLU	-	EXPRESSION TAG	UNP Q5HLM5
D	724	HIS	-	EXPRESSION TAG	UNP Q5HLM5
D	725	HIS	-	EXPRESSION TAG	UNP Q5HLM5
D	726	HIS	-	EXPRESSION TAG	UNP Q5HLM5
D	727	HIS	-	EXPRESSION TAG	UNP Q5HLM5
D	728	HIS	-	EXPRESSION TAG	UNP Q5HLM5
D	729	HIS	-	EXPRESSION TAG	UNP Q5HLM5

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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	A	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	C	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

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	8	Total Cl 8 8	0	0
3	A	11	Total Cl 11 11	0	0
3	D	6	Total Cl 6 6	0	0
3	C	7	Total Cl 7 7	0	0

- Molecule 4 is water.

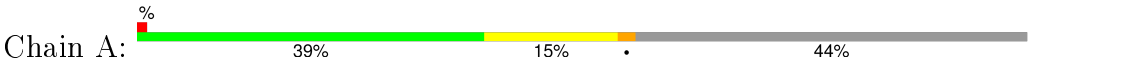
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	8	Total O 8 8	0	0
4	B	5	Total O 5 5	0	0
4	C	3	Total O 3 3	0	0
4	D	4	Total O 4 4	0	0

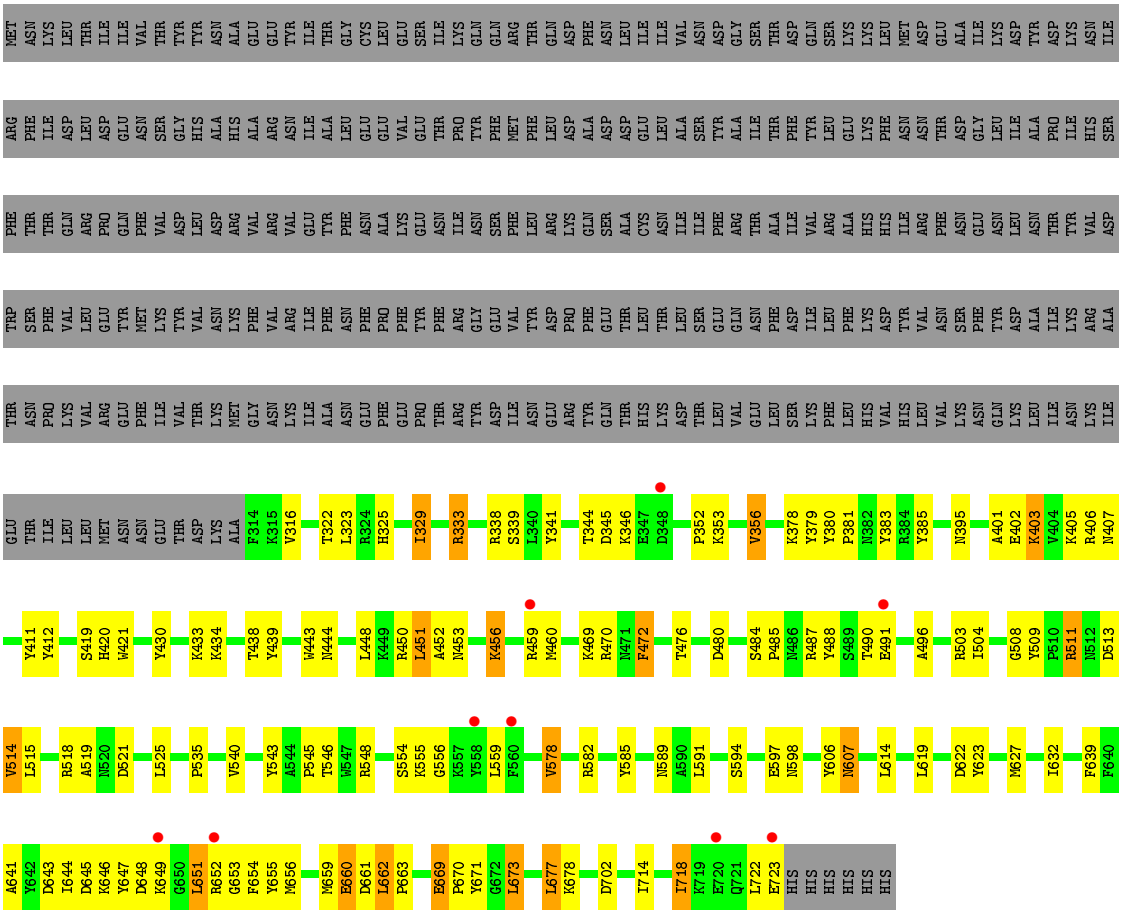
3 Residue-property plots (i)

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

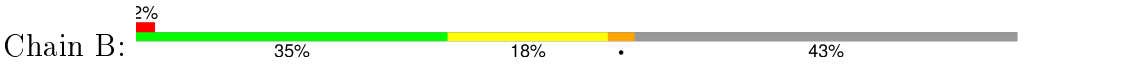
Chain A:

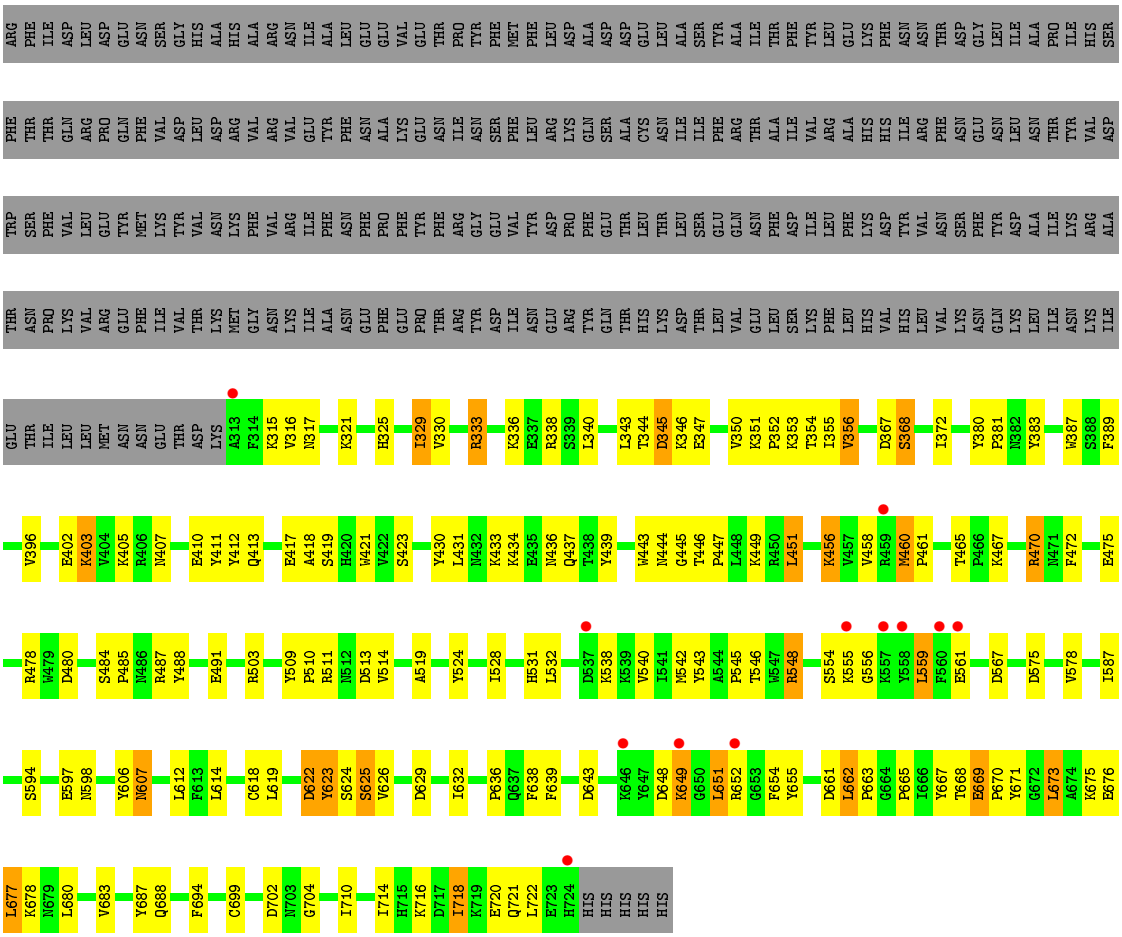




- Molecule 1: Teichoic acid biosynthesis protein F

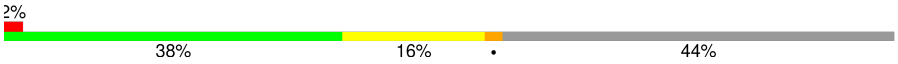
Chain B:

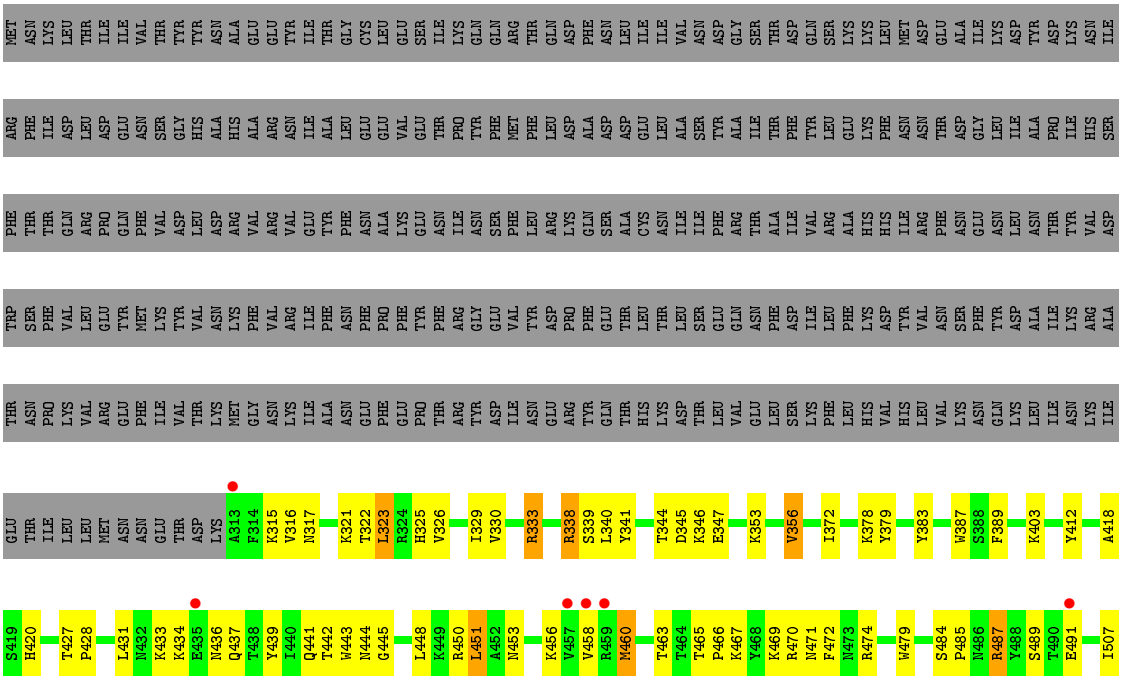


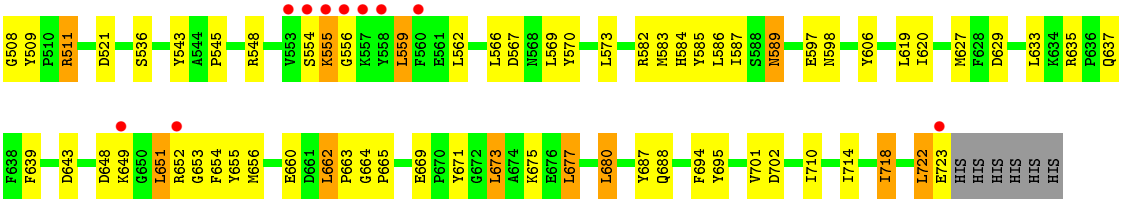


• Molecule 1: Teichoic acid biosynthesis protein F

Chain C:

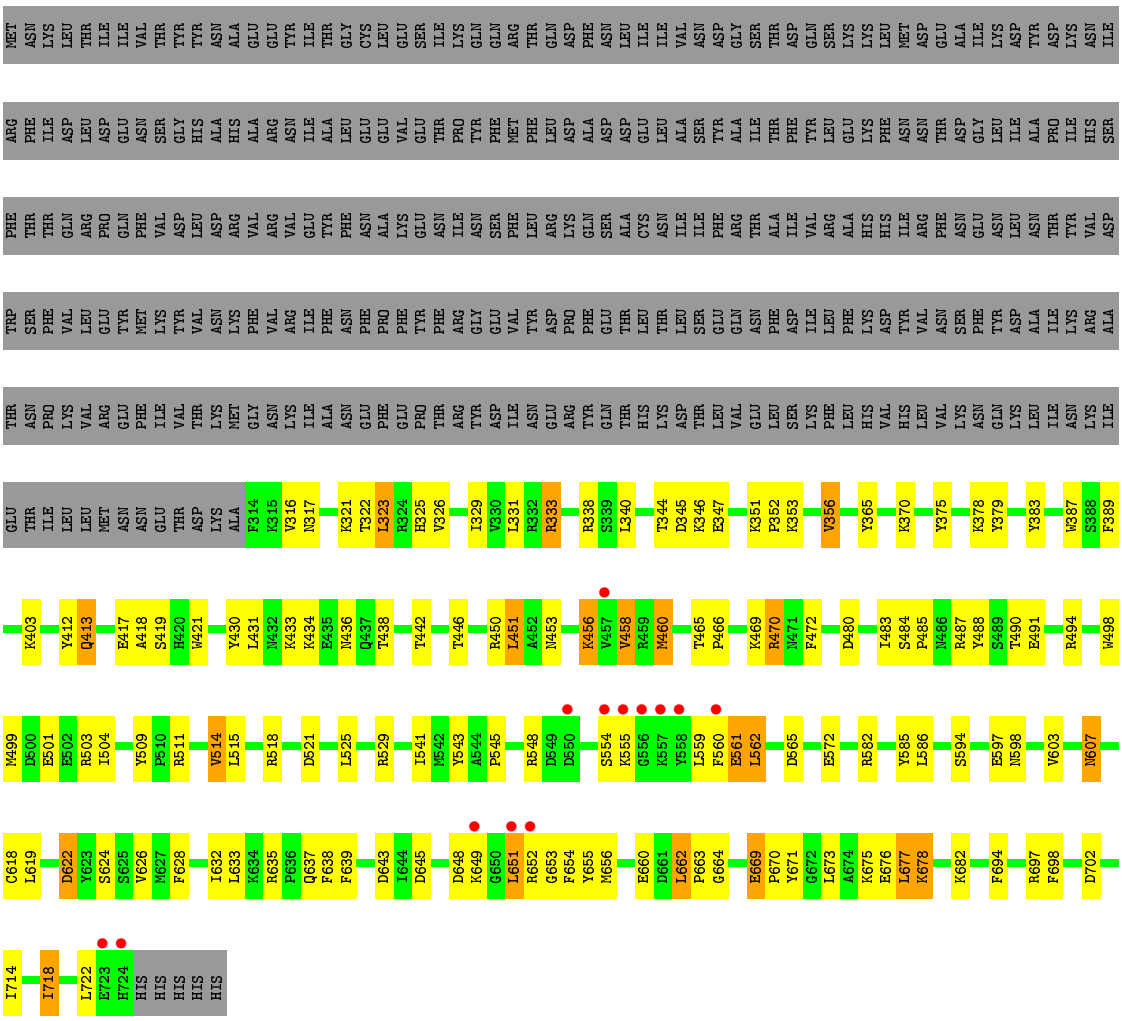






• Molecule 1: Teichoic acid biosynthesis protein F

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α, β, γ	223.60Å 223.60Å 100.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.80 – 2.81 19.80 – 2.81	Depositor EDS
% Data completeness (in resolution range)	97.9 (19.80-2.81) 97.9 (19.80-2.81)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	1.64 (at 2.79Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.197 , 0.246 0.189 , 0.237	Depositor DCC
R_{free} test set	3045 reflections (5.00%)	DCC
Wilson B-factor (Å^2)	72.4	Xtriage
Anisotropy	0.048	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.27 , 29.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	0 of 60941 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	13970	wwPDB-VP
Average B, all atoms (Å^2)	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.40% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| > 5$	RMSZ	$\#\|Z\| > 5$
1	A	0.48	0/3551	0.62	0/4798
1	B	0.49	0/3567	0.65	0/4820
1	C	0.46	0/3556	0.61	0/4805
1	D	0.46	0/3562	0.61	0/4813
All	All	0.47	0/14236	0.62	0/19236

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3462	0	3369	86	0
1	B	3477	0	3381	105	0
1	C	3467	0	3374	104	0
1	D	3472	0	3376	96	0
2	A	15	0	0	0	0
2	B	5	0	0	2	0
2	C	15	0	0	0	0
2	D	5	0	0	0	0
3	A	11	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	8	0	0	6	0
3	C	7	0	0	2	0
3	D	6	0	0	4	0
4	A	8	0	0	1	0
4	B	5	0	0	0	0
4	C	3	0	0	0	0
4	D	4	0	0	0	0
All	All	13970	0	13500	389	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 14.

All (389) 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:A:333:ARG:HG2	1:A:333:ARG:HH11	1.02	1.19
1:D:470:ARG:HG2	1:D:470:ARG:HH11	1.04	1.10
1:C:333:ARG:HG2	1:C:333:ARG:HH11	1.28	0.99
1:B:333:ARG:HH11	1:B:333:ARG:HG2	1.28	0.98
1:C:470:ARG:HG2	1:C:470:ARG:HH11	1.27	0.97
1:B:470:ARG:HG2	1:B:470:ARG:HH11	1.25	0.97
1:D:333:ARG:HH11	1:D:333:ARG:HG2	1.35	0.91
1:A:470:ARG:HH11	1:A:470:ARG:HG2	1.35	0.91
1:D:470:ARG:HG2	1:D:470:ARG:NH1	1.85	0.89
1:A:333:ARG:NH1	1:A:333:ARG:HG2	1.84	0.86
1:A:333:ARG:CG	1:A:333:ARG:HH11	1.87	0.84
1:A:648:ASP:HB2	1:A:651:LEU:HB2	1.64	0.80
1:B:663:PRO:HG3	1:B:694:PHE:CG	2.17	0.79
1:B:413:GLN:O	1:B:417:GLU:HG3	1.83	0.79
1:C:648:ASP:HB2	1:C:651:LEU:HB2	1.63	0.79
1:C:333:ARG:CG	1:C:333:ARG:HH11	1.95	0.78
1:B:548:ARG:HD3	1:B:643:ASP:OD2	1.83	0.78
1:C:383:TYR:CE2	1:C:718:ILE:HD11	2.20	0.77
1:A:383:TYR:CE2	1:A:718:ILE:HD11	2.20	0.77
1:C:662:LEU:H	1:C:662:LEU:HD23	1.51	0.75
1:A:485:PRO:HG3	1:A:509:TYR:CE2	2.20	0.75
1:B:561:GLU:HB2	3:B:734:CL:CL	2.24	0.74
1:B:470:ARG:CG	1:B:470:ARG:HH11	2.00	0.74
1:A:662:LEU:HB2	1:A:663:PRO:HD3	1.68	0.74
1:C:383:TYR:CZ	1:C:718:ILE:HD11	2.24	0.73
1:A:353:LYS:HG2	3:A:733:CL:CL	2.26	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:543:TYR:CZ	1:C:545:PRO:HG3	2.24	0.72
1:A:514:VAL:HG12	1:A:518:ARG:HG3	1.69	0.72
1:D:639:PHE:CZ	1:D:663:PRO:HD2	2.25	0.72
1:D:572:GLU:OE1	1:D:678:LYS:HE3	1.90	0.72
1:C:470:ARG:HG2	1:C:470:ARG:NH1	2.03	0.71
1:C:662:LEU:H	1:C:662:LEU:CD2	2.03	0.71
1:C:372:ILE:HD13	1:C:710:ILE:HG21	1.72	0.71
1:D:485:PRO:HG3	1:D:509:TYR:CE2	2.27	0.70
1:A:338:ARG:HD2	1:A:430:TYR:CD1	2.25	0.70
1:D:648:ASP:HB2	1:D:651:LEU:HB2	1.72	0.70
1:B:467:LYS:HE3	1:B:470:ARG:NH2	2.07	0.69
1:A:714:ILE:O	1:A:718:ILE:HG23	1.92	0.69
1:C:620:ILE:HD11	1:C:677:LEU:HD21	1.73	0.69
1:D:548:ARG:NH2	1:D:622:ASP:OD1	2.23	0.68
1:C:548:ARG:HD3	1:C:643:ASP:OD2	1.94	0.68
1:D:333:ARG:NH1	1:D:333:ARG:HG2	2.08	0.67
1:C:467:LYS:HE3	1:C:470:ARG:HH22	1.60	0.67
1:D:543:TYR:CZ	1:D:545:PRO:HG3	2.29	0.67
1:A:325:HIS:CE1	1:A:329:ILE:HD13	2.30	0.67
1:B:403:LYS:HE2	3:B:736:CL:CL	2.32	0.66
1:C:673:LEU:HD22	1:C:677:LEU:HD22	1.76	0.66
1:D:325:HIS:CE1	1:D:329:ILE:HD11	2.30	0.66
1:C:444:ASN:O	1:C:511:ARG:NH1	2.28	0.66
1:B:383:TYR:CE2	1:B:718:ILE:HD11	2.31	0.66
1:D:451:LEU:H	1:D:451:LEU:HD12	1.60	0.66
1:B:333:ARG:HH11	1:B:333:ARG:CG	2.08	0.65
1:A:470:ARG:NH1	1:A:470:ARG:HG2	2.07	0.65
1:C:487:ARG:HH11	1:C:487:ARG:HB2	1.61	0.65
1:A:508:GLY:HA3	1:A:702:ASP:OD1	1.95	0.65
1:B:470:ARG:NH1	1:B:470:ARG:HG2	2.03	0.65
1:C:465:THR:HB	1:C:466:PRO:HD3	1.77	0.65
1:B:456:LYS:HA	1:B:456:LYS:HE2	1.78	0.64
1:B:538:LYS:HE2	1:B:575:ASP:O	1.97	0.64
1:C:333:ARG:HG2	1:C:333:ARG:NH1	2.02	0.64
1:B:511:ARG:HH22	1:B:629:ASP:CG	2.02	0.64
1:C:714:ILE:O	1:C:718:ILE:HG23	1.98	0.63
1:B:542:MET:CE	1:B:612:LEU:HB3	2.29	0.63
1:C:569:LEU:HD22	1:C:573:LEU:HD12	1.80	0.63
1:B:372:ILE:HD13	1:B:710:ILE:HG21	1.79	0.63
1:D:548:ARG:HH22	1:D:622:ASP:CG	2.01	0.63
1:A:383:TYR:CZ	1:A:718:ILE:HD11	2.33	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:633:LEU:HB3	1:C:635:ARG:HD3	1.82	0.62
1:D:413:GLN:O	1:D:417:GLU:HG3	2.00	0.61
1:D:353:LYS:HG2	3:D:731:CL:CL	2.37	0.61
1:D:338:ARG:HD2	1:D:430:TYR:CD1	2.35	0.61
1:D:662:LEU:HB2	1:D:663:PRO:HD3	1.82	0.61
1:A:662:LEU:HD23	1:A:662:LEU:H	1.64	0.61
1:B:353:LYS:HG2	3:B:732:CL:CL	2.38	0.61
1:A:673:LEU:HD22	1:A:677:LEU:HD22	1.83	0.61
1:C:487:ARG:NH1	1:C:487:ARG:HB2	2.16	0.60
1:C:665:PRO:HG3	1:C:687:TYR:CZ	2.35	0.60
1:D:317:ASN:O	1:D:321:LYS:HG3	2.01	0.60
1:B:451:LEU:H	1:B:451:LEU:HD12	1.64	0.60
1:B:662:LEU:HD23	1:B:662:LEU:H	1.67	0.60
1:B:325:HIS:CE1	1:B:329:ILE:HD13	2.37	0.60
1:B:542:MET:HE1	1:B:612:LEU:CB	2.32	0.60
1:D:470:ARG:CG	1:D:470:ARG:HH11	1.95	0.59
1:B:559:LEU:HD23	1:B:587:ILE:HG23	1.84	0.59
1:D:347:GLU:HG3	1:D:436:ASN:HB2	1.82	0.59
1:D:438:THR:HA	1:D:480:ASP:OD2	2.02	0.59
1:C:701:VAL:HG12	1:C:701:VAL:O	2.02	0.59
1:D:670:PRO:HD2	1:D:671:TYR:CD2	2.38	0.59
1:D:554:SER:O	1:D:555:LYS:HB2	2.02	0.59
1:A:338:ARG:HD3	1:A:412:TYR:OH	2.03	0.59
1:B:480:ASP:O	1:B:503:ARG:HG2	2.03	0.59
1:D:465:THR:HB	1:D:466:PRO:HD3	1.84	0.58
1:A:641:ALA:HB1	1:A:644:ILE:HB	1.84	0.58
1:D:453:ASN:HB2	1:D:498:TRP:CZ2	2.38	0.58
1:D:582:ARG:HD3	3:D:733:CL:CL	2.40	0.58
1:A:378:LYS:HD3	1:A:379:TYR:CE2	2.39	0.58
1:C:347:GLU:O	1:C:436:ASN:ND2	2.37	0.58
1:A:546:THR:HG21	1:A:623:TYR:O	2.04	0.58
1:B:546:THR:HG21	1:B:623:TYR:O	2.04	0.58
1:D:714:ILE:O	1:D:718:ILE:HG23	2.04	0.58
1:D:585:TYR:CZ	1:D:586:LEU:HG	2.38	0.57
1:C:665:PRO:HG3	1:C:687:TYR:CE1	2.39	0.57
1:B:434:LYS:C	1:B:436:ASN:H	2.07	0.57
1:A:662:LEU:HB2	1:A:663:PRO:CD	2.35	0.57
1:A:513:ASP:HB2	4:A:748:HOH:O	2.04	0.57
1:B:624:SER:OG	1:B:626:VAL:HG22	2.03	0.57
1:B:648:ASP:HB2	1:B:651:LEU:HB2	1.84	0.57
1:A:420:HIS:CD2	1:A:438:THR:HB	2.40	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:548:ARG:HD3	1:A:643:ASP:OD2	2.05	0.57
1:B:338:ARG:HD2	1:B:430:TYR:CD1	2.40	0.56
1:C:434:LYS:C	1:C:436:ASN:H	2.07	0.56
1:C:387:TRP:HB3	1:C:389:PHE:CE2	2.41	0.56
1:B:347:GLU:O	1:B:436:ASN:ND2	2.39	0.56
1:C:559:LEU:HD23	1:C:587:ILE:HG23	1.88	0.56
1:B:675:LYS:O	1:B:678:LYS:HB2	2.06	0.56
1:B:368:SER:HB3	1:B:510:PRO:HD2	1.89	0.56
1:B:652:ARG:C	1:B:654:PHE:H	2.08	0.55
1:B:542:MET:HE1	1:B:612:LEU:HB3	1.86	0.55
1:D:325:HIS:CE1	1:D:340:LEU:HB2	2.42	0.55
1:C:511:ARG:HH22	1:C:629:ASP:CG	2.10	0.55
1:B:548:ARG:HH22	1:B:622:ASP:CG	2.08	0.55
1:B:625:SER:N	2:B:730:SO4:O2	2.39	0.55
1:C:662:LEU:CD2	1:C:662:LEU:N	2.68	0.55
1:B:632:ILE:HG12	1:B:699:CYS:HB3	1.89	0.54
1:A:325:HIS:CE1	1:A:329:ILE:CD1	2.90	0.54
1:A:451:LEU:HD12	1:A:451:LEU:H	1.72	0.54
1:C:340:LEU:O	1:C:344:THR:HG23	2.08	0.54
1:D:597:GLU:O	1:D:598:ASN:HB2	2.07	0.54
1:B:548:ARG:NH2	1:B:622:ASP:OD1	2.30	0.54
1:C:554:SER:C	1:C:556:GLY:H	2.11	0.53
1:B:638:PHE:HE1	1:B:676:GLU:HG2	1.73	0.53
1:A:548:ARG:NH2	1:A:622:ASP:OD1	2.35	0.53
1:A:405:LYS:O	1:A:411:TYR:HB2	2.08	0.53
1:C:582:ARG:NH2	1:C:606:TYR:O	2.42	0.53
1:C:485:PRO:HD2	1:C:489:SER:HB2	1.90	0.53
1:A:485:PRO:HG3	1:A:509:TYR:CZ	2.45	0.52
1:C:485:PRO:HG3	1:C:509:TYR:CE2	2.44	0.52
1:D:383:TYR:CE2	1:D:718:ILE:HD11	2.44	0.52
1:B:475:GLU:OE1	1:B:478:ARG:NH1	2.41	0.52
1:B:543:TYR:CZ	1:B:545:PRO:HG3	2.45	0.52
1:D:514:VAL:HG12	1:D:518:ARG:HG3	1.92	0.52
1:D:352:PRO:HD2	3:D:731:CL:CL	2.47	0.52
1:B:716:LYS:O	1:B:720:GLU:HG3	2.09	0.52
1:C:418:ALA:O	1:C:437:GLN:HG2	2.09	0.52
1:C:453:ASN:HA	1:C:469:LYS:HD3	1.92	0.52
1:B:559:LEU:CD2	1:B:587:ILE:HG23	2.40	0.52
1:C:663:PRO:HG3	1:C:694:PHE:CG	2.45	0.52
1:D:442:THR:HG22	1:D:483:ILE:HD12	1.92	0.52
1:B:325:HIS:NE2	1:B:329:ILE:HD13	2.24	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:648:ASP:HB2	1:A:651:LEU:CB	2.37	0.51
1:C:718:ILE:HG13	1:C:718:ILE:O	2.10	0.51
1:D:662:LEU:HB2	1:D:663:PRO:CD	2.40	0.51
1:C:445:GLY:HA2	1:C:509:TYR:CE2	2.45	0.51
1:D:338:ARG:HD3	1:D:412:TYR:OH	2.10	0.51
1:B:387:TRP:HB3	1:B:389:PHE:CE2	2.44	0.51
1:A:395:ASN:O	1:A:403:LYS:HE2	2.11	0.51
1:C:316:VAL:HG13	1:C:317:ASN:N	2.25	0.51
1:A:662:LEU:CD2	1:A:662:LEU:H	2.24	0.51
1:B:513:ASP:HA	1:B:704:GLY:HA2	1.92	0.51
1:D:352:PRO:HA	1:D:419:SER:HB3	1.92	0.51
1:D:697:ARG:HD3	1:D:698:PHE:CZ	2.45	0.51
1:C:637:GLN:O	1:C:664:GLY:HA3	2.10	0.51
1:D:638:PHE:HE1	1:D:676:GLU:HG2	1.76	0.51
1:B:467:LYS:HE3	1:B:470:ARG:HH22	1.76	0.51
1:A:521:ASP:O	1:A:525:LEU:HG	2.10	0.51
1:B:329:ILE:HG22	1:B:330:VAL:N	2.26	0.50
1:D:639:PHE:CE2	1:D:663:PRO:HD2	2.46	0.50
1:D:451:LEU:N	1:D:451:LEU:HD12	2.23	0.50
1:A:341:TYR:CD1	1:A:412:TYR:HB3	2.46	0.50
1:D:669:GLU:HG2	1:D:671:TYR:H	1.76	0.50
1:A:543:TYR:CZ	1:A:545:PRO:HG3	2.46	0.50
1:C:467:LYS:CE	1:C:470:ARG:HH22	2.25	0.50
1:C:344:THR:O	1:C:346:LYS:HG3	2.12	0.50
1:A:718:ILE:O	1:A:722:LEU:HD13	2.11	0.50
1:C:507:ILE:HG13	1:C:508:GLY:O	2.12	0.50
1:C:322:THR:O	1:C:326:VAL:HG23	2.12	0.50
1:C:639:PHE:CZ	1:C:663:PRO:HD2	2.46	0.50
1:A:456:LYS:HA	1:A:456:LYS:HE2	1.93	0.50
1:D:456:LYS:HE2	1:D:456:LYS:HA	1.94	0.50
1:C:652:ARG:C	1:C:654:PHE:H	2.15	0.49
1:C:450:ARG:HA	1:C:655:TYR:CE2	2.47	0.49
1:C:347:GLU:HA	1:C:434:LYS:HD3	1.95	0.49
1:D:322:THR:O	1:D:326:VAL:HG23	2.12	0.49
1:C:341:TYR:O	1:C:345:ASP:HB2	2.12	0.49
1:C:442:THR:O	1:C:509:TYR:HE1	1.96	0.49
1:A:420:HIS:CE1	1:A:718:ILE:HG22	2.48	0.49
1:D:490:THR:HG23	1:D:504:ILE:HD13	1.93	0.49
1:A:597:GLU:O	1:A:598:ASN:HB2	2.13	0.49
1:B:519:ALA:HA	1:B:614:LEU:HD22	1.95	0.49
1:A:333:ARG:NH1	1:A:333:ARG:CG	2.57	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:607:ASN:N	1:D:607:ASN:OD1	2.45	0.48
1:A:448:LEU:HD13	1:A:627:MET:HE1	1.94	0.48
1:C:545:PRO:HG2	1:C:583:MET:HE1	1.95	0.48
1:C:448:LEU:HD13	1:C:627:MET:HE1	1.95	0.48
1:A:722:LEU:O	1:A:723:GLU:HG3	2.13	0.48
1:A:325:HIS:NE2	1:A:329:ILE:HD13	2.28	0.48
1:D:353:LYS:CG	3:D:731:CL:CL	2.99	0.48
1:B:344:THR:O	1:B:346:LYS:HG2	2.13	0.48
1:C:584:HIS:ND1	1:C:586:LEU:HB2	2.28	0.48
1:D:387:TRP:HB3	1:D:389:PHE:CE2	2.48	0.48
1:A:646:LYS:O	1:A:648:ASP:N	2.41	0.48
1:A:662:LEU:N	1:A:662:LEU:CD2	2.77	0.48
1:A:540:VAL:HG22	1:A:578:VAL:HG12	1.95	0.48
1:D:637:GLN:O	1:D:664:GLY:HA3	2.14	0.47
1:B:663:PRO:HG3	1:B:694:PHE:CD2	2.49	0.47
1:C:485:PRO:HA	1:C:508:GLY:HA2	1.95	0.47
1:D:652:ARG:C	1:D:654:PHE:H	2.16	0.47
1:D:325:HIS:NE2	1:D:329:ILE:HD11	2.28	0.47
1:D:347:GLU:O	1:D:436:ASN:ND2	2.47	0.47
1:B:356:VAL:HG23	1:B:421:TRP:HA	1.96	0.47
1:B:542:MET:HE2	1:B:612:LEU:HB3	1.94	0.47
1:B:651:LEU:O	1:B:651:LEU:HD23	2.15	0.47
1:C:316:VAL:HG23	1:D:331:LEU:HD22	1.96	0.47
1:D:378:LYS:HD3	1:D:379:TYR:CE2	2.49	0.47
1:C:420:HIS:CE1	1:C:718:ILE:HG22	2.49	0.47
1:C:338:ARG:HD3	1:C:412:TYR:OH	2.13	0.47
1:A:453:ASN:HA	1:A:469:LYS:HD3	1.96	0.47
1:B:670:PRO:HD2	1:B:671:TYR:CD2	2.49	0.47
1:D:663:PRO:HG3	1:D:694:PHE:CG	2.49	0.47
1:D:365:TYR:CZ	1:D:370:LYS:HG3	2.50	0.47
1:A:515:LEU:HD12	1:A:632:ILE:HD12	1.96	0.47
1:D:697:ARG:HD3	1:D:698:PHE:CE2	2.49	0.47
1:B:524:TYR:O	1:B:528:ILE:HG13	2.15	0.47
1:D:465:THR:CG2	1:D:469:LYS:HE3	2.44	0.47
1:B:338:ARG:HD3	1:B:412:TYR:OH	2.14	0.47
1:B:669:GLU:HG2	1:B:671:TYR:H	1.79	0.47
1:A:480:ASP:O	1:A:503:ARG:HG2	2.15	0.47
1:B:625:SER:HB3	2:B:730:SO4:O2	2.14	0.47
1:B:434:LYS:C	1:B:436:ASN:N	2.67	0.47
1:D:356:VAL:HG23	1:D:421:TRP:HA	1.97	0.47
1:A:344:THR:O	1:A:346:LYS:HG3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:406:ARG:O	1:A:407:ASN:HB2	2.14	0.46
1:A:484:SER:HA	1:A:485:PRO:HD3	1.73	0.46
1:D:651:LEU:HG	1:D:653:GLY:H	1.81	0.46
1:C:450:ARG:HG3	1:C:653:GLY:O	2.15	0.46
1:D:340:LEU:O	1:D:344:THR:HG23	2.15	0.46
1:B:353:LYS:HA	1:B:718:ILE:HD12	1.98	0.46
1:B:354:THR:C	1:B:355:ILE:HG13	2.35	0.46
1:B:639:PHE:CD2	1:B:639:PHE:N	2.83	0.46
1:C:470:ARG:NH1	1:C:471:ASN:OD1	2.48	0.46
1:B:445:GLY:HA2	1:B:509:TYR:CE2	2.50	0.46
1:D:494:ARG:HA	1:D:499:MET:HB2	1.96	0.46
1:C:353:LYS:HA	1:C:718:ILE:HD12	1.97	0.46
1:D:345:ASP:OD2	1:D:434:LYS:HE3	2.15	0.46
1:A:490:THR:HG23	1:A:504:ILE:HD13	1.98	0.46
1:B:532:LEU:HD21	1:B:606:TYR:CE2	2.50	0.46
1:D:639:PHE:HZ	1:D:663:PRO:HD2	1.73	0.46
1:B:662:LEU:HB2	1:B:663:PRO:CD	2.46	0.46
1:B:714:ILE:O	1:B:718:ILE:HG23	2.15	0.46
1:B:673:LEU:HD22	1:B:677:LEU:HD22	1.97	0.46
1:A:352:PRO:HA	1:A:419:SER:HB3	1.97	0.45
1:B:317:ASN:O	1:B:321:LYS:HG3	2.17	0.45
1:A:450:ARG:HA	1:A:655:TYR:CE2	2.52	0.45
1:A:509:TYR:HB3	1:A:511:ARG:HG2	1.96	0.45
1:B:451:LEU:N	1:B:451:LEU:HD12	2.32	0.45
1:C:323:LEU:HA	1:C:323:LEU:HD12	1.75	0.45
1:D:560:PHE:O	1:D:561:GLU:C	2.54	0.45
1:D:675:LYS:O	1:D:678:LYS:HB2	2.17	0.45
1:D:325:HIS:CE1	1:D:329:ILE:CD1	2.98	0.45
1:B:638:PHE:CZ	1:B:683:VAL:HG11	2.52	0.45
1:C:470:ARG:CG	1:C:470:ARG:NH1	2.75	0.45
1:D:434:LYS:C	1:D:436:ASN:H	2.19	0.45
1:B:380:TYR:N	1:B:381:PRO:HD3	2.31	0.45
1:C:663:PRO:HG3	1:C:694:PHE:CD2	2.52	0.45
1:A:652:ARG:C	1:A:654:PHE:H	2.19	0.45
1:B:531:HIS:CD2	1:B:531:HIS:O	2.70	0.45
1:B:488:TYR:OH	1:B:655:TYR:HB2	2.16	0.45
1:A:519:ALA:HA	1:A:614:LEU:HD22	1.98	0.45
1:B:418:ALA:O	1:B:437:GLN:HG2	2.17	0.45
1:B:687:TYR:O	1:B:688:GLN:C	2.54	0.45
1:D:325:HIS:HE1	1:D:340:LEU:HB2	1.81	0.44
1:B:329:ILE:HD12	1:B:336:LYS:HB2	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:347:GLU:HG3	1:B:436:ASN:HB2	1.98	0.44
1:C:316:VAL:HG23	1:D:331:LEU:CD2	2.47	0.44
1:A:385:TYR:O	1:A:401:ALA:HA	2.17	0.44
1:D:548:ARG:NH1	1:D:643:ASP:OD1	2.50	0.44
1:B:554:SER:C	1:B:556:GLY:H	2.20	0.44
1:C:680:LEU:HD23	1:C:680:LEU:HA	1.80	0.44
1:B:607:ASN:N	1:B:607:ASN:OD1	2.50	0.44
1:A:345:ASP:OD1	1:A:434:LYS:CE	2.66	0.44
1:D:450:ARG:HA	1:D:655:TYR:CE2	2.52	0.44
1:A:607:ASN:N	1:A:607:ASN:OD1	2.48	0.44
1:B:405:LYS:O	1:B:411:TYR:HB2	2.18	0.44
1:B:403:LYS:CE	3:B:736:CL:CL	3.01	0.44
1:D:525:LEU:O	1:D:529:ARG:HG3	2.18	0.44
1:C:470:ARG:HD2	1:C:474:ARG:HH21	1.83	0.44
1:C:353:LYS:HG2	3:C:733:CL:CL	2.54	0.44
1:A:472:PHE:O	1:A:476:THR:HG23	2.18	0.44
1:B:458:VAL:C	1:B:460:MET:H	2.21	0.44
1:D:325:HIS:O	1:D:329:ILE:HD13	2.17	0.43
1:B:340:LEU:O	1:B:344:THR:HG23	2.18	0.43
1:C:585:TYR:O	1:C:589:ASN:OD1	2.36	0.43
1:C:356:VAL:HG22	1:C:418:ALA:CB	2.48	0.43
1:A:450:ARG:HD2	1:A:653:GLY:HA2	2.01	0.43
1:B:352:PRO:HA	1:B:419:SER:HB3	2.00	0.43
1:C:718:ILE:O	1:C:722:LEU:HD22	2.19	0.43
1:C:347:GLU:CG	1:C:436:ASN:HB2	2.49	0.43
1:C:325:HIS:CE1	1:C:340:LEU:HB2	2.53	0.43
1:C:431:LEU:HA	1:C:431:LEU:HD23	1.60	0.43
1:C:687:TYR:O	1:C:688:GLN:C	2.56	0.43
1:C:441:GLN:HG2	1:C:479:TRP:CE2	2.53	0.43
1:A:443:TRP:CG	1:A:444:ASN:N	2.86	0.43
1:B:663:PRO:HG3	1:B:694:PHE:CD1	2.53	0.43
1:C:722:LEU:O	1:C:723:GLU:HG3	2.19	0.43
1:C:341:TYR:CD1	1:C:412:TYR:HB3	2.53	0.43
1:C:316:VAL:CG1	1:C:317:ASN:N	2.81	0.43
1:D:638:PHE:CD2	1:D:638:PHE:N	2.87	0.43
1:D:446:THR:HB	1:D:628:PHE:CD2	2.54	0.43
1:B:345:ASP:OD2	1:B:434:LYS:HE3	2.19	0.43
1:D:515:LEU:HD12	1:D:632:ILE:HD12	2.01	0.43
1:D:624:SER:OG	1:D:626:VAL:HG22	2.19	0.43
1:A:356:VAL:HG23	1:A:421:TRP:HA	2.01	0.43
1:C:545:PRO:HG2	1:C:583:MET:CE	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:378:LYS:HD3	1:C:379:TYR:CE2	2.53	0.42
1:C:554:SER:C	1:C:556:GLY:N	2.72	0.42
1:B:484:SER:HA	1:B:485:PRO:HD3	1.78	0.42
1:D:375:TYR:C	1:D:375:TYR:CD2	2.93	0.42
1:D:458:VAL:HG12	1:D:460:MET:HB2	2.00	0.42
1:C:695:TYR:CD2	1:C:695:TYR:C	2.92	0.42
1:C:458:VAL:HG13	1:C:460:MET:HE2	2.01	0.42
1:C:722:LEU:HD12	1:C:722:LEU:HA	1.70	0.42
1:C:586:LEU:HD23	1:C:586:LEU:HA	1.88	0.42
1:C:458:VAL:HG13	1:C:460:MET:CE	2.48	0.42
1:B:618:CYS:HB2	1:B:636:PRO:O	2.18	0.42
1:A:380:TYR:N	1:A:381:PRO:HD3	2.34	0.42
1:C:671:TYR:O	1:C:675:LYS:HG2	2.20	0.42
1:D:652:ARG:C	1:D:654:PHE:N	2.72	0.42
1:D:633:LEU:HB3	1:D:635:ARG:HD3	2.01	0.42
1:B:662:LEU:CD2	1:B:662:LEU:H	2.32	0.42
1:A:554:SER:C	1:A:556:GLY:H	2.21	0.42
1:A:651:LEU:HG	1:A:653:GLY:H	1.84	0.42
1:A:591:LEU:HD23	1:A:591:LEU:HA	1.91	0.42
1:D:562:LEU:H	1:D:562:LEU:HG	1.67	0.42
1:D:480:ASP:O	1:D:503:ARG:HG2	2.20	0.42
1:B:460:MET:HA	1:B:461:PRO:HD2	1.77	0.42
1:A:459:ARG:HG3	1:A:459:ARG:O	2.19	0.42
1:B:597:GLU:O	1:B:598:ASN:HB2	2.19	0.42
1:C:569:LEU:HD23	1:C:569:LEU:HA	1.83	0.42
1:B:648:ASP:O	1:B:649:LYS:HB2	2.20	0.42
1:B:652:ARG:C	1:B:654:PHE:N	2.71	0.42
1:C:467:LYS:HE3	1:C:470:ARG:NH2	2.33	0.42
1:A:450:ARG:HG3	1:A:653:GLY:O	2.20	0.42
1:A:639:PHE:CZ	1:A:663:PRO:HD2	2.55	0.42
1:D:356:VAL:HG22	1:D:418:ALA:CB	2.50	0.42
1:A:669:GLU:HA	1:A:670:PRO:HD3	1.89	0.42
1:A:353:LYS:HB2	1:A:383:TYR:CD2	2.55	0.41
1:B:325:HIS:CE1	1:B:340:LEU:HB2	2.55	0.41
1:B:540:VAL:HG22	1:B:578:VAL:HG12	2.01	0.41
1:B:665:PRO:HB2	1:B:667:TYR:CE2	2.55	0.41
1:D:682:LYS:HB2	1:D:682:LYS:HE3	1.93	0.41
1:D:501:GLU:HA	1:D:504:ILE:HD12	2.03	0.41
1:C:329:ILE:HG22	1:C:330:VAL:N	2.32	0.41
1:A:660:GLU:HB3	1:A:661:ASP:H	1.53	0.41
1:A:582:ARG:NH2	1:A:606:TYR:O	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:431:LEU:HD23	1:D:431:LEU:HA	1.88	0.41
1:C:451:LEU:H	1:C:451:LEU:HD12	1.85	0.41
1:C:597:GLU:O	1:C:598:ASN:HB2	2.21	0.41
1:C:353:LYS:CG	3:C:733:CL:CL	3.06	0.41
1:D:586:LEU:HD23	1:D:586:LEU:HA	1.65	0.41
1:B:431:LEU:HA	1:B:431:LEU:HD23	1.68	0.41
1:A:352:PRO:HB2	1:A:722:LEU:HD11	2.02	0.41
1:A:353:LYS:HA	1:A:718:ILE:HD12	2.02	0.41
1:A:470:ARG:NH1	1:A:470:ARG:CG	2.79	0.41
1:A:325:HIS:HB2	1:A:339:SER:HB2	2.02	0.41
1:C:427:THR:HA	1:C:428:PRO:HD3	1.91	0.41
1:B:410:GLU:N	1:B:410:GLU:OE1	2.47	0.41
1:D:677:LEU:HA	1:D:677:LEU:HD12	1.87	0.41
1:A:438:THR:HA	1:A:480:ASP:OD2	2.21	0.41
1:D:351:LYS:HA	1:D:352:PRO:HD3	1.87	0.41
1:C:652:ARG:C	1:C:654:PHE:N	2.73	0.41
1:B:350:VAL:HA	1:B:417:GLU:O	2.20	0.41
1:B:351:LYS:HB3	3:B:732:CL:CL	2.57	0.41
1:B:338:ARG:HD2	1:B:430:TYR:CG	2.56	0.41
1:C:554:SER:O	1:C:555:LYS:HB2	2.20	0.41
1:C:566:LEU:HA	1:C:566:LEU:HD23	1.86	0.41
1:C:484:SER:HA	1:C:485:PRO:HD3	1.86	0.41
1:A:543:TYR:CE2	1:A:545:PRO:HG3	2.55	0.41
1:D:458:VAL:O	1:D:458:VAL:HG12	2.21	0.41
1:A:669:GLU:HG2	1:A:671:TYR:H	1.86	0.40
1:B:446:THR:HA	1:B:447:PRO:HD3	1.90	0.40
1:D:453:ASN:HA	1:D:469:LYS:HD3	2.02	0.40
1:C:434:LYS:C	1:C:436:ASN:N	2.71	0.40
1:A:452:ALA:CB	1:A:496:ALA:O	2.70	0.40
1:B:396:VAL:HA	3:B:736:CL:CL	2.59	0.40
1:B:443:TRP:CG	1:B:444:ASN:N	2.89	0.40
1:D:323:LEU:HA	1:D:323:LEU:HD12	1.75	0.40
1:D:484:SER:HA	1:D:485:PRO:HD3	1.84	0.40
1:C:317:ASN:O	1:C:321:LYS:HG3	2.21	0.40
1:D:541:ILE:HG12	1:D:618:CYS:SG	2.61	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	408/729 (56%)	384 (94%)	21 (5%)	3 (1%)	26	60
1	B	410/729 (56%)	378 (92%)	30 (7%)	2 (0%)	34	68
1	C	409/729 (56%)	370 (90%)	36 (9%)	3 (1%)	26	60
1	D	409/729 (56%)	376 (92%)	26 (6%)	7 (2%)	11	34
All	All	1636/2916 (56%)	1508 (92%)	113 (7%)	15 (1%)	21	53

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	647	TYR
1	D	561	GLU
1	A	535	PRO
1	B	345	ASP
1	B	649	LYS
1	D	649	LYS
1	C	521	ASP
1	C	649	LYS
1	D	521	ASP
1	D	562	LEU
1	D	565	ASP
1	A	649	LYS
1	C	443	TRP
1	D	458	VAL
1	D	603	VAL

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	379/675 (56%)	341 (90%)	38 (10%)	9 27
1	B	380/675 (56%)	334 (88%)	46 (12%)	6 18
1	C	379/675 (56%)	343 (90%)	36 (10%)	11 29
1	D	380/675 (56%)	345 (91%)	35 (9%)	11 31
All	All	1518/2700 (56%)	1363 (90%)	155 (10%)	9 26

All (155) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	316	VAL
1	A	322	THR
1	A	323	LEU
1	A	329	ILE
1	A	333	ARG
1	A	356	VAL
1	A	402	GLU
1	A	403	LYS
1	A	433	LYS
1	A	439	TYR
1	A	451	LEU
1	A	456	LYS
1	A	460	MET
1	A	472	PHE
1	A	487	ARG
1	A	488	TYR
1	A	491	GLU
1	A	511	ARG
1	A	514	VAL
1	A	555	LYS
1	A	559	LEU
1	A	578	VAL
1	A	585	TYR
1	A	589	ASN
1	A	594	SER
1	A	607	ASN
1	A	619	LEU
1	A	645	ASP
1	A	651	LEU
1	A	656	MET
1	A	659	MET

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Mol	Chain	Res	Type
1	A	660	GLU
1	A	662	LEU
1	A	669	GLU
1	A	673	LEU
1	A	677	LEU
1	A	678	LYS
1	A	718	ILE
1	B	315	LYS
1	B	316	VAL
1	B	329	ILE
1	B	333	ARG
1	B	343	LEU
1	B	356	VAL
1	B	367	ASP
1	B	368	SER
1	B	402	GLU
1	B	403	LYS
1	B	407	ASN
1	B	423	SER
1	B	433	LYS
1	B	439	TYR
1	B	449	LYS
1	B	451	LEU
1	B	456	LYS
1	B	460	MET
1	B	465	THR
1	B	470	ARG
1	B	472	PHE
1	B	487	ARG
1	B	491	GLU
1	B	514	VAL
1	B	548	ARG
1	B	555	LYS
1	B	559	LEU
1	B	567	ASP
1	B	594	SER
1	B	607	ASN
1	B	619	LEU
1	B	622	ASP
1	B	623	TYR
1	B	625	SER
1	B	651	LEU

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Mol	Chain	Res	Type
1	B	661	ASP
1	B	662	LEU
1	B	668	THR
1	B	669	GLU
1	B	673	LEU
1	B	677	LEU
1	B	680	LEU
1	B	702	ASP
1	B	718	ILE
1	B	721	GLN
1	B	722	LEU
1	C	315	LYS
1	C	323	LEU
1	C	333	ARG
1	C	338	ARG
1	C	339	SER
1	C	356	VAL
1	C	403	LYS
1	C	433	LYS
1	C	439	TYR
1	C	451	LEU
1	C	456	LYS
1	C	460	MET
1	C	463	THR
1	C	472	PHE
1	C	487	ARG
1	C	491	GLU
1	C	511	ARG
1	C	536	SER
1	C	555	LYS
1	C	559	LEU
1	C	562	LEU
1	C	567	ASP
1	C	570	TYR
1	C	589	ASN
1	C	619	LEU
1	C	651	LEU
1	C	656	MET
1	C	660	GLU
1	C	662	LEU
1	C	669	GLU
1	C	673	LEU

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Mol	Chain	Res	Type
1	C	677	LEU
1	C	680	LEU
1	C	702	ASP
1	C	718	ILE
1	C	722	LEU
1	D	316	VAL
1	D	323	LEU
1	D	333	ARG
1	D	346	LYS
1	D	356	VAL
1	D	403	LYS
1	D	413	GLN
1	D	433	LYS
1	D	451	LEU
1	D	456	LYS
1	D	460	MET
1	D	470	ARG
1	D	472	PHE
1	D	487	ARG
1	D	488	TYR
1	D	491	GLU
1	D	511	ARG
1	D	514	VAL
1	D	559	LEU
1	D	594	SER
1	D	607	ASN
1	D	619	LEU
1	D	622	ASP
1	D	645	ASP
1	D	651	LEU
1	D	656	MET
1	D	660	GLU
1	D	662	LEU
1	D	669	GLU
1	D	673	LEU
1	D	677	LEU
1	D	678	LYS
1	D	702	ASP
1	D	718	ILE
1	D	722	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	349	ASN
1	A	708	GLN
1	B	349	ASN
1	B	436	ASN
1	B	531	HIS
1	B	686	GLN
1	C	349	ASN
1	C	436	ASN
1	C	531	HIS
1	C	686	GLN
1	D	349	ASN
1	D	413	GLN
1	D	531	HIS
1	D	686	GLN

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

Of 40 ligands modelled in this entry, 32 are monoatomic - leaving 8 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	730	-	4,4,4	0.24	0	6,6,6	0.17	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	SO4	A	731	-	4,4,4	0.16	0	6,6,6	0.35	0
2	SO4	A	732	-	4,4,4	0.24	0	6,6,6	0.06	0
2	SO4	B	730	-	4,4,4	0.23	0	6,6,6	0.22	0
2	SO4	C	730	-	4,4,4	0.26	0	6,6,6	0.17	0
2	SO4	C	731	-	4,4,4	0.20	0	6,6,6	0.15	0
2	SO4	C	732	-	4,4,4	0.21	0	6,6,6	0.14	0
2	SO4	D	730	-	4,4,4	0.20	0	6,6,6	0.09	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	730	-	-	0/0/0/0	0/0/0/0
2	SO4	A	731	-	-	0/0/0/0	0/0/0/0
2	SO4	A	732	-	-	0/0/0/0	0/0/0/0
2	SO4	B	730	-	-	0/0/0/0	0/0/0/0
2	SO4	C	730	-	-	0/0/0/0	0/0/0/0
2	SO4	C	731	-	-	0/0/0/0	0/0/0/0
2	SO4	C	732	-	-	0/0/0/0	0/0/0/0
2	SO4	D	730	-	-	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.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	730	SO4	2	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

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, 95^{th} 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	410/729 (56%)	-0.27	9 (2%) 65 54	50, 73, 143, 221	0
1	B	412/729 (56%)	-0.24	12 (2%) 55 43	51, 72, 158, 239	0
1	C	411/729 (56%)	-0.19	16 (3%) 43 31	56, 79, 158, 237	0
1	D	411/729 (56%)	-0.17	13 (3%) 51 39	57, 78, 159, 221	0
All	All	1644/2916 (56%)	-0.22	50 (3%) 54 41	50, 76, 154, 239	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	724	HIS	5.5
1	C	558	TYR	5.3
1	D	556	GLY	5.3
1	D	652	ARG	5.1
1	B	652	ARG	5.1
1	B	560	PHE	4.2
1	B	558	TYR	4.2
1	D	558	TYR	4.1
1	C	556	GLY	4.0
1	C	652	ARG	4.0
1	A	652	ARG	3.8
1	D	560	PHE	3.8
1	C	555	LYS	3.7
1	B	537	ASP	3.5
1	D	555	LYS	3.4
1	C	560	PHE	3.3
1	A	560	PHE	3.3
1	B	313	ALA	3.2
1	C	313	ALA	3.2
1	C	557	LYS	3.2
1	D	557	LYS	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	557	LYS	3.1
1	A	459	ARG	3.1
1	D	649	LYS	3.0
1	D	651	LEU	2.9
1	A	558	TYR	2.9
1	A	723	GLU	2.8
1	B	561	GLU	2.8
1	D	554	SER	2.7
1	D	550	ASP	2.7
1	A	649	LYS	2.7
1	B	646	LYS	2.6
1	C	554	SER	2.6
1	C	723	GLU	2.5
1	C	458	VAL	2.5
1	B	459	ARG	2.5
1	D	457	VAL	2.5
1	C	459	ARG	2.4
1	B	649	LYS	2.4
1	B	724	HIS	2.4
1	C	649	LYS	2.3
1	C	435	GLU	2.3
1	B	555	LYS	2.2
1	A	491	GLU	2.2
1	A	348	ASP	2.1
1	D	723	GLU	2.1
1	C	457	VAL	2.1
1	C	491	GLU	2.1
1	C	553	VAL	2.1
1	A	720	GLU	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, 95^{th} 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	SO4	A	732	5/5	0.85	0.41	4.25	155,155,156,157	0
3	CL	C	738	1/1	0.94	0.34	1.96	87,87,87,87	0
2	SO4	D	730	5/5	0.80	0.29	1.88	188,188,189,190	0
2	SO4	B	730	5/5	0.84	0.27	1.63	174,174,176,176	0
3	CL	A	736	1/1	0.96	0.13	0.11	82,82,82,82	0
3	CL	D	734	1/1	0.92	0.13	-0.01	93,93,93,93	0
3	CL	A	739	1/1	0.98	0.18	-0.34	68,68,68,68	0
3	CL	A	740	1/1	0.91	0.10	-0.79	95,95,95,95	0
2	SO4	A	730	5/5	0.98	0.12	-0.98	99,100,102,103	0
3	CL	C	739	1/1	0.92	0.08	-1.14	99,99,99,99	0
2	SO4	C	730	5/5	0.97	0.11	-1.33	113,114,115,116	0
3	CL	B	737	1/1	0.98	0.10	-1.70	86,86,86,86	0
3	CL	B	731	1/1	0.99	0.08	-2.55	50,50,50,50	0
3	CL	D	735	1/1	0.86	0.14	-	103,103,103,103	0
3	CL	B	738	1/1	0.97	0.08	-	97,97,97,97	0
3	CL	A	734	1/1	0.88	0.34	-	108,108,108,108	0
3	CL	D	732	1/1	0.94	0.27	-	86,86,86,86	0
3	CL	B	733	1/1	0.90	0.41	-	111,111,111,111	0
3	CL	C	737	1/1	0.87	0.30	-	92,92,92,92	0
3	CL	A	738	1/1	0.98	0.28	-	98,98,98,98	0
3	CL	A	735	1/1	0.95	0.60	-	98,98,98,98	0
3	CL	A	741	1/1	0.92	0.17	-	91,91,91,91	0
3	CL	B	736	1/1	0.96	0.33	-	103,103,103,103	0
2	SO4	C	731	5/5	0.95	0.12	-	118,119,121,121	0
3	CL	C	735	1/1	0.96	0.22	-	81,81,81,81	0
3	CL	B	735	1/1	0.79	0.44	-	103,103,103,103	0
3	CL	B	734	1/1	0.77	0.18	-	104,104,104,104	0
2	SO4	C	732	5/5	0.62	0.48	-	177,177,178,178	0
2	SO4	A	731	5/5	0.98	0.15	-	87,87,88,91	0
3	CL	A	737	1/1	0.96	0.50	-	93,93,93,93	0
3	CL	A	743	1/1	0.81	0.14	-	99,99,99,99	0
3	CL	A	742	1/1	0.94	0.17	-	95,95,95,95	0
3	CL	D	731	1/1	0.84	0.08	-	90,90,90,90	0
3	CL	B	732	1/1	0.82	0.30	-	99,99,99,99	0
3	CL	C	734	1/1	0.97	0.21	-	93,93,93,93	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
3	CL	C	736	1/1	0.95	0.10	-	92,92,92,92	0
3	CL	C	733	1/1	0.94	0.16	-	105,105,105,105	0
3	CL	A	733	1/1	0.80	0.43	-	97,97,97,97	0
3	CL	D	733	1/1	0.87	0.28	-	96,96,96,96	0
3	CL	D	736	1/1	0.92	0.08	-	93,93,93,93	0

6.5 Other polymers

There are no such residues in this entry.