Full wwPDB X-ray Structure Validation Report (i)

Nov 13, 2023 – 07:01 PM JST

PDB ID	:	5XSV
Title	:	Crystal structure of an archaeal chitinase in the ligand-free form
Authors	:	Nishitani, Y.; Miki, K.
Deposited on	:	2017-06-15
Resolution	:	3.72 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

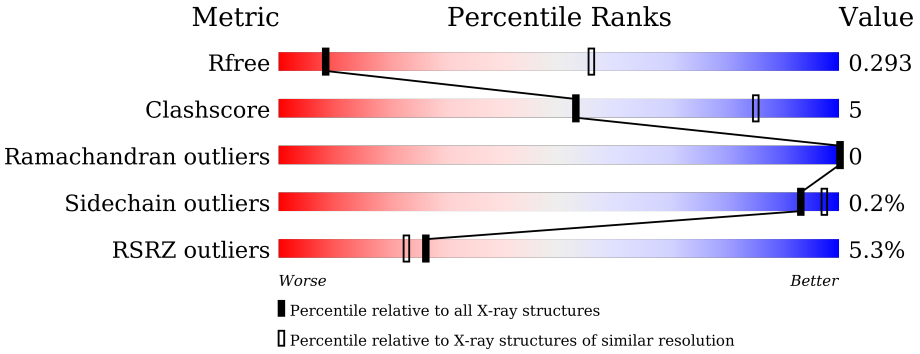
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 3.72 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1089 (3.90-3.54)
Clashscore	141614	1012 (3.88-3.56)
Ramachandran outliers	138981	1114 (3.90-3.54)
Sidechain outliers	138945	1110 (3.90-3.54)
RSRZ outliers	127900	1020 (3.90-3.54)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	486	.% 86% 14%
1	B	486	.% 86% 14%
1	C	486	6% 86% 13%
1	D	486	13% 86% 14%

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
3	SO4	D	1004	-	-	-	X

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 15938 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 Chitinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	486	3975	2583	647	731	14	0	0	0
1	B	485	3967	2578	646	730	13	0	0	0
1	C	485	3967	2578	646	730	13	0	0	0
1	D	486	3975	2583	647	731	14	0	0	0

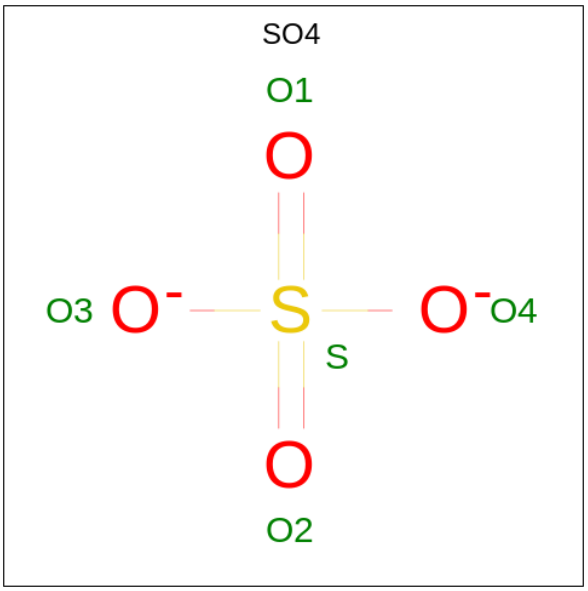
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	320	MET	-	initiating methionine	UNP A0A161KIT4
B	320	MET	-	initiating methionine	UNP A0A161KIT4
C	320	MET	-	initiating methionine	UNP A0A161KIT4
D	320	MET	-	initiating methionine	UNP A0A161KIT4

- Molecule 2 is COBALT (II) ION (three-letter code: CO) (formula: Co).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Co		
2	A	5	5	5	0	0
2	B	4	4	4	0	0
2	C	2	2	2	0	0
2	D	3	3	3	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



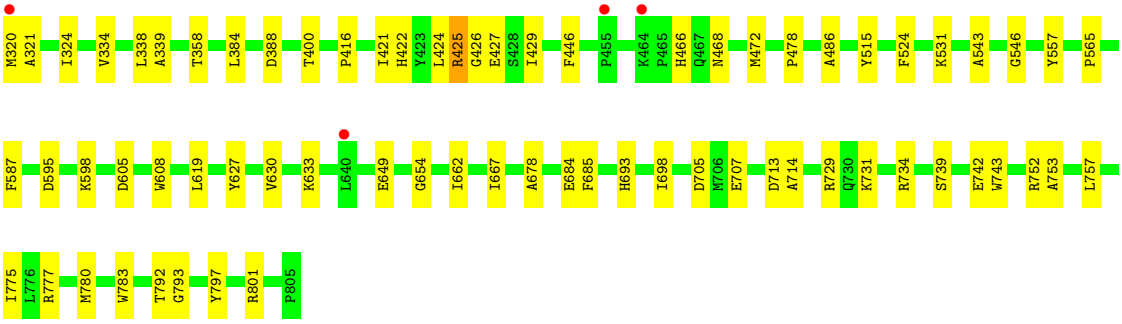
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

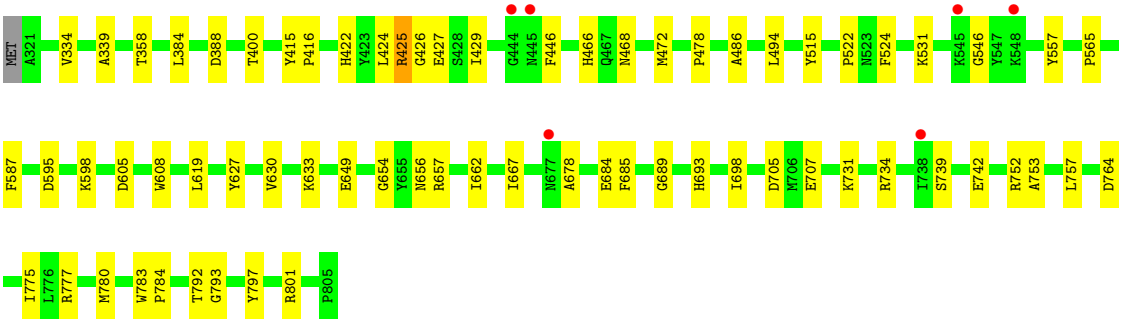
- Molecule 1: Chitinase

Chain A:



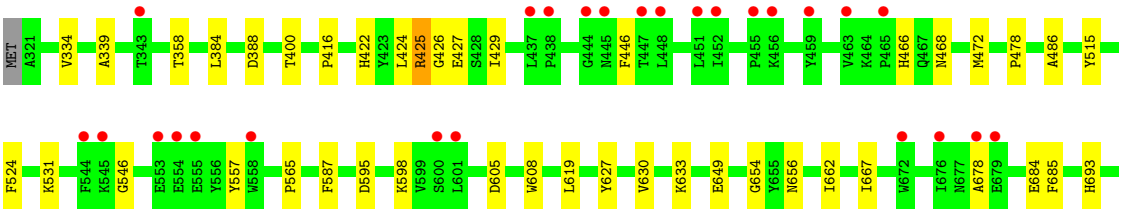
- Molecule 1: Chitinase

Chain B:



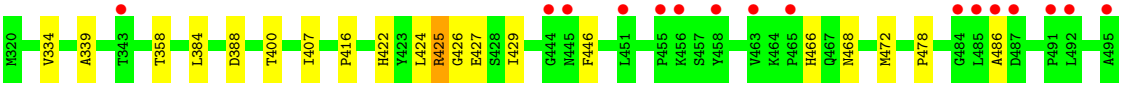
- Molecule 1: Chitinase

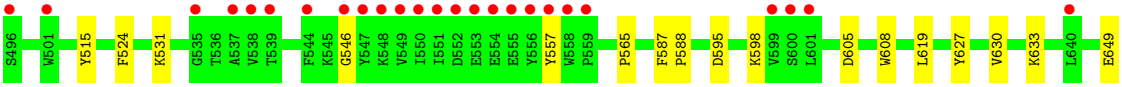
Chain C:

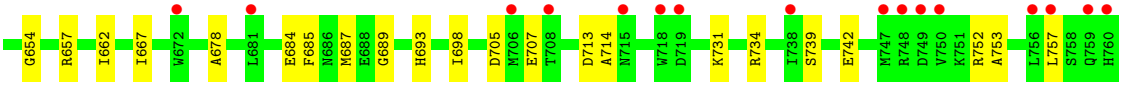


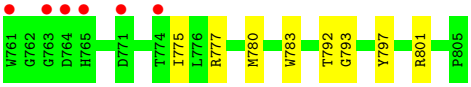
- Molecule 1: Chitinase

Chain D:









4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	126.92Å 104.73Å 156.46Å 90.00° 113.43° 90.00°	Depositor
Resolution (Å)	41.80 – 3.72 41.80 – 3.72	Depositor EDS
% Data completeness (in resolution range)	98.9 (41.80-3.72) 99.1 (41.80-3.72)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.60 (at 3.76Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.278 , 0.298 0.277 , 0.293	Depositor DCC
R_{free} test set	1976 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	108.7	Xtriage
Anisotropy	0.339	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 64.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.40$, $< L^2 > = 0.23$	Xtriage
Estimated twinning fraction	0.105 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.86	EDS
Total number of atoms	15938	wwPDB-VP
Average B, all atoms (Å^2)	118.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.77% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: CO, 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	0.22	0/4106	0.38	0/5606
1	B	0.22	0/4098	0.38	0/5596
1	C	0.22	0/4098	0.38	0/5596
1	D	0.21	0/4106	0.38	0/5606
All	All	0.22	0/16408	0.38	0/22404

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 (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	3975	0	3800	54	0
1	B	3967	0	3791	39	0
1	C	3967	0	3791	38	0
1	D	3975	0	3800	37	0
2	A	5	0	0	0	0
2	B	4	0	0	0	0
2	C	2	0	0	0	0
2	D	3	0	0	0	0
3	A	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	10	0	0	0	0
3	C	10	0	0	0	0
3	D	10	0	0	0	0
All	All	15938	0	15182	167	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 (167) 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:320:MET:CE	1:A:338:LEU:HD23	1.55	1.36
1:A:320:MET:HE3	1:A:338:LEU:HD23	1.12	1.05
1:A:320:MET:CE	1:A:338:LEU:CD2	2.36	1.03
1:A:320:MET:HE2	1:A:338:LEU:CD2	1.88	1.02
1:A:320:MET:HE2	1:A:338:LEU:HD23	1.42	0.89
1:A:320:MET:HG3	1:A:321:ALA:H	1.36	0.87
1:A:320:MET:HE3	1:A:338:LEU:CD2	2.05	0.80
1:A:630:VAL:HG12	1:A:633:LYS:HB3	1.66	0.78
1:B:630:VAL:HG12	1:B:633:LYS:HB3	1.66	0.77
1:C:630:VAL:HG12	1:C:633:LYS:HB3	1.66	0.76
1:D:630:VAL:HG12	1:D:633:LYS:HB3	1.67	0.75
1:A:320:MET:HE2	1:A:338:LEU:HD21	1.69	0.74
1:A:320:MET:HG3	1:A:321:ALA:N	2.04	0.72
1:A:472:MET:HG3	1:A:515:TYR:HE1	1.54	0.72
1:B:472:MET:HG3	1:B:515:TYR:HE1	1.54	0.72
1:C:472:MET:HG3	1:C:515:TYR:HE1	1.54	0.72
1:D:472:MET:HG3	1:D:515:TYR:HE1	1.54	0.71
1:B:334:VAL:HG21	1:B:416:PRO:HG2	1.72	0.71
1:D:334:VAL:HG21	1:D:416:PRO:HG2	1.72	0.71
1:A:334:VAL:HG21	1:A:416:PRO:HG2	1.72	0.70
1:C:334:VAL:HG21	1:C:416:PRO:HG2	1.73	0.70
1:B:446:PHE:O	1:B:752:ARG:NH2	2.26	0.69
1:D:446:PHE:O	1:D:752:ARG:NH2	2.27	0.68
1:A:446:PHE:O	1:A:752:ARG:NH2	2.26	0.68
1:C:446:PHE:O	1:C:752:ARG:NH2	2.27	0.68
1:A:320:MET:HE2	1:A:421:ILE:HD11	1.75	0.68
1:A:486:ALA:HB2	1:A:707:GLU:HB2	1.79	0.65
1:B:486:ALA:HB2	1:B:707:GLU:HB2	1.79	0.65
1:A:320:MET:CE	1:A:421:ILE:HD11	2.27	0.64
1:C:486:ALA:HB2	1:C:707:GLU:HB2	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:486:ALA:HB2	1:D:707:GLU:HB2	1.79	0.63
1:D:358:THR:HB	1:D:384:LEU:HD11	1.81	0.62
1:B:358:THR:HB	1:B:384:LEU:HD11	1.81	0.62
1:B:649:GLU:HG2	1:B:698:ILE:HD11	1.82	0.61
1:C:358:THR:HB	1:C:384:LEU:HD11	1.80	0.61
1:A:649:GLU:HG2	1:A:698:ILE:HD11	1.82	0.61
1:A:358:THR:HB	1:A:384:LEU:HD11	1.81	0.61
1:D:649:GLU:HG2	1:D:698:ILE:HD11	1.83	0.60
1:C:649:GLU:HG2	1:C:698:ILE:HD11	1.82	0.59
1:C:468:ASN:ND2	1:C:472:MET:SD	2.75	0.59
1:D:468:ASN:ND2	1:D:472:MET:SD	2.75	0.59
1:A:468:ASN:ND2	1:A:472:MET:SD	2.75	0.59
1:A:320:MET:CE	1:A:421:ILE:CD1	2.81	0.58
1:B:468:ASN:ND2	1:B:472:MET:SD	2.75	0.58
1:C:757:LEU:HD11	1:C:775:ILE:HD11	1.86	0.57
1:A:320:MET:CG	1:A:321:ALA:H	2.12	0.57
1:B:757:LEU:HD11	1:B:775:ILE:HD11	1.87	0.56
1:A:757:LEU:HD11	1:A:775:ILE:HD11	1.87	0.56
1:A:678:ALA:HB1	1:A:684:GLU:HG3	1.88	0.56
1:D:757:LEU:HD11	1:D:775:ILE:HD11	1.86	0.56
1:C:678:ALA:HB1	1:C:684:GLU:HG3	1.88	0.56
1:A:320:MET:HE1	1:A:421:ILE:CD1	2.35	0.56
1:B:678:ALA:HB1	1:B:684:GLU:HG3	1.87	0.56
1:D:565:PRO:HA	1:D:598:LYS:HB2	1.88	0.55
1:C:426:GLY:HA3	1:C:627:TYR:HB3	1.89	0.55
1:D:422:HIS:ND1	1:D:427:GLU:OE2	2.40	0.55
1:C:565:PRO:HA	1:C:598:LYS:HB2	1.88	0.54
1:D:678:ALA:HB1	1:D:684:GLU:HG3	1.88	0.54
1:B:422:HIS:ND1	1:B:427:GLU:OE2	2.40	0.54
1:B:565:PRO:HA	1:B:598:LYS:HB2	1.88	0.54
1:B:654:GLY:HA3	1:B:662:ILE:HD11	1.89	0.54
1:A:426:GLY:HA3	1:A:627:TYR:HB3	1.89	0.54
1:D:654:GLY:HA3	1:D:662:ILE:HD11	1.89	0.54
1:A:565:PRO:HA	1:A:598:LYS:HB2	1.88	0.54
1:B:426:GLY:HA3	1:B:627:TYR:HB3	1.89	0.53
1:C:792:THR:OG1	1:C:793:GLY:N	2.42	0.53
1:D:426:GLY:HA3	1:D:627:TYR:HB3	1.89	0.53
1:C:654:GLY:HA3	1:C:662:ILE:HD11	1.89	0.53
1:A:654:GLY:HA3	1:A:662:ILE:HD11	1.90	0.53
1:B:595:ASP:OD2	1:B:731:LYS:NZ	2.38	0.52
1:C:422:HIS:ND1	1:C:427:GLU:OE2	2.41	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:792:THR:OG1	1:A:793:GLY:N	2.42	0.52
1:A:422:HIS:ND1	1:A:427:GLU:OE2	2.41	0.52
1:B:792:THR:OG1	1:B:793:GLY:N	2.43	0.52
1:A:524:PHE:HE2	1:A:619:LEU:HD13	1.75	0.51
1:A:546:GLY:H	1:A:557:TYR:HD2	1.59	0.51
1:C:546:GLY:H	1:C:557:TYR:HD2	1.59	0.51
1:C:595:ASP:OD2	1:C:731:LYS:NZ	2.39	0.51
1:D:792:THR:OG1	1:D:793:GLY:N	2.41	0.51
1:D:546:GLY:H	1:D:557:TYR:HD2	1.59	0.51
1:C:524:PHE:HE2	1:C:619:LEU:HD13	1.76	0.51
1:D:595:ASP:OD2	1:D:731:LYS:NZ	2.39	0.50
1:B:524:PHE:HE2	1:B:619:LEU:HD13	1.76	0.50
1:B:546:GLY:H	1:B:557:TYR:HD2	1.59	0.49
1:C:734:ARG:HG3	1:C:735:PRO:HD2	1.94	0.49
1:D:524:PHE:HE2	1:D:619:LEU:HD13	1.76	0.48
1:B:388:ASP:CG	1:B:801:ARG:HH22	2.18	0.47
1:C:425:ARG:HD3	1:C:426:GLY:N	2.30	0.47
1:D:425:ARG:HD3	1:D:426:GLY:N	2.30	0.47
1:B:425:ARG:HD3	1:B:426:GLY:N	2.30	0.47
1:C:339:ALA:HB2	1:C:400:THR:HG23	1.97	0.47
1:A:425:ARG:HD3	1:A:426:GLY:N	2.30	0.46
1:D:388:ASP:CG	1:D:801:ARG:HH22	2.18	0.46
1:B:780:MET:HG2	1:B:783:TRP:CE3	2.50	0.46
1:C:388:ASP:CG	1:C:801:ARG:HH22	2.18	0.46
1:C:725:PHE:HB3	1:C:743:TRP:CZ2	2.50	0.46
1:C:780:MET:HG2	1:C:783:TRP:CE3	2.51	0.46
1:A:729:ARG:HD2	1:A:743:TRP:NE1	2.30	0.46
1:C:478:PRO:O	1:C:777:ARG:NH2	2.46	0.46
1:A:320:MET:HE2	1:A:421:ILE:CD1	2.42	0.46
1:A:339:ALA:HB2	1:A:400:THR:HG23	1.97	0.46
1:D:739:SER:OG	1:D:742:GLU:OE2	2.34	0.46
1:B:478:PRO:O	1:B:777:ARG:NH2	2.46	0.46
1:A:780:MET:HG2	1:A:783:TRP:CE3	2.51	0.46
1:D:339:ALA:HB2	1:D:400:THR:HG23	1.98	0.45
1:B:753:ALA:O	1:B:757:LEU:HG	2.16	0.45
1:D:780:MET:HG2	1:D:783:TRP:CE3	2.51	0.45
1:A:595:ASP:OD2	1:A:731:LYS:NZ	2.39	0.45
1:A:320:MET:HB2	1:A:421:ILE:HD11	1.97	0.45
1:A:388:ASP:CG	1:A:801:ARG:HH22	2.18	0.45
1:A:424:LEU:HD23	1:A:797:TYR:HA	1.98	0.45
1:D:424:LEU:HD23	1:D:797:TYR:HA	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:478:PRO:O	1:D:777:ARG:NH2	2.46	0.45
1:D:667:ILE:HG12	1:D:685:PHE:HE1	1.82	0.45
1:C:667:ILE:HG12	1:C:685:PHE:HE1	1.82	0.45
1:B:339:ALA:HB2	1:B:400:THR:HG23	1.97	0.45
1:B:424:LEU:HD23	1:B:797:TYR:HA	1.98	0.45
1:B:667:ILE:HG12	1:B:685:PHE:HE1	1.82	0.45
1:A:753:ALA:O	1:A:757:LEU:HG	2.17	0.44
1:C:725:PHE:HB3	1:C:743:TRP:HZ2	1.82	0.44
1:C:424:LEU:HD23	1:C:797:TYR:HA	1.99	0.44
1:C:753:ALA:O	1:C:757:LEU:HG	2.17	0.44
1:A:478:PRO:O	1:A:777:ARG:NH2	2.46	0.44
1:D:531:LYS:HZ2	1:D:693:HIS:HA	1.83	0.44
1:D:753:ALA:O	1:D:757:LEU:HG	2.17	0.44
1:B:605:ASP:HB3	1:B:608:TRP:HD1	1.83	0.44
1:A:531:LYS:HZ2	1:A:693:HIS:HA	1.83	0.43
1:C:466:HIS:O	1:C:705:ASP:HB3	2.18	0.43
1:A:466:HIS:O	1:A:705:ASP:HB3	2.18	0.43
1:B:587:PHE:O	1:B:734:ARG:NH2	2.52	0.43
1:A:605:ASP:HB3	1:A:608:TRP:HD1	1.83	0.43
1:D:466:HIS:O	1:D:705:ASP:HB3	2.19	0.43
1:A:667:ILE:HG12	1:A:685:PHE:HE1	1.82	0.43
1:A:739:SER:HG	1:A:742:GLU:CD	2.23	0.42
1:D:605:ASP:HB3	1:D:608:TRP:HD1	1.83	0.42
1:B:531:LYS:HZ2	1:B:693:HIS:HA	1.84	0.42
1:C:605:ASP:HB3	1:C:608:TRP:HD1	1.83	0.42
1:C:531:LYS:HZ2	1:C:693:HIS:HA	1.84	0.42
1:D:687:MET:H	1:D:687:MET:HG2	1.66	0.42
1:A:739:SER:N	1:A:742:GLU:OE1	2.52	0.42
1:B:466:HIS:O	1:B:705:ASP:HB3	2.19	0.42
1:D:587:PHE:O	1:D:734:ARG:NH2	2.52	0.42
1:C:734:ARG:CG	1:C:735:PRO:HD2	2.49	0.42
1:A:587:PHE:O	1:A:734:ARG:NH2	2.53	0.42
1:A:429:ILE:HA	1:A:627:TYR:CE2	2.55	0.41
1:B:739:SER:N	1:B:742:GLU:OE1	2.53	0.41
1:C:587:PHE:O	1:C:734:ARG:NH2	2.53	0.41
1:C:429:ILE:HA	1:C:627:TYR:CE2	2.55	0.41
1:B:429:ILE:HA	1:B:627:TYR:CE2	2.55	0.41
1:B:415:TYR:HA	1:B:416:PRO:HD3	1.96	0.41
1:B:657:ARG:HH11	1:B:689:GLY:HA2	1.85	0.41
1:C:744:ASN:O	1:C:748:ARG:HG3	2.20	0.41
1:D:657:ARG:HH11	1:D:689:GLY:HA2	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:320:MET:CG	1:A:321:ALA:N	2.75	0.41
1:B:764:ASP:OD1	1:B:764:ASP:N	2.51	0.41
1:D:429:ILE:HA	1:D:627:TYR:CE2	2.55	0.41
1:D:587:PHE:HA	1:D:588:PRO:HD2	1.96	0.41
1:D:713:ASP:OD1	1:D:714:ALA:N	2.54	0.41
1:B:656:ASN:HD22	1:B:693:HIS:CD2	2.39	0.41
1:C:656:ASN:HD22	1:C:693:HIS:CD2	2.39	0.41
1:D:407:ILE:HD12	1:D:416:PRO:HG3	2.03	0.41
1:A:324:ILE:HD13	1:A:334:VAL:HG22	2.03	0.40
1:B:783:TRP:HA	1:B:784:PRO:HD3	1.96	0.40
1:A:713:ASP:OD1	1:A:714:ALA:N	2.55	0.40
1:A:543:ALA:HA	1:B:494:LEU:HD13	2.04	0.40
1:B:515:TYR:CD1	1:B:522:PRO:HD3	2.57	0.40
1:C:713:ASP:OD1	1:C:714:ALA:N	2.55	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	484/486 (100%)	468 (97%)	16 (3%)	0	100	100
1	B	483/486 (99%)	465 (96%)	18 (4%)	0	100	100
1	C	483/486 (99%)	467 (97%)	16 (3%)	0	100	100
1	D	484/486 (100%)	467 (96%)	17 (4%)	0	100	100
All	All	1934/1944 (100%)	1867 (96%)	67 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	421/421 (100%)	420 (100%)	1 (0%)	93	97
1	B	420/421 (100%)	419 (100%)	1 (0%)	93	97
1	C	420/421 (100%)	419 (100%)	1 (0%)	93	97
1	D	421/421 (100%)	420 (100%)	1 (0%)	93	97
All	All	1682/1684 (100%)	1678 (100%)	4 (0%)	93	97

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	425	ARG
1	B	425	ARG
1	C	425	ARG
1	D	425	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	656	ASN
1	B	656	ASN
1	C	656	ASN
1	D	656	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 22 ligands modelled in this entry, 14 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	$\#\|Z\| > 2$
3	SO4	C	1003	-	4,4,4	0.14	0	6,6,6	0.04	0
3	SO4	B	1005	-	4,4,4	0.14	0	6,6,6	0.05	0
3	SO4	A	1005	-	4,4,4	0.14	0	6,6,6	0.06	0
3	SO4	A	1006	-	4,4,4	0.15	0	6,6,6	0.05	0
3	SO4	C	1004	-	4,4,4	0.14	0	6,6,6	0.06	0
3	SO4	B	1006	-	4,4,4	0.14	0	6,6,6	0.05	0
3	SO4	D	1004	-	4,4,4	0.14	0	6,6,6	0.05	0
3	SO4	D	1005	-	4,4,4	0.14	0	6,6,6	0.04	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	486/486 (100%)	0.03	4 (0%) 86 83	71, 92, 110, 129	0
1	B	485/486 (99%)	0.06	6 (1%) 79 74	77, 93, 111, 128	0
1	C	485/486 (99%)	0.47	29 (5%) 21 16	77, 140, 168, 178	0
1	D	486/486 (100%)	0.75	63 (12%) 3 4	77, 167, 204, 229	0
All	All	1942/1944 (99%)	0.33	102 (5%) 26 23	71, 101, 196, 229	0

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	547	TYR	4.9
1	D	553	GLU	4.9
1	D	451	LEU	4.8
1	D	555	GLU	4.6
1	D	600	SER	4.5
1	D	537	ALA	4.3
1	D	496	SER	4.2
1	C	555	GLU	3.9
1	D	444	GLY	3.9
1	D	708	THR	3.9
1	D	484	GLY	3.9
1	D	640	LEU	3.7
1	D	601	LEU	3.6
1	D	544	PHE	3.5
1	D	765	HIS	3.5
1	C	437	LEU	3.5
1	A	320	MET	3.5
1	D	756	LEU	3.4
1	D	554	GLU	3.4
1	D	546	GLY	3.4
1	D	718	TRP	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	455	PRO	3.3
1	D	485	LEU	3.2
1	D	558	TRP	3.2
1	D	486	ALA	3.2
1	D	465	PRO	3.2
1	C	600	SER	3.2
1	C	451	LEU	3.2
1	D	445	ASN	3.1
1	C	553	GLU	3.1
1	D	487	ASP	3.1
1	B	444	GLY	3.1
1	C	444	GLY	3.1
1	D	715	ASN	3.0
1	D	759	GLN	3.0
1	D	539	THR	3.0
1	D	551	ILE	3.0
1	C	438	PRO	2.9
1	D	749	ASP	2.9
1	D	761	TRP	2.9
1	C	672	TRP	2.9
1	D	548	LYS	2.8
1	D	491	PRO	2.8
1	C	715	ASN	2.8
1	D	747	MET	2.7
1	D	535	GLY	2.7
1	B	545	LYS	2.7
1	C	456	LYS	2.6
1	D	764	ASP	2.6
1	C	448	LEU	2.6
1	C	764	ASP	2.6
1	D	552	ASP	2.6
1	D	760	HIS	2.5
1	C	445	ASN	2.5
1	B	548	LYS	2.5
1	D	763	GLY	2.5
1	D	706	MET	2.5
1	D	455	PRO	2.5
1	C	465	PRO	2.5
1	D	538	VAL	2.5
1	D	748	ARG	2.5
1	D	343	THR	2.5
1	D	495	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	549	VAL	2.4
1	C	554	GLU	2.4
1	C	679	GLU	2.4
1	B	445	ASN	2.3
1	D	556	TYR	2.3
1	C	757	LEU	2.3
1	D	557	TYR	2.3
1	D	463	VAL	2.3
1	C	459	TYR	2.3
1	C	463	VAL	2.3
1	C	447	THR	2.3
1	C	558	TRP	2.3
1	C	678	ALA	2.3
1	D	757	LEU	2.2
1	A	640	LEU	2.2
1	C	544	PHE	2.2
1	D	492	LEU	2.2
1	D	681	LEU	2.2
1	D	550	ILE	2.2
1	D	719	ASP	2.2
1	D	456	LYS	2.1
1	D	750	VAL	2.1
1	D	599	VAL	2.1
1	B	677	ASN	2.1
1	C	452	ILE	2.1
1	B	738	ILE	2.1
1	C	545	LYS	2.1
1	C	676	ILE	2.1
1	D	458	TYR	2.1
1	C	343	THR	2.1
1	D	672	TRP	2.1
1	D	501	TRP	2.1
1	D	771	ASP	2.1
1	D	559	PRO	2.0
1	C	601	LEU	2.0
1	D	774	THR	2.0
1	D	738	ILE	2.0
1	A	464	LYS	2.0
1	A	455	PRO	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 monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	CO	B	1003	1/1	0.26	0.17	114,114,114,114	1
2	CO	D	1003	1/1	0.67	0.21	128,128,128,128	1
3	SO4	D	1005	5/5	0.69	0.23	175,175,175,175	0
3	SO4	C	1004	5/5	0.72	0.27	156,156,156,156	0
2	CO	C	1002	1/1	0.75	0.16	126,126,126,126	1
3	SO4	D	1004	5/5	0.80	0.41	178,178,178,178	0
2	CO	A	1004	1/1	0.81	0.26	112,112,112,112	1
2	CO	A	1003	1/1	0.84	0.27	98,98,98,98	1
2	CO	D	1001	1/1	0.84	0.42	150,150,150,150	1
2	CO	C	1001	1/1	0.84	0.18	127,127,127,127	1
3	SO4	C	1003	5/5	0.89	0.31	128,128,128,128	0
3	SO4	A	1006	5/5	0.89	0.16	105,105,105,105	5
2	CO	D	1002	1/1	0.92	0.21	88,88,88,88	0
3	SO4	B	1006	5/5	0.92	0.16	107,107,107,107	0
2	CO	B	1004	1/1	0.92	0.14	121,121,121,121	0
2	CO	A	1001	1/1	0.93	0.12	85,85,85,85	0
2	CO	A	1002	1/1	0.94	0.17	73,73,73,73	0
3	SO4	A	1005	5/5	0.96	0.31	99,99,99,99	0
2	CO	B	1002	1/1	0.96	0.19	79,79,79,79	0
3	SO4	B	1005	5/5	0.96	0.31	94,94,94,94	0
2	CO	B	1001	1/1	0.96	0.10	84,84,84,84	0
2	CO	A	1007	1/1	0.97	0.19	83,83,83,83	0

6.5 Other polymers (i)

There are no such residues in this entry.