Full wwPDB X-ray Structure Validation Report (i)

Jun 18, 2024 – 05:41 PM EDT

PDB ID	:	4MHB
Title	:	Structure of a putative reductase from Yersinia pestis
Authors	:	Anderson, S.M.; Wawrzak, Z.; Kudritska, M.; Kwon, K.; Rembert, P.; Savchenko, A.; Anderson, W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on	:	2013-08-29
Resolution	:	1.75 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:

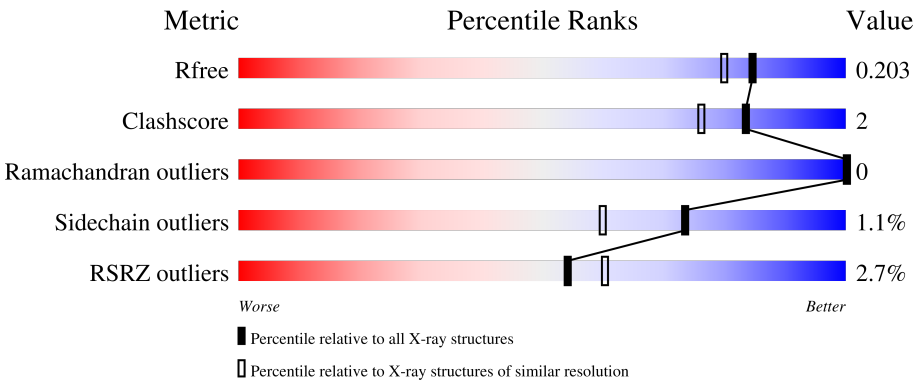
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 1.75 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	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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	297	% 89% 6% 5%
1	B	297	3% 91% • 5%
1	C	297	2% 92% • • 5%
1	D	297	2% 89% 6% 5%
1	E	297	3% 90% • • 5%

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Mol	Chain	Length	Quality of chain
1	F	297	3% 89% 5% 5%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	D	304	-	-	X	X

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 15426 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.

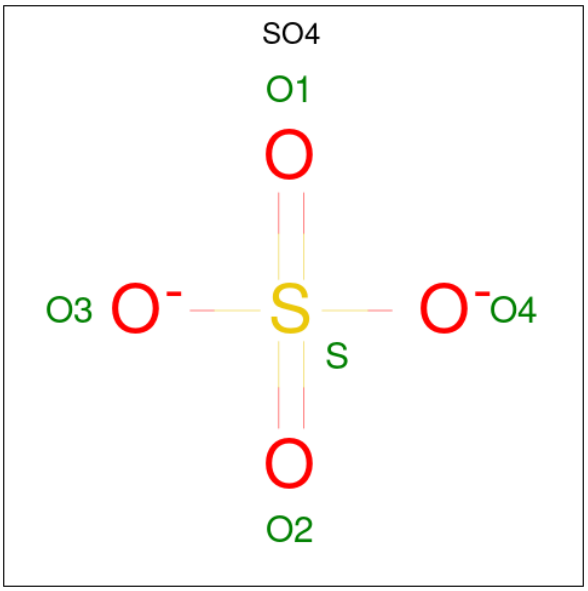
- Molecule 1 is a protein called Putative aldo/keto reductase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	283	2253	1435	394	413	1	10	0	1	0
1	B	283	2255	1436	394	414	1	10	0	1	0
1	C	283	2253	1435	394	413	1	10	0	1	0
1	D	282	2246	1431	392	412	1	10	0	1	0
1	E	282	2254	1436	395	412	1	10	0	2	0
1	F	282	2246	1431	392	412	1	10	0	1	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	EXPRESSION TAG	UNP Q7CJX8
B	1	MSE	-	EXPRESSION TAG	UNP Q7CJX8
C	1	MSE	-	EXPRESSION TAG	UNP Q7CJX8
D	1	MSE	-	EXPRESSION TAG	UNP Q7CJX8
E	1	MSE	-	EXPRESSION TAG	UNP Q7CJX8
F	1	MSE	-	EXPRESSION TAG	UNP Q7CJX8

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



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

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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	F	1	5	4	1	0	0
			Total	O	S		
2	F	1	5	4	1	0	0
			Total	O	S		
2	F	1	5	4	1	0	0
			Total	O	S		
2	F	1	5	4	1	0	0
			Total	O	S		
2	F	1	5	4	1	0	0
			Total	O	S		
2	F	1	5	4	1	0	0
			Total	O	S		
2	F	1	5	4	1	0	0

- Molecule 3 is water.

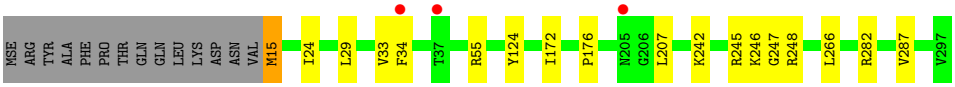
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	335	338	338	0	3
			Total	O		
3	B	262	266	266	0	4
			Total	O		
3	C	293	299	299	0	6
			Total	O		
3	D	279	285	285	0	6
			Total	O		
3	E	268	271	271	0	3
			Total	O		
3	F	245	250	250	0	5

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: Putative aldo/keto reductase

Chain A:

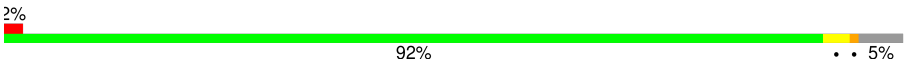


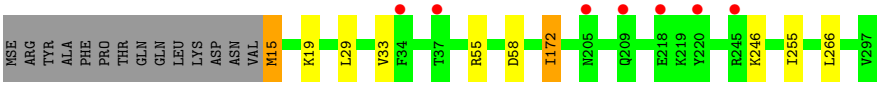
• Molecule 1: Putative aldo/keto reductase

Chain B:

• Molecule 1: Putative aldo/keto reductase

Chain C:





• Molecule 1: Putative aldo/keto reductase

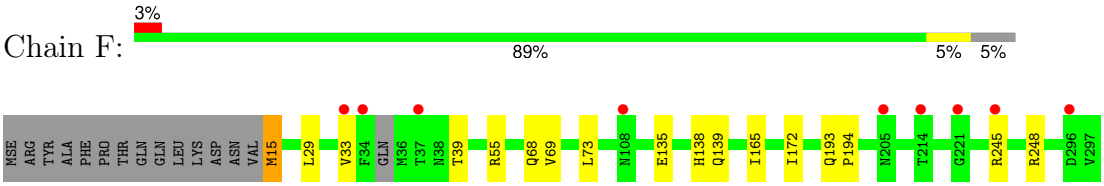
Chain D:

• Molecule 1: Putative aldo/keto reductase

Chain E:

• Molecule 1: Putative aldo/keto reductase

Chain F:


3%
89%
5%
5%
MSE
ARG
TYR
ALA
PHE
PRO
THR
GLN
GLN
LEU
LYS
ASP
ASN
VAL
M15
L29
V33
F34
GLN
M36
T37
N38
T39
R55
Q68
V69
L73
N108
E135
H138
Q139
I165
I172
Q193
P194
N205
T214
G221
R245
R248
D296
V297

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α, β, γ	128.75Å 128.75Å 277.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.40 – 1.75 36.36 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.1 (36.40-1.75) 86.0 (36.36-1.75)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.10 (at 1.75Å)	Xtriage
Refinement program	PHENIX dev_1101	Depositor
R, R_{free}	0.178 , 0.202 0.179 , 0.203	Depositor DCC
R_{free} test set	11600 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	20.0	Xtriage
Anisotropy	0.348	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 53.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.53$, $< L^2 > = 0.37$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	15426	wwPDB-VP
Average B, all atoms (Å^2)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 65.21 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 7.3469e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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: 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	0.86	0/2293	0.76	2/3092 (0.1%)
1	B	0.72	0/2295	0.69	0/3095
1	C	0.76	0/2293	0.70	0/3092
1	D	0.76	0/2286	0.71	0/3083
1	E	0.72	0/2296	0.68	0/3094
1	F	0.67	0/2286	0.68	0/3083
All	All	0.75	0/13749	0.70	2/18539 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	282	ARG	NE-CZ-NH1	-5.39	117.60	120.30
1	A	266	LEU	CB-CG-CD1	-5.32	101.96	111.00

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	2253	0	2227	13	0
1	B	2255	0	2232	8	0
1	C	2253	0	2227	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2246	0	2223	15	0
1	E	2254	0	2236	8	0
1	F	2246	0	2223	13	0
2	A	40	0	0	2	0
2	B	30	0	0	1	0
2	C	40	0	0	1	0
2	D	30	0	0	4	0
2	E	35	0	0	1	0
2	F	35	0	0	0	0
3	A	338	0	0	0	0
3	B	266	0	0	1	0
3	C	299	0	0	1	0
3	D	285	0	0	1	0
3	E	271	0	0	1	0
3	F	250	0	0	0	0
All	All	15426	0	13368	66	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 2.

All (66) 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:34:PHE:HD2	1:A:242:LYS:HD2	1.41	0.84
1:F:245:ARG:HG2	1:F:248:ARG:HG3	1.71	0.72
1:B:38:ASN:OD1	1:B:41:GLU:N	2.23	0.69
1:D:210:ASN:O	1:D:214:THR:HG23	1.93	0.68
1:E:38:ASN:HB3	1:E:41:GLU:HB3	1.76	0.67
1:D:184:VAL:HB	1:D:185:PRO:HD3	1.77	0.67
1:B:210:ASN:O	1:B:214:THR:HG23	1.97	0.65
1:E:36:MSE:HG2	1:E:242:LYS:NZ	2.12	0.64
1:A:15[A]:MSE:HE1	1:A:29:LEU:HD12	1.80	0.63
1:E:15[A]:MSE:HE1	1:E:29:LEU:HD12	1.83	0.60
1:E:36:MSE:HG2	1:E:242:LYS:HZ1	1.65	0.59
1:B:15[A]:MSE:HE1	1:B:29:LEU:HD12	1.83	0.58
1:D:135:GLU:HG2	1:D:165:ILE:HD12	1.86	0.58
1:D:138:HIS:HD2	1:D:165:ILE:HD13	1.68	0.57
1:A:246:LYS:NZ	2:A:307:SO4:O3	2.40	0.55
1:F:138:HIS:HD2	1:F:165:ILE:HD13	1.70	0.55
1:D:15[A]:MSE:HE1	1:D:29:LEU:HD12	1.87	0.55
1:F:39:THR:HG23	1:F:68:GLN:CG	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:PHE:CD2	1:A:242:LYS:HD2	2.31	0.55
1:D:152:HIS:CE1	2:D:304:SO4:O1	2.61	0.54
1:D:74:LYS:HD3	2:D:303:SO4:O3	2.07	0.54
1:F:135:GLU:HG2	1:F:165:ILE:HD12	1.92	0.52
1:F:39:THR:HG23	1:F:68:GLN:HG3	1.92	0.52
1:A:248:ARG:NH1	2:A:308:SO4:O1	2.43	0.52
1:C:19:LYS:NZ	3:C:655:HOH:O	2.42	0.52
1:F:15[B]:MSE:HE1	1:F:55:ARG:HG3	1.91	0.51
1:C:246:LYS:NZ	2:C:306:SO4:O2	2.44	0.51
1:A:34:PHE:HB2	1:A:242:LYS:NZ	2.25	0.51
1:D:152:HIS:CD2	2:D:304:SO4:O1	2.64	0.51
1:D:184:VAL:HG13	1:D:194:PRO:HG3	1.95	0.48
1:C:29:LEU:HD21	1:C:255:ILE:HD11	1.96	0.48
1:C:15[A]:MSE:HE1	1:C:29:LEU:HB2	1.97	0.47
1:D:36:MSE:O	1:D:36:MSE:HG3	2.13	0.47
1:D:33:VAL:HG23	1:D:33:VAL:O	2.13	0.47
1:D:104:GLU:HG2	3:D:658:HOH:O	2.13	0.47
1:A:33:VAL:O	1:A:34:PHE:CG	2.67	0.47
1:A:245:ARG:NH1	1:A:247:GLY:H	2.12	0.47
1:B:15[B]:MSE:HE1	1:B:55:ARG:HG3	1.96	0.47
1:C:15[B]:MSE:HE1	1:C:55:ARG:HG3	1.97	0.47
1:F:33:VAL:O	1:F:33:VAL:HG23	2.14	0.47
1:D:181:LEU:O	1:D:185:PRO:HG2	2.15	0.46
1:F:135:GLU:O	1:F:139:GLN:HG3	2.16	0.45
1:F:15[A]:MSE:HE1	1:F:29:LEU:HD12	1.97	0.45
1:A:15[B]:MSE:HE1	1:A:55:ARG:HG3	1.99	0.45
1:B:47:ILE:HD11	1:B:75:LEU:HD12	1.98	0.45
1:C:33:VAL:HG22	1:C:58:ASP:O	2.16	0.44
1:D:152:HIS:NE2	2:D:304:SO4:O1	2.50	0.44
1:F:15[A]:MSE:HE1	1:F:29:LEU:HB2	2.00	0.44
1:D:210:ASN:HA	1:D:211:PRO:HD3	1.85	0.44
1:C:266:LEU:HD12	1:C:266:LEU:HA	1.75	0.44
1:A:176:PRO:HD2	1:A:207:LEU:HD11	2.01	0.43
1:E:33:VAL:HG22	1:E:58:ASP:O	2.19	0.43
1:C:15[A]:MSE:HE1	1:C:29:LEU:HD12	2.00	0.43
1:C:172:ILE:H	1:C:172:ILE:HD13	1.83	0.42
1:E:74:LYS:HD3	2:E:303:SO4:O3	2.20	0.42
1:B:246:LYS:NZ	2:B:303:SO4:O2	2.53	0.42
1:B:225:GLY:HA3	3:B:575:HOH:O	2.20	0.41
1:F:69:VAL:O	1:F:73:LEU:HG	2.20	0.41
1:E:248[B]:ARG:NH1	3:E:616:HOH:O	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:ARG:HH11	1:A:247:GLY:H	1.68	0.41
1:A:124:TYR:CE1	1:A:287:VAL:HG22	2.56	0.41
1:B:232:ILE:HG23	1:B:237:ILE:HB	2.03	0.41
1:F:15[B]:MSE:HE1	1:F:55:ARG:CG	2.50	0.41
1:F:193:GLN:NE2	1:F:194:PRO:HD2	2.35	0.41
1:A:24:ILE:HG21	1:A:24:ILE:HD13	1.92	0.40
1:C:266:LEU:HD21	1:E:188:GLN:HB3	2.02	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	281/297 (95%)	274 (98%)	7 (2%)	0	100	100
1	B	281/297 (95%)	276 (98%)	5 (2%)	0	100	100
1	C	281/297 (95%)	276 (98%)	5 (2%)	0	100	100
1	D	278/297 (94%)	273 (98%)	5 (2%)	0	100	100
1	E	279/297 (94%)	275 (99%)	4 (1%)	0	100	100
1	F	278/297 (94%)	273 (98%)	5 (2%)	0	100	100
All	All	1678/1782 (94%)	1647 (98%)	31 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	234/237 (99%)	231 (99%)	3 (1%)	69	54
1	B	235/237 (99%)	232 (99%)	3 (1%)	69	54
1	C	234/237 (99%)	231 (99%)	3 (1%)	69	54
1	D	234/237 (99%)	230 (98%)	4 (2%)	60	42
1	E	235/237 (99%)	229 (97%)	6 (3%)	46	23
1	F	234/237 (99%)	231 (99%)	3 (1%)	69	54
All	All	1406/1422 (99%)	1384 (98%)	22 (2%)	73	45

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15[A]	MSE
1	A	15[B]	MSE
1	A	172	ILE
1	B	15[A]	MSE
1	B	15[B]	MSE
1	B	172	ILE
1	C	15[A]	MSE
1	C	15[B]	MSE
1	C	172	ILE
1	D	15[A]	MSE
1	D	15[B]	MSE
1	D	172	ILE
1	D	207	LEU
1	E	15[A]	MSE
1	E	15[B]	MSE
1	E	36	MSE
1	E	172	ILE
1	E	207	LEU
1	E	245	ARG
1	F	15[A]	MSE
1	F	15[B]	MSE
1	F	172	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	122	GLN

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)

42 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	SO4	D	303	-	4,4,4	0.36	0	6,6,6	0.53	0
2	SO4	F	306	-	4,4,4	0.26	0	6,6,6	0.63	0
2	SO4	C	308	-	4,4,4	0.25	0	6,6,6	0.26	0
2	SO4	A	307	-	4,4,4	0.34	0	6,6,6	0.41	0
2	SO4	B	303	-	4,4,4	0.30	0	6,6,6	0.28	0
2	SO4	F	307	-	4,4,4	0.25	0	6,6,6	0.34	0
2	SO4	A	306	-	4,4,4	0.23	0	6,6,6	0.27	0
2	SO4	B	305	-	4,4,4	0.31	0	6,6,6	0.74	0
2	SO4	C	305	-	4,4,4	0.54	0	6,6,6	0.67	0
2	SO4	A	305	-	4,4,4	0.37	0	6,6,6	0.35	0
2	SO4	E	306	-	4,4,4	0.33	0	6,6,6	0.20	0
2	SO4	F	305	-	4,4,4	0.36	0	6,6,6	0.60	0
2	SO4	B	301	-	4,4,4	0.37	0	6,6,6	0.29	0
2	SO4	C	306	-	4,4,4	0.29	0	6,6,6	0.40	0
2	SO4	D	305	-	4,4,4	0.28	0	6,6,6	0.41	0
2	SO4	D	302	-	4,4,4	0.29	0	6,6,6	0.16	0
2	SO4	C	302	-	4,4,4	0.28	0	6,6,6	0.36	0
2	SO4	E	307	-	4,4,4	0.31	0	6,6,6	0.12	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	SO4	E	301	-	4,4,4	0.24	0	6,6,6	0.44	0
2	SO4	C	301	-	4,4,4	0.40	0	6,6,6	0.58	0
2	SO4	F	301	-	4,4,4	0.04	0	6,6,6	0.58	0
2	SO4	D	301	-	4,4,4	0.20	0	6,6,6	0.46	0
2	SO4	C	303	-	4,4,4	0.32	0	6,6,6	0.12	0
2	SO4	B	306	-	4,4,4	0.29	0	6,6,6	0.19	0
2	SO4	F	303	-	4,4,4	0.27	0	6,6,6	0.17	0
2	SO4	A	301	-	4,4,4	0.18	0	6,6,6	0.45	0
2	SO4	B	304	-	4,4,4	0.17	0	6,6,6	0.21	0
2	SO4	E	303	-	4,4,4	0.41	0	6,6,6	0.44	0
2	SO4	A	303	-	4,4,4	0.32	0	6,6,6	0.43	0
2	SO4	A	302	-	4,4,4	0.28	0	6,6,6	0.29	0
2	SO4	C	304	-	4,4,4	0.40	0	6,6,6	0.50	0
2	SO4	E	304	-	4,4,4	0.29	0	6,6,6	0.15	0
2	SO4	A	304	-	4,4,4	0.31	0	6,6,6	0.14	0
2	SO4	E	305	-	4,4,4	0.31	0	6,6,6	0.19	0
2	SO4	D	304	-	4,4,4	0.49	0	6,6,6	0.61	0
2	SO4	A	308	-	4,4,4	0.29	0	6,6,6	0.24	0
2	SO4	B	302	-	4,4,4	0.34	0	6,6,6	0.30	0
2	SO4	F	302	-	4,4,4	0.19	0	6,6,6	0.90	0
2	SO4	F	304	-	4,4,4	0.28	0	6,6,6	0.14	0
2	SO4	E	302	-	4,4,4	0.35	0	6,6,6	0.29	0
2	SO4	C	307	-	4,4,4	0.28	0	6,6,6	0.16	0
2	SO4	D	306	-	4,4,4	0.31	0	6,6,6	0.16	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.

7 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	303	SO4	1	0
2	A	307	SO4	1	0
2	B	303	SO4	1	0
2	C	306	SO4	1	0
2	E	303	SO4	1	0
2	D	304	SO4	3	0
2	A	308	SO4	1	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (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	274/297 (92%)	-0.21	3 (1%)	80	86	12, 19, 38, 77	0
1	B	274/297 (92%)	0.04	10 (3%)	42	49	15, 25, 44, 77	0
1	C	274/297 (92%)	-0.18	7 (2%)	56	62	14, 23, 42, 86	0
1	D	273/297 (91%)	-0.12	6 (2%)	62	69	14, 21, 37, 80	0
1	E	273/297 (91%)	-0.02	9 (3%)	46	53	15, 25, 41, 71	0
1	F	273/297 (91%)	-0.07	9 (3%)	46	53	17, 27, 44, 91	0
All	All	1641/1782 (92%)	-0.09	44 (2%)	54	60	12, 23, 42, 91	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	37	THR	8.5
1	C	37	THR	8.2
1	F	37	THR	7.7
1	C	34	PHE	7.1
1	E	37	THR	6.6
1	E	34	PHE	5.6
1	D	37	THR	5.5
1	A	34	PHE	5.1
1	D	34	PHE	4.8
1	A	37	THR	4.7
1	B	34	PHE	4.3
1	F	34	PHE	4.1
1	B	139	GLN	3.9
1	D	33	VAL	3.7
1	E	205	ASN	3.7
1	B	35	GLN	3.5
1	F	33	VAL	3.5
1	B	140	ALA	3.1
1	F	108	ASN	3.1

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Mol	Chain	Res	Type	RSRZ
1	E	139	GLN	3.1
1	E	38	ASN	2.9
1	B	205	ASN	2.9
1	E	140	ALA	2.9
1	B	33	VAL	2.8
1	E	33	VAL	2.8
1	C	205	ASN	2.8
1	D	31	PHE	2.7
1	C	245	ARG	2.5
1	F	205	ASN	2.5
1	C	209	GLN	2.4
1	F	214	THR	2.4
1	F	296	ASP	2.4
1	D	38	ASN	2.3
1	D	205	ASN	2.3
1	E	148	VAL	2.3
1	C	220	TYR	2.3
1	E	108	ASN	2.3
1	F	245	ARG	2.2
1	F	221	GLY	2.2
1	A	205	ASN	2.1
1	B	296	ASP	2.1
1	B	258	PHE	2.1
1	C	218	GLU	2.1
1	B	209	GLN	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	SO4	C	306	5/5	0.44	0.37	86,92,94,94	0
2	SO4	F	306	5/5	0.52	0.32	70,72,80,82	0
2	SO4	E	303	5/5	0.53	0.27	76,79,84,85	0
2	SO4	D	303	5/5	0.55	0.26	31,37,44,45	5
2	SO4	D	304	5/5	0.58	0.42	23,28,39,46	5
2	SO4	A	307	5/5	0.59	0.28	70,77,80,82	0
2	SO4	D	302	5/5	0.68	0.21	95,95,97,98	0
2	SO4	B	306	5/5	0.69	0.26	93,94,96,96	0
2	SO4	E	305	5/5	0.70	0.32	69,73,79,81	0
2	SO4	F	307	5/5	0.71	0.33	86,87,90,92	0
2	SO4	F	303	5/5	0.72	0.26	89,91,92,94	0
2	SO4	D	306	5/5	0.73	0.31	67,71,78,78	0
2	SO4	B	303	5/5	0.74	0.25	78,83,86,86	0
2	SO4	D	305	5/5	0.75	0.19	75,77,80,83	0
2	SO4	C	304	5/5	0.77	0.22	79,84,85,85	0
2	SO4	F	304	5/5	0.79	0.29	91,94,95,96	0
2	SO4	E	307	5/5	0.79	0.21	93,93,95,95	0
2	SO4	A	306	5/5	0.79	0.21	34,34,36,40	5
2	SO4	B	305	5/5	0.80	0.21	38,39,42,43	5
2	SO4	A	304	5/5	0.81	0.31	78,79,81,83	0
2	SO4	E	302	5/5	0.82	0.19	78,79,81,84	0
2	SO4	E	306	5/5	0.82	0.29	86,88,89,90	0
2	SO4	F	305	5/5	0.82	0.34	19,23,44,44	5
2	SO4	B	302	5/5	0.82	0.19	66,69,71,73	0
2	SO4	F	301	5/5	0.82	0.23	82,84,89,90	0
2	SO4	A	302	5/5	0.83	0.28	83,85,86,87	0
2	SO4	C	308	5/5	0.84	0.28	70,76,80,81	0
2	SO4	C	305	5/5	0.84	0.32	18,18,39,41	5
2	SO4	E	304	5/5	0.84	0.22	93,94,96,96	0
2	SO4	A	301	5/5	0.85	0.18	70,76,77,77	0
2	SO4	A	305	5/5	0.85	0.20	71,75,77,77	0
2	SO4	C	302	5/5	0.85	0.26	76,76,77,79	0
2	SO4	A	308	5/5	0.87	0.24	59,67,72,74	0
2	SO4	C	307	5/5	0.87	0.37	90,91,92,93	0
2	SO4	B	304	5/5	0.89	0.16	37,40,43,45	5
2	SO4	C	303	5/5	0.90	0.35	64,70,71,73	0
2	SO4	F	302	5/5	0.91	0.26	33,53,59,59	0
2	SO4	C	301	5/5	0.95	0.21	36,50,52,52	0
2	SO4	B	301	5/5	0.95	0.16	29,49,50,54	0
2	SO4	A	303	5/5	0.96	0.17	24,42,45,46	0
2	SO4	D	301	5/5	0.97	0.15	24,40,45,45	0
2	SO4	E	301	5/5	0.98	0.14	30,45,47,51	0

6.5 Other polymers (i)

There are no such residues in this entry.