Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 06:22 PM GMT

PDB ID	:	1AFW
Title	:	THE 1.8 ANGSTROM CRYSTAL STRUCTURE OF THE DIMERIC PEROXISOMAL THIOLASE OF SACCHAROMYCES CEREVISIAE
Authors	:	Mathieu, M.; Wierenga, R.K.
Deposited on	:	1997-03-15
Resolution	:	1.80 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:

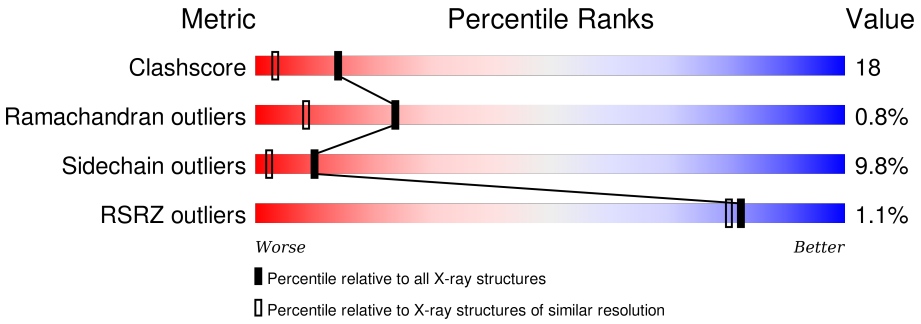
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 1.80 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))
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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	393	% 64% 30% 5%
1	B	393	% 64% 30% 6%

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	MRD	A	501	-	-	-	X

2 Entry composition

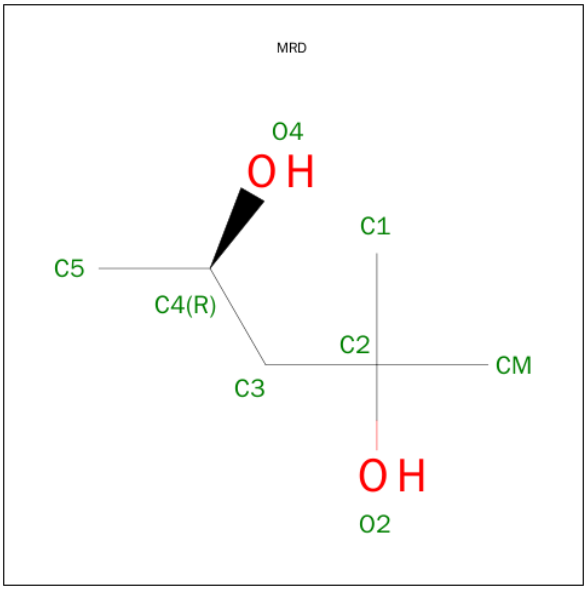
There are 3 unique types of molecules in this entry. The entry contains 6338 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 3-KETOACETYL-COA THIOLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	390	Total	C	N	O	S	0	0	0
			2923	1841	509	559	14			
1	B	393	Total	C	N	O	S	0	0	0
			2951	1857	516	564	14			

- Molecule 2 is (4R)-2-METHYLPENTANE-2,4-DIOL (three-letter code: MRD) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			8	6	2		
2	B	1	Total	C	O	0	0
			8	6	2		

- Molecule 3 is water.

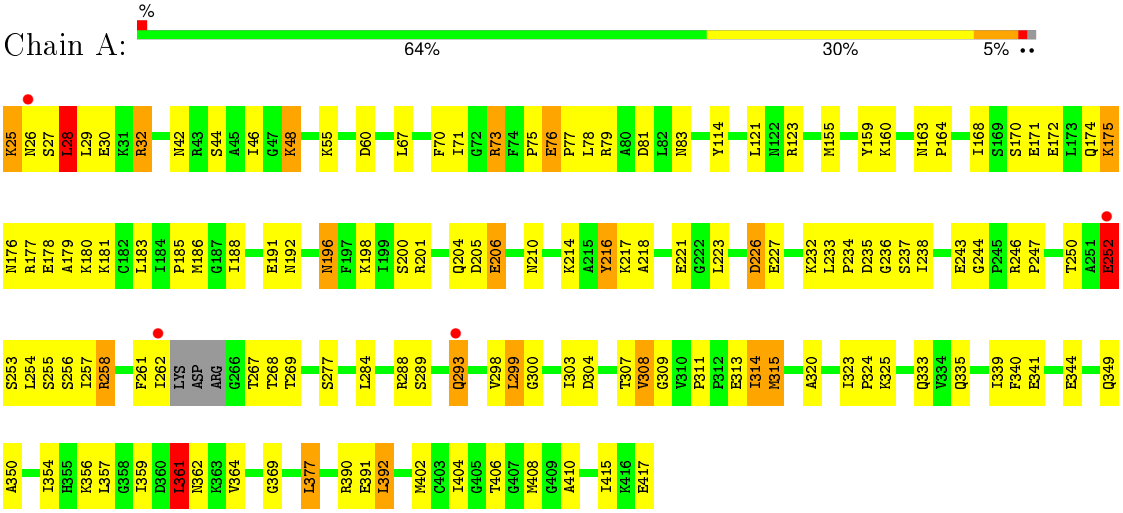
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	208	Total O 208 208	0	0
3	B	240	Total O 240 240	0	0

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.

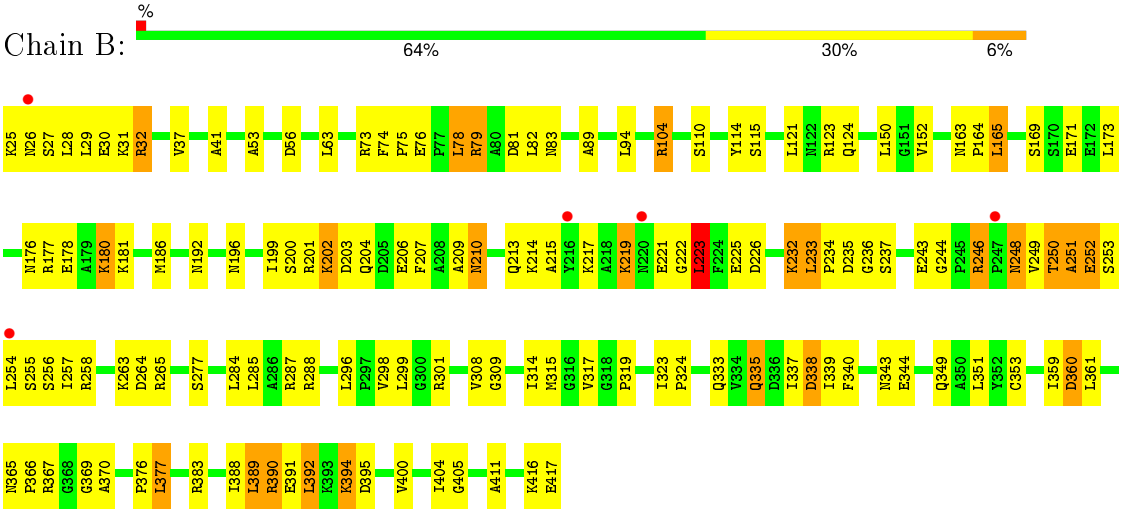
- Molecule 1: 3-KETOACETYL-COA THIOLASE

Chain A:



- Molecule 1: 3-KETOACETYL-COA THIOLASE

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	71.17Å 92.65Å 116.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 1.80 9.96 – 1.80	Depositor EDS
% Data completeness (in resolution range)	98.9 (10.00-1.80) 99.6 (9.96-1.80)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$< I/\sigma(I) >$ [1]	12.34 (at 1.80Å)	Xtriage
Refinement program	TNT V. 5-E	Depositor
R, R_{free}	0.203 , 0.240 0.187 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å^2)	17.1	Xtriage
Anisotropy	0.074	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 118.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.45$, $< L^2 > = 0.27$	Xtriage
Outliers	0 of 71365 reflections	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	6338	wwPDB-VP
Average B, all atoms (Å^2)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.40% 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: MRD

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	0.93	0/2966	1.34	18/4013 (0.4%)
1	B	0.95	1/2995 (0.0%)	1.39	25/4052 (0.6%)
All	All	0.94	1/5961 (0.0%)	1.37	43/8065 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	1	0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	41	ALA	CA-CB	-5.33	1.41	1.52

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	56	ASP	CB-CG-OD2	-8.05	111.05	118.30
1	B	123	ARG	NE-CZ-NH2	-7.74	116.43	120.30
1	B	244	GLY	N-CA-C	7.24	131.19	113.10
1	B	308	VAL	CG1-CB-CG2	-7.09	99.56	110.90
1	A	252	GLU	CB-CA-C	7.02	124.43	110.40
1	A	123	ARG	NE-CZ-NH1	-6.89	116.86	120.30
1	B	299	LEU	CB-CG-CD2	-6.62	99.74	111.00
1	A	60	ASP	CB-CG-OD1	6.49	124.14	118.30
1	A	32	ARG	NE-CZ-NH2	-6.46	117.07	120.30
1	B	308	VAL	CB-CA-C	-6.45	99.15	111.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	304	ASP	CB-CG-OD1	6.38	124.04	118.30
1	B	123	ARG	NE-CZ-NH1	6.30	123.45	120.30
1	B	377	LEU	CB-CG-CD2	-6.29	100.30	111.00
1	A	55	LYS	CD-CE-NZ	-6.17	97.50	111.70
1	A	315	MET	CG-SD-CE	6.16	110.06	100.20
1	B	31	LYS	CD-CE-NZ	-6.03	97.83	111.70
1	A	226	ASP	CB-CG-OD2	6.00	123.70	118.30
1	B	288	ARG	NE-CZ-NH1	5.97	123.28	120.30
1	B	78	LEU	CB-CG-CD2	-5.97	100.86	111.00
1	B	152	VAL	CG1-CB-CG2	-5.93	101.42	110.90
1	A	298	VAL	CB-CA-C	-5.91	100.17	111.40
1	A	392	LEU	CB-CG-CD1	-5.91	100.95	111.00
1	B	223	LEU	CB-CG-CD2	-5.87	101.02	111.00
1	B	360	ASP	CB-CG-OD1	5.75	123.47	118.30
1	B	186	MET	CG-SD-CE	-5.72	91.05	100.20
1	B	246	ARG	CG-CD-NE	5.71	123.80	111.80
1	A	123	ARG	NE-CZ-NH2	5.70	123.15	120.30
1	B	392	LEU	CB-CG-CD1	-5.62	101.45	111.00
1	A	28	LEU	CB-CG-CD2	-5.46	101.71	111.00
1	A	216	TYR	CA-CB-CG	-5.42	103.11	113.40
1	B	94	LEU	N-CA-C	5.31	125.33	111.00
1	B	173	LEU	CA-CB-CG	-5.31	103.10	115.30
1	A	254	LEU	CB-CG-CD1	5.29	120.00	111.00
1	A	32	ARG	NE-CZ-NH1	5.28	122.94	120.30
1	A	356	LYS	CD-CE-NZ	5.28	123.83	111.70
1	A	44	SER	N-CA-C	-5.25	96.83	111.00
1	B	32	ARG	NE-CZ-NH1	5.23	122.92	120.30
1	B	104	ARG	NE-CZ-NH2	5.18	122.89	120.30
1	B	376	PRO	N-CA-C	-5.16	98.68	112.10
1	B	338	ASP	CB-CG-OD2	5.13	122.92	118.30
1	A	361	LEU	CA-CB-CG	-5.12	103.54	115.30
1	B	199	ILE	CG1-CB-CG2	5.10	122.62	111.40
1	B	177	ARG	NE-CZ-NH2	-5.02	117.79	120.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	252	GLU	CA

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	2923	0	2969	113	0
1	B	2951	0	3000	109	0
2	A	8	0	14	0	0
2	B	8	0	14	0	0
3	A	208	0	0	7	0
3	B	240	0	0	17	0
All	All	6338	0	5997	217	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 18.

All (217) 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:B:207:PHE:HB2	1:B:351:LEU:HD21	1.34	1.08
1:A:354:ILE:HD12	1:A:364:VAL:HG21	1.43	1.00
1:A:188:ILE:HD13	1:A:262:ILE:HD11	1.45	0.95
1:A:201:ARG:HB2	1:A:267:THR:HG22	1.56	0.86
1:B:73:ARG:HG3	1:B:73:ARG:HH11	1.44	0.81
1:B:201:ARG:HD3	3:B:615:HOH:O	1.83	0.78
1:A:48:LYS:NZ	1:A:244:GLY:HA2	1.98	0.78
1:B:32:ARG:HD2	3:B:545:HOH:O	1.84	0.76
1:B:344:GLU:CD	1:B:369:GLY:HA3	2.06	0.75
1:A:354:ILE:CD1	1:A:364:VAL:HG21	2.16	0.74
1:A:191:GLU:HG3	1:A:262:ILE:CD1	2.17	0.74
1:A:121:LEU:CD2	1:B:121:LEU:HD23	2.18	0.73
1:A:81:ASP:OD1	1:A:83:ASN:HB2	1.90	0.72
1:B:226:ASP:HB3	3:B:568:HOH:O	1.91	0.70
1:A:350:ALA:O	1:A:354:ILE:HG12	1.92	0.70
1:B:207:PHE:CB	1:B:351:LEU:HD21	2.17	0.69
1:B:207:PHE:CD2	1:B:351:LEU:HD23	2.28	0.69
1:B:165:LEU:HD22	3:B:716:HOH:O	1.92	0.68
1:B:366:PRO:HA	3:B:742:HOH:O	1.92	0.68
1:A:175:LYS:N	1:A:175:LYS:HD2	2.06	0.68
1:A:252:GLU:HA	1:A:255:SER:OG	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:394:LYS:HD3	1:B:394:LYS:N	2.08	0.67
1:B:217:LYS:O	1:B:221:GLU:HG3	1.94	0.67
1:A:78:LEU:HD23	1:A:284:LEU:HD21	1.77	0.65
1:A:26:ASN:OD1	1:A:27:SER:N	2.30	0.65
1:A:361:LEU:HD12	3:A:590:HOH:O	1.97	0.64
1:A:210:ASN:OD1	1:A:214:LYS:HE2	1.97	0.64
1:B:78:LEU:HD21	1:B:284:LEU:HD21	1.81	0.63
1:A:206:GLU:OE1	1:A:206:GLU:HA	1.98	0.63
1:A:311:PRO:HB2	1:A:314:ILE:HD11	1.80	0.62
1:B:210:ASN:ND2	3:B:631:HOH:O	2.31	0.62
1:A:155:MET:O	1:A:159:TYR:HB2	2.00	0.62
1:B:209:ALA:O	1:B:213:GLN:N	2.30	0.62
1:B:203:ASP:O	1:B:206:GLU:HB3	1.99	0.62
1:A:78:LEU:CD2	1:A:284:LEU:HD21	2.29	0.62
1:A:27:SER:O	1:A:32:ARG:NH2	2.30	0.62
1:A:46:ILE:HG12	3:A:628:HOH:O	1.98	0.62
1:A:155:MET:HG3	1:A:277:SER:O	2.00	0.61
1:A:192:ASN:O	1:A:196:ASN:HB2	2.00	0.61
1:B:222:GLY:O	1:B:225:GLU:HB2	2.00	0.61
1:A:27:SER:HA	1:A:30:GLU:HG3	1.82	0.61
1:B:74:PHE:HB2	1:B:79:ARG:HG2	1.83	0.60
1:A:377:LEU:HD13	3:A:685:HOH:O	2.00	0.60
1:B:390:ARG:HD2	1:B:390:ARG:H	1.65	0.60
1:B:232:LYS:NZ	1:B:236:GLY:O	2.30	0.59
1:B:25:LYS:O	1:B:29:LEU:HD23	2.02	0.59
1:A:210:ASN:O	1:A:214:LYS:HG3	2.03	0.59
1:A:300:GLY:HA2	1:A:417:GLU:HG3	1.85	0.59
1:B:390:ARG:HD2	1:B:390:ARG:N	2.17	0.59
1:B:314:ILE:O	1:B:317:VAL:HG22	2.03	0.59
1:B:63:LEU:HD11	1:B:150:LEU:HD13	1.86	0.58
1:A:188:ILE:HD13	1:A:262:ILE:CD1	2.29	0.57
1:B:252:GLU:N	3:B:715:HOH:O	2.37	0.57
1:B:27:SER:O	1:B:30:GLU:HB3	2.05	0.57
1:A:188:ILE:CD1	1:A:262:ILE:HD11	2.27	0.56
1:A:121:LEU:HD23	1:B:121:LEU:HD23	1.88	0.56
1:B:176:ASN:O	1:B:180:LYS:HG2	2.06	0.56
1:A:218:ALA:HA	1:A:223:LEU:HD12	1.87	0.56
1:A:324:PRO:HD3	1:A:357:LEU:HD21	1.88	0.56
1:A:25:LYS:O	1:A:28:LEU:HB2	2.06	0.56
1:B:251:ALA:O	1:B:254:LEU:HB2	2.06	0.55
1:B:163:ASN:OD1	1:B:164:PRO:HD2	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:79:ARG:HH11	1:A:79:ARG:HG2	1.70	0.55
1:A:232:LYS:HG3	1:A:237:SER:O	2.07	0.55
1:B:27:SER:O	1:B:32:ARG:NH1	2.37	0.55
1:B:73:ARG:NH1	1:B:73:ARG:HG3	2.18	0.55
1:B:215:ALA:O	1:B:219:LYS:HB2	2.06	0.55
1:B:223:LEU:HB3	1:B:367:ARG:NH1	2.23	0.55
1:A:252:GLU:HB3	1:A:255:SER:OG	2.06	0.54
1:B:207:PHE:HB2	1:B:351:LEU:CD2	2.24	0.54
1:B:344:GLU:OE2	1:B:369:GLY:HA3	2.07	0.54
1:B:323:ILE:HB	1:B:324:PRO:HD3	1.89	0.54
1:A:73:ARG:O	1:A:73:ARG:HG3	2.08	0.54
1:B:335:GLN:HB2	3:B:612:HOH:O	2.06	0.53
1:A:73:ARG:NH1	1:A:226:ASP:O	2.35	0.53
1:A:191:GLU:OE2	1:A:268:THR:HB	2.08	0.53
1:A:75:PRO:C	1:A:77:PRO:HD2	2.30	0.53
1:B:37:VAL:HB	1:B:298:VAL:HG13	1.91	0.53
1:A:233:LEU:N	1:A:233:LEU:HD22	2.24	0.52
1:A:289:SER:O	1:A:293:GLN:NE2	2.42	0.52
1:A:308:VAL:CG1	1:A:325:LYS:HD2	2.39	0.52
1:B:248:ASN:HA	3:B:671:HOH:O	2.09	0.52
1:A:339:ILE:CG1	1:A:392:LEU:HD23	2.40	0.52
1:A:344:GLU:CD	1:A:369:GLY:HA3	2.31	0.52
1:A:48:LYS:HZ3	1:A:244:GLY:HA2	1.74	0.51
1:A:48:LYS:HZ2	1:A:244:GLY:HA2	1.74	0.51
1:A:339:ILE:HG12	1:A:392:LEU:HD23	1.91	0.51
1:A:79:ARG:HH11	1:A:79:ARG:CG	2.23	0.50
1:A:308:VAL:HG13	1:A:325:LYS:HD3	1.93	0.50
1:A:32:ARG:NH1	3:A:597:HOH:O	2.45	0.50
1:B:73:ARG:HH11	1:B:73:ARG:CG	2.20	0.50
1:B:83:ASN:ND2	1:B:115:SER:OG	2.41	0.50
1:A:313:GLU:HG2	1:A:314:ILE:HG13	1.93	0.50
1:B:383:ARG:HD2	1:B:383:ARG:C	2.32	0.50
1:B:213:GLN:NE2	1:B:249:VAL:O	2.41	0.50
1:B:233:LEU:HB3	1:B:234:PRO:CD	2.41	0.50
1:A:201:ARG:HB2	1:A:267:THR:CG2	2.34	0.49
1:A:261:PHE:HB2	1:A:262:ILE:HD12	1.93	0.49
1:A:79:ARG:HB3	1:A:79:ARG:CZ	2.42	0.49
1:A:325:LYS:HE3	3:A:598:HOH:O	2.12	0.49
1:B:81:ASP:OD1	1:B:83:ASN:HB2	2.12	0.49
1:A:178:GLU:OE1	1:A:178:GLU:HA	2.12	0.49
1:A:250:THR:N	1:A:253:SER:OG	2.44	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:303:ILE:HD11	1:A:415:ILE:HD12	1.94	0.49
1:A:308:VAL:HG13	1:A:325:LYS:CD	2.42	0.49
1:B:390:ARG:H	1:B:390:ARG:CD	2.23	0.49
1:B:53:ALA:HB2	1:B:243:GLU:OE2	2.13	0.49
1:B:213:GLN:HG3	3:B:729:HOH:O	2.13	0.48
1:A:250:THR:HG1	1:A:253:SER:CB	2.22	0.48
1:A:114:TYR:O	1:B:309:GLY:HA3	2.12	0.48
1:B:264:ASP:O	1:B:265:ARG:HD3	2.12	0.48
1:A:191:GLU:HG3	1:A:262:ILE:HD13	1.94	0.48
1:A:261:PHE:C	1:A:262:ILE:HG13	2.33	0.48
1:A:76:GLU:OE2	1:A:79:ARG:NE	2.43	0.48
1:A:404:ILE:HB	1:A:408:MET:HB2	1.95	0.48
1:A:309:GLY:HA3	1:B:114:TYR:O	2.14	0.48
1:B:251:ALA:N	3:B:715:HOH:O	2.47	0.48
1:A:200:SER:O	1:A:204:GLN:HG3	2.14	0.48
1:A:344:GLU:OE2	1:A:369:GLY:HA3	2.14	0.48
1:B:338:ASP:O	1:B:339:ILE:HG13	2.13	0.48
1:A:176:ASN:HB3	1:A:179:ALA:HB3	1.95	0.48
1:B:89:ALA:HB1	1:B:121:LEU:HD11	1.95	0.47
1:B:249:VAL:HG12	1:B:250:THR:N	2.29	0.47
1:B:178:GLU:HA	1:B:178:GLU:OE1	2.15	0.47
1:B:251:ALA:O	1:B:254:LEU:N	2.44	0.47
1:A:237:SER:O	1:A:238:ILE:HG13	2.14	0.47
1:B:124:GLN:HB2	1:B:404:ILE:HG23	1.97	0.47
1:A:311:PRO:CB	1:A:314:ILE:HD11	2.45	0.47
1:A:180:LYS:NZ	3:A:687:HOH:O	2.48	0.47
1:B:76:GLU:HG3	3:B:708:HOH:O	2.15	0.46
1:B:400:VAL:O	1:B:411:ALA:HA	2.15	0.46
1:B:235:ASP:OD1	1:B:237:SER:N	2.31	0.46
1:A:320:ALA:O	1:A:324:PRO:HG2	2.16	0.46
1:A:354:ILE:HD11	1:A:364:VAL:HG11	1.98	0.46
1:A:42:ASN:HD21	1:A:73:ARG:NH1	2.13	0.46
1:A:261:PHE:O	1:A:262:ILE:HG13	2.16	0.46
1:B:361:LEU:H	1:B:361:LEU:CD2	2.29	0.46
1:A:258:ARG:O	1:A:269:THR:HB	2.15	0.45
1:A:252:GLU:O	1:A:255:SER:OG	2.30	0.45
1:B:232:LYS:HE3	1:B:236:GLY:HA2	1.98	0.45
1:A:79:ARG:HB3	1:A:79:ARG:NH1	2.30	0.45
1:A:216:TYR:CE2	1:A:247:PRO:HB3	2.50	0.45
1:A:76:GLU:N	1:A:77:PRO:CD	2.80	0.45
1:A:233:LEU:N	1:A:233:LEU:CD2	2.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:GLU:HA	1:B:76:GLU:OE1	2.16	0.45
1:B:249:VAL:CG1	1:B:250:THR:N	2.80	0.45
1:B:389:LEU:O	1:B:391:GLU:N	2.50	0.45
1:B:73:ARG:NH1	1:B:73:ARG:CG	2.80	0.45
1:B:178:GLU:HA	1:B:181:LYS:HD2	1.99	0.45
1:B:75:PRO:HG2	1:B:296:LEU:HD22	1.99	0.45
1:A:171:GLU:OE1	1:A:171:GLU:HA	2.17	0.44
1:B:361:LEU:N	1:B:361:LEU:CD2	2.79	0.44
1:B:192:ASN:O	1:B:196:ASN:HB2	2.17	0.44
1:A:406:THR:HB	1:B:104:ARG:HG2	1.98	0.44
1:B:365:ASN:N	1:B:366:PRO:CD	2.80	0.44
1:A:183:LEU:O	1:A:185:PRO:HD3	2.18	0.44
1:B:344:GLU:OE1	1:B:369:GLY:HA3	2.17	0.44
1:A:311:PRO:CG	1:A:314:ILE:HD11	2.48	0.44
1:B:339:ILE:HD12	1:B:392:LEU:HD23	1.98	0.44
1:B:361:LEU:HD22	1:B:361:LEU:N	2.33	0.44
1:B:210:ASN:OD1	1:B:214:LYS:HE2	2.18	0.44
1:A:340:PHE:CE2	1:A:359:ILE:HG12	2.52	0.44
1:B:394:LYS:O	1:B:395:ASP:HB2	2.18	0.44
1:A:178:GLU:HA	1:A:181:LYS:HD2	2.00	0.44
1:B:343:ASN:HD21	1:B:370:ALA:HB2	1.83	0.43
1:A:234:PRO:C	1:A:236:GLY:H	2.22	0.43
1:A:341:GLU:OE2	1:A:391:GLU:OE1	2.37	0.43
1:B:255:SER:HB3	3:B:650:HOH:O	2.18	0.43
1:B:390:ARG:N	1:B:390:ARG:CD	2.81	0.43
1:A:198:LYS:HD2	1:A:198:LYS:HA	1.71	0.43
1:A:250:THR:O	1:A:253:SER:OG	2.30	0.43
1:A:323:ILE:HB	1:A:324:PRO:HD3	2.00	0.43
1:A:67:LEU:O	1:A:71:ILE:HG13	2.19	0.43
1:A:201:ARG:HE	1:A:205:ASP:CG	2.22	0.43
1:A:361:LEU:HA	1:A:361:LEU:HD13	1.35	0.43
1:A:168:ILE:HG13	1:A:183:LEU:HD21	2.01	0.43
1:B:181:LYS:NZ	3:B:733:HOH:O	2.51	0.43
1:B:257:ILE:HG22	1:B:258:ARG:N	2.33	0.43
1:A:171:GLU:O	1:A:174:GLN:HB2	2.19	0.43
1:B:200:SER:O	1:B:204:GLN:HG3	2.19	0.43
1:A:324:PRO:HG3	1:A:357:LEU:CD2	2.49	0.43
1:B:82:LEU:HA	1:B:82:LEU:HD23	1.78	0.43
1:A:79:ARG:NH1	1:A:79:ARG:CG	2.80	0.43
1:A:299:LEU:HD23	1:A:299:LEU:N	2.34	0.43
1:B:388:ILE:CG2	1:B:389:LEU:N	2.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:221:GLU:HB2	1:B:223:LEU:HG	2.01	0.42
1:A:186:MET:HG3	3:A:645:HOH:O	2.18	0.42
1:A:174:GLN:C	1:A:175:LYS:HD2	2.40	0.42
1:B:233:LEU:HB3	1:B:234:PRO:HD3	2.01	0.42
1:B:257:ILE:CG2	1:B:258:ARG:N	2.82	0.42
1:A:307:THR:HG22	1:A:410:ALA:HB2	2.01	0.42
1:B:359:ILE:CG2	1:B:360:ASP:N	2.81	0.42
1:A:227:GLU:HB3	1:A:390:ARG:HD3	2.01	0.42
1:A:48:LYS:HE2	1:A:48:LYS:HB2	1.50	0.42
1:B:226:ASP:N	3:B:618:HOH:O	2.35	0.42
1:B:359:ILE:HG22	1:B:360:ASP:N	2.34	0.42
1:B:319:PRO:HD2	1:B:353:CYS:SG	2.60	0.42
1:B:394:LYS:HD2	1:B:417:GLU:O	2.20	0.42
1:B:200:SER:O	1:B:203:ASP:HB2	2.19	0.42
1:A:181:LYS:C	1:A:183:LEU:H	2.23	0.41
1:B:207:PHE:CD2	1:B:351:LEU:CD2	3.00	0.41
1:B:223:LEU:O	1:B:367:ARG:NH1	2.52	0.41
1:A:32:ARG:O	1:A:288:ARG:HD2	2.20	0.41
1:B:202:LYS:H	1:B:202:LYS:HG2	1.26	0.41
1:B:124:GLN:HG3	1:B:405:GLY:O	2.21	0.41
1:B:263:LYS:H	1:B:263:LYS:HG2	1.70	0.41
1:B:83:ASN:ND2	3:B:613:HOH:O	2.50	0.41
1:B:235:ASP:OD1	1:B:237:SER:OG	2.37	0.41
1:A:163:ASN:HA	1:A:164:PRO:HD2	1.48	0.41
1:B:287:ARG:NH1	3:B:555:HOH:O	2.52	0.41
1:B:215:ALA:O	1:B:219:LYS:N	2.49	0.40
1:B:301:ARG:HD3	1:B:417:GLU:OE2	2.21	0.40
1:A:324:PRO:HG3	1:A:357:LEU:HD21	2.03	0.40
1:B:124:GLN:HB2	1:B:404:ILE:CG2	2.51	0.40
1:A:79:ARG:CB	1:A:79:ARG:CZ	3.00	0.40
1:B:337:ILE:HB	1:B:340:PHE:CZ	2.56	0.40
1:A:67:LEU:HA	1:A:67:LEU:HD23	1.86	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	386/393 (98%)	366 (95%)	18 (5%)	2 (0%)	34 17
1	B	391/393 (100%)	364 (93%)	23 (6%)	4 (1%)	19 5
All	All	777/786 (99%)	730 (94%)	41 (5%)	6 (1%)	24 8

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	377	LEU
1	B	377	LEU
1	B	390	ARG
1	B	223	LEU
1	B	251	ALA
1	A	235	ASP

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	311/314 (99%)	278 (89%)	33 (11%)	8 2
1	B	314/314 (100%)	286 (91%)	28 (9%)	12 3
All	All	625/628 (100%)	564 (90%)	61 (10%)	10 2

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	LYS
1	A	28	LEU
1	A	29	LEU
1	A	48	LYS
1	A	70	PHE
1	A	73	ARG

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Mol	Chain	Res	Type
1	A	76	GLU
1	A	160	LYS
1	A	170	SER
1	A	172	GLU
1	A	175	LYS
1	A	177	ARG
1	A	196	ASN
1	A	206	GLU
1	A	217	LYS
1	A	221	GLU
1	A	243	GLU
1	A	246	ARG
1	A	252	GLU
1	A	256	SER
1	A	257	ILE
1	A	258	ARG
1	A	293	GLN
1	A	299	LEU
1	A	308	VAL
1	A	314	ILE
1	A	315	MET
1	A	333	GLN
1	A	335	GLN
1	A	349	GLN
1	A	361	LEU
1	A	362	ASN
1	A	402	MET
1	B	26	ASN
1	B	28	LEU
1	B	79	ARG
1	B	110	SER
1	B	165	LEU
1	B	169	SER
1	B	171	GLU
1	B	180	LYS
1	B	202	LYS
1	B	210	ASN
1	B	219	LYS
1	B	232	LYS
1	B	233	LEU
1	B	246	ARG
1	B	248	ASN

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Mol	Chain	Res	Type
1	B	250	THR
1	B	252	GLU
1	B	253	SER
1	B	256	SER
1	B	277	SER
1	B	285	LEU
1	B	315	MET
1	B	333	GLN
1	B	335	GLN
1	B	349	GLN
1	B	389	LEU
1	B	394	LYS
1	B	416	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	124	GLN
1	A	157	ASN
1	A	292	ASN
1	A	293	GLN
1	B	68	ASN
1	B	83	ASN
1	B	240	GLN
1	B	292	ASN
1	B	343	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 carbohydrates in this entry.

5.6 Ligand geometry (i)

2 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	MRD	A	501	-	6,7,7	0.55	0	7,10,10	2.41	2 (28%)
2	MRD	B	502	-	6,7,7	0.46	0	7,10,10	1.30	1 (14%)

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	MRD	A	501	-	-	0/5/5/5	0/0/0/0
2	MRD	B	502	-	-	0/5/5/5	0/0/0/0

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	501	MRD	CM-C2-C1	-5.07	99.19	110.24
2	B	502	MRD	CM-C2-C1	-2.84	104.05	110.24
2	A	501	MRD	O2-C2-C1	3.45	120.74	108.09

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

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	390/393 (99%)	-0.44	4 (1%)	84	82	7, 22, 56, 92	0
1	B	393/393 (100%)	-0.44	5 (1%)	79	76	6, 21, 51, 84	0
All	All	783/786 (99%)	-0.44	9 (1%)	82	80	6, 22, 53, 92	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	26	ASN	4.4
1	A	26	ASN	3.7
1	A	262	ILE	3.5
1	B	220	ASN	3.3
1	A	293	GLN	2.9
1	B	247	PRO	2.9
1	B	216	TYR	2.4
1	B	254	LEU	2.2
1	A	252	GLU	2.1

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

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	MRD	A	501	8/8	0.90	0.13	5.48	28,34,43,44	0
2	MRD	B	502	8/8	0.96	0.08	1.95	21,24,30,34	0

6.5 Other polymers (i)

There are no such residues in this entry.