



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:26 AM GMT

PDB ID : 2QQ6
Title : Crystal structure of mandelate racemase/muconate lactonizing enzyme-like protein from *Rubrobacter xylanophilus* DSM 9941
Authors : Eswaramoorthy, S.; Madegowda, M.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2007-07-26
Resolution : 2.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

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

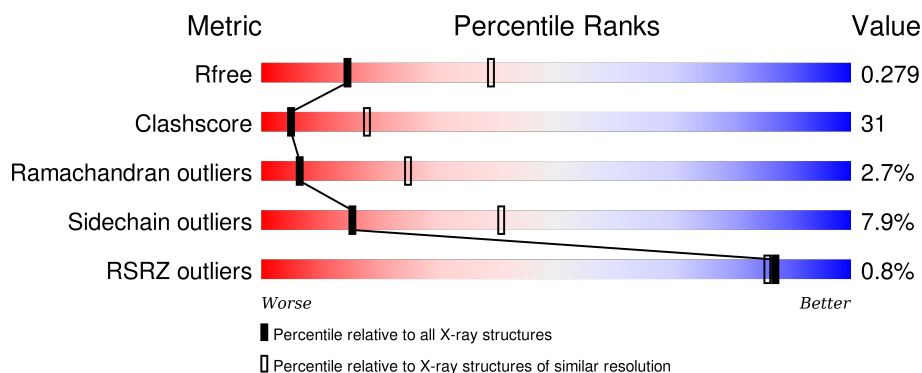
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	410	<div> <div></div> <div>50% 41% 6% .</div> </div>
1	B	410	<div> <div></div> <div>48% 43% . . .</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6051 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 Mandelate racemase/muconate lactonizing enzyme-like protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	394	Total	C	N	O	S	Se	0	0	0
			3000	1902	522	563	5	8			
1	B	396	Total	C	N	O	S	Se	0	0	0
			3021	1913	530	565	5	8			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MSE	-	EXPRESSION TAG	UNP Q1AYK7
A	0	SER	-	EXPRESSION TAG	UNP Q1AYK7
A	1	LEU	-	EXPRESSION TAG	UNP Q1AYK7
A	401	GLU	-	EXPRESSION TAG	UNP Q1AYK7
A	402	GLY	-	EXPRESSION TAG	UNP Q1AYK7
A	403	HIS	-	EXPRESSION TAG	UNP Q1AYK7
A	404	HIS	-	EXPRESSION TAG	UNP Q1AYK7
A	405	HIS	-	EXPRESSION TAG	UNP Q1AYK7
A	406	HIS	-	EXPRESSION TAG	UNP Q1AYK7
A	407	HIS	-	EXPRESSION TAG	UNP Q1AYK7
A	408	HIS	-	EXPRESSION TAG	UNP Q1AYK7
B	-1	MSE	-	EXPRESSION TAG	UNP Q1AYK7
B	0	SER	-	EXPRESSION TAG	UNP Q1AYK7
B	1	LEU	-	EXPRESSION TAG	UNP Q1AYK7
B	401	GLU	-	EXPRESSION TAG	UNP Q1AYK7
B	402	GLY	-	EXPRESSION TAG	UNP Q1AYK7
B	403	HIS	-	EXPRESSION TAG	UNP Q1AYK7
B	404	HIS	-	EXPRESSION TAG	UNP Q1AYK7
B	405	HIS	-	EXPRESSION TAG	UNP Q1AYK7
B	406	HIS	-	EXPRESSION TAG	UNP Q1AYK7
B	407	HIS	-	EXPRESSION TAG	UNP Q1AYK7
B	408	HIS	-	EXPRESSION TAG	UNP Q1AYK7

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total 1	Mg 1	0	0
2	A	1	Total 1	Mg 1	0	0

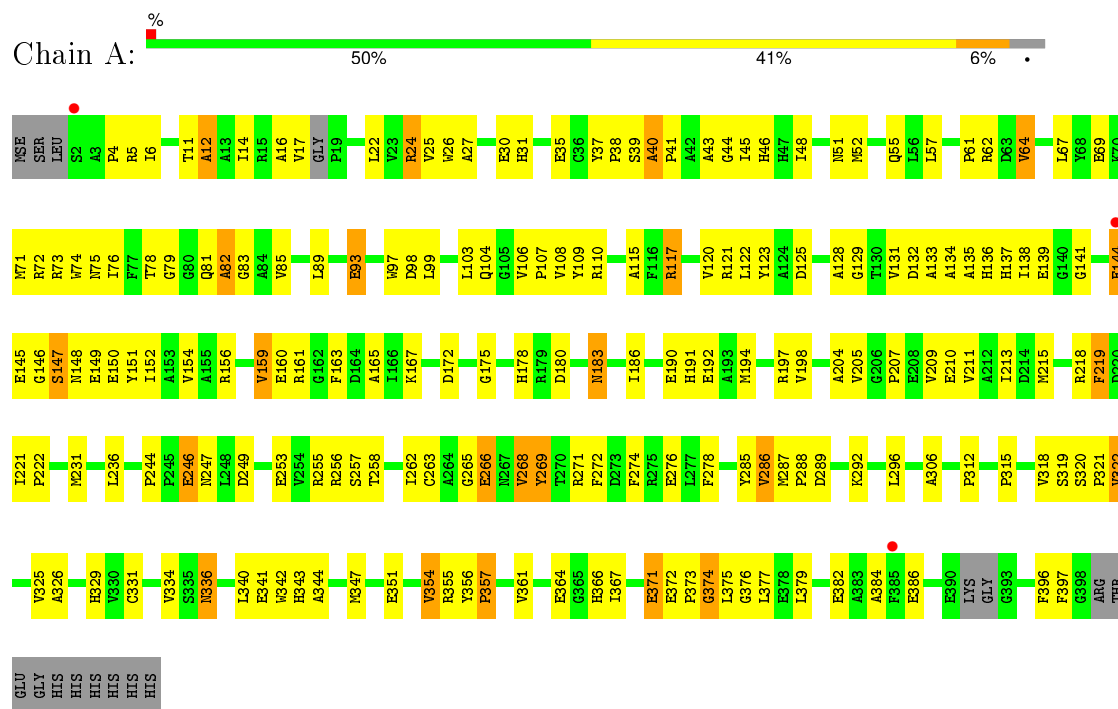
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	14	Total 14	O 14	0	0
3	B	14	Total 14	O 14	0	0

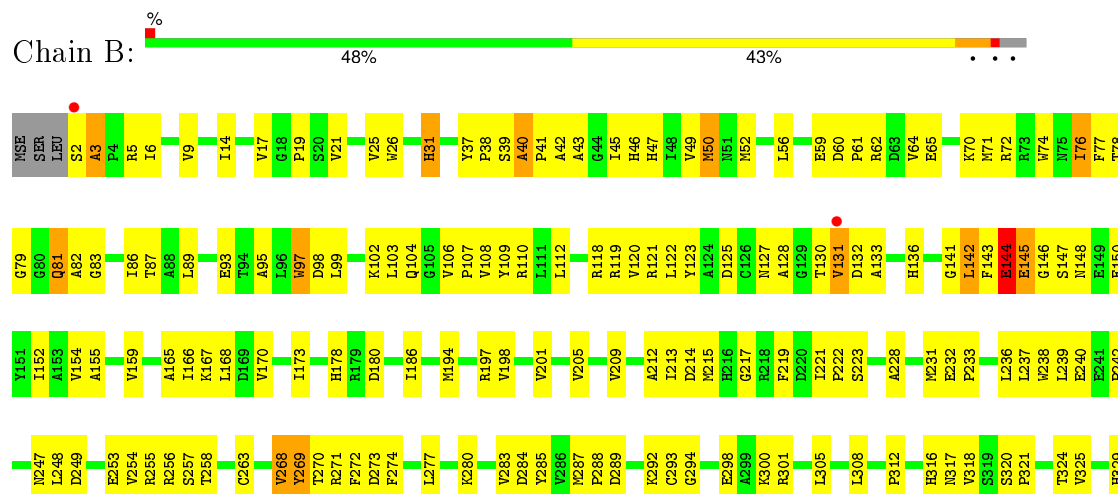
3 Residue-property plots

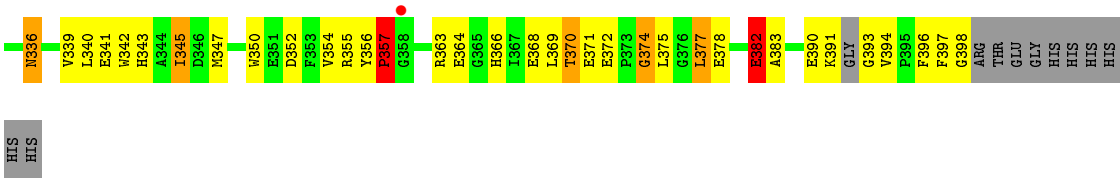
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Mandelate racemase/muconate lactonizing enzyme-like protein



- Molecule 1: Mandelate racemase/muconate lactonizing enzyme-like protein





4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	184.37Å 184.37Å 118.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.90 49.73 – 2.90	Depositor EDS
% Data completeness (in resolution range)	94.4 (50.00-2.90) 94.4 (49.73-2.90)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.35 (at 2.91Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.222 , 0.282 0.221 , 0.279	Depositor DCC
R_{free} test set	829 reflections (3.85%)	DCC
Wilson B-factor (Å ²)	45.0	Xtriage
Anisotropy	0.749	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 61.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 22823 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6051	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.10% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.43	0/3067	0.69	1/4165 (0.0%)
1	B	0.41	0/3089	0.69	0/4195
All	All	0.42	0/6156	0.69	1/8360 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	269	TYR	N-CA-C	5.24	125.15	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	3000	0	2858	183	0
1	B	3021	0	2885	196	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	14	0	0	7	0
3	B	14	0	0	1	0
All	All	6051	0	5743	363	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 31.

All (363) 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:17:VAL:H	1:A:347:MSE:HE1	1.11	1.12
1:B:17:VAL:HG21	1:B:318:VAL:HG22	1.33	1.05
1:A:82:ALA:CB	1:B:83:GLY:HA3	1.88	1.04
1:A:82:ALA:HB2	1:B:83:GLY:HA3	1.42	1.01
1:B:40:ALA:HB1	1:B:41:PRO:HD2	1.40	0.99
1:B:336:ASN:HD22	1:B:336:ASN:H	0.99	0.94
1:A:336:ASN:HD22	1:A:336:ASN:H	1.17	0.92
1:A:40:ALA:HB1	1:A:41:PRO:HD2	1.51	0.92
1:B:347:MSE:HE2	1:B:350:TRP:HA	1.52	0.89
1:B:391:LYS:C	1:B:393:GLY:HA2	1.94	0.88
1:A:321:PRO:O	1:A:325:VAL:HG23	1.74	0.87
1:A:190:GLU:HG2	1:A:194:MSE:HE2	1.55	0.87
1:B:40:ALA:HB1	1:B:41:PRO:CD	2.04	0.87
1:B:215:MSE:HE2	1:B:239:LEU:HG	1.57	0.86
1:B:215:MSE:HE2	1:B:239:LEU:CG	2.05	0.86
1:B:336:ASN:HD22	1:B:336:ASN:N	1.75	0.84
1:A:40:ALA:HB1	1:A:41:PRO:CD	2.08	0.83
1:A:52:MSE:HE1	1:A:74:TRP:CD1	2.15	0.82
1:B:132:ASP:HB3	1:B:136:HIS:H	1.46	0.81
1:A:191:HIS:HA	1:A:194:MSE:HE3	1.63	0.81
1:B:50:MSE:HA	1:B:50:MSE:HE3	1.63	0.80
1:B:248:LEU:HD13	1:B:277:LEU:HD12	1.60	0.80
1:B:215:MSE:CE	1:B:242:PRO:HB3	2.13	0.78
1:A:17:VAL:HG11	1:A:318:VAL:HG22	1.63	0.78
1:A:17:VAL:N	1:A:347:MSE:HE1	1.95	0.77
1:B:336:ASN:ND2	1:B:336:ASN:H	1.78	0.77
1:A:82:ALA:HB1	1:B:83:GLY:HA3	1.66	0.76
1:A:82:ALA:HB2	1:B:83:GLY:CA	2.16	0.75
1:A:221:ILE:HB	1:A:222:PRO:HD3	1.68	0.75
1:A:319:SER:HB3	3:A:415:HOH:O	1.88	0.73
1:B:321:PRO:O	1:B:325:VAL:HG23	1.89	0.73
1:B:186:ILE:HG13	1:B:223:SER:HB3	1.71	0.73
1:A:129:GLY:HA3	1:A:139:GLU:O	1.88	0.73
1:B:123:TYR:CZ	1:B:341:GLU:HB2	2.25	0.72
1:A:40:ALA:CB	1:A:41:PRO:HD2	2.19	0.72
1:A:4:PRO:HD2	1:A:62:ARG:NH2	2.05	0.71
1:B:42:ALA:HA	1:B:45:ILE:HD12	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:VAL:HG11	1:A:209:VAL:HG11	1.72	0.71
1:B:214:ASP:HA	1:B:240:GLU:HB3	1.73	0.71
1:A:6:ILE:CD1	1:A:61:PRO:HG3	2.20	0.71
1:B:40:ALA:CB	1:B:41:PRO:HD2	2.20	0.70
1:A:40:ALA:H	1:B:78:THR:HB	1.57	0.70
1:B:317:ASN:HD21	1:B:324:THR:HA	1.56	0.70
1:A:183:ASN:HD22	1:A:183:ASN:C	1.93	0.70
1:B:248:LEU:HD13	1:B:277:LEU:CD1	2.20	0.70
1:B:274:PHE:CZ	1:B:288:PRO:HB3	2.26	0.70
1:A:17:VAL:H	1:A:347:MSE:CE	1.97	0.69
1:A:172:ASP:OD2	1:A:175:GLY:HA3	1.93	0.69
1:B:329:HIS:NE2	1:B:375:LEU:HA	2.07	0.69
1:B:39:SER:O	1:B:45:ILE:HD11	1.93	0.69
1:A:6:ILE:HD12	1:A:61:PRO:HG3	1.75	0.68
1:B:391:LYS:O	1:B:393:GLY:HA2	1.94	0.68
1:B:320:SER:HB2	1:B:321:PRO:CD	2.23	0.68
1:A:78:THR:HB	1:B:40:ALA:H	1.58	0.68
1:B:119:ARG:HB3	1:B:368:GLU:OE2	1.94	0.68
1:A:356:TYR:CD1	1:A:357:PRO:HD2	2.30	0.67
1:B:382:GLU:HA	1:B:382:GLU:OE1	1.95	0.67
1:B:72:ARG:HH12	1:B:271:ARG:NH2	1.91	0.67
1:A:148:ASN:ND2	1:A:197:ARG:HA	2.11	0.66
1:B:205:VAL:HG12	1:B:209:VAL:HG21	1.76	0.66
1:B:62:ARG:HD2	1:B:104:GLN:HE22	1.61	0.66
1:B:215:MSE:HE3	1:B:242:PRO:HB3	1.78	0.66
1:A:148:ASN:HD21	1:A:197:ARG:HA	1.60	0.66
1:A:334:VAL:HG13	1:A:336:ASN:ND2	2.11	0.66
1:B:347:MSE:HE2	1:B:350:TRP:CA	2.26	0.66
1:B:132:ASP:HB3	1:B:136:HIS:N	2.10	0.65
1:A:40:ALA:CB	1:A:41:PRO:CD	2.75	0.65
1:A:35:GLU:OE2	1:A:319:SER:HB2	1.97	0.65
1:B:144:GLU:HG2	1:B:145:GLU:H	1.63	0.64
1:B:173:ILE:HA	1:B:178:HIS:CD2	2.34	0.62
1:A:379:LEU:HD11	1:A:384:ALA:HB2	1.81	0.62
1:A:246:GLU:HG3	1:B:272:PHE:CE1	2.35	0.62
1:B:50:MSE:HA	1:B:50:MSE:CE	2.30	0.61
1:B:125:ASP:OD2	1:B:167:LYS:HD3	2.00	0.61
1:B:255:ARG:HH21	1:B:284:ASP:CG	2.04	0.61
1:A:69:GLU:O	1:A:73:ARG:HG3	2.00	0.61
1:A:17:VAL:HG11	1:A:318:VAL:CG2	2.30	0.60
1:B:49:VAL:O	1:B:52:MSE:HB2	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:278:PHE:CZ	1:A:306:ALA:HB2	2.36	0.60
1:B:123:TYR:HE1	1:B:167:LYS:HD2	1.66	0.60
1:A:351:GLU:OE1	1:A:361:VAL:HG23	2.02	0.60
1:A:117:ARG:HH11	1:A:117:ARG:HB3	1.67	0.60
1:B:64:VAL:HG11	1:B:97:TRP:CE2	2.37	0.60
1:A:265:GLY:HA2	1:A:268:VAL:CG1	2.31	0.60
1:A:315:PRO:HG2	1:A:340:LEU:HD13	1.83	0.59
1:A:354:VAL:HG11	1:A:377:LEU:HD22	1.83	0.59
1:B:215:MSE:CE	1:B:239:LEU:HG	2.28	0.59
1:B:72:ARG:HH12	1:B:271:ARG:HH22	1.48	0.59
1:A:274:PHE:CZ	1:A:288:PRO:HB3	2.38	0.59
1:A:146:GLY:HA2	1:A:149:GLU:OE2	2.02	0.59
1:A:186:ILE:CG2	1:A:194:MSE:HE1	2.32	0.58
1:B:215:MSE:HE2	1:B:239:LEU:CD2	2.31	0.58
1:B:320:SER:HA	1:B:350:TRP:HE1	1.66	0.58
1:B:268:VAL:HG13	1:B:274:PHE:CZ	2.37	0.58
1:B:354:VAL:HG11	1:B:377:LEU:HB2	1.85	0.58
1:A:52:MSE:HE1	1:A:74:TRP:HD1	1.63	0.58
1:A:266:GLU:HB3	3:A:411:HOH:O	2.03	0.58
1:A:289:ASP:HB3	1:A:292:LYS:HB3	1.86	0.58
1:B:374:GLY:O	1:B:375:LEU:HB2	2.04	0.58
1:A:99:LEU:HG	1:A:103:LEU:HD12	1.86	0.58
1:A:231:MSE:SE	1:A:236:LEU:HD11	2.54	0.58
1:A:334:VAL:HG13	1:A:336:ASN:HD22	1.68	0.58
1:A:122:LEU:HB2	1:A:163:PHE:HE2	1.68	0.57
1:B:144:GLU:HB3	1:B:147:SER:OG	2.03	0.57
1:B:155:ALA:O	1:B:159:VAL:HG23	2.04	0.57
1:A:244:PRO:HB3	1:A:246:GLU:OE2	2.04	0.57
1:A:356:TYR:CG	1:A:357:PRO:HD2	2.40	0.57
1:B:232:GLU:HB2	1:B:233:PRO:HD3	1.86	0.57
1:B:356:TYR:CD1	1:B:357:PRO:HD2	2.40	0.57
1:A:22:LEU:N	1:A:22:LEU:HD22	2.20	0.57
1:A:336:ASN:HD22	1:A:336:ASN:N	1.88	0.56
1:A:109:TYR:CD1	1:A:115:ALA:HB2	2.39	0.56
1:A:336:ASN:ND2	1:A:336:ASN:H	1.95	0.56
1:B:370:THR:OG1	1:B:371:GLU:N	2.39	0.56
1:B:112:LEU:O	1:B:300:LYS:HE3	2.05	0.56
1:A:183:ASN:ND2	1:A:183:ASN:C	2.58	0.56
1:B:215:MSE:HE1	1:B:242:PRO:HB3	1.84	0.56
1:A:14:ILE:HG12	1:A:22:LEU:HD21	1.88	0.56
1:A:186:ILE:HG23	1:A:194:MSE:HE1	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:TYR:CE1	1:A:341:GLU:HG3	2.41	0.56
1:B:17:VAL:HG21	1:B:318:VAL:CG2	2.22	0.56
1:A:108:VAL:HG23	1:A:375:LEU:N	2.21	0.55
1:A:107:PRO:HB3	1:A:372:GLU:O	2.06	0.55
1:A:107:PRO:HG2	1:A:110:ARG:CG	2.37	0.55
1:A:11:THR:HG1	1:A:46:HIS:CE1	2.24	0.55
1:B:231:MSE:SE	1:B:236:LEU:HD11	2.56	0.55
1:B:45:ILE:O	1:B:49:VAL:HG23	2.06	0.54
1:A:253:GLU:OE1	1:A:256:ARG:NH1	2.40	0.54
1:B:128:ALA:HA	1:B:343:HIS:CE1	2.42	0.54
1:A:342:TRP:CE2	1:A:344:ALA:HB3	2.42	0.54
1:A:73:ARG:O	1:A:76:ILE:HG22	2.08	0.54
1:B:356:TYR:CG	1:B:357:PRO:HD2	2.43	0.54
1:B:396:PHE:O	1:B:398:GLY:N	2.41	0.54
1:B:82:ALA:HA	1:B:86:ILE:HG13	1.89	0.54
1:B:215:MSE:HE2	1:B:239:LEU:HD21	1.90	0.54
1:B:283:VAL:HB	1:B:285:TYR:O	2.07	0.54
1:A:128:ALA:HB3	1:A:137:HIS:CE1	2.43	0.54
1:A:52:MSE:CE	1:A:55:GLN:HG3	2.38	0.53
1:B:221:ILE:HB	1:B:222:PRO:HD3	1.89	0.53
1:B:287:MSE:HE3	1:B:316:HIS:HB2	1.89	0.53
1:A:150:GLU:O	1:A:154:VAL:HG23	2.08	0.53
1:B:377:LEU:HD12	1:B:377:LEU:C	2.28	0.53
1:B:285:TYR:CD2	1:B:312:PRO:HB2	2.43	0.53
1:B:370:THR:OG1	1:B:372:GLU:HG2	2.08	0.53
1:B:342:TRP:CH2	1:B:345:ILE:HA	2.44	0.53
1:A:82:ALA:HB3	3:A:417:HOH:O	2.08	0.53
1:B:125:ASP:HA	1:B:167:LYS:HB3	1.89	0.53
1:A:106:VAL:CG1	1:A:107:PRO:HD2	2.39	0.53
1:B:317:ASN:HB3	1:B:341:GLU:O	2.09	0.53
1:A:329:HIS:NE2	1:A:375:LEU:HA	2.23	0.53
1:B:6:ILE:HG13	1:B:59:GLU:O	2.09	0.53
1:A:106:VAL:HG12	1:A:110:ARG:HB2	1.90	0.52
1:B:2:SER:O	1:B:3:ALA:HB3	2.09	0.52
1:A:51:ASN:ND2	1:B:41:PRO:HB3	2.25	0.52
1:B:6:ILE:HD12	1:B:59:GLU:HB2	1.91	0.52
1:A:133:ALA:HB3	1:B:74:TRP:CZ2	2.44	0.52
1:A:285:TYR:CD2	1:A:312:PRO:HB2	2.45	0.52
1:B:178:HIS:CE1	1:B:180:ASP:O	2.63	0.52
1:B:121:ARG:HB2	1:B:366:HIS:CE1	2.45	0.52
3:A:410:HOH:O	1:B:40:ALA:HB2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:205:VAL:CG1	1:B:209:VAL:HG21	2.38	0.51
1:A:296:LEU:HD11	1:A:326:ALA:HB1	1.92	0.51
1:B:320:SER:HA	1:B:350:TRP:NE1	2.25	0.51
1:A:71:MSE:O	1:A:75:ASN:ND2	2.40	0.51
1:B:123:TYR:CE2	1:B:341:GLU:HB2	2.46	0.51
1:B:396:PHE:C	1:B:398:GLY:H	2.14	0.51
1:B:43:ALA:O	1:B:46:HIS:HB3	2.11	0.51
1:B:363:ARG:O	1:B:364:GLU:HB2	2.10	0.51
1:A:82:ALA:CB	3:A:417:HOH:O	2.58	0.51
1:B:52:MSE:HE1	1:B:74:TRP:HD1	1.76	0.51
1:A:120:VAL:CG2	1:A:367:ILE:HB	2.41	0.51
1:B:148:ASN:O	1:B:152:ILE:HG13	2.11	0.51
1:B:347:MSE:HG3	1:B:347:MSE:O	2.11	0.51
1:A:246:GLU:HB2	1:B:272:PHE:CG	2.45	0.51
1:B:43:ALA:HA	1:B:394:VAL:CG2	2.41	0.50
1:A:272:PHE:O	1:A:276:GLU:HG3	2.10	0.50
1:B:71:MSE:SE	1:B:89:LEU:HD22	2.62	0.50
1:A:6:ILE:HD11	1:A:61:PRO:HG3	1.92	0.50
1:B:26:TRP:CE3	1:B:31:HIS:HB3	2.46	0.50
1:B:40:ALA:CB	1:B:41:PRO:CD	2.77	0.50
1:B:197:ARG:O	1:B:201:VAL:HG23	2.11	0.50
1:A:128:ALA:HB2	1:A:151:TYR:CZ	2.46	0.50
1:A:131:VAL:HA	1:A:136:HIS:O	2.12	0.50
1:A:144:GLU:HB3	1:A:147:SER:HB2	1.94	0.50
1:B:198:VAL:HG13	1:B:213:ILE:HD11	1.94	0.50
1:A:247:ASN:OD1	1:A:249:ASP:HB2	2.11	0.50
1:A:255:ARG:HA	1:A:262:ILE:HG13	1.93	0.50
1:A:98:ASP:HB2	1:A:375:LEU:HD22	1.93	0.49
1:A:135:ALA:HB3	1:A:218:ARG:HH22	1.77	0.49
1:B:320:SER:HB2	1:B:321:PRO:HD2	1.93	0.49
1:A:37:TYR:CB	1:A:292:LYS:HB2	2.43	0.49
1:B:215:MSE:HE2	1:B:239:LEU:CD1	2.43	0.49
1:A:315:PRO:HG2	1:A:340:LEU:CD1	2.42	0.49
1:B:25:VAL:O	1:B:31:HIS:HA	2.13	0.49
1:B:72:ARG:NH1	1:B:271:ARG:HH22	2.10	0.49
1:B:289:ASP:HB3	1:B:292:LYS:HB3	1.94	0.49
1:B:194:MSE:O	1:B:198:VAL:HG23	2.12	0.49
1:A:186:ILE:HG23	1:A:194:MSE:CE	2.42	0.49
1:B:354:VAL:CG1	1:B:377:LEU:HB2	2.41	0.48
1:B:336:ASN:ND2	1:B:336:ASN:N	2.48	0.48
1:B:146:GLY:O	1:B:150:GLU:OE1	2.32	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:228:ALA:HB1	1:B:258:THR:HG21	1.94	0.48
1:B:143:PHE:O	1:B:144:GLU:HB2	2.14	0.48
1:A:43:ALA:O	1:A:46:HIS:HB3	2.13	0.48
1:B:31:HIS:HD2	1:B:102:LYS:NZ	2.10	0.48
1:A:134:ALA:HB3	1:A:136:HIS:CD2	2.48	0.48
1:B:355:ARG:NH1	1:B:378:GLU:HG2	2.28	0.48
1:A:4:PRO:O	1:A:5:ARG:NE	2.47	0.48
1:A:205:VAL:CG1	1:A:209:VAL:HG11	2.40	0.48
1:A:107:PRO:HG2	1:A:110:ARG:HG2	1.95	0.48
1:B:318:VAL:HG13	1:B:318:VAL:O	2.14	0.48
1:B:37:TYR:HB3	1:B:87:THR:HG23	1.96	0.48
1:A:246:GLU:HG3	1:B:272:PHE:CZ	2.48	0.48
1:B:238:TRP:CD1	1:B:263:CYS:HB3	2.49	0.48
1:A:141:GLY:O	1:A:144:GLU:HB2	2.14	0.48
1:A:198:VAL:HG22	1:A:213:ILE:HD13	1.95	0.47
1:A:135:ALA:HB3	1:A:218:ARG:NH2	2.30	0.47
1:B:141:GLY:O	1:B:142:LEU:C	2.52	0.47
1:A:44:GLY:O	1:A:48:ILE:HG13	2.14	0.47
1:A:278:PHE:CZ	1:A:286:VAL:HG11	2.49	0.47
1:B:81:GLN:O	1:B:82:ALA:HB2	2.14	0.47
1:B:83:GLY:HA2	1:B:269:TYR:CZ	2.49	0.47
1:A:22:LEU:CD2	1:A:22:LEU:N	2.77	0.47
1:B:31:HIS:CD2	1:B:102:LYS:NZ	2.82	0.47
1:B:123:TYR:CE1	1:B:167:LYS:HD2	2.47	0.47
1:B:110:ARG:HG2	1:B:110:ARG:HH11	1.79	0.47
1:A:45:ILE:HD12	3:A:418:HOH:O	2.15	0.47
1:B:64:VAL:HG11	1:B:97:TRP:CZ2	2.50	0.47
1:A:26:TRP:CD2	1:A:31:HIS:HB3	2.50	0.47
1:B:357:PRO:HG2	1:B:370:THR:HG21	1.96	0.47
1:A:107:PRO:HG2	1:A:110:ARG:HG3	1.97	0.47
1:B:106:VAL:HG12	1:B:107:PRO:HD2	1.97	0.47
1:B:212:ALA:HB2	1:B:237:LEU:HB3	1.97	0.47
1:A:51:ASN:HD22	1:B:41:PRO:HB3	1.79	0.46
1:A:221:ILE:CB	1:A:222:PRO:HD3	2.39	0.46
1:A:106:VAL:HG12	1:A:107:PRO:HD2	1.97	0.46
1:A:374:GLY:C	1:A:376:GLY:H	2.19	0.46
1:B:238:TRP:C	1:B:238:TRP:CD1	2.88	0.46
1:A:269:TYR:OH	1:B:79:GLY:HA2	2.15	0.46
1:A:364:GLU:HA	1:A:364:GLU:OE1	2.14	0.46
1:A:178:HIS:CD2	1:A:180:ASP:O	2.68	0.46
1:A:108:VAL:HG23	1:A:374:GLY:HA2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:ILE:HG22	1:B:77:PHE:N	2.29	0.46
1:B:144:GLU:O	1:B:145:GLU:C	2.54	0.46
1:A:79:GLY:O	1:A:85:VAL:HB	2.16	0.46
1:A:354:VAL:HG13	1:A:355:ARG:N	2.31	0.46
1:A:278:PHE:CE2	1:A:286:VAL:HG11	2.51	0.46
1:A:5:ARG:O	1:A:27:ALA:HA	2.16	0.46
1:B:64:VAL:HG13	1:B:65:GLU:N	2.31	0.46
1:A:108:VAL:HG21	1:A:375:LEU:HD13	1.96	0.46
1:A:156:ARG:HG2	1:A:156:ARG:HH11	1.81	0.46
1:A:26:TRP:CE3	1:A:31:HIS:HB3	2.52	0.45
1:B:37:TYR:CE1	1:B:318:VAL:HG11	2.52	0.45
1:B:9:VAL:O	1:B:50:MSE:HE1	2.16	0.45
1:A:161:ARG:HG2	1:A:161:ARG:HH11	1.82	0.45
1:B:352:ASP:O	1:B:383:ALA:CB	2.65	0.45
1:B:298:GLU:HB2	3:B:416:HOH:O	2.16	0.45
1:A:41:PRO:HA	3:A:420:HOH:O	2.16	0.45
1:A:215:MSE:HE2	1:A:219:PHE:CD1	2.51	0.45
1:B:6:ILE:HD13	1:B:56:LEU:HB3	1.99	0.45
1:A:25:VAL:O	1:A:31:HIS:HA	2.17	0.45
1:B:263:CYS:SG	1:B:287:MSE:SE	3.25	0.45
1:B:342:TRP:HZ3	1:B:345:ILE:HG12	1.82	0.45
1:A:326:ALA:O	1:A:329:HIS:HB2	2.16	0.45
1:B:81:GLN:CD	1:B:81:GLN:H	2.19	0.45
1:B:21:VAL:HG21	1:B:45:ILE:HB	1.98	0.45
1:A:148:ASN:O	1:A:151:TYR:HB2	2.16	0.45
1:B:37:TYR:CD1	1:B:38:PRO:HD2	2.51	0.45
1:A:79:GLY:O	1:A:82:ALA:HB3	2.17	0.45
1:A:6:ILE:HG22	1:A:57:LEU:HA	1.98	0.45
1:B:305:LEU:O	1:B:308:LEU:HB2	2.17	0.45
1:A:35:GLU:CD	1:A:319:SER:HB2	2.38	0.44
1:A:128:ALA:HB2	1:A:151:TYR:CE1	2.52	0.44
1:A:315:PRO:HG3	1:A:331:CYS:SG	2.58	0.44
1:A:372:GLU:HB3	1:A:373:PRO:HD2	1.98	0.44
1:A:207:PRO:O	1:A:209:VAL:HG23	2.17	0.44
1:B:205:VAL:HG12	1:B:209:VAL:CG2	2.47	0.44
1:B:64:VAL:HG11	1:B:97:TRP:CD2	2.52	0.44
1:B:219:PHE:N	1:B:219:PHE:CD1	2.85	0.44
1:B:254:VAL:O	1:B:257:SER:HB2	2.16	0.44
1:B:52:MSE:HE1	1:B:74:TRP:CD1	2.52	0.44
1:B:263:CYS:HB2	1:B:285:TYR:HB2	1.98	0.44
1:A:24:ARG:HD3	1:A:31:HIS:CD2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:371:GLU:O	1:A:371:GLU:HG3	2.16	0.44
1:B:145:GLU:HG3	1:B:146:GLY:N	2.33	0.44
1:B:390:GLU:O	1:B:391:LYS:CB	2.65	0.44
1:A:16:ALA:O	1:A:17:VAL:O	2.35	0.44
1:A:325:VAL:HG21	1:A:377:LEU:HD21	1.98	0.44
1:A:268:VAL:HG13	1:A:274:PHE:CZ	2.53	0.44
1:A:318:VAL:HG13	1:A:318:VAL:O	2.16	0.44
1:B:128:ALA:O	1:B:131:VAL:HG23	2.18	0.43
1:A:30:GLU:OE1	1:A:30:GLU:HA	2.19	0.43
1:A:132:ASP:HB3	1:A:138:ILE:HD11	2.00	0.43
1:A:125:ASP:OD1	1:A:167:LYS:HD3	2.17	0.43
1:A:320:SER:CB	1:A:321:PRO:CD	2.96	0.43
1:A:107:PRO:HA	1:A:374:GLY:N	2.32	0.43
1:B:106:VAL:CG1	1:B:107:PRO:HD2	2.48	0.43
1:B:72:ARG:NH1	1:B:271:ARG:NH2	2.62	0.43
1:A:83:GLY:HA2	1:A:269:TYR:CE2	2.53	0.43
1:A:12:ALA:HB2	1:A:397:PHE:CE1	2.53	0.43
1:B:253:GLU:HA	1:B:256:ARG:NH2	2.33	0.43
1:A:121:ARG:HB2	1:A:366:HIS:CE1	2.54	0.43
1:A:85:VAL:O	1:A:89:LEU:HG	2.19	0.43
1:B:144:GLU:HG2	1:B:145:GLU:N	2.31	0.43
1:B:108:VAL:HG12	1:B:329:HIS:CG	2.53	0.43
1:B:106:VAL:HG12	1:B:110:ARG:HB2	2.01	0.43
1:A:93:GLU:OE1	1:A:97:TRP:CD1	2.71	0.43
1:A:148:ASN:O	1:A:152:ILE:HG13	2.19	0.43
1:B:342:TRP:CZ3	1:B:345:ILE:HG12	2.53	0.43
1:A:343:HIS:O	1:A:343:HIS:ND1	2.52	0.43
1:A:74:TRP:CZ2	1:B:133:ALA:HB3	2.54	0.43
1:A:62:ARG:HG2	1:A:104:GLN:CD	2.39	0.43
1:B:247:ASN:OD1	1:B:249:ASP:HB2	2.19	0.42
1:B:144:GLU:OE1	1:B:145:GLU:HB3	2.18	0.42
1:A:152:ILE:HG23	1:A:204:ALA:CB	2.49	0.42
1:B:369:LEU:HA	1:B:369:LEU:HD23	1.81	0.42
1:A:35:GLU:OE1	1:A:322:VAL:HB	2.20	0.42
1:B:122:LEU:HD23	1:B:340:LEU:HB3	2.01	0.42
1:B:269:TYR:CD1	1:B:269:TYR:N	2.88	0.42
1:A:52:MSE:HE2	1:A:55:GLN:HG3	2.01	0.42
1:B:147:SER:HA	1:B:150:GLU:OE1	2.19	0.42
1:B:119:ARG:HB3	1:B:368:GLU:CD	2.40	0.42
1:B:154:VAL:O	1:B:155:ALA:C	2.58	0.42
1:B:345:ILE:O	1:B:345:ILE:HD13	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:170:VAL:O	1:B:194:MSE:HG2	2.20	0.42
1:A:263:CYS:SG	1:A:287:MSE:SE	3.28	0.42
1:A:14:ILE:HG23	1:A:22:LEU:CD2	2.50	0.42
1:B:173:ILE:HD12	1:B:173:ILE:N	2.34	0.41
1:A:43:ALA:HB1	1:B:47:HIS:CE1	2.54	0.41
1:A:271:ARG:HG3	1:A:272:PHE:N	2.35	0.41
1:A:382:GLU:O	1:A:386:GLU:HG3	2.20	0.41
1:A:40:ALA:N	1:B:78:THR:HB	2.30	0.41
1:A:205:VAL:HG11	1:A:209:VAL:CG1	2.44	0.41
1:B:76:ILE:HG22	1:B:77:PHE:CD1	2.55	0.41
1:B:76:ILE:HG22	1:B:77:PHE:CG	2.54	0.41
1:A:396:PHE:O	1:A:397:PHE:HB2	2.20	0.41
1:A:64:VAL:HG21	1:A:97:TRP:CE2	2.55	0.41
1:B:197:ARG:HH11	1:B:197:ARG:HG2	1.85	0.41
1:B:108:VAL:HG12	1:B:109:TYR:N	2.35	0.41
1:B:14:ILE:O	1:B:19:PRO:HA	2.20	0.41
1:B:70:LYS:O	1:B:74:TRP:HB2	2.20	0.41
1:A:156:ARG:O	1:A:160:GLU:HB2	2.20	0.41
1:A:165:ALA:HA	1:A:210:GLU:O	2.20	0.41
1:B:60:ASP:OD2	1:B:62:ARG:NH2	2.53	0.41
1:A:11:THR:OG1	1:A:46:HIS:CE1	2.74	0.41
1:B:354:VAL:HG11	1:B:377:LEU:HD22	2.02	0.41
1:A:108:VAL:HG23	1:A:375:LEU:H	1.83	0.41
1:B:144:GLU:CG	1:B:145:GLU:N	2.84	0.41
1:B:165:ALA:C	1:B:166:ILE:HD12	2.41	0.41
1:A:106:VAL:CG1	1:A:110:ARG:HB2	2.51	0.41
1:B:99:LEU:O	1:B:103:LEU:HG	2.21	0.41
1:B:396:PHE:C	1:B:398:GLY:N	2.74	0.40
1:A:145:GLU:C	1:A:147:SER:H	2.24	0.40
1:A:37:TYR:O	1:A:39:SER:N	2.54	0.40
1:A:186:ILE:HD13	1:A:194:MSE:HE1	2.04	0.40
1:A:39:SER:O	1:A:40:ALA:C	2.60	0.40
1:B:215:MSE:CE	1:B:239:LEU:HD21	2.51	0.40
1:A:52:MSE:HG2	1:A:71:MSE:SE	2.72	0.40
1:B:268:VAL:HG23	1:B:273:ASP:HB3	2.02	0.40
1:B:86:ILE:CG2	1:B:293:CYS:O	2.69	0.40
1:B:95:ALA:O	1:B:98:ASP:HB3	2.22	0.40
1:B:9:VAL:HG12	1:B:50:MSE:HE1	2.03	0.40
1:A:159:VAL:CG2	1:A:160:GLU:N	2.84	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	388/410 (95%)	346 (89%)	35 (9%)	7 (2%)	11	37
1	B	392/410 (96%)	336 (86%)	42 (11%)	14 (4%)	4	18
All	All	780/820 (95%)	682 (87%)	77 (10%)	21 (3%)	6	25

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	40	ALA
1	A	82	ALA
1	B	142	LEU
1	A	374	GLY
1	B	40	ALA
1	B	145	GLU
1	B	357	PRO
1	B	397	PHE
1	A	357	PRO
1	B	270	THR
1	B	374	GLY
1	A	12	ALA
1	B	144	GLU
1	B	382	GLU
1	B	131	VAL
1	A	322	VAL
1	B	61	PRO
1	B	217	GLY
1	B	294	GLY
1	A	38	PRO
1	B	3	ALA

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	298/312 (96%)	275 (92%)	23 (8%)	16	42
1	B	300/312 (96%)	276 (92%)	24 (8%)	15	40
All	All	598/624 (96%)	551 (92%)	47 (8%)	15	41

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	ARG
1	A	64	VAL
1	A	67	LEU
1	A	72	ARG
1	A	81	GLN
1	A	93	GLU
1	A	117	ARG
1	A	144	GLU
1	A	147	SER
1	A	159	VAL
1	A	183	ASN
1	A	192	GLU
1	A	211	VAL
1	A	219	PHE
1	A	246	GLU
1	A	257	SER
1	A	258	THR
1	A	266	GLU
1	A	268	VAL
1	A	286	VAL
1	A	336	ASN
1	A	354	VAL
1	A	371	GLU
1	B	5	ARG
1	B	31	HIS
1	B	50	MSE
1	B	76	ILE

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Mol	Chain	Res	Type
1	B	81	GLN
1	B	93	GLU
1	B	97	TRP
1	B	118	ARG
1	B	120	VAL
1	B	127	ASN
1	B	130	THR
1	B	144	GLU
1	B	168	LEU
1	B	268	VAL
1	B	269	TYR
1	B	280	LYS
1	B	301	ARG
1	B	336	ASN
1	B	339	VAL
1	B	345	ILE
1	B	357	PRO
1	B	370	THR
1	B	377	LEU
1	B	382	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	HIS
1	A	51	ASN
1	A	136	HIS
1	A	137	HIS
1	A	148	ASN
1	A	183	ASN
1	A	191	HIS
1	A	336	ASN
1	B	31	HIS
1	B	127	ASN
1	B	178	HIS
1	B	336	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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	386/410 (94%)	-0.29	3 (0%) 87 86	20, 39, 59, 69	0
1	B	388/410 (94%)	-0.14	3 (0%) 87 86	18, 43, 62, 72	0
All	All	774/820 (94%)	-0.21	6 (0%) 87 86	18, 42, 61, 72	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	2	SER	3.6
1	B	131	VAL	2.6
1	A	2	SER	2.5
1	B	358	GLY	2.2
1	A	144	GLU	2.1
1	A	385	PHE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors

of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	MG	B	409	1/1	0.87	0.21	1.68	45,45,45,45	0
2	MG	A	409	1/1	0.98	0.17	1.17	18,18,18,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.