



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:54 AM UTC

PDB ID : 4GBG / pdb_00004gbg
Title : Crystal structure of Ethyl acetoacetate treated lipase from *Thermomyces lanuginosa* at 2.9 Å resolution
Authors : Yamini, S.; Mukherjee, J.; Gupta, M.N.; Sinha, M.; Kaur, P.; Sharma, S.; Singh, T.P.
Deposited on : 2012-07-27
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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

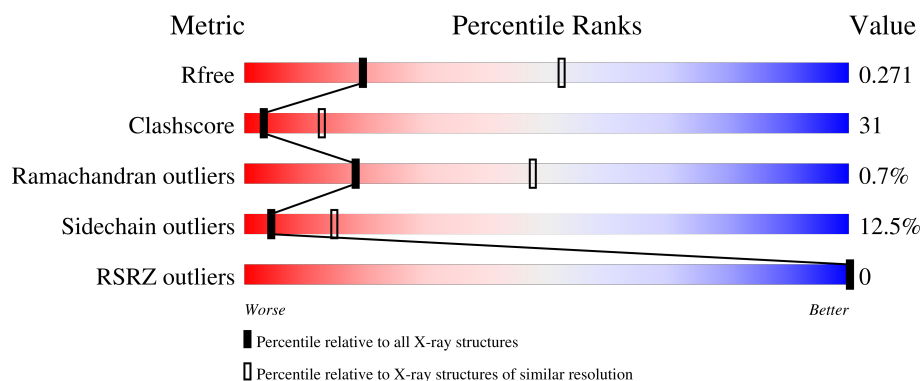
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

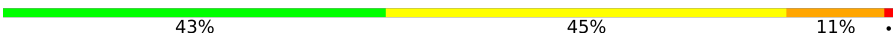
The reported resolution of this entry is 2.90 Å.

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}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (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 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 $\leq 5\%$. 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	269	
1	B	269	

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	EAC	A	301	-	X	-	-

2 Entry composition [i](#)

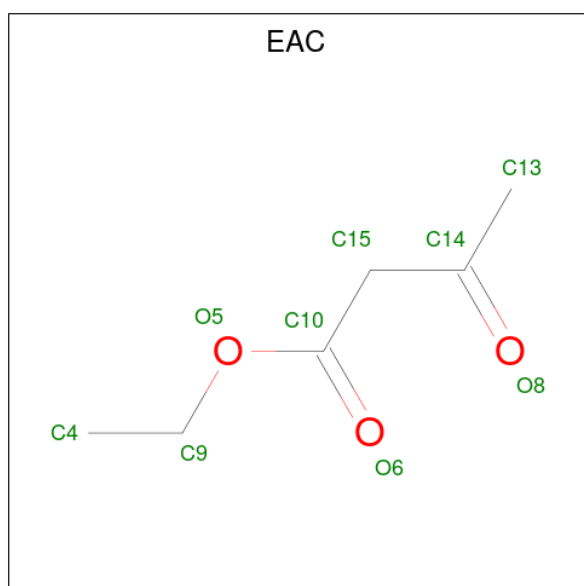
There are 4 unique types of molecules in this entry. The entry contains 4527 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Lipase.

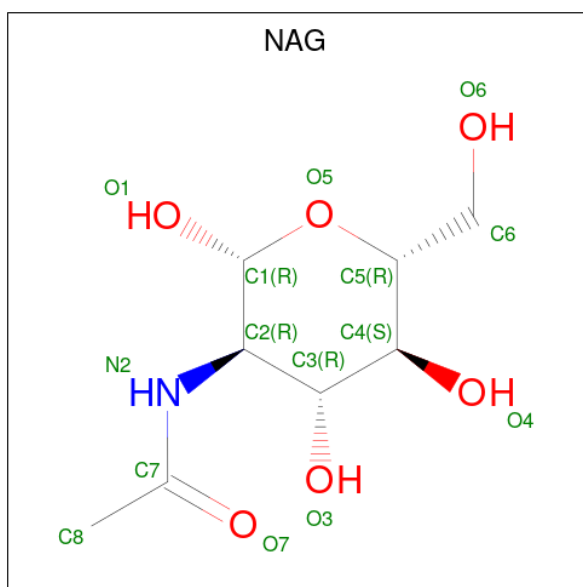
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	269	Total	C	N	O	S	0	0	0
			2070	1303	359	402	6			
1	B	269	Total	C	N	O	S	0	0	0
			2070	1303	359	402	6			

- Molecule 2 is ethyl 3-oxobutanoate (CCD ID: EAC) (formula: $C_6H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			9	6	3		

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		

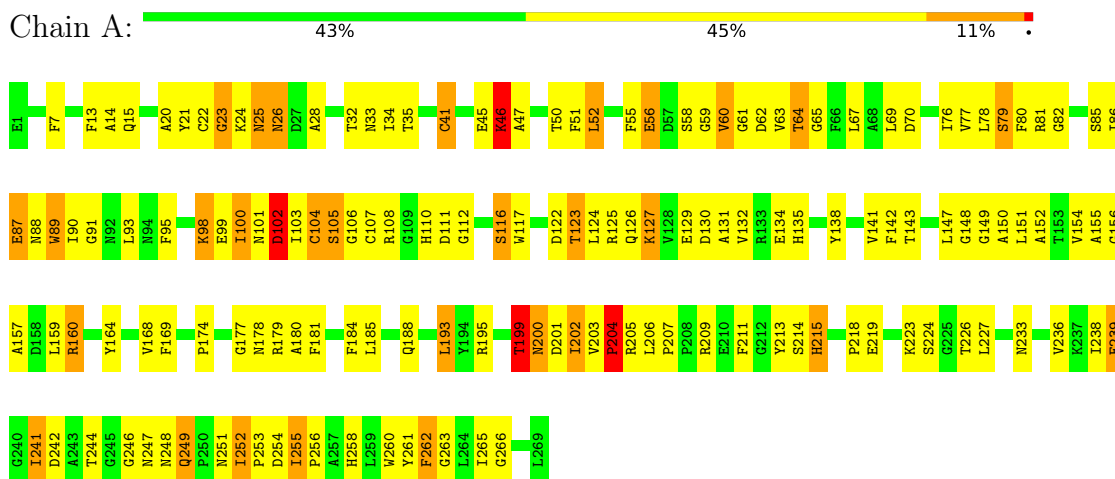
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	175	Total	O	0	0
			175	175		
4	B	175	Total	O	0	0
			175	175		

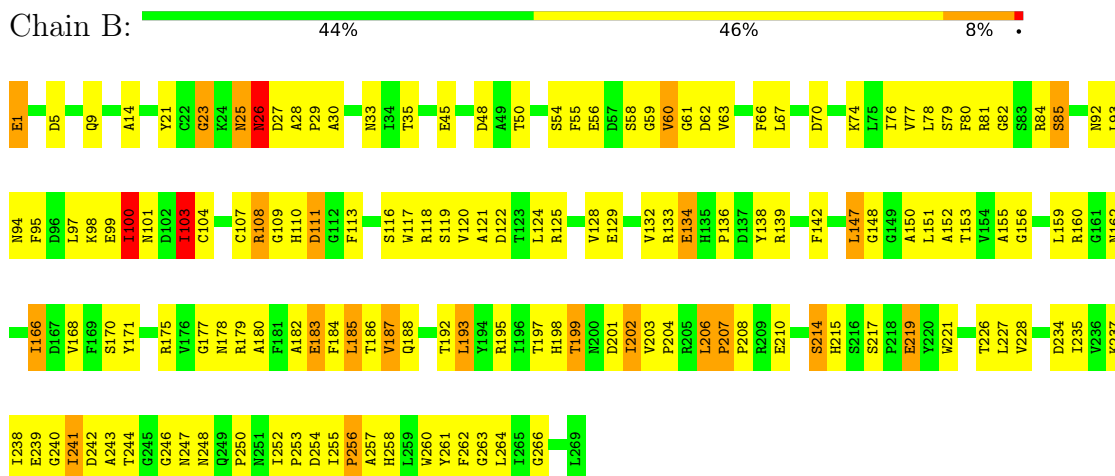
3 Residue-property plots

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: Lipase



• Molecule 1: Lipase



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	140.49Å 140.49Å 80.69Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.56 – 2.90 40.56 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.56-2.90) 100.0 (40.56-2.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.83 (at 2.90Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.226 , 0.247 0.238 , 0.271	Depositor DCC
R_{free} test set	1036 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	53.5	Xtriage
Anisotropy	0.063	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 56.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.084 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4527	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: EAC, NAG

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	0.73	2/2120 (0.1%)	1.18	25/2887 (0.9%)
1	B	0.72	1/2120 (0.0%)	1.29	23/2887 (0.8%)
All	All	0.73	3/4240 (0.1%)	1.24	48/5774 (0.8%)

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	B	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	258	HIS	CG-ND1	-5.95	1.31	1.38
1	B	26	ASN	CA-C	-5.33	1.48	1.53
1	A	258	HIS	CD2-NE2	-5.17	1.32	1.37

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	187	VAL	CB-CA-C	-12.91	90.12	111.29
1	B	103	ILE	N-CA-C	-11.91	101.32	111.56
1	B	155	ALA	N-CA-C	-11.58	99.15	113.28
1	B	203	VAL	CA-C-N	-10.21	110.42	120.83
1	B	203	VAL	C-N-CA	-10.21	110.42	120.83
1	A	60	VAL	N-CA-C	9.73	119.77	110.42
1	B	60	VAL	N-CA-C	9.73	120.58	110.36
1	A	102	ASP	N-CA-C	-9.54	101.06	112.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	122	ASP	N-CA-C	9.07	120.78	111.07
1	A	76	ILE	N-CA-C	-7.88	94.75	107.28
1	B	30	ALA	N-CA-C	-7.70	99.88	110.35
1	B	179	ARG	N-CA-C	7.18	119.97	111.71
1	B	76	ILE	N-CA-C	-7.11	95.97	107.28
1	A	52	LEU	N-CA-C	-6.96	104.66	113.43
1	A	262	PHE	CB-CA-C	-6.89	96.71	110.42
1	B	180	ALA	N-CA-C	-6.85	103.81	111.28
1	A	199	THR	CB-CA-C	-6.63	108.92	116.54
1	A	93	LEU	N-CA-C	6.62	119.62	108.23
1	A	56	GLU	N-CA-C	6.46	119.67	108.76
1	A	41	CYS	CA-C-N	6.42	126.18	119.05
1	A	41	CYS	C-N-CA	6.42	126.18	119.05
1	B	153	THR	N-CA-C	-6.33	105.71	113.50
1	B	177	GLY	N-CA-C	6.22	118.96	110.69
1	A	14	ALA	N-CA-C	-6.21	104.62	111.82
1	B	111	ASP	N-CA-C	6.12	118.48	111.02
1	B	23	GLY	N-CA-C	-6.04	103.41	113.02
1	A	23	GLY	N-CA-C	-6.00	101.80	112.54
1	B	85	SER	N-CA-C	-5.96	97.48	107.80
1	B	14	ALA	N-CA-C	-5.96	104.86	111.36
1	B	234	ASP	N-CA-C	-5.81	105.52	113.30
1	B	59	GLY	N-CA-C	5.63	122.50	112.02
1	B	82	GLY	N-CA-C	-5.62	104.86	112.57
1	A	244	THR	N-CA-C	-5.54	101.20	109.96
1	A	46	LYS	N-CA-C	-5.49	105.39	111.71
1	A	209	ARG	N-CA-C	-5.49	105.85	112.54
1	A	265	ILE	N-CA-C	5.47	115.07	108.06
1	B	26	ASN	N-CA-C	-5.31	100.23	108.31
1	A	104	CYS	N-CA-C	5.27	117.08	108.76
1	A	89	TRP	N-CA-C	-5.24	105.04	111.75
1	A	103	ILE	N-CA-C	-5.23	106.59	111.45
1	A	98	LYS	N-CA-C	-5.22	101.65	109.95
1	A	22	CYS	N-CA-C	5.22	119.61	112.88
1	A	25	ASN	N-CA-C	5.21	118.75	111.56
1	A	77	VAL	N-CA-C	5.12	115.28	108.11
1	A	239	GLU	N-CA-C	5.07	118.47	109.96
1	A	105	SER	CB-CA-C	5.05	117.44	110.16
1	B	228	VAL	CA-C-N	-5.01	115.40	120.31
1	B	228	VAL	C-N-CA	-5.01	115.40	120.31

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	171	TYR	Sidechain

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	2070	0	1963	135	0
1	B	2070	0	1964	117	0
2	A	9	0	10	1	0
3	A	14	0	13	0	0
3	B	14	0	13	0	0
4	A	175	0	0	7	0
4	B	175	0	0	4	0
All	All	4527	0	3963	252	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 (252) 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:101:ASN:OD1	1:A:104:CYS:O	1.80	0.99
1:B:195:ARG:NH2	1:B:219:GLU:HG3	1.78	0.97
1:A:202:ILE:HG13	1:A:253:PRO:HB2	1.48	0.91
1:A:252:ILE:HD12	1:A:253:PRO:HD2	1.53	0.90
1:A:201:ASP:OD1	1:A:202:ILE:N	2.05	0.89
1:A:61:GLY:HA3	1:A:116:SER:HB3	1.54	0.88
1:B:254:ASP:OD1	1:B:256:PRO:HD2	1.75	0.86
1:A:56:GLU:HA	1:A:64:THR:HG23	1.57	0.86
1:B:182:ALA:HB2	1:B:214:SER:HB3	1.58	0.85
1:A:101:ASN:OD1	1:A:106:GLY:N	2.09	0.85
1:A:131:ALA:HA	1:A:134:GLU:HG3	1.58	0.84
1:B:23:GLY:HA3	1:B:25:ASN:HD21	1.46	0.81
1:B:195:ARG:HH22	1:B:219:GLU:HG3	1.44	0.79
1:B:26:ASN:N	1:B:26:ASN:HD22	1.81	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:117:TRP:HZ3	1:B:124:LEU:HD12	1.48	0.77
1:B:101:ASN:HA	1:B:104:CYS:O	1.85	0.77
1:A:261:TYR:O	1:A:262:PHE:HB2	1.85	0.76
1:B:254:ASP:CG	1:B:256:PRO:HD2	2.11	0.75
1:A:248:ASN:HB2	4:A:515:HOH:O	1.86	0.74
1:A:32:THR:HG22	1:A:33:ASN:H	1.54	0.73
1:B:156:GLY:O	1:B:160:ARG:HG3	1.89	0.73
1:A:195:ARG:NH2	1:A:219:GLU:HB2	2.04	0.72
1:A:254:ASP:OD2	1:A:256:PRO:HD2	1.89	0.72
1:A:223:LYS:HD2	1:A:236:VAL:CG2	2.21	0.70
1:B:175:ARG:HG3	1:B:215:HIS:CD2	2.26	0.70
1:A:201:ASP:OD1	1:A:201:ASP:C	2.30	0.70
1:A:254:ASP:CG	1:A:256:PRO:HD2	2.17	0.69
1:A:15:GLN:HG2	1:A:41:CYS:HA	1.74	0.69
1:B:226:THR:O	1:B:227:LEU:HB2	1.91	0.68
1:A:81:ARG:HG2	1:A:82:GLY:N	2.09	0.68
1:B:128:VAL:O	1:B:132:VAL:HG23	1.94	0.68
1:A:20:ALA:HB1	1:A:81:ARG:HB2	1.74	0.68
1:A:32:THR:HG22	1:A:33:ASN:N	2.09	0.68
1:B:133:ARG:O	1:B:136:PRO:HD3	1.94	0.68
1:B:95:PHE:HE2	1:B:208:PRO:HD3	1.59	0.67
1:B:1:GLU:O	1:B:1:GLU:HG2	1.93	0.67
1:B:81:ARG:NH2	1:B:84:ARG:HG2	2.10	0.67
1:A:87:GLU:HB2	2:A:301:EAC:O6	1.94	0.66
1:A:101:ASN:HA	1:A:104:CYS:O	1.94	0.66
1:A:200:ASN:O	1:A:201:ASP:C	2.39	0.66
1:A:101:ASN:CG	1:A:106:GLY:H	2.03	0.66
1:B:206:LEU:HA	1:B:207:PRO:C	2.21	0.64
1:B:67:LEU:HD12	1:B:77:VAL:O	1.97	0.64
1:B:219:GLU:HB3	1:B:238:ILE:HB	1.80	0.64
1:A:168:VAL:HG12	1:A:193:LEU:HD12	1.80	0.64
1:A:202:ILE:CG1	1:A:253:PRO:HB2	2.23	0.64
1:A:104:CYS:SG	1:A:184:PHE:CD2	2.91	0.64
1:A:226:THR:O	1:A:227:LEU:HB2	1.96	0.63
1:B:61:GLY:HA3	1:B:116:SER:OG	1.99	0.63
1:A:200:ASN:O	1:A:201:ASP:O	2.16	0.62
1:A:205:ARG:NH2	1:A:251:ASN:OD1	2.20	0.62
1:A:135:HIS:ND1	4:A:517:HOH:O	2.31	0.62
1:A:202:ILE:HG13	1:A:253:PRO:CB	2.26	0.62
1:A:203:VAL:C	1:A:205:ARG:H	2.08	0.61
1:A:117:TRP:CD1	1:A:154:VAL:HG12	2.36	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:TYR:CZ	1:A:266:GLY:HA2	2.35	0.61
1:B:109:GLY:O	1:B:110:HIS:C	2.39	0.61
1:B:110:HIS:HB3	1:B:113:PHE:CD2	2.35	0.61
1:B:80:PHE:HD2	1:B:152:ALA:HB2	1.66	0.60
1:B:85:SER:HB3	4:B:434:HOH:O	2.01	0.60
1:B:175:ARG:HD2	1:B:207:PRO:O	2.02	0.59
1:B:100:ILE:HG13	1:B:107:CYS:O	2.03	0.59
1:B:183:GLU:HA	1:B:241:ILE:HD11	1.83	0.59
1:B:182:ALA:CB	1:B:214:SER:HB3	2.31	0.59
1:A:95:PHE:HD2	1:A:211:PHE:CE1	2.20	0.58
1:A:227:LEU:H	1:A:260:TRP:CG	2.22	0.58
1:A:107:CYS:SG	1:A:181:PHE:HA	2.45	0.57
1:B:201:ASP:CG	1:B:258:HIS:HD1	2.12	0.57
1:A:88:ASN:O	1:A:89:TRP:C	2.47	0.57
1:A:168:VAL:CG1	1:A:193:LEU:HD12	2.34	0.57
1:B:103:ILE:HG22	1:B:104:CYS:SG	2.45	0.56
1:B:195:ARG:CZ	1:B:219:GLU:HG3	2.35	0.56
1:B:98:LYS:HB2	1:B:111:ASP:OD1	2.05	0.56
1:B:237:LYS:O	1:B:238:ILE:HD13	2.06	0.56
1:A:59:GLY:HA2	4:A:411:HOH:O	2.06	0.56
1:A:206:LEU:HA	1:A:207:PRO:C	2.31	0.56
1:B:254:ASP:OD2	1:B:256:PRO:HB2	2.06	0.56
1:B:125:ARG:HD3	1:B:159:LEU:HD22	1.88	0.55
1:A:100:ILE:HD12	1:A:107:CYS:O	2.06	0.55
1:B:95:PHE:CE2	1:B:208:PRO:HD3	2.40	0.55
1:B:93:LEU:HB3	4:B:512:HOH:O	2.07	0.55
1:A:88:ASN:O	1:A:91:GLY:N	2.40	0.54
1:A:80:PHE:O	1:A:148:GLY:HA3	2.07	0.54
1:B:26:ASN:N	1:B:26:ASN:ND2	2.52	0.54
1:B:261:TYR:O	1:B:262:PHE:HB2	2.08	0.54
1:A:203:VAL:HB	1:A:204:PRO:HD3	1.90	0.54
1:B:80:PHE:O	1:B:148:GLY:HA3	2.07	0.54
1:B:117:TRP:CZ3	1:B:124:LEU:HD12	2.37	0.54
1:A:107:CYS:HA	1:A:180:ALA:HB3	1.90	0.54
1:A:179:ARG:HG3	1:A:179:ARG:O	2.08	0.54
1:A:157:ALA:HB2	1:A:185:LEU:CD2	2.38	0.54
1:A:34:ILE:HG12	1:A:51:PHE:CZ	2.43	0.53
1:A:101:ASN:OD1	1:A:105:SER:CA	2.56	0.53
1:A:130:ASP:O	1:A:134:GLU:HG3	2.08	0.53
1:A:223:LYS:HD2	1:A:236:VAL:HG23	1.89	0.53
1:B:25:ASN:O	1:B:26:ASN:C	2.49	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:23:GLY:HA3	1:B:25:ASN:ND2	2.21	0.53
1:A:102:ASP:OD2	1:A:102:ASP:N	2.30	0.53
1:B:61:GLY:O	1:B:62:ASP:C	2.52	0.53
1:A:58:SER:HA	4:A:497:HOH:O	2.09	0.53
1:B:204:PRO:HB2	1:B:247:ASN:OD1	2.08	0.52
1:A:34:ILE:HG12	1:A:51:PHE:CE2	2.45	0.52
1:A:101:ASN:OD1	1:A:105:SER:HA	2.09	0.52
1:A:203:VAL:C	1:A:205:ARG:N	2.68	0.52
1:A:204:PRO:HB2	1:A:247:ASN:OD1	2.09	0.52
1:A:238:ILE:HD12	1:A:246:GLY:HA3	1.92	0.52
1:A:238:ILE:HD12	1:A:246:GLY:CA	2.40	0.52
1:A:156:GLY:O	1:A:160:ARG:HG3	2.10	0.52
1:B:80:PHE:HB2	1:B:148:GLY:O	2.10	0.51
1:A:178:ASN:OD1	1:A:180:ALA:HB3	2.10	0.51
1:A:223:LYS:HD2	1:A:236:VAL:HG21	1.91	0.51
1:A:25:ASN:HA	1:A:28:ALA:HB2	1.93	0.51
1:B:184:PHE:O	1:B:185:LEU:C	2.51	0.51
1:A:101:ASN:ND2	1:A:106:GLY:H	2.09	0.51
1:B:108:ARG:HH11	1:B:108:ARG:CG	2.22	0.51
1:A:215:HIS:HB2	4:A:522:HOH:O	2.11	0.51
1:A:78:LEU:HB3	1:A:142:PHE:CD2	2.46	0.50
1:A:98:LYS:HB2	1:A:111:ASP:OD1	2.12	0.50
1:A:242:ASP:HA	4:A:416:HOH:O	2.10	0.50
1:A:227:LEU:H	1:A:260:TRP:CD1	2.29	0.50
1:B:26:ASN:HA	1:B:56:GLU:OE2	2.12	0.50
1:A:45:GLU:O	1:A:46:LYS:C	2.54	0.50
1:B:227:LEU:H	1:B:260:TRP:CD1	2.29	0.50
1:B:78:LEU:HB3	1:B:142:PHE:CD1	2.47	0.49
1:B:125:ARG:O	1:B:129:GLU:HG3	2.13	0.49
1:A:150:ALA:HB2	1:A:174:PRO:HD2	1.95	0.49
1:A:52:LEU:HB2	1:A:67:LEU:O	2.12	0.49
1:A:81:ARG:HG2	1:A:82:GLY:O	2.12	0.49
1:A:143:THR:HA	1:A:169:PHE:O	2.12	0.49
1:A:101:ASN:OD1	1:A:105:SER:C	2.56	0.49
1:B:26:ASN:O	1:B:27:ASP:OD2	2.30	0.49
1:B:99:GLU:HB2	1:B:108:ARG:NH1	2.27	0.49
1:A:130:ASP:O	1:A:134:GLU:CG	2.60	0.49
1:A:261:TYR:C	1:A:263:GLY:H	2.20	0.49
1:B:104:CYS:O	1:B:107:CYS:HB2	2.12	0.49
1:A:32:THR:O	1:A:50:THR:HB	2.13	0.48
1:A:122:ASP:O	1:A:123:THR:C	2.53	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:254:ASP:OD2	1:B:257:ALA:N	2.43	0.48
1:A:255:ILE:N	1:A:256:PRO:CD	2.76	0.48
1:B:100:ILE:HD12	1:B:107:CYS:HB3	1.95	0.48
1:B:244:THR:HG22	1:B:248:ASN:O	2.13	0.48
1:B:201:ASP:OD2	1:B:258:HIS:HB2	2.14	0.47
1:A:7:PHE:CE1	1:A:262:PHE:HE1	2.32	0.47
1:B:175:ARG:NH1	4:B:404:HOH:O	2.36	0.47
1:B:97:LEU:HD12	1:B:97:LEU:N	2.30	0.47
1:B:186:THR:HG23	1:B:217:SER:HB3	1.96	0.47
1:A:85:SER:O	1:A:88:ASN:HB2	2.15	0.47
1:B:184:PHE:CE1	1:B:188:GLN:HG3	2.50	0.47
1:A:81:ARG:CG	1:A:82:GLY:N	2.77	0.47
1:B:208:PRO:HB2	1:B:210:GLU:OE2	2.14	0.47
1:B:94:ASN:O	1:B:110:HIS:NE2	2.42	0.46
1:B:263:GLY:O	1:B:264:LEU:C	2.58	0.46
1:A:61:GLY:O	1:A:62:ASP:C	2.58	0.46
1:A:124:LEU:O	1:A:127:LYS:N	2.49	0.46
1:B:26:ASN:HD22	1:B:26:ASN:H	1.60	0.46
1:B:197:THR:HB	1:B:247:ASN:HB2	1.98	0.46
1:A:88:ASN:O	1:A:90:ILE:N	2.48	0.46
1:A:25:ASN:HB3	1:A:51:PHE:CZ	2.51	0.46
1:A:252:ILE:CD1	1:A:253:PRO:HD2	2.37	0.46
1:A:184:PHE:O	1:A:188:GLN:HB2	2.16	0.46
1:A:203:VAL:O	1:A:205:ARG:N	2.49	0.46
1:B:48:ASP:CG	1:B:48:ASP:O	2.53	0.46
1:B:100:ILE:CD1	1:B:107:CYS:HB3	2.46	0.46
1:B:175:ARG:HG3	1:B:215:HIS:NE2	2.31	0.46
1:A:127:LYS:HA	1:A:130:ASP:OD1	2.17	0.45
1:B:21:TYR:CZ	1:B:266:GLY:HA2	2.50	0.45
1:B:70:ASP:C	1:B:70:ASP:OD1	2.59	0.45
1:B:183:GLU:O	1:B:187:VAL:HG22	2.17	0.45
1:B:160:ARG:HG2	1:B:166:ILE:HG13	1.99	0.45
1:A:69:LEU:HD11	1:A:138:TYR:CE2	2.51	0.45
1:A:174:PRO:HB3	1:A:203:VAL:HG12	1.99	0.45
1:A:88:ASN:C	1:A:90:ILE:N	2.72	0.45
1:A:117:TRP:NE1	1:A:155:ALA:HA	2.32	0.45
1:B:97:LEU:HD12	1:B:97:LEU:H	1.82	0.45
1:B:33:ASN:HA	1:B:50:THR:HG22	1.99	0.44
1:B:128:VAL:O	1:B:129:GLU:C	2.60	0.44
1:B:84:ARG:HB2	4:B:503:HOH:O	2.16	0.44
1:B:198:HIS:O	1:B:199:THR:C	2.59	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:255:ILE:N	1:B:256:PRO:CD	2.80	0.44
1:B:134:GLU:H	1:B:134:GLU:HG2	1.42	0.44
1:A:110:HIS:CE1	1:A:112:GLY:HA3	2.52	0.44
1:B:110:HIS:CB	1:B:113:PHE:CD2	3.00	0.44
1:B:120:VAL:O	1:B:121:ALA:C	2.59	0.44
1:A:108:ARG:H	1:A:178:ASN:CG	2.26	0.44
1:B:99:GLU:HG2	1:B:100:ILE:N	2.33	0.44
1:A:56:GLU:HG2	1:A:64:THR:CG2	2.48	0.44
1:A:110:HIS:HE1	1:A:112:GLY:HA3	1.83	0.44
1:A:125:ARG:O	1:A:129:GLU:HG3	2.17	0.44
1:A:131:ALA:CA	1:A:134:GLU:HG3	2.39	0.44
1:A:23:GLY:HA2	1:A:35:THR:O	2.18	0.43
1:B:118:ARG:HD3	1:B:118:ARG:HA	1.57	0.43
1:B:192:THR:HG22	1:B:193:LEU:N	2.33	0.43
1:B:35:THR:HA	1:B:45:GLU:HG2	2.01	0.43
1:A:25:ASN:ND2	1:A:51:PHE:CZ	2.87	0.43
1:A:124:LEU:O	1:A:125:ARG:C	2.58	0.43
1:B:104:CYS:SG	1:B:184:PHE:CD2	3.11	0.43
1:A:241:ILE:HG22	1:A:242:ASP:OD1	2.19	0.43
1:B:255:ILE:O	1:B:258:HIS:HB3	2.19	0.43
1:A:85:SER:O	1:A:86:ILE:C	2.60	0.43
1:B:240:GLY:O	1:B:243:ALA:HB2	2.19	0.43
1:A:155:ALA:O	1:A:159:LEU:HG	2.19	0.42
1:B:108:ARG:CG	1:B:108:ARG:NH1	2.80	0.42
1:B:55:PHE:HE1	1:B:63:VAL:HG12	1.85	0.42
1:A:56:GLU:HG2	1:A:64:THR:HG23	2.01	0.42
1:A:199:THR:HB	1:A:200:ASN:H	1.62	0.42
1:A:21:TYR:CE1	1:A:266:GLY:HA2	2.55	0.42
1:A:70:ASP:OD1	1:A:70:ASP:C	2.62	0.42
1:A:86:ILE:O	1:A:90:ILE:HG13	2.18	0.42
1:B:9:GLN:OE1	1:B:139:ARG:NH2	2.45	0.42
1:B:74:LYS:HD2	1:B:138:TYR:CE1	2.54	0.42
1:A:151:LEU:O	1:A:154:VAL:N	2.53	0.42
1:B:21:TYR:CE2	1:B:266:GLY:HA2	2.54	0.42
1:B:185:LEU:H	1:B:185:LEU:HG	1.66	0.42
1:B:221:TRP:CD2	1:B:246:GLY:HA2	2.55	0.42
1:A:132:VAL:HG11	1:A:164:TYR:CD2	2.55	0.42
1:A:126:GLN:O	1:A:130:ASP:OD1	2.37	0.42
1:B:5:ASP:O	1:B:9:GLN:HG3	2.19	0.42
1:A:151:LEU:O	1:A:152:ALA:C	2.62	0.42
1:A:246:GLY:O	1:A:249:GLN:NE2	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:186:THR:O	1:B:186:THR:HG22	2.19	0.42
1:B:210:GLU:H	1:B:210:GLU:CD	2.27	0.42
1:A:65:GLY:HA3	1:A:79:SER:O	2.20	0.42
1:B:28:ALA:O	1:B:29:PRO:C	2.62	0.42
1:A:147:LEU:C	1:A:149:GLY:N	2.76	0.41
1:B:142:PHE:O	1:B:168:VAL:HA	2.19	0.41
1:A:52:LEU:HD23	1:A:52:LEU:HA	1.74	0.41
1:B:201:ASP:OD2	1:B:258:HIS:ND1	2.54	0.41
1:B:214:SER:OG	1:B:242:ASP:OD1	2.23	0.41
1:A:55:PHE:HE1	1:A:63:VAL:HG12	1.86	0.41
1:A:45:GLU:C	1:A:47:ALA:N	2.76	0.41
1:A:101:ASN:HD21	1:A:106:GLY:H	1.69	0.41
1:B:66:PHE:CE2	1:B:79:SER:HB3	2.56	0.41
1:B:95:PHE:CD1	1:B:95:PHE:N	2.88	0.41
1:B:118:ARG:O	1:B:119:SER:C	2.63	0.41
1:B:227:LEU:H	1:B:260:TRP:CG	2.38	0.41
1:A:95:PHE:HD2	1:A:211:PHE:CD1	2.39	0.41
1:B:150:ALA:C	1:B:152:ALA:H	2.29	0.41
1:B:182:ALA:HB1	1:B:241:ILE:HD13	2.03	0.41
1:A:108:ARG:O	1:A:178:ASN:N	2.51	0.41
1:A:218:PRO:HA	4:A:412:HOH:O	2.20	0.41
1:B:170:SER:OG	1:B:195:ARG:HA	2.21	0.41
1:A:25:ASN:CG	1:A:51:PHE:CE1	2.99	0.40
1:A:52:LEU:HB2	1:A:67:LEU:HD23	2.04	0.40
1:A:13:PHE:CE1	1:A:141:VAL:HG11	2.56	0.40
1:A:25:ASN:O	1:A:26:ASN:C	2.64	0.40
1:A:177:GLY:O	1:A:213:TYR:HA	2.22	0.40
1:A:178:ASN:OD1	1:A:178:ASN:C	2.63	0.40
1:B:92:ASN:O	1:B:93:LEU:C	2.63	0.40
1:B:125:ARG:CD	1:B:159:LEU:HD22	2.52	0.40
1:B:147:LEU:O	1:B:147:LEU:HG	2.20	0.40
1:B:202:ILE:HA	1:B:253:PRO:HB2	2.04	0.40
1:A:98:LYS:HG2	1:A:99:GLU:N	2.37	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	267/269 (99%)	239 (90%)	27 (10%)	1 (0%)	30	59
1	B	267/269 (99%)	245 (92%)	19 (7%)	3 (1%)	11	36
All	All	534/538 (99%)	484 (91%)	46 (9%)	4 (1%)	18	48

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	199	THR
1	B	100	ILE
1	A	204	PRO
1	B	250	PRO

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	220/220 (100%)	193 (88%)	27 (12%)	4	15
1	B	220/220 (100%)	192 (87%)	28 (13%)	4	14
All	All	440/440 (100%)	385 (88%)	55 (12%)	4	15

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	LYS
1	A	26	ASN

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Mol	Chain	Res	Type
1	A	46	LYS
1	A	60	VAL
1	A	64	THR
1	A	79	SER
1	A	87	GLU
1	A	100	ILE
1	A	102	ASP
1	A	116	SER
1	A	123	THR
1	A	127	LYS
1	A	160	ARG
1	A	193	LEU
1	A	199	THR
1	A	200	ASN
1	A	202	ILE
1	A	204	PRO
1	A	214	SER
1	A	215	HIS
1	A	224	SER
1	A	233	ASN
1	A	239	GLU
1	A	241	ILE
1	A	249	GLN
1	A	252	ILE
1	A	255	ILE
1	B	1	GLU
1	B	25	ASN
1	B	26	ASN
1	B	54	SER
1	B	58	SER
1	B	60	VAL
1	B	100	ILE
1	B	103	ILE
1	B	108	ARG
1	B	134	GLU
1	B	147	LEU
1	B	151	LEU
1	B	162	ASN
1	B	166	ILE
1	B	178	ASN
1	B	183	GLU
1	B	185	LEU

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Mol	Chain	Res	Type
1	B	193	LEU
1	B	202	ILE
1	B	206	LEU
1	B	207	PRO
1	B	214	SER
1	B	219	GLU
1	B	235	ILE
1	B	239	GLU
1	B	241	ILE
1	B	252	ILE
1	B	256	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	25	ASN
1	A	88	ASN
1	B	25	ASN
1	B	26	ASN
1	B	71	ASN
1	B	88	ASN
1	B	94	ASN
1	B	135	HIS
1	B	188	GLN
1	B	249	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 oligosaccharides in this entry.

5.6 Ligand geometry

3 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	EAC	A	301	-	8,8,8	3.05	2 (25%)	9,9,9	1.99	4 (44%)
3	NAG	A	302	1	14,14,15	0.42	0	17,19,21	1.15	2 (11%)
3	NAG	B	301	1	14,14,15	0.43	0	17,19,21	1.16	2 (11%)

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	EAC	A	301	-	-	4/7/7/7	-
3	NAG	A	302	1	-	0/6/23/26	0/1/1/1
3	NAG	B	301	1	-	2/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	EAC	O5-C10	7.06	1.53	1.33
2	A	301	EAC	C15-C10	4.17	1.58	1.51

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	EAC	C9-O5-C10	3.94	127.85	116.67
2	A	301	EAC	C14-C15-C10	2.55	118.51	112.78
2	A	301	EAC	O5-C10-C15	2.50	118.90	111.38
3	A	302	NAG	C8-C7-N2	2.37	120.05	116.12
3	B	301	NAG	C8-C7-N2	2.36	120.03	116.12
3	B	301	NAG	C2-N2-C7	-2.13	120.04	122.90
3	A	302	NAG	C2-N2-C7	-2.10	120.08	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	EAC	O5-C9-C4	2.00	115.58	108.40

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	EAC	C13-C14-C15-C10
2	A	301	EAC	O8-C14-C15-C10
3	B	301	NAG	O5-C5-C6-O6
3	B	301	NAG	C4-C5-C6-O6
2	A	301	EAC	C15-C10-O5-C9
2	A	301	EAC	O6-C10-O5-C9

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	EAC	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	269/269 (100%)	-1.63	0 100 100	22, 48, 74, 88	1 (0%)
1	B	269/269 (100%)	-1.61	0 100 100	22, 52, 72, 92	1 (0%)
All	All	538/538 (100%)	-1.62	0 100 100	22, 50, 74, 92	2 (0%)

There are no RSRZ outliers to report.

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 oligosaccharides 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	EAC	A	301	9/9	0.98	0.08	50,51,54,58	0
3	NAG	A	302	14/15	0.98	0.04	41,43,46,47	0
3	NAG	B	301	14/15	0.99	0.04	40,42,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.