



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 08:26 PM UTC

PDB ID : 4S0X / pdb_00004s0x
Title : Structure of three phase partition - treated lipase from *Thermomyces lanuginosa* in complex with lauric acid at 2.1 Å resolution
Authors : Kumar, M.; Mukherjee, J.; Gupta, M.N.; Sinha, M.; Kaur, P.; Sharma, S.; Singh, T.P.
Deposited on : 2015-01-07
Resolution : 2.10 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

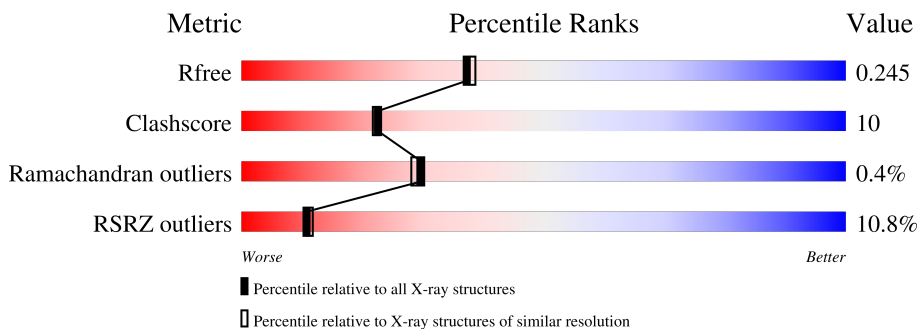
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	8% 77% 22%
1	B	269	14% 80% 20%

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
4	ACT	B	305	-	-	X	-
5	DAO	A	307	-	-	X	-

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 4520 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 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



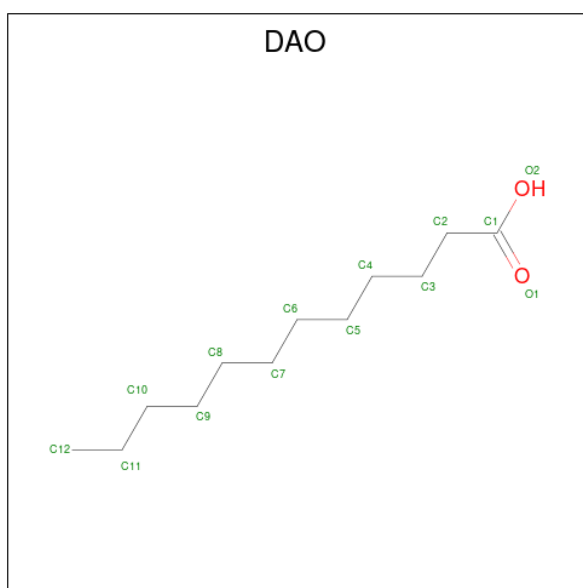
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



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

- Molecule 5 is LAURIC ACID (CCD ID: DAO) (formula: $C_{12}H_{24}O_2$).



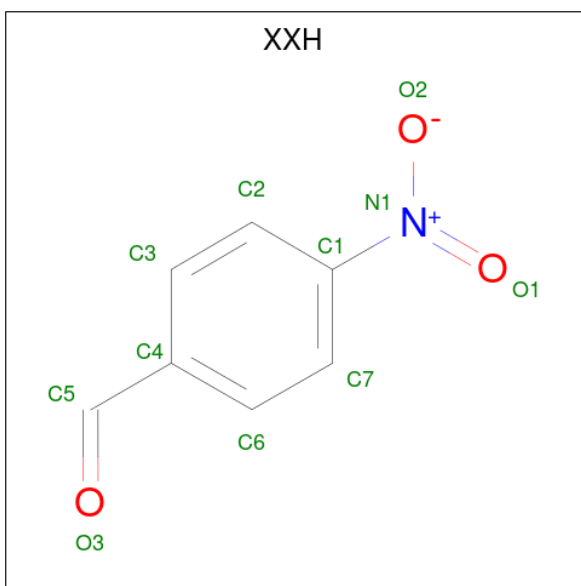
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			14	12	2		

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is 4-nitrobenzaldehyde (CCD ID: XXH) (formula: $C_7H_5NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	B	1	Total	C	N	O	0	0
			11	7	1	3		

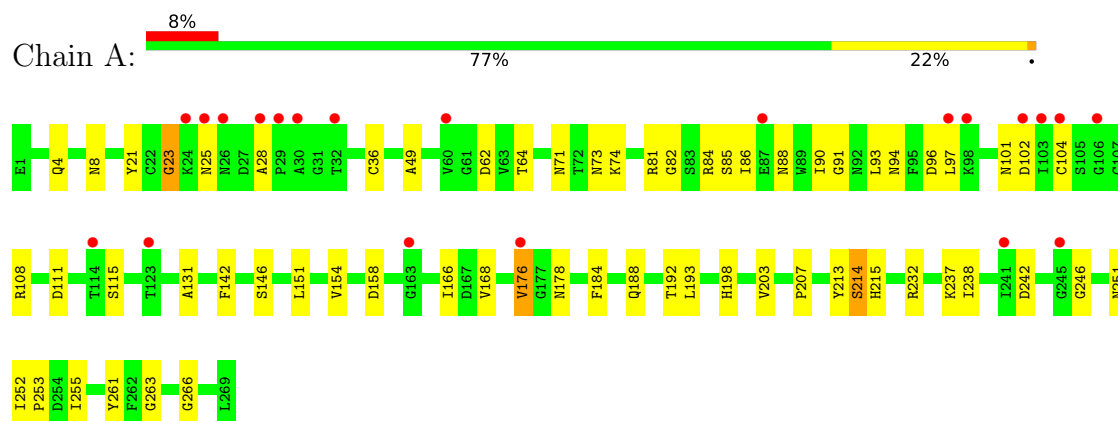
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	140	Total 140	O 140	0	0
8	B	139	Total 139	O 139	0	0

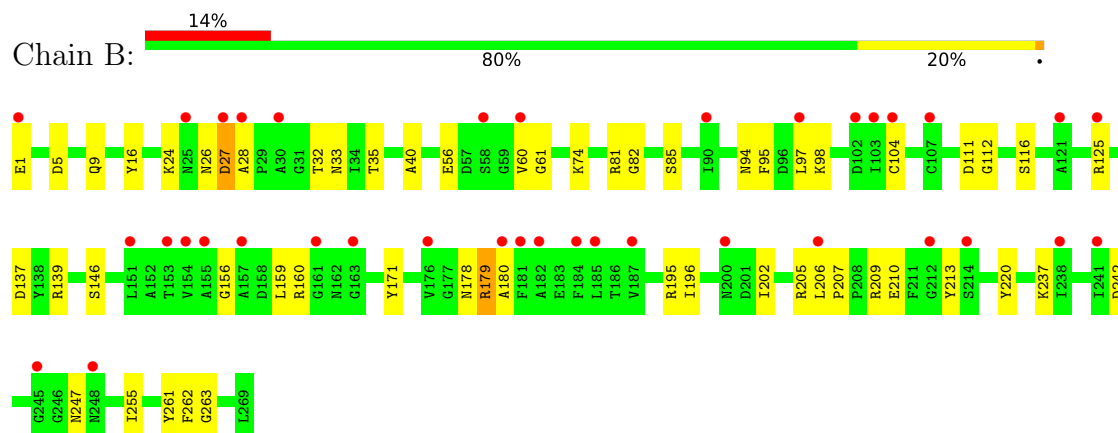
3 Residue-property plots

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	140.00Å 140.00Å 80.49Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.10 50.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	98.0 (50.00-2.10) 98.1 (50.00-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.61 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.209 , 0.260 (Not available) , 0.245	Depositor DCC
R_{free} test set	2643 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	32.1	Xtriage
Anisotropy	0.741	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 42.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.039 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4520	wwPDB-VP
Average B, all atoms (Å ²)	46.0	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: DAO, NAG, ACT, GOL, EDO, XXH

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	1.27	6/2120 (0.3%)	1.15	9/2887 (0.3%)
1	B	1.29	4/2120 (0.2%)	1.22	7/2887 (0.2%)
All	All	1.28	10/4240 (0.2%)	1.18	16/5774 (0.3%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	146	SER	C-O	-6.38	1.18	1.24
1	A	146	SER	C-O	-6.03	1.19	1.24
1	A	166	ILE	CA-C	-5.85	1.45	1.52
1	B	262	PHE	C-O	-5.72	1.16	1.24
1	A	214	SER	CA-C	5.68	1.59	1.52
1	A	71	ASN	CA-C	-5.35	1.45	1.52
1	B	255	ILE	CA-C	5.21	1.58	1.52
1	B	16	TYR	C-O	-5.13	1.18	1.24
1	A	198	HIS	C-O	-5.11	1.18	1.24
1	A	237	LYS	C-O	-5.02	1.18	1.24

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	179	ARG	N-CA-C	-11.53	96.23	110.41
1	B	202	ILE	N-CA-C	6.34	117.09	110.62
1	A	255	ILE	CA-C-N	-6.31	113.13	119.56
1	A	255	ILE	C-N-CA	-6.31	113.13	119.56
1	B	27	ASP	N-CA-C	-6.14	105.37	112.86
1	B	40	ALA	N-CA-C	5.89	118.65	111.82
1	A	176	VAL	CB-CA-C	5.85	116.94	110.62
1	A	207	PRO	CA-C-N	-5.80	113.99	119.85
1	A	207	PRO	C-N-CA	-5.80	113.99	119.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	95	PHE	N-CA-C	5.77	119.30	112.72
1	B	60	VAL	N-CA-C	5.76	116.54	110.72
1	A	158	ASP	N-CA-C	5.51	118.21	111.82
1	A	8	ASN	N-CA-C	5.34	117.10	111.28
1	A	131	ALA	N-CA-C	-5.19	105.31	111.69
1	B	195	ARG	NE-CZ-NH1	-5.07	116.43	121.50
1	A	49	ALA	N-CA-C	5.04	116.61	108.34

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	1964	47	0
1	B	2070	0	1964	40	0
2	A	14	0	13	0	0
2	B	14	0	13	0	0
3	A	24	0	32	2	0
3	B	12	0	16	0	0
4	A	4	0	3	1	0
4	B	4	0	3	2	0
5	A	14	0	23	8	0
6	B	4	0	6	0	0
7	B	11	0	4	1	0
8	A	140	0	0	3	0
8	B	139	0	0	6	0
All	All	4520	0	4041	87	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 10.

All (87) 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:23:GLY:HA3	1:A:36:CYS:HA	1.51	0.93
1:A:23:GLY:CA	1:A:36:CYS:HA	2.01	0.90
1:B:104:CYS:HB2	8:B:487:HOH:O	1.70	0.90
1:B:85:SER:HB2	8:B:525:HOH:O	1.78	0.84
1:A:252:ILE:HD12	1:A:253:PRO:HD2	1.63	0.78
1:B:26:ASN:HD21	1:B:56:GLU:HB2	1.48	0.77
1:B:26:ASN:ND2	1:B:56:GLU:HB2	2.01	0.76
1:A:238:ILE:HD12	1:A:246:GLY:HA3	1.72	0.72
1:A:97:LEU:HD11	1:A:213:TYR:CE1	2.25	0.72
1:A:193:LEU:H	3:A:302:GOL:H11	1.58	0.68
1:A:64:THR:H	5:A:307:DAO:H112	1.57	0.68
1:A:85:SER:HB2	8:A:536:HOH:O	1.94	0.66
1:B:24:LYS:HD2	1:B:35:THR:OG1	1.97	0.65
1:A:62:ASP:O	5:A:307:DAO:H82	1.96	0.64
1:A:90:ILE:HD11	1:A:203:VAL:HG22	1.79	0.64
1:B:210:GLU:HG3	8:B:467:HOH:O	1.99	0.63
1:A:64:THR:N	5:A:307:DAO:H112	2.13	0.63
1:B:26:ASN:O	1:B:27:ASP:HB3	1.96	0.63
1:A:252:ILE:HG22	8:A:469:HOH:O	1.99	0.61
1:A:86:ILE:O	1:A:90:ILE:HG12	2.02	0.60
1:B:210:GLU:HA	8:B:467:HOH:O	2.00	0.60
1:A:101:ASN:HA	1:A:104:CYS:O	2.04	0.58
1:A:4:GLN:HE21	1:A:232:ARG:HD3	1.69	0.58
1:A:111:ASP:O	1:A:115:SER:OG	2.22	0.57
1:A:215:HIS:HD2	8:A:508:HOH:O	1.86	0.57
1:A:23:GLY:HA2	1:A:36:CYS:HA	1.85	0.56
1:B:5:ASP:O	1:B:9:GLN:HG3	2.06	0.56
1:A:97:LEU:CD1	1:A:213:TYR:CE1	2.88	0.56
1:B:9:GLN:NE2	1:B:139:ARG:HH12	2.05	0.55
1:A:81:ARG:HD2	1:A:82:GLY:O	2.07	0.53
1:B:97:LEU:HD11	1:B:213:TYR:CE1	2.44	0.53
1:A:151:LEU:HD12	5:A:307:DAO:H123	1.92	0.52
1:A:90:ILE:CD1	1:A:203:VAL:HG22	2.39	0.51
1:B:179:ARG:O	1:B:180:ALA:HB3	2.09	0.51
1:A:192:THR:HA	3:A:302:GOL:H2	1.92	0.51
1:A:84:ARG:H	5:A:307:DAO:H52	1.75	0.51
1:A:154:VAL:HG23	1:A:176:VAL:HG21	1.93	0.51
1:A:238:ILE:CD1	1:A:246:GLY:CA	2.89	0.50
1:A:81:ARG:CD	1:A:82:GLY:O	2.60	0.50
1:A:238:ILE:HD12	1:A:246:GLY:CA	2.40	0.50
1:B:98:LYS:HB2	1:B:111:ASP:OD1	2.12	0.50
1:B:81:ARG:CD	1:B:82:GLY:O	2.60	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:81:ARG:HD2	1:B:82:GLY:O	2.11	0.50
1:A:4:GLN:NE2	1:A:232:ARG:HD3	2.28	0.49
1:A:25:ASN:HA	1:A:28:ALA:HB2	1.96	0.48
1:B:26:ASN:H	1:B:28:ALA:CB	2.27	0.48
1:A:151:LEU:CD1	5:A:307:DAO:C12	2.92	0.47
1:B:61:GLY:HA3	1:B:116:SER:OG	2.14	0.47
1:B:9:GLN:HE22	1:B:139:ARG:HH22	1.63	0.47
1:B:9:GLN:HG2	7:B:306:XXH:C2	2.45	0.47
1:A:94:ASN:HD21	1:A:96:ASP:CG	2.23	0.46
1:A:108:ARG:HG3	1:A:178:ASN:ND2	2.29	0.46
1:A:214:SER:OG	1:A:242:ASP:OD2	2.33	0.46
1:A:25:ASN:C	1:A:25:ASN:ND2	2.73	0.46
1:B:178:ASN:CG	1:B:179:ARG:O	2.59	0.46
1:B:137:ASP:HB3	8:B:409:HOH:O	2.17	0.44
1:A:25:ASN:HD22	1:A:25:ASN:H	1.65	0.44
1:A:151:LEU:HD12	5:A:307:DAO:C12	2.47	0.44
1:A:142:PHE:O	1:A:168:VAL:HA	2.18	0.44
1:B:178:ASN:C	1:B:179:ARG:O	2.46	0.44
1:B:220:TYR:CE1	1:B:237:LYS:HG3	2.52	0.43
1:B:112:GLY:HA3	4:B:305:ACT:H3	1.99	0.43
1:A:93:LEU:HD12	4:A:306:ACT:H2	2.00	0.43
1:A:261:TYR:C	1:A:263:GLY:H	2.27	0.43
1:B:32:THR:HG22	1:B:33:ASN:O	2.18	0.43
1:A:101:ASN:O	1:A:102:ASP:C	2.61	0.43
1:B:1:GLU:OE1	1:B:1:GLU:C	2.62	0.42
1:B:74:LYS:NZ	8:B:488:HOH:O	2.53	0.42
1:A:88:ASN:O	1:A:91:GLY:N	2.53	0.42
1:A:90:ILE:HD11	1:A:203:VAL:CG2	2.48	0.42
1:B:178:ASN:OD1	1:B:179:ARG:O	2.37	0.42
1:B:97:LEU:HD12	1:B:97:LEU:N	2.35	0.42
1:B:206:LEU:HA	1:B:207:PRO:C	2.45	0.42
1:A:64:THR:H	5:A:307:DAO:C11	2.30	0.41
1:A:73:ASN:O	1:A:74:LYS:HB2	2.20	0.41
1:B:26:ASN:H	1:B:28:ALA:HB2	1.86	0.41
1:B:26:ASN:C	1:B:28:ALA:N	2.72	0.41
1:B:125:ARG:HD3	1:B:159:LEU:HD22	2.02	0.41
1:B:156:GLY:O	1:B:160:ARG:HG3	2.20	0.41
1:B:261:TYR:C	1:B:263:GLY:H	2.28	0.41
1:A:21:TYR:CZ	1:A:266:GLY:HA2	2.56	0.41
1:B:171:TYR:CE2	1:B:196:ILE:HD13	2.55	0.41
1:B:205:ARG:HD2	1:B:247:ASN:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:209:ARG:NH1	1:B:242:ASP:HB3	2.35	0.41
1:B:94:ASN:HB3	4:B:305:ACT:O	2.19	0.41
1:B:179:ARG:O	1:B:180:ALA:CB	2.69	0.40
1:A:184:PHE:O	1:A:188:GLN:HB2	2.21	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	267/269 (99%)	254 (95%)	11 (4%)	2 (1%)	18	15
1	B	267/269 (99%)	259 (97%)	8 (3%)	0	100	100
All	All	534/538 (99%)	513 (96%)	19 (4%)	2 (0%)	30	28

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	23	GLY
1	A	251	ASN

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

13 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$
3	GOL	B	303	-	5,5,5	0.80	0	5,5,5	0.67	0
3	GOL	A	304	-	5,5,5	0.98	0	5,5,5	1.05	0
4	ACT	B	305	-	3,3,3	1.30	1 (33%)	3,3,3	0.36	0
3	GOL	A	302	-	5,5,5	1.06	0	5,5,5	0.65	0
3	GOL	B	302	-	5,5,5	0.89	0	5,5,5	0.94	0
2	NAG	A	301	1	14,14,15	0.83	0	17,19,21	2.51	8 (47%)
4	ACT	A	306	-	3,3,3	1.10	0	3,3,3	0.55	0
6	EDO	B	304	-	3,3,3	0.84	0	2,2,2	0.42	0
5	DAO	A	307	-	13,13,13	0.75	0	13,13,13	0.72	0
7	XXH	B	306	-	11,11,11	3.41	4 (36%)	12,14,14	5.18	7 (58%)
3	GOL	A	303	-	5,5,5	1.19	0	5,5,5	1.14	0
2	NAG	B	301	1	14,14,15	0.85	0	17,19,21	2.31	8 (47%)
3	GOL	A	305	-	5,5,5	0.29	0	5,5,5	1.15	0

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
3	GOL	B	303	-	-	0/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	A	304	-	-	2/4/4/4	-
3	GOL	A	302	-	-	2/4/4/4	-
3	GOL	B	302	-	-	0/4/4/4	-
2	NAG	A	301	1	-	2/6/23/26	0/1/1/1
6	EDO	B	304	-	-	1/1/1/1	-
5	DAO	A	307	-	-	9/11/11/11	-
7	XXH	B	306	-	-	0/4/6/6	0/1/1/1
3	GOL	A	303	-	-	2/4/4/4	-
2	NAG	B	301	1	-	2/6/23/26	0/1/1/1
3	GOL	A	305	-	-	3/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	306	XXH	O1-N1	7.49	1.35	1.22
7	B	306	XXH	C4-C5	-7.26	1.26	1.47
7	B	306	XXH	C2-C3	2.69	1.43	1.38
4	B	305	ACT	OXT-C	-2.24	1.20	1.30
7	B	306	XXH	C1-N1	-2.13	1.40	1.45

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	306	XXH	O3-C5-C4	13.30	169.64	124.56
7	B	306	XXH	C2-C1-N1	7.14	125.57	119.34
7	B	306	XXH	C7-C1-N1	-5.38	114.64	119.34
7	B	306	XXH	C6-C4-C5	-5.29	106.59	120.33
2	A	301	NAG	O5-C1-C2	-5.24	103.18	111.29
7	B	306	XXH	C3-C4-C5	4.42	131.83	120.33
2	B	301	NAG	O3-C3-C2	-4.27	100.54	109.40
2	A	301	NAG	O5-C5-C6	4.05	115.55	107.66
2	A	301	NAG	O5-C5-C4	-4.01	101.08	110.83
2	B	301	NAG	C4-C3-C2	3.82	116.62	111.02
2	B	301	NAG	C1-C2-N2	-3.26	105.30	110.43
2	A	301	NAG	O4-C4-C5	-3.24	101.33	109.32
2	A	301	NAG	O3-C3-C2	-3.23	102.70	109.40
2	B	301	NAG	C2-N2-C7	3.04	126.97	122.90
2	B	301	NAG	O5-C1-C2	-2.90	106.80	111.29
2	B	301	NAG	O5-C5-C4	-2.59	104.53	110.83
7	B	306	XXH	C3-C4-C6	2.49	121.35	117.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	NAG	O5-C5-C6	2.49	112.50	107.66
2	A	301	NAG	O3-C3-C4	2.35	115.92	110.38
2	A	301	NAG	C1-O5-C5	2.35	115.34	112.19
2	B	301	NAG	C3-C4-C5	-2.27	106.12	110.23
2	A	301	NAG	C2-N2-C7	-2.23	119.91	122.90
7	B	306	XXH	C2-C3-C4	-2.02	118.61	121.22

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	302	GOL	C1-C2-C3-O3
3	A	303	GOL	C1-C2-C3-O3
3	A	304	GOL	C1-C2-C3-O3
2	B	301	NAG	O5-C5-C6-O6
2	A	301	NAG	O5-C5-C6-O6
2	A	301	NAG	C4-C5-C6-O6
2	B	301	NAG	C4-C5-C6-O6
5	A	307	DAO	C1-C2-C3-C4
3	A	302	GOL	O2-C2-C3-O3
3	A	305	GOL	C1-C2-C3-O3
3	A	304	GOL	O2-C2-C3-O3
5	A	307	DAO	C5-C6-C7-C8
5	A	307	DAO	C2-C3-C4-C5
5	A	307	DAO	C11-C10-C9-C8
3	A	303	GOL	O2-C2-C3-O3
3	A	305	GOL	O2-C2-C3-O3
5	A	307	DAO	C3-C4-C5-C6
5	A	307	DAO	C7-C8-C9-C10
5	A	307	DAO	C6-C7-C8-C9
6	B	304	EDO	O1-C1-C2-O2
5	A	307	DAO	O2-C1-C2-C3
3	A	305	GOL	O1-C1-C2-O2
5	A	307	DAO	O1-C1-C2-C3

There are no ring outliers.

5 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	305	ACT	2	0
3	A	302	GOL	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	306	ACT	1	0
5	A	307	DAO	8	0
7	B	306	XXH	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

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, 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%)	0.62	21 (7%)	19 20	25, 42, 68, 120	1 (0%)
1	B	269/269 (100%)	0.88	37 (13%)	6 6	25, 45, 72, 116	1 (0%)
All	All	538/538 (100%)	0.75	58 (10%)	11 11	25, 44, 71, 120	2 (0%)

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	176	VAL	5.3
1	B	180	ALA	4.1
1	A	106	GLY	4.0
1	A	60	VAL	3.8
1	A	103	ILE	3.8
1	B	155	ALA	3.4
1	A	26	ASN	3.2
1	A	176	VAL	3.2
1	B	27	ASP	3.1
1	A	102	ASP	2.8
1	A	24	LYS	2.8
1	B	184	PHE	2.8
1	B	25	ASN	2.8
1	B	161	GLY	2.8
1	B	245	GLY	2.8
1	A	241	ILE	2.8
1	A	97	LEU	2.8
1	B	206	LEU	2.8
1	B	154	VAL	2.7
1	B	181	PHE	2.7
1	B	60	VAL	2.7
1	B	125	ARG	2.6
1	B	241	ILE	2.6
1	A	25	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	163	GLY	2.5
1	A	114	THR	2.5
1	A	29	PRO	2.5
1	B	97	LEU	2.5
1	B	163	GLY	2.5
1	B	187	VAL	2.4
1	B	151	LEU	2.4
1	B	90	ILE	2.4
1	B	102	ASP	2.4
1	B	107	CYS	2.4
1	B	28	ALA	2.4
1	B	185	LEU	2.4
1	B	103	ILE	2.4
1	B	121	ALA	2.3
1	B	238	ILE	2.3
1	B	30	ALA	2.3
1	B	153	THR	2.3
1	A	32	THR	2.2
1	A	28	ALA	2.2
1	A	30	ALA	2.2
1	B	214	SER	2.1
1	A	98	LYS	2.1
1	A	245	GLY	2.1
1	B	157	ALA	2.1
1	B	182	ALA	2.1
1	B	104	CYS	2.1
1	B	248	ASN	2.1
1	B	212	GLY	2.1
1	B	1	GLU	2.1
1	A	87	GLU	2.0
1	B	200	ASN	2.0
1	A	104	CYS	2.0
1	B	58	SER	2.0
1	A	123	THR	2.0

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 [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($\text{\AA}^2$)	Q<0.9
2	NAG	B	301	14/15	0.77	0.13	54,57,72,73	0
3	GOL	B	302	6/6	0.79	0.17	48,55,60,60	0
5	DAO	A	307	14/14	0.79	0.23	36,43,56,59	14
2	NAG	A	301	14/15	0.80	0.13	48,57,63,66	0
4	ACT	B	305	4/4	0.86	0.16	47,49,51,58	0
3	GOL	A	303	6/6	0.86	0.14	43,56,58,60	0
3	GOL	B	303	6/6	0.87	0.12	50,56,59,65	0
6	EDO	B	304	4/4	0.87	0.16	41,45,45,46	0
3	GOL	A	302	6/6	0.88	0.12	42,45,54,59	0
3	GOL	A	305	6/6	0.89	0.16	43,46,49,53	0
3	GOL	A	304	6/6	0.92	0.13	33,45,49,52	0
7	XXH	B	306	11/11	0.92	0.12	30,34,42,51	11
4	ACT	A	306	4/4	0.94	0.12	52,52,54,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.