Full wwPDB X-ray Structure Validation Report (i)

Jun 18, 2024 – 04:03 PM EDT

PDB ID	:	4KJX
Title	:	Crystal Structure of the complex of three phase partition treated lipase from Thermomyces lanuginosa with Lauric acid and P-nitrobenzaldehyde (PNB) at 2.1 resolution
Authors	:	Kumar, M.; Mukherjee, J.; Gupta, M.N.; Sinha, M.; Kaur, P.; Sharma, S.; Singh, T.P.
Deposited on	:	2013-05-04
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 (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

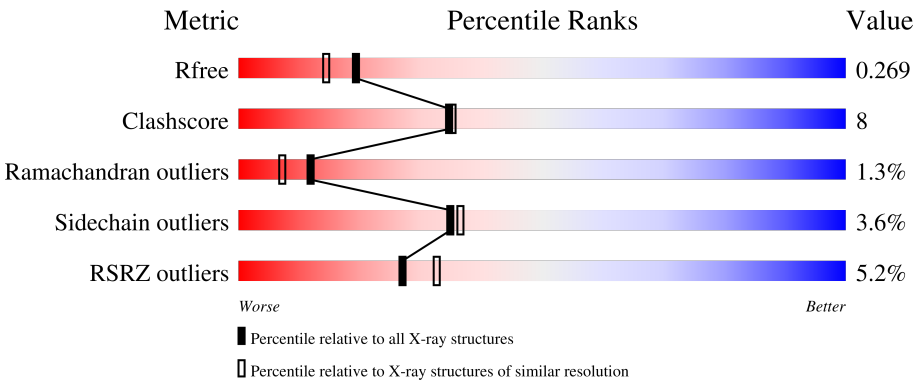
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA,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

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 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}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (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 <=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	4% 82% 16% •
1	B	269	7% 83% 14% •

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
5	DAO	A	307	-	-	X	-

2 Entry composition (i)

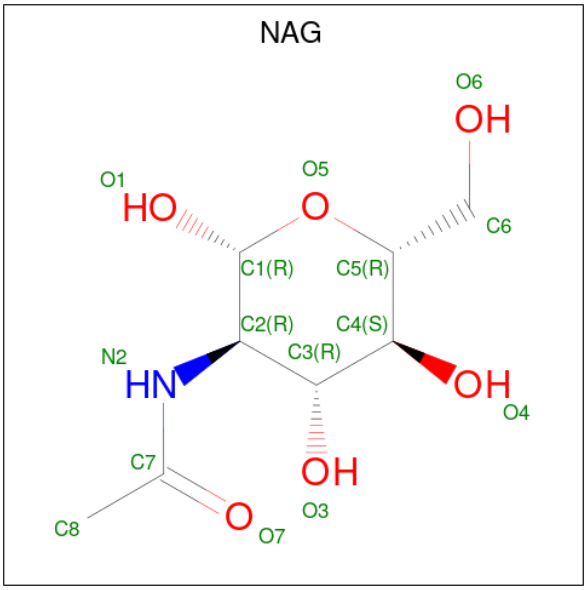
There are 8 unique types of molecules in this entry. The entry contains 4520 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

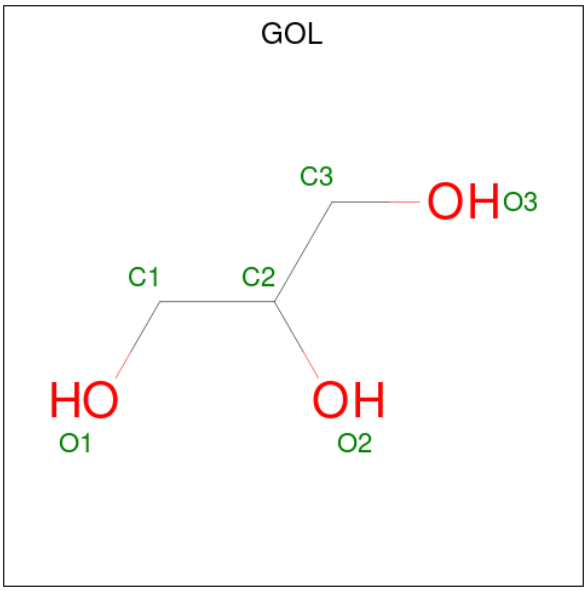
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	269	2070	1303	359	402	6	0	0	0
1	B	269	2070	1303	359	402	6	0	0	0

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



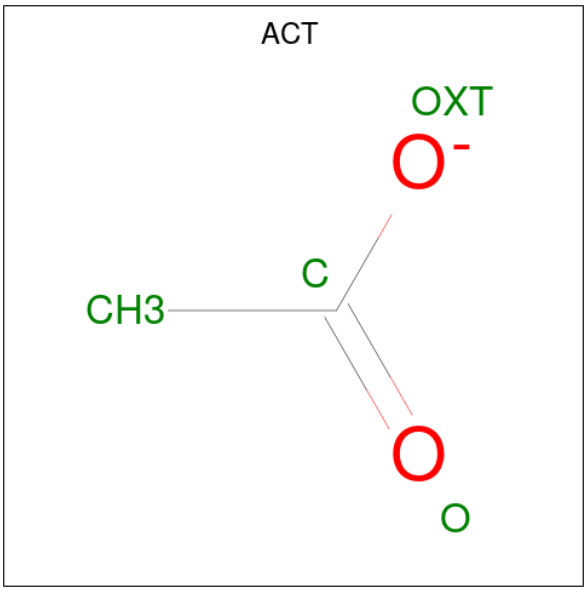
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	14	8	1	5	0	0
2	B	1	14	8	1	5	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



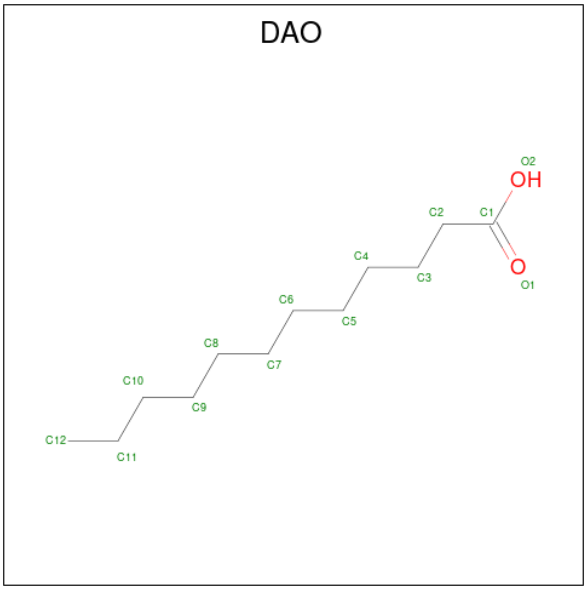
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	6	3	3	0	0
			Total	C	O		
3	A	1	6	3	3	0	0
			Total	C	O		
3	A	1	6	3	3	0	0
			Total	C	O		
3	A	1	6	3	3	0	0
			Total	C	O		
3	B	1	6	3	3	0	0
			Total	C	O		
3	B	1	6	3	3	0	0

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



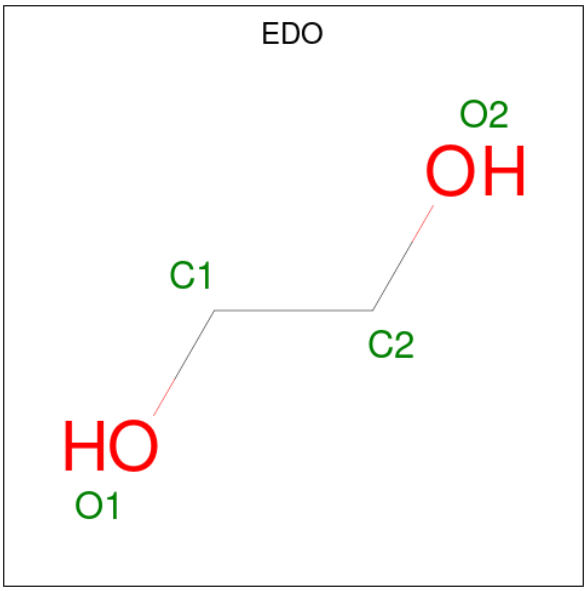
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
4	A	1	4	2	2	0	0

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



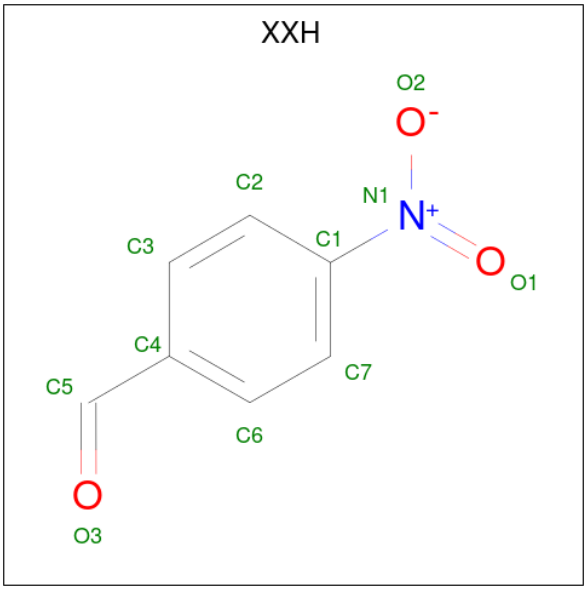
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
5	A	1	14	12	2	0	0
			Total	C	O		
5	B	1	14	12	2	0	0

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

- Molecule 7 is 4-nitrobenzaldehyde (three-letter code: XXH) (formula: $C_7H_5NO_3$).



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

- Molecule 8 is water.

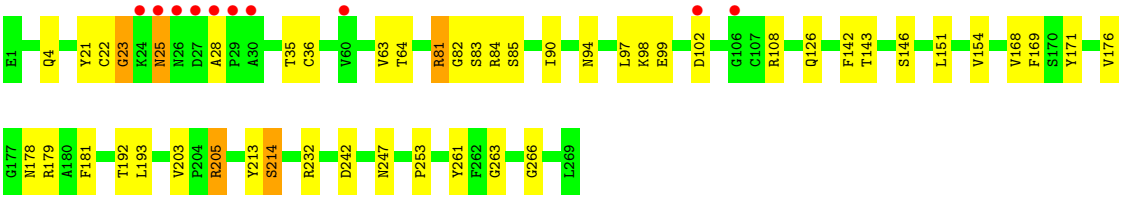
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
8	A	135	135	135	0	0
			Total	O		
8	B	134	134	134	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

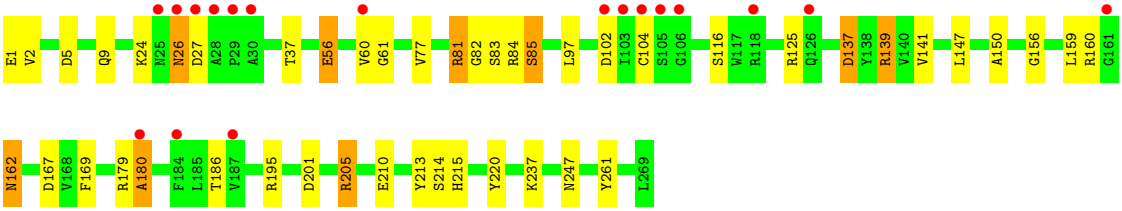
- Molecule 1: Lipase

Chain A:



- Molecule 1: Lipase

Chain B:



4 Data and refinement statistics (i)

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)	45.83 – 2.10 45.82 – 2.10	Depositor EDS
% Data completeness (in resolution range)	98.0 (45.83-2.10) 98.0 (45.82-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	2.61 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.214 , 0.265 0.217 , 0.269	Depositor DCC
R_{free} test set	2642 reflections (5.11%)	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 , 46.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	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)	41.0	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.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, DAO, NAG, XXH, EDO, ACT

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	1.06	1/2120 (0.0%)	1.04	3/2887 (0.1%)
1	B	1.12	3/2120 (0.1%)	1.10	9/2887 (0.3%)
All	All	1.09	4/4240 (0.1%)	1.07	12/5774 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	137	ASP	CB-CG	5.79	1.64	1.51
1	A	171	TYR	CE1-CZ	-5.40	1.31	1.38
1	B	261	TYR	CB-CG	-5.33	1.43	1.51
1	B	169	PHE	CG-CD2	-5.20	1.30	1.38

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	205	ARG	NE-CZ-NH1	10.56	125.58	120.30
1	A	205	ARG	NE-CZ-NH1	8.80	124.70	120.30
1	B	205	ARG	NE-CZ-NH2	-7.59	116.51	120.30
1	B	195	ARG	NE-CZ-NH1	-7.34	116.63	120.30
1	A	81	ARG	NE-CZ-NH2	-7.17	116.72	120.30
1	B	137	ASP	CB-CG-OD1	6.93	124.54	118.30
1	B	139	ARG	NE-CZ-NH1	6.73	123.67	120.30
1	A	81	ARG	NE-CZ-NH1	6.04	123.32	120.30
1	B	139	ARG	NE-CZ-NH2	-5.79	117.40	120.30
1	B	201	ASP	CB-CG-OD2	5.63	123.37	118.30
1	B	81	ARG	NE-CZ-NH2	-5.56	117.52	120.30
1	B	167	ASP	CB-CG-OD2	5.02	122.82	118.30

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	2070	0	1964	35	0
1	B	2070	0	1964	32	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
5	A	14	0	23	10	0
5	B	14	0	23	2	0
6	B	4	0	6	0	0
7	B	11	0	5	1	0
8	A	135	0	0	4	0
8	B	134	0	0	5	0
All	All	4520	0	4062	69	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 8.

All (69) 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:26:ASN:O	1:B:27:ASP:HB3	1.61	1.00
1:A:90:ILE:HD11	1:A:203:VAL:HG22	1.51	0.92
1:A:23:GLY:CA	1:A:36:CYS:HA	2.04	0.88
1:A:23:GLY:HA3	1:A:36:CYS:HA	1.57	0.86
1:B:26:ASN:HD21	1:B:56:GLU:HB2	1.41	0.85
1:B:104:CYS:HB2	8:B:489:HOH:O	1.81	0.78
1:A:97:LEU:HD11	1:A:213:TYR:CE1	2.20	0.76
1:A:64:THR:H	5:A:307:DAO:H112	1.52	0.74
1:B:137:ASP:HB3	8:B:409:HOH:O	1.88	0.73
1:B:210:GLU:HG3	8:B:469:HOH:O	1.90	0.71
1:B:26:ASN:ND2	1:B:56:GLU:HB2	2.05	0.71
1:B:210:GLU:HA	8:B:469:HOH:O	1.89	0.71
1:B:9:GLN:NE2	1:B:139:ARG:HH12	1.90	0.70
1:B:179:ARG:O	1:B:180:ALA:HB3	1.93	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:307:DAO:H41	5:A:307:DAO:O2	1.95	0.67
1:A:83:SER:HA	5:A:307:DAO:H71	1.76	0.65
4:A:306:ACT:H1	8:A:457:HOH:O	1.95	0.65
1:A:64:THR:N	5:A:307:DAO:H112	2.13	0.63
1:A:85:SER:HB2	8:A:471:HOH:O	1.99	0.62
1:A:192:THR:HA	3:A:302:GOL:H2	1.83	0.61
1:A:4:GLN:HE21	1:A:232:ARG:HD3	1.66	0.61
1:B:179:ARG:O	1:B:180:ALA:CB	2.53	0.57
1:A:4:GLN:NE2	1:A:232:ARG:HD3	2.20	0.57
1:B:9:GLN:HE22	1:B:139:ARG:HH22	1.54	0.56
1:A:151:LEU:CD1	5:A:307:DAO:C12	2.84	0.55
1:A:25:ASN:HA	1:A:28:ALA:HB2	1.87	0.55
1:A:97:LEU:CD1	1:A:213:TYR:CE1	2.90	0.55
1:B:26:ASN:O	1:B:27:ASP:CB	2.41	0.54
1:A:154:VAL:HG23	1:A:176:VAL:HG21	1.88	0.54
1:B:5:ASP:O	1:B:9:GLN:HG3	2.08	0.54
1:B:83:SER:OG	5:B:306:DAO:H111	2.07	0.54
1:B:125:ARG:HD3	1:B:159:LEU:HD22	1.90	0.54
1:A:81:ARG:CD	1:A:82:GLY:O	2.57	0.52
1:A:81:ARG:HD2	1:A:82:GLY:O	2.09	0.52
1:B:97:LEU:HD11	1:B:213:TYR:CE1	2.44	0.52
1:A:151:LEU:CD1	5:A:307:DAO:H122	2.40	0.52
1:A:22:CYS:O	1:A:23:GLY:O	2.28	0.51
1:A:151:LEU:HD12	5:A:307:DAO:C12	2.40	0.51
1:A:97:LEU:HD13	8:A:531:HOH:O	2.09	0.51
1:B:162:ASN:ND2	1:B:162:ASN:O	2.44	0.51
1:A:64:THR:H	5:A:307:DAO:C11	2.21	0.51
1:A:179:ARG:NH1	8:A:472:HOH:O	2.44	0.50
1:B:220:TYR:CE1	1:B:237:LYS:HG3	2.46	0.50
1:A:205:ARG:HD2	1:A:247:ASN:O	2.11	0.50
1:A:193:LEU:H	3:A:302:GOL:H11	1.75	0.50
1:B:81:ARG:CD	1:B:82:GLY:O	2.60	0.49
1:B:156:GLY:O	1:B:160:ARG:HG3	2.12	0.49
1:A:214:SER:OG	1:A:242:ASP:OD2	2.31	0.49
1:A:151:LEU:HD12	5:A:307:DAO:H123	1.94	0.49
1:A:23:GLY:HA2	1:A:36:CYS:HA	1.92	0.48
1:B:9:GLN:HG2	7:B:305:XXH:C2	2.43	0.48
1:B:84:ARG:HB2	5:B:306:DAO:H121	1.95	0.48
1:B:61:GLY:HA3	1:B:116:SER:OG	2.13	0.47
1:B:186:THR:HG22	1:B:186:THR:O	2.15	0.47
1:B:85:SER:HB3	8:B:528:HOH:O	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:77:VAL:HA	1:B:141:VAL:O	2.16	0.45
1:A:142:PHE:O	1:A:168:VAL:HA	2.16	0.45
1:A:63:VAL:HA	5:A:307:DAO:H112	1.98	0.44
1:B:205:ARG:HD2	1:B:247:ASN:O	2.18	0.44
1:A:261:TYR:C	1:A:263:GLY:H	2.21	0.43
1:B:97:LEU:HD12	1:B:97:LEU:N	2.34	0.43
1:B:24:LYS:HG2	1:B:37:THR:HG23	2.00	0.42
1:A:23:GLY:HA3	1:A:35:THR:O	2.20	0.42
1:A:108:ARG:HG3	1:A:178:ASN:ND2	2.33	0.42
1:B:81:ARG:HD2	1:B:82:GLY:O	2.19	0.42
1:B:27:ASP:O	1:B:27:ASP:CG	2.59	0.40
1:B:147:LEU:O	1:B:150:ALA:HB3	2.20	0.40
1:A:143:THR:HA	1:A:169:PHE:O	2.21	0.40
1:A:21:TYR:CZ	1:A:266:GLY:HA2	2.57	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%)	251 (94%)	12 (4%)	4 (2%)	10 5
1	B	267/269 (99%)	251 (94%)	13 (5%)	3 (1%)	14 9
All	All	534/538 (99%)	502 (94%)	25 (5%)	7 (1%)	12 7

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	23	GLY
1	A	102	ASP
1	B	26	ASN

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Mol	Chain	Res	Type
1	B	102	ASP
1	A	146	SER
1	B	180	ALA

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	220/220 (100%)	212 (96%)	8 (4%)	35	36
1	B	220/220 (100%)	212 (96%)	8 (4%)	35	36
All	All	440/440 (100%)	424 (96%)	16 (4%)	35	36

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	ARG
1	A	94	ASN
1	A	98	LYS
1	A	99	GLU
1	A	126	GLN
1	A	181	PHE
1	A	214	SER
1	A	253	PRO
1	B	1	GLU
1	B	2	VAL
1	B	56	GLU
1	B	60	VAL
1	B	85	SER
1	B	162	ASN
1	B	214	SER
1	B	215	HIS

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	4	GLN
1	A	88	ASN
1	A	249	GLN
1	B	4	GLN
1	B	9	GLN
1	B	26	ASN
1	B	88	ASN
1	B	92	ASN
1	B	200	ASN
1	B	233	ASN
1	B	248	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

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	$\#\|Z\| > 2$
4	ACT	A	306	-	3,3,3	0.72	0	3,3,3	1.36	0
3	GOL	B	302	-	5,5,5	0.52	0	5,5,5	0.47	0
6	EDO	B	303	-	3,3,3	0.74	0	2,2,2	0.23	0
3	GOL	A	303	-	5,5,5	0.97	0	5,5,5	0.99	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	GOL	A	302	-	5,5,5	0.80	0	5,5,5	0.81	0
3	GOL	B	304	-	5,5,5	0.53	0	5,5,5	0.68	0
2	NAG	B	301	1	14,14,15	0.75	0	17,19,21	1.54	4 (23%)
5	DAO	B	306	-	13,13,13	0.96	1 (7%)	13,13,13	1.66	1 (7%)
3	GOL	A	304	-	5,5,5	0.63	0	5,5,5	0.70	0
5	DAO	A	307	-	13,13,13	0.78	1 (7%)	13,13,13	1.48	3 (23%)
7	XXH	B	305	-	11,11,11	1.26	1 (9%)	12,14,14	2.28	5 (41%)
3	GOL	A	305	-	5,5,5	0.18	0	5,5,5	0.77	0
2	NAG	A	301	1	14,14,15	0.67	0	17,19,21	1.62	3 (17%)

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	302	-	-	0/4/4/4	-
6	EDO	B	303	-	-	1/1/1/1	-
3	GOL	A	303	-	-	2/4/4/4	-
3	GOL	A	302	-	-	2/4/4/4	-
3	GOL	B	304	-	-	0/4/4/4	-
2	NAG	B	301	1	-	2/6/23/26	0/1/1/1
5	DAO	B	306	-	-	8/11/11/11	-
3	GOL	A	304	-	-	2/4/4/4	-
5	DAO	A	307	-	-	8/11/11/11	-
7	XXH	B	305	-	-	0/4/6/6	0/1/1/1
3	GOL	A	305	-	-	2/4/4/4	-
2	NAG	A	301	1	-	2/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	305	XXH	C2-C3	2.65	1.43	1.38
5	B	306	DAO	O2-C1	-2.33	1.23	1.30
5	A	307	DAO	O1-C1	2.04	1.28	1.22

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	B	306	DAO	C3-C2-C1	-5.62	99.84	114.51
7	B	305	XXH	C2-C1-N1	5.06	123.75	119.34
2	A	301	NAG	C1-O5-C5	3.51	116.89	112.19
2	A	301	NAG	O5-C5-C6	3.43	114.35	107.66
2	B	301	NAG	C2-N2-C7	-3.36	118.40	122.90
7	B	305	XXH	C6-C7-C1	3.02	124.08	120.08
2	B	301	NAG	O5-C5-C4	-3.02	103.48	110.83
7	B	305	XXH	C7-C6-C4	-2.92	117.44	121.22
5	A	307	DAO	O2-C1-C2	2.89	123.13	114.00
2	A	301	NAG	O5-C5-C4	-2.63	104.43	110.83
2	B	301	NAG	O5-C1-C2	-2.38	107.62	111.29
5	A	307	DAO	O1-C1-C2	-2.24	115.99	123.09
2	B	301	NAG	O5-C5-C6	2.14	111.82	107.66
5	A	307	DAO	C6-C5-C4	-2.09	103.80	114.37
7	B	305	XXH	O3-C5-C4	-2.06	117.58	124.56
7	B	305	XXH	C3-C4-C6	2.01	120.63	117.65

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	303	GOL	C1-C2-C3-O3
3	A	304	GOL	C1-C2-C3-O3
3	A	305	GOL	C1-C2-C3-O3
2	A	301	NAG	O5-C5-C6-O6
2	B	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
5	B	306	DAO	C5-C6-C7-C8
3	A	302	GOL	C1-C2-C3-O3
3	A	302	GOL	O2-C2-C3-O3
3	A	303	GOL	O2-C2-C3-O3
3	A	304	GOL	O2-C2-C3-O3
3	A	305	GOL	O2-C2-C3-O3
5	B	306	DAO	C2-C3-C4-C5
5	B	306	DAO	C9-C10-C11-C12
5	B	306	DAO	C3-C4-C5-C6
5	A	307	DAO	C4-C5-C6-C7
5	B	306	DAO	C7-C8-C9-C10
5	A	307	DAO	C3-C4-C5-C6
5	A	307	DAO	C7-C8-C9-C10
5	A	307	DAO	C11-C10-C9-C8

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Mol	Chain	Res	Type	Atoms
5	B	306	DAO	C6-C7-C8-C9
6	B	303	EDO	O1-C1-C2-O2
5	A	307	DAO	C2-C3-C4-C5
5	A	307	DAO	O1-C1-C2-C3
5	A	307	DAO	O2-C1-C2-C3
5	B	306	DAO	O1-C1-C2-C3
5	B	306	DAO	O2-C1-C2-C3

There are no ring outliers.

5 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	306	ACT	1	0
3	A	302	GOL	2	0
5	B	306	DAO	2	0
5	A	307	DAO	10	0
7	B	305	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 (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%)	0.12	10 (3%)	41	48	21, 37, 61, 118	1 (0%)
1	B	269/269 (100%)	0.17	18 (6%)	17	22	20, 40, 66, 113	1 (0%)
All	All	538/538 (100%)	0.15	28 (5%)	27	32	20, 38, 64, 118	2 (0%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	26	ASN	6.0
1	A	25	ASN	5.0
1	B	27	ASP	4.6
1	B	26	ASN	4.6
1	A	60	VAL	3.8
1	A	106	GLY	3.8
1	A	27	ASP	3.3
1	B	28	ALA	3.1
1	B	184	PHE	3.0
1	B	102	ASP	2.8
1	B	103	ILE	2.8
1	B	105	SER	2.8
1	B	60	VAL	2.7
1	B	29	PRO	2.6
1	B	25	ASN	2.6
1	A	24	LYS	2.5
1	B	106	GLY	2.4
1	B	180	ALA	2.4
1	A	29	PRO	2.3
1	B	187	VAL	2.3
1	B	104	CYS	2.3
1	A	102	ASP	2.2
1	B	161	GLY	2.2
1	B	30	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	28	ALA	2.2
1	A	30	ALA	2.2
1	B	126	GLN	2.0
1	B	118	ARG	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	GOL	B	302	6/6	0.68	0.27	58,68,70,73	0
2	NAG	B	301	14/15	0.77	0.32	61,69,72,72	0
5	DAO	B	306	14/14	0.81	0.25	33,42,46,53	0
5	DAO	A	307	14/14	0.82	0.33	30,36,49,54	0
3	GOL	A	303	6/6	0.85	0.15	44,63,66,67	0
3	GOL	A	302	6/6	0.85	0.18	47,48,60,69	0
6	EDO	B	303	4/4	0.85	0.17	41,41,42,45	0
3	GOL	B	304	6/6	0.86	0.18	53,61,67,78	0
2	NAG	A	301	14/15	0.87	0.30	58,71,76,82	0
3	GOL	A	304	6/6	0.91	0.16	32,53,58,64	0
3	GOL	A	305	6/6	0.92	0.18	49,55,63,65	0
7	XXH	B	305	11/11	0.95	0.11	34,40,54,56	0
4	ACT	A	306	4/4	0.96	0.09	47,47,52,62	0

6.5 Other polymers (i)

There are no such residues in this entry.