



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:44 PM GMT

PDB ID : 1GZ7
Title : CRYSTAL STRUCTURE OF THE CLOSED STATE OF LIPASE 2 FROM CANDIDA RUGOSA
Authors : Mancheno, J.M.; Hermoso, J.A.
Deposited on : 2002-05-17
Resolution : 1.97 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

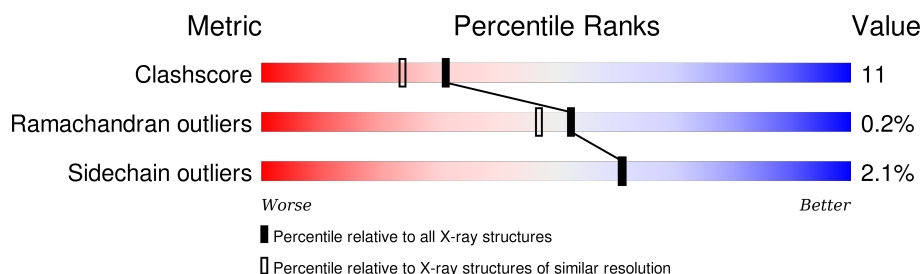
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.97 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	9905 (2.00-1.96)
Ramachandran outliers	100387	9792 (2.00-1.96)
Sidechain outliers	100360	9791 (2.00-1.96)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	534	 81% 18% •
1	B	534	 79% 19% •
1	C	534	 79% 20% •
1	D	534	 78% 21% •

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	NAG	B	1535	X	-	-	-
3	GOL	B	1537	-	-	X	-
3	GOL	C	1537	-	-	X	-
3	GOL	C	1538	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 17843 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 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	534	Total	C	N	O	S	0	0	0
			4053	2585	668	781	19			
1	B	534	Total	C	N	O	S	0	0	0
			4053	2585	668	781	19			
1	C	534	Total	C	N	O	S	0	0	0
			4053	2585	668	781	19			
1	D	534	Total	C	N	O	S	0	0	0
			4053	2585	668	781	19			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	78	LEU	ARG	CONFLICT	UNP P32946
A	79	ASP	HIS	CONFLICT	UNP P32946
B	78	LEU	ARG	CONFLICT	UNP P32946
B	79	ASP	HIS	CONFLICT	UNP P32946
C	78	LEU	ARG	CONFLICT	UNP P32946
C	79	ASP	HIS	CONFLICT	UNP P32946
D	78	LEU	ARG	CONFLICT	UNP P32946
D	79	ASP	HIS	CONFLICT	UNP P32946

- Molecule 2 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	2	Total	C	N	O	0	0
			28	16	2	10		
2	B	2	Total	C	N	O	0	0
			28	16	2	10		
2	C	2	Total	C	N	O	0	0
			28	16	2	10		
2	D	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 3 is GLYCEROL (three-letter code: 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	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	462	Total	O	0	0
			462	462		
4	B	352	Total	O	0	0
			352	352		
4	C	312	Total	O	0	0
			312	312		

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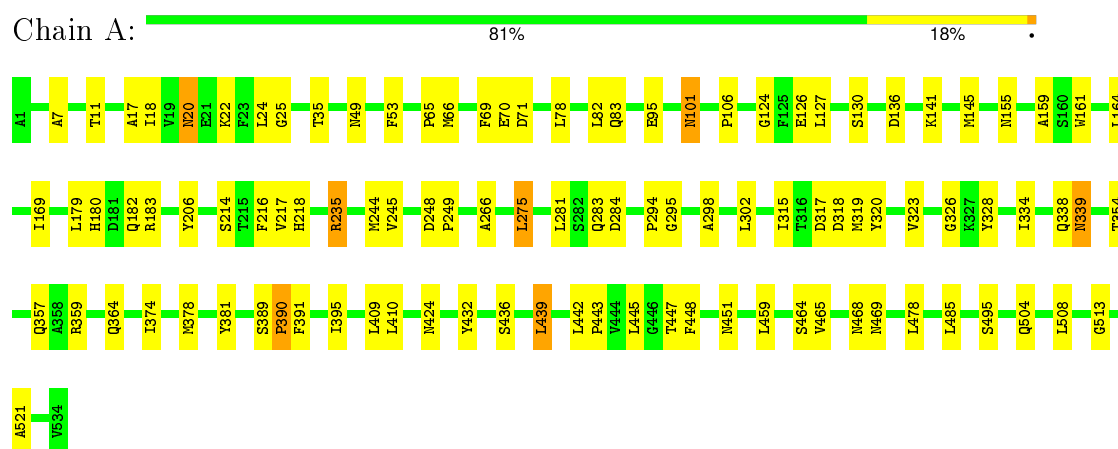
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	345	Total	O	0	0
			345	345		

3 Residue-property plots

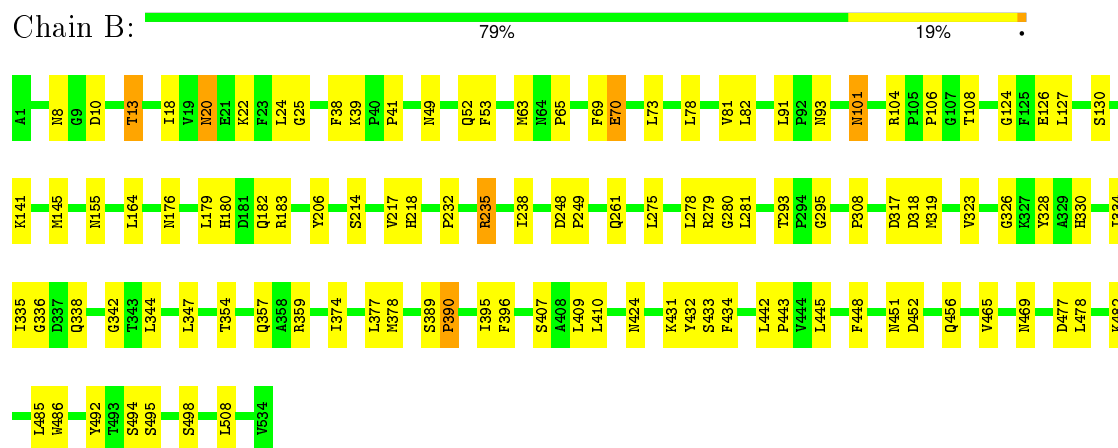
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

Note EDS was not executed.

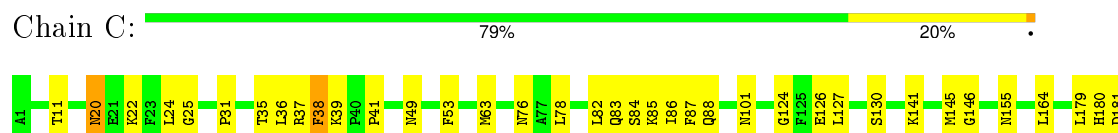
• Molecule 1: LIPASE 2

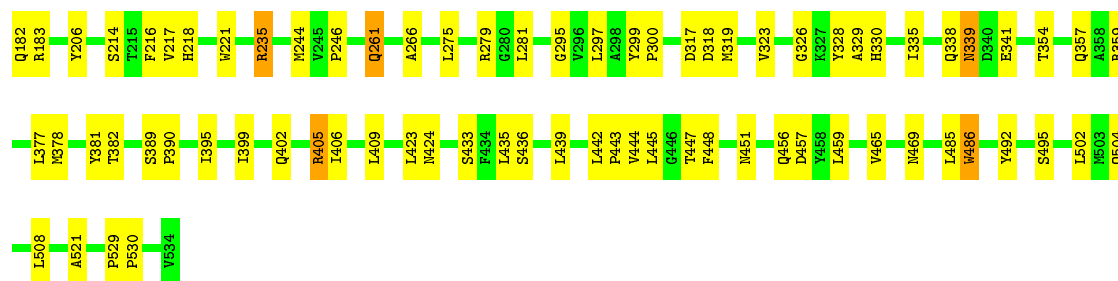


• Molecule 1: LIPASE 2



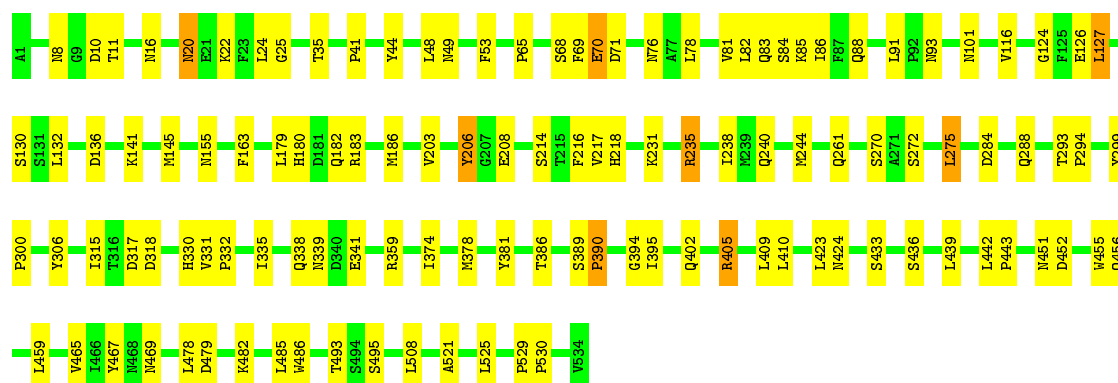
• Molecule 1: LIPASE 2





• Molecule 1: LIPASE 2

Chain D: 78% 21% .



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	61.15Å 91.14Å 108.46Å 90.78° 106.31° 86.91°	Depositor
Resolution (Å)	11.98 – 1.97	Depositor
% Data completeness (in resolution range)	94.4 (11.98-1.97)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.200 , 0.236	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	17843	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

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, 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	0.36	0/4159	0.63	1/5668 (0.0%)
1	B	0.35	0/4159	0.62	0/5668
1	C	0.35	1/4159 (0.0%)	0.63	2/5668 (0.0%)
1	D	0.34	0/4159	0.62	1/5668 (0.0%)
All	All	0.35	1/16636 (0.0%)	0.63	4/22672 (0.0%)

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	C	0	1
2	B	1	0
All	All	1	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	39	LYS	N-CA	5.60	1.57	1.46

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	39	LYS	N-CA-C	-7.31	91.27	111.00
1	C	39	LYS	N-CA-CB	5.72	120.90	110.60
1	A	315	ILE	N-CA-C	-5.41	96.41	111.00
1	D	315	ILE	N-CA-C	-5.04	97.39	111.00

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	1535	NAG	C1

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	38	PHE	Peptide

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	4053	0	3913	74	0
1	B	4053	0	3912	85	1
1	C	4053	0	3913	93	1
1	D	4053	0	3913	97	1
2	A	28	0	25	0	1
2	B	28	0	25	2	0
2	C	28	0	25	2	0
2	D	28	0	25	0	0
3	A	12	0	16	1	0
3	B	12	0	16	5	0
3	C	12	0	16	9	0
3	D	12	0	16	2	0
4	A	462	0	0	5	1
4	B	352	0	0	6	1
4	C	312	0	0	6	2
4	D	345	0	0	10	4
All	All	17843	0	15815	347	8

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 11.

All (347) 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:C:146:GLY:HA2	4:C:2099:HOH:O	1.59	1.02
1:A:359:ARG:HH11	1:A:378:MET:HE1	1.26	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:11:THR:H	1:D:49:ASN:HD22	1.11	0.95
1:C:20:ASN:HD21	1:C:22:LYS:HE3	1.31	0.93
1:D:180:HIS:ND1	3:D:1538:GOL:H12	1.82	0.93
1:A:66:MET:CE	1:A:295:GLY:H	1.83	0.92
1:A:66:MET:HE2	1:A:295:GLY:H	1.38	0.88
1:C:180:HIS:ND1	3:C:1538:GOL:H32	1.88	0.87
1:B:182:GLN:HE22	1:B:214:SER:HB3	1.43	0.84
1:A:20:ASN:HD21	1:A:22:LYS:HE3	1.42	0.84
1:C:183:ARG:HG2	3:C:1538:GOL:H11	1.60	0.82
1:B:498:SER:OG	4:B:2318:HOH:O	1.63	0.82
1:C:146:GLY:O	4:C:2098:HOH:O	1.97	0.82
1:B:424:ASN:HD21	1:B:495:SER:H	1.23	0.81
1:A:424:ASN:HD21	1:A:495:SER:H	1.25	0.81
1:D:11:THR:H	1:D:49:ASN:ND2	1.77	0.81
1:D:424:ASN:HD21	1:D:495:SER:H	1.26	0.80
1:D:182:GLN:HE22	1:D:214:SER:HB3	1.47	0.79
1:D:331:VAL:HG12	3:D:1537:GOL:H11	1.64	0.79
1:D:359:ARG:HH11	1:D:378:MET:HE1	1.48	0.78
1:B:13:THR:CG2	1:B:52:GLN:HG2	2.14	0.78
1:A:182:GLN:HE22	1:A:214:SER:HB3	1.47	0.78
1:B:359:ARG:HH11	1:B:378:MET:CE	1.97	0.78
1:D:359:ARG:HH11	1:D:378:MET:CE	1.97	0.77
1:C:182:GLN:HE22	1:C:214:SER:HB3	1.49	0.76
1:A:338:GLN:HE21	1:A:451:ASN:HD21	1.33	0.76
1:A:359:ARG:HH11	1:A:378:MET:CE	1.99	0.73
1:A:465:VAL:O	1:A:469:ASN:HB2	1.88	0.73
1:A:66:MET:HE1	1:A:294:PRO:HA	1.72	0.72
1:B:280:GLY:HA2	4:B:2182:HOH:O	1.90	0.72
1:B:180:HIS:ND1	3:B:1538:GOL:H12	2.05	0.72
1:D:130:SER:H	1:D:155:ASN:ND2	1.88	0.72
1:C:359:ARG:HH11	1:C:378:MET:CE	2.03	0.71
1:D:272:SER:HA	4:D:2177:HOH:O	1.89	0.71
1:C:11:THR:H	1:C:49:ASN:HD22	1.38	0.71
1:B:330:HIS:H	3:B:1537:GOL:H32	1.55	0.71
1:D:183:ARG:HH11	1:D:218:HIS:HD1	1.38	0.71
1:C:130:SER:H	1:C:155:ASN:ND2	1.89	0.71
1:B:330:HIS:H	3:B:1537:GOL:C3	2.03	0.71
1:D:70:GLU:H	1:D:70:GLU:CD	1.94	0.70
1:B:141:LYS:O	1:B:145:MET:HG2	1.91	0.70
1:C:424:ASN:HD21	1:C:495:SER:H	1.38	0.69
1:D:402:GLN:OE1	1:D:405:ARG:HD2	1.94	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:83:GLN:HE22	1:D:459:LEU:HD11	1.59	0.67
1:C:359:ARG:HG3	1:C:378:MET:HE1	1.76	0.67
1:B:338:GLN:HE21	1:B:451:ASN:HD21	1.41	0.67
1:C:465:VAL:O	1:C:469:ASN:HB2	1.95	0.67
1:D:130:SER:H	1:D:155:ASN:HD21	1.41	0.67
1:A:20:ASN:HD21	1:A:22:LYS:CE	2.07	0.67
1:C:330:HIS:H	3:C:1537:GOL:C3	2.08	0.67
1:D:493:THR:HG23	4:D:2312:HOH:O	1.95	0.66
2:B:1535:NAG:H61	2:B:1536:NAG:H82	1.77	0.66
1:A:25:GLY:H	1:A:101:ASN:HD22	1.44	0.66
1:C:261:GLN:HG3	4:C:2170:HOH:O	1.94	0.66
1:B:359:ARG:HH11	1:B:378:MET:HE1	1.60	0.65
1:A:130:SER:H	1:A:155:ASN:ND2	1.94	0.65
1:D:25:GLY:H	1:D:101:ASN:ND2	1.93	0.65
1:B:25:GLY:H	1:B:101:ASN:HD22	1.43	0.65
1:D:91:LEU:HD23	1:D:93:ASN:ND2	2.11	0.65
1:A:7:ALA:O	4:A:2005:HOH:O	2.13	0.65
1:B:179:LEU:HD13	1:B:217:VAL:HG11	1.79	0.65
1:D:25:GLY:H	1:D:101:ASN:HD22	1.43	0.64
1:C:486:TRP:CE3	1:C:486:TRP:HA	2.32	0.64
1:C:24:LEU:O	1:C:53:PHE:HB3	1.98	0.63
1:A:183:ARG:HH11	1:A:218:HIS:HD1	1.47	0.62
1:B:13:THR:HG21	1:B:52:GLN:HG2	1.81	0.62
1:A:395:ILE:O	4:A:2362:HOH:O	2.15	0.62
1:D:84:SER:OG	1:D:86:ILE:HG12	2.00	0.62
1:A:179:LEU:HD13	1:A:217:VAL:HG11	1.81	0.62
1:B:124:GLY:O	1:B:126:GLU:HG3	2.01	0.61
1:B:65:PRO:HG3	1:B:127:LEU:CD2	2.30	0.61
1:C:124:GLY:O	1:C:126:GLU:HG3	1.99	0.61
1:A:338:GLN:NE2	1:A:451:ASN:HD21	1.99	0.61
1:C:299:TYR:HB3	2:C:1536:NAG:H81	1.83	0.61
1:C:486:TRP:HA	1:C:486:TRP:HE3	1.66	0.60
1:A:66:MET:HE1	1:A:295:GLY:H	1.65	0.60
1:C:216:PHE:CZ	1:C:244:MET:HG3	2.36	0.60
1:D:465:VAL:O	1:D:469:ASN:HB2	2.00	0.60
1:D:65:PRO:HG3	1:D:127:LEU:HD21	1.81	0.60
1:C:508:LEU:CG	1:D:70:GLU:HG3	2.32	0.60
1:D:83:GLN:NE2	1:D:459:LEU:HD11	2.15	0.60
1:B:442:LEU:HD12	1:B:443:PRO:HD2	1.84	0.60
1:B:20:ASN:HD22	1:B:20:ASN:C	2.05	0.60
1:C:330:HIS:H	3:C:1537:GOL:H32	1.65	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:76:ASN:HD21	1:D:443:PRO:HD2	1.66	0.60
1:C:465:VAL:CG1	1:C:485:LEU:HD22	2.32	0.59
1:C:354:THR:OG1	1:C:357:GLN:HG3	2.03	0.59
1:B:465:VAL:O	1:B:469:ASN:HB2	2.02	0.59
1:D:235:ARG:HD2	4:D:2135:HOH:O	2.02	0.59
1:D:338:GLN:HE21	1:D:451:ASN:HD21	1.49	0.59
1:B:63:MET:HE2	1:B:126:GLU:HB3	1.83	0.59
1:C:359:ARG:HH11	1:C:378:MET:HE1	1.67	0.59
1:A:11:THR:H	1:A:49:ASN:ND2	2.01	0.58
1:D:235:ARG:HD3	1:D:478:LEU:HD21	1.84	0.58
1:D:124:GLY:O	1:D:126:GLU:HG3	2.03	0.58
1:C:399:ILE:HD11	1:C:444:VAL:HG13	1.87	0.57
1:B:13:THR:HG23	1:B:52:GLN:HG2	1.86	0.57
1:C:141:LYS:O	1:C:145:MET:HG2	2.05	0.57
1:B:130:SER:H	1:B:155:ASN:ND2	2.03	0.57
1:B:338:GLN:NE2	1:B:451:ASN:HD21	2.02	0.57
1:A:124:GLY:O	1:A:126:GLU:HG3	2.05	0.57
1:D:76:ASN:HD21	1:D:443:PRO:CD	2.17	0.56
1:C:326:GLY:HA2	1:C:328:TYR:CE2	2.41	0.56
1:A:20:ASN:HA	1:A:106:PRO:HD3	1.87	0.56
1:C:508:LEU:HG	1:D:70:GLU:HG3	1.87	0.56
1:A:65:PRO:HG3	1:A:127:LEU:CD2	2.36	0.56
1:C:20:ASN:HD22	1:C:20:ASN:C	2.09	0.56
1:A:83:GLN:HE22	1:A:459:LEU:HD11	1.69	0.56
1:D:20:ASN:HD22	1:D:20:ASN:C	2.09	0.56
1:B:359:ARG:HH11	1:B:378:MET:HE2	1.70	0.56
1:A:465:VAL:CG1	1:A:485:LEU:HD22	2.36	0.56
1:B:330:HIS:N	3:B:1537:GOL:H32	2.20	0.56
1:B:183:ARG:HH11	1:B:218:HIS:HD1	1.52	0.56
1:A:235:ARG:HD2	4:A:2400:HOH:O	2.05	0.56
1:B:25:GLY:H	1:B:101:ASN:ND2	2.05	0.55
1:D:389:SER:HA	1:D:390:PRO:C	2.26	0.55
1:C:179:LEU:HD13	1:C:217:VAL:HG11	1.88	0.55
1:B:261:GLN:OE1	1:B:293:THR:HG22	2.06	0.55
1:D:270:SER:HA	4:D:2176:HOH:O	2.06	0.55
1:A:248:ASP:HB3	1:A:249:PRO:HD2	1.88	0.55
1:B:38:PHE:CE1	1:B:164:LEU:HD22	2.42	0.55
1:D:442:LEU:HD12	1:D:443:PRO:HD2	1.89	0.54
1:D:24:LEU:O	1:D:53:PHE:HB3	2.08	0.54
1:D:272:SER:CA	4:D:2177:HOH:O	2.52	0.54
1:C:338:GLN:HE21	1:C:451:ASN:HD21	1.55	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:266:ALA:HB1	1:A:281:LEU:HD11	1.90	0.54
1:D:65:PRO:HG3	1:D:127:LEU:CD2	2.37	0.54
1:B:20:ASN:HA	1:B:106:PRO:HD3	1.90	0.54
1:C:38:PHE:CE1	1:C:164:LEU:HD22	2.42	0.53
1:D:68:SER:HB2	1:D:71:ASP:OD2	2.09	0.53
1:D:486:TRP:CE3	1:D:486:TRP:HA	2.43	0.53
1:A:354:THR:OG1	1:A:357:GLN:HG3	2.09	0.53
1:C:183:ARG:CG	3:C:1538:GOL:H11	2.36	0.53
1:D:20:ASN:HD21	1:D:22:LYS:HE3	1.72	0.53
1:B:70:GLU:CD	1:B:70:GLU:H	2.12	0.53
1:B:20:ASN:HD21	1:B:22:LYS:HE3	1.73	0.53
1:A:25:GLY:H	1:A:101:ASN:ND2	2.05	0.53
1:A:141:LYS:O	1:A:145:MET:HG2	2.09	0.53
1:C:78:LEU:O	1:C:82:LEU:HG	2.08	0.53
1:C:330:HIS:H	3:C:1537:GOL:H31	1.74	0.52
1:B:24:LEU:HD23	1:B:101:ASN:HB3	1.91	0.52
1:A:127:LEU:C	1:A:127:LEU:HD12	2.30	0.52
1:C:235:ARG:HD2	4:C:2135:HOH:O	2.10	0.52
1:A:130:SER:H	1:A:155:ASN:HD21	1.57	0.52
1:A:11:THR:H	1:A:49:ASN:HD22	1.56	0.52
1:D:386:THR:HA	1:D:394:GLY:O	2.10	0.52
1:D:395:ILE:C	1:D:395:ILE:HD12	2.29	0.52
1:D:338:GLN:NE2	1:D:451:ASN:HD21	2.06	0.52
1:A:180:HIS:ND1	3:A:1538:GOL:H32	2.25	0.52
1:D:11:THR:N	1:D:49:ASN:HD22	1.95	0.51
1:B:508:LEU:HD22	1:B:508:LEU:N	2.25	0.51
1:B:465:VAL:CG1	1:B:485:LEU:HD22	2.39	0.51
1:B:41:PRO:HD2	1:B:180:HIS:HB3	1.92	0.51
1:A:445:LEU:HB3	1:A:448:PHE:HB3	1.92	0.51
1:A:66:MET:CE	1:A:298:ALA:HB2	2.41	0.51
1:B:78:LEU:O	1:B:81:VAL:HG22	2.10	0.51
1:B:359:ARG:NH1	1:B:378:MET:HE2	2.26	0.51
1:C:130:SER:H	1:C:155:ASN:HD21	1.58	0.51
1:D:335:ILE:O	1:D:433:SER:HA	2.10	0.51
1:B:395:ILE:HD12	1:B:395:ILE:C	2.31	0.51
1:A:66:MET:HE2	1:A:295:GLY:N	2.18	0.50
1:B:63:MET:CE	1:B:295:GLY:HA2	2.41	0.50
1:D:76:ASN:ND2	1:D:443:PRO:HD2	2.26	0.50
1:D:338:GLN:NE2	1:D:436:SER:HB3	2.27	0.50
1:C:409:LEU:HD23	1:C:409:LEU:C	2.31	0.50
1:A:439:LEU:HB2	1:A:447:THR:HA	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:MET:HE3	1:A:298:ALA:HB2	1.92	0.50
1:C:508:LEU:N	1:C:508:LEU:HD22	2.26	0.50
1:D:85:LYS:HA	1:D:88:GLN:NE2	2.26	0.50
1:C:389:SER:HA	1:C:390:PRO:C	2.31	0.50
1:D:409:LEU:C	1:D:409:LEU:HD23	2.32	0.50
1:C:338:GLN:NE2	1:C:436:SER:HB3	2.26	0.49
1:D:141:LYS:O	1:D:145:MET:HG2	2.12	0.49
1:B:389:SER:HA	1:B:390:PRO:C	2.32	0.49
2:C:1535:NAG:H61	2:C:1536:NAG:H82	1.94	0.49
1:B:232:PRO:HB3	3:B:1537:GOL:H11	1.95	0.49
1:D:455:TRP:HA	1:D:459:LEU:O	2.12	0.49
1:B:73:LEU:HD21	1:B:347:LEU:HD11	1.95	0.49
1:C:37:ARG:HB2	1:C:279:ARG:HG2	1.95	0.49
1:A:317:ASP:O	1:A:318:ASP:C	2.50	0.49
1:C:319:MET:O	1:C:323:VAL:HG23	2.13	0.49
1:C:299:TYR:N	1:C:300:PRO:HD2	2.28	0.49
1:A:216:PHE:CZ	1:A:244:MET:HG3	2.47	0.49
1:D:132:LEU:HD12	4:D:2081:HOH:O	2.13	0.48
1:B:465:VAL:HG13	1:B:485:LEU:HD22	1.95	0.48
1:C:63:MET:CE	1:C:295:GLY:HA2	2.44	0.48
1:D:76:ASN:HD21	1:D:443:PRO:CG	2.26	0.48
1:B:395:ILE:HG13	4:B:2272:HOH:O	2.12	0.48
1:B:354:THR:OG1	1:B:357:GLN:HG3	2.14	0.48
1:B:10:ASP:HA	1:B:49:ASN:HD22	1.79	0.48
1:C:183:ARG:HH11	1:C:218:HIS:HD1	1.62	0.48
1:B:8:ASN:ND2	1:B:10:ASP:OD2	2.44	0.48
1:B:359:ARG:NH1	1:B:378:MET:CE	2.73	0.47
1:D:299:TYR:N	1:D:300:PRO:HD2	2.29	0.47
1:A:374:ILE:O	1:A:378:MET:HG3	2.14	0.47
1:B:63:MET:HE3	1:B:295:GLY:HA2	1.95	0.47
1:D:395:ILE:HG13	4:D:2271:HOH:O	2.14	0.47
1:C:127:LEU:C	1:C:127:LEU:HD12	2.35	0.47
1:A:326:GLY:HA2	1:A:328:TYR:CE2	2.49	0.47
1:A:24:LEU:O	1:A:53:PHE:HB3	2.14	0.47
1:A:508:LEU:HD22	1:A:508:LEU:N	2.30	0.47
1:C:87:PHE:CD1	1:C:297:LEU:HD21	2.49	0.47
1:D:208:GLU:HA	1:D:240:GLN:O	2.14	0.47
1:C:359:ARG:NH1	1:C:378:MET:CE	2.77	0.47
1:D:317:ASP:CG	1:D:318:ASP:N	2.68	0.47
1:A:465:VAL:HG13	1:A:485:LEU:HD22	1.97	0.47
1:B:176:ASN:OD1	1:B:308:PRO:HA	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:221:TRP:HZ2	3:C:1538:GOL:H12	1.79	0.46
1:D:529:PRO:HB2	1:D:530:PRO:HD3	1.96	0.46
1:C:25:GLY:H	1:C:101:ASN:HD22	1.63	0.46
1:B:104:ARG:HD2	1:B:108:THR:HB	1.98	0.46
1:D:359:ARG:NH1	1:D:378:MET:HE2	2.30	0.46
1:B:275:LEU:HB2	4:B:2181:HOH:O	2.15	0.46
1:A:389:SER:HA	1:A:390:PRO:C	2.36	0.46
1:D:479:ASP:HB3	1:D:482:LYS:HG2	1.97	0.46
1:C:85:LYS:HA	1:C:88:GLN:NE2	2.31	0.46
1:A:235:ARG:HD3	1:A:478:LEU:HD21	1.97	0.46
1:C:338:GLN:NE2	1:C:451:ASN:HD21	2.13	0.46
1:D:465:VAL:CG1	1:D:485:LEU:HD22	2.45	0.46
1:D:216:PHE:CZ	1:D:244:MET:HG3	2.51	0.46
1:C:35:THR:HG23	1:C:36:LEU:HG	1.96	0.46
1:D:275:LEU:HA	1:D:275:LEU:HD12	1.85	0.46
1:C:335:ILE:O	1:C:433:SER:HA	2.15	0.46
1:A:20:ASN:C	1:A:20:ASN:HD22	2.19	0.46
1:B:477:ASP:HB2	1:B:482:LYS:HG3	1.98	0.46
1:A:294:PRO:HG2	4:A:2239:HOH:O	2.14	0.46
1:D:317:ASP:CG	1:D:318:ASP:H	2.19	0.46
1:C:317:ASP:CG	1:C:318:ASP:H	2.19	0.46
1:C:317:ASP:CG	1:C:318:ASP:N	2.69	0.45
1:B:248:ASP:HB3	1:B:249:PRO:HD2	1.97	0.45
1:B:486:TRP:CE3	1:B:486:TRP:HA	2.50	0.45
1:A:95:GLU:OE2	1:A:283:GLN:HG3	2.16	0.45
1:C:266:ALA:HB1	1:C:281:LEU:HD11	1.97	0.45
1:B:18:ILE:HD12	1:B:18:ILE:N	2.32	0.45
1:C:423:LEU:HB3	1:C:492:TYR:CE1	2.51	0.45
1:B:326:GLY:HA2	1:B:328:TYR:CE2	2.51	0.45
1:B:261:GLN:HG3	4:B:2178:HOH:O	2.16	0.45
1:B:395:ILE:HD12	1:B:396:PHE:N	2.32	0.45
1:B:69:PHE:CZ	1:B:344:LEU:HG	2.51	0.45
1:C:329:ALA:HB1	3:C:1537:GOL:H11	1.99	0.45
1:C:339:ASN:HB2	1:C:436:SER:O	2.17	0.45
1:C:338:GLN:O	1:C:341:GLU:HG2	2.16	0.45
1:C:435:LEU:HB3	1:C:504:GLN:HG2	1.99	0.45
1:D:41:PRO:HD2	1:D:180:HIS:HB3	1.99	0.44
1:C:395:ILE:HD12	1:C:395:ILE:C	2.38	0.44
1:A:359:ARG:NH1	1:A:378:MET:CE	2.74	0.44
1:B:24:LEU:O	1:B:53:PHE:HB3	2.18	0.44
1:D:78:LEU:O	1:D:82:LEU:HG	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:442:LEU:HD12	1:C:443:PRO:HD2	2.00	0.44
1:D:374:ILE:O	1:D:378:MET:HG3	2.18	0.44
1:C:246:PRO:HD3	4:C:2202:HOH:O	2.16	0.44
1:D:294:PRO:HG2	4:D:2194:HOH:O	2.18	0.44
1:A:334:ILE:HG12	1:A:432:TYR:HB2	1.99	0.44
1:A:164:LEU:HD23	1:A:169:ILE:HD11	2.00	0.44
1:C:402:GLN:O	1:C:406:ILE:HG13	2.18	0.44
1:B:334:ILE:HG12	1:B:432:TYR:HB2	1.99	0.44
1:B:275:LEU:HA	1:B:275:LEU:HD12	1.84	0.43
1:B:39:LYS:HE3	1:B:279:ARG:CZ	2.48	0.43
1:C:508:LEU:HD21	1:D:70:GLU:HG3	2.00	0.43
1:A:339:ASN:HB2	1:A:436:SER:O	2.18	0.43
1:B:319:MET:O	1:B:323:VAL:HG23	2.18	0.43
1:B:127:LEU:HD12	1:B:127:LEU:C	2.38	0.43
1:A:245:VAL:HB	1:A:319:MET:CE	2.49	0.43
1:D:179:LEU:HD13	1:D:217:VAL:HG11	1.99	0.43
1:C:83:GLN:HE22	1:C:459:LEU:HD11	1.82	0.43
1:A:35:THR:HG21	1:D:35:THR:HB	2.01	0.43
1:A:442:LEU:HD12	1:A:443:PRO:HD2	1.99	0.43
1:D:261:GLN:OE1	1:D:293:THR:HG22	2.18	0.43
1:B:206:TYR:CB	1:B:238:ILE:HB	2.49	0.43
1:D:91:LEU:HD23	1:D:93:ASN:HD22	1.84	0.43
1:C:377:LEU:HD21	1:C:409:LEU:HD11	2.00	0.43
1:C:382:THR:C	1:C:405:ARG:HH21	2.22	0.43
1:A:381:TYR:CZ	1:A:521:ALA:HB1	2.54	0.43
1:C:41:PRO:HD2	1:C:180:HIS:HB3	2.00	0.43
1:B:13:THR:HG23	1:B:52:GLN:HA	2.01	0.43
1:C:76:ASN:HD21	1:C:443:PRO:HD2	1.84	0.43
1:B:456:GLN:NE2	4:B:2297:HOH:O	2.52	0.43
1:A:266:ALA:HB1	1:A:281:LEU:CD1	2.48	0.43
1:D:163:PHE:HB2	1:D:306:TYR:HB3	2.01	0.42
1:C:445:LEU:HB3	1:C:448:PHE:HB3	2.01	0.42
1:D:508:LEU:N	1:D:508:LEU:HD22	2.33	0.42
1:B:235:ARG:HD3	1:B:478:LEU:HD21	2.02	0.42
1:C:84:SER:OG	1:C:86:ILE:HG12	2.19	0.42
1:B:342:GLY:HA3	1:B:407:SER:O	2.18	0.42
1:C:439:LEU:HB2	1:C:447:THR:HA	2.00	0.42
1:A:320:TYR:O	1:A:323:VAL:HG12	2.20	0.42
1:C:330:HIS:HB2	3:C:1537:GOL:H32	2.01	0.42
1:C:63:MET:HE3	1:C:295:GLY:HA2	2.02	0.42
1:D:330:HIS:HE1	4:D:2215:HOH:O	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:69:PHE:C	1:D:71:ASP:H	2.23	0.42
1:A:78:LEU:O	1:A:82:LEU:HG	2.19	0.42
1:C:11:THR:H	1:C:49:ASN:ND2	2.12	0.42
1:C:409:LEU:O	1:C:409:LEU:HD23	2.18	0.42
1:A:409:LEU:C	1:A:409:LEU:HD23	2.40	0.42
1:D:44:TYR:CE2	1:D:48:LEU:HD21	2.55	0.42
2:B:1535:NAG:H61	2:B:1536:NAG:C8	2.49	0.42
1:D:409:LEU:O	1:D:409:LEU:HD23	2.19	0.42
1:C:25:GLY:H	1:C:101:ASN:ND2	2.18	0.42
1:D:206:TYR:HB2	1:D:238:ILE:HB	2.01	0.42
1:B:377:LEU:HD21	1:B:409:LEU:HD11	2.00	0.42
1:D:284:ASP:O	1:D:288:GLN:HG2	2.20	0.42
1:B:445:LEU:HB3	1:B:448:PHE:HB3	2.01	0.42
1:D:335:ILE:CG2	1:D:423:LEU:HD21	2.50	0.42
1:D:81:VAL:HG23	1:D:82:LEU:N	2.35	0.42
1:B:336:GLY:HA3	1:B:434:PHE:CZ	2.55	0.42
1:A:504:GLN:NE2	1:A:513:GLY:HA3	2.34	0.42
1:A:275:LEU:HD12	1:A:275:LEU:HA	1.78	0.41
1:D:231:LYS:HE3	1:D:231:LYS:HB2	1.80	0.41
1:B:465:VAL:HG13	1:B:485:LEU:CD2	2.50	0.41
1:C:317:ASP:O	1:C:318:ASP:C	2.59	0.41
1:C:402:GLN:OE1	1:C:405:ARG:HD2	2.19	0.41
1:A:464:SER:O	1:A:468:ASN:HB2	2.19	0.41
1:A:69:PHE:C	1:A:71:ASP:H	2.22	0.41
1:B:91:LEU:HD23	1:B:93:ASN:ND2	2.35	0.41
1:D:456:GLN:HB2	1:D:467:TYR:HB3	2.02	0.41
1:D:116:VAL:HB	1:D:203:VAL:HG22	2.02	0.41
1:C:330:HIS:HE1	4:C:2208:HOH:O	2.02	0.41
1:D:206:TYR:CB	1:D:238:ILE:HB	2.50	0.41
1:D:381:TYR:CZ	1:D:521:ALA:HB1	2.55	0.41
1:C:456:GLN:HG2	1:C:457:ASP:OD1	2.20	0.41
1:D:452:ASP:OD1	1:D:452:ASP:N	2.53	0.41
1:C:433:SER:HB3	1:C:502:LEU:HD23	2.02	0.41
1:C:381:TYR:CZ	1:C:521:ALA:HB1	2.56	0.41
1:C:508:LEU:CD2	1:D:70:GLU:HG3	2.51	0.41
1:D:16:ASN:ND2	4:D:2009:HOH:O	2.51	0.41
1:A:159:ALA:HB3	1:A:161:TRP:CE3	2.56	0.41
1:D:235:ARG:O	1:D:332:PRO:HD2	2.21	0.41
1:D:318:ASP:OD2	1:D:530:PRO:HG3	2.20	0.41
1:C:529:PRO:HB2	1:C:530:PRO:HD3	2.03	0.41
1:B:452:ASP:N	1:B:452:ASP:OD1	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:317:ASP:O	1:B:318:ASP:C	2.59	0.40
1:B:278:LEU:HA	1:B:281:LEU:CD1	2.51	0.40
1:A:302:LEU:HA	1:A:302:LEU:HD23	1.79	0.40
1:A:17:ALA:O	1:A:18:ILE:HB	2.21	0.40
1:D:182:GLN:O	1:D:186:MET:HG3	2.20	0.40
1:B:81:VAL:HG23	1:B:82:LEU:N	2.37	0.40
1:B:431:LYS:HE3	1:B:492:TYR:O	2.22	0.40
1:B:335:ILE:O	1:B:433:SER:HA	2.22	0.40
1:C:182:GLN:NE2	1:C:218:HIS:NE2	2.65	0.40
1:A:395:ILE:C	1:A:395:ILE:HD12	2.41	0.40
1:D:338:GLN:O	1:D:341:GLU:HG2	2.22	0.40
1:C:339:ASN:HA	1:C:339:ASN:HD22	1.71	0.40
1:A:389:SER:HA	1:A:391:PHE:N	2.36	0.40
1:C:41:PRO:HG3	1:C:181:ASP:HA	2.03	0.40
1:B:374:ILE:HG22	1:B:378:MET:HE2	2.03	0.40
1:A:284:ASP:HB3	4:A:2268:HOH:O	2.22	0.40
1:D:8:ASN:ND2	1:D:10:ASP:OD2	2.53	0.40

All (8) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:272:SER:OG	2:A:1535:NAG:O7[1_655]	1.51	0.69
4:B:2179:HOH:O	4:C:2033:HOH:O[1_645]	1.59	0.61
4:A:2041:HOH:O	4:A:2151:HOH:O[1_655]	1.95	0.25
4:C:2080:HOH:O	4:D:2230:HOH:O[1_455]	2.01	0.19
4:D:2008:HOH:O	4:D:2251:HOH:O[1_455]	2.03	0.17
1:B:280:GLY:O	1:C:31:PRO:CD[1_645]	2.07	0.13
4:D:2013:HOH:O	4:D:2338:HOH:O[1_455]	2.09	0.11
4:D:2016:HOH:O	4:D:2095:HOH:O[1_455]	2.12	0.08

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	532/534 (100%)	507 (95%)	23 (4%)	2 (0%)	39	31
1	B	532/534 (100%)	509 (96%)	22 (4%)	1 (0%)	52	47
1	C	532/534 (100%)	500 (94%)	32 (6%)	0	100	100
1	D	532/534 (100%)	506 (95%)	25 (5%)	1 (0%)	52	47
All	All	2128/2136 (100%)	2022 (95%)	102 (5%)	4 (0%)	52	47

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	70	GLU
1	B	390	PRO
1	D	390	PRO
1	A	390	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	435/435 (100%)	425 (98%)	10 (2%)	58	57
1	B	435/435 (100%)	428 (98%)	7 (2%)	70	71
1	C	435/435 (100%)	427 (98%)	8 (2%)	66	67
1	D	435/435 (100%)	423 (97%)	12 (3%)	51	48
All	All	1740/1740 (100%)	1703 (98%)	37 (2%)	61	61

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	ASN
1	A	101	ASN
1	A	136	ASP
1	A	206	TYR
1	A	235	ARG

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Mol	Chain	Res	Type
1	A	275	LEU
1	A	339	ASN
1	A	364	GLN
1	A	410	LEU
1	A	439	LEU
1	B	13	THR
1	B	20	ASN
1	B	70	GLU
1	B	101	ASN
1	B	235	ARG
1	B	410	LEU
1	B	494	SER
1	C	20	ASN
1	C	206	TYR
1	C	235	ARG
1	C	261	GLN
1	C	275	LEU
1	C	339	ASN
1	C	405	ARG
1	C	486	TRP
1	D	20	ASN
1	D	70	GLU
1	D	127	LEU
1	D	136	ASP
1	D	206	TYR
1	D	235	ARG
1	D	275	LEU
1	D	339	ASN
1	D	405	ARG
1	D	410	LEU
1	D	439	LEU
1	D	525	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (74) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	ASN
1	A	49	ASN
1	A	51	GLN
1	A	83	GLN
1	A	101	ASN
1	A	137	GLN

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Mol	Chain	Res	Type
1	A	155	ASN
1	A	182	GLN
1	A	240	GLN
1	A	283	GLN
1	A	330	HIS
1	A	338	GLN
1	A	339	ASN
1	A	424	ASN
1	A	456	GLN
1	A	500	ASN
1	A	528	ASN
1	B	20	ASN
1	B	49	ASN
1	B	51	GLN
1	B	93	ASN
1	B	101	ASN
1	B	137	GLN
1	B	155	ASN
1	B	182	GLN
1	B	187	GLN
1	B	240	GLN
1	B	283	GLN
1	B	330	HIS
1	B	338	GLN
1	B	339	ASN
1	B	424	ASN
1	B	456	GLN
1	B	500	ASN
1	B	528	ASN
1	C	20	ASN
1	C	49	ASN
1	C	51	GLN
1	C	83	GLN
1	C	101	ASN
1	C	137	GLN
1	C	155	ASN
1	C	182	GLN
1	C	187	GLN
1	C	222	ASN
1	C	240	GLN
1	C	330	HIS
1	C	338	GLN

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Mol	Chain	Res	Type
1	C	339	ASN
1	C	424	ASN
1	C	456	GLN
1	C	488	ASN
1	C	500	ASN
1	C	504	GLN
1	C	528	ASN
1	D	20	ASN
1	D	49	ASN
1	D	51	GLN
1	D	76	ASN
1	D	83	GLN
1	D	93	ASN
1	D	101	ASN
1	D	137	GLN
1	D	155	ASN
1	D	182	GLN
1	D	240	GLN
1	D	283	GLN
1	D	330	HIS
1	D	338	GLN
1	D	339	ASN
1	D	424	ASN
1	D	456	GLN
1	D	500	ASN
1	D	528	ASN

5.3.3 RNA ⓘ

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 ⓘ

8 carbohydrates 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	NAG	A	1535	1,2	14,14,15	0.58	0	15,19,21	0.69	0
2	NAG	A	1536	2	14,14,15	0.54	0	15,19,21	0.81	1 (6%)
2	NAG	B	1535	1,2	14,14,15	0.76	1 (7%)	15,19,21	0.88	1 (6%)
2	NAG	B	1536	2	14,14,15	0.54	0	15,19,21	1.04	1 (6%)
2	NAG	C	1535	1,2	14,14,15	0.55	0	15,19,21	0.79	1 (6%)
2	NAG	C	1536	2	14,14,15	0.50	0	15,19,21	0.68	0
2	NAG	D	1535	1,2	14,14,15	0.52	0	15,19,21	0.75	1 (6%)
2	NAG	D	1536	2	14,14,15	0.47	0	15,19,21	0.74	1 (6%)

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	NAG	A	1535	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	1536	2	-	0/6/23/26	0/1/1/1
2	NAG	B	1535	1,2	1/1/5/7	0/6/23/26	0/1/1/1
2	NAG	B	1536	2	-	0/6/23/26	0/1/1/1
2	NAG	C	1535	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	1536	2	-	0/6/23/26	0/1/1/1
2	NAG	D	1535	1,2	-	0/6/23/26	0/1/1/1
2	NAG	D	1536	2	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1535	NAG	C1-C2	2.18	1.55	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1536	NAG	C2-N2-C7	-2.67	119.61	123.04
2	B	1535	NAG	C2-N2-C7	-2.45	119.90	123.04
2	D	1535	NAG	C2-N2-C7	-2.42	119.93	123.04

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1536	NAG	C2-N2-C7	-2.33	120.05	123.04
2	A	1536	NAG	C2-N2-C7	-2.29	120.09	123.04
2	C	1535	NAG	C2-N2-C7	-2.29	120.10	123.04

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	1535	NAG	C1

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1535	NAG	0	1
2	B	1535	NAG	2	0
2	B	1536	NAG	2	0
2	C	1535	NAG	1	0
2	C	1536	NAG	2	0

5.6 Ligand geometry [i](#)

8 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	A	1537	-	5,5,5	0.12	0	5,5,5	0.40	0
3	GOL	A	1538	-	5,5,5	0.36	0	5,5,5	0.50	0
3	GOL	B	1537	-	5,5,5	0.23	0	5,5,5	0.43	0
3	GOL	B	1538	-	5,5,5	0.33	0	5,5,5	0.35	0
3	GOL	C	1537	-	5,5,5	0.15	0	5,5,5	0.54	0
3	GOL	C	1538	-	5,5,5	0.33	0	5,5,5	0.57	0
3	GOL	D	1537	-	5,5,5	0.40	0	5,5,5	0.28	0
3	GOL	D	1538	-	5,5,5	0.32	0	5,5,5	0.55	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	A	1537	-	-	0/4/4/4	0/0/0/0
3	GOL	A	1538	-	-	0/4/4/4	0/0/0/0
3	GOL	B	1537	-	-	0/4/4/4	0/0/0/0
3	GOL	B	1538	-	-	0/4/4/4	0/0/0/0
3	GOL	C	1537	-	-	0/4/4/4	0/0/0/0
3	GOL	C	1538	-	-	0/4/4/4	0/0/0/0
3	GOL	D	1537	-	-	0/4/4/4	0/0/0/0
3	GOL	D	1538	-	-	0/4/4/4	0/0/0/0

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.

7 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1538	GOL	1	0
3	B	1537	GOL	4	0
3	B	1538	GOL	1	0
3	C	1537	GOL	5	0
3	C	1538	GOL	4	0
3	D	1537	GOL	1	0
3	D	1538	GOL	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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.