



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2014 – 02:46 AM GMT

PDB ID : 2W2D
Title : Crystal Structure of a Catalytically Active, Non-toxic Endopeptidase Derivative of Clostridium botulinum Toxin A
Authors : Masuyer, G.; Thiagarajan, N.; James, P.L.; Marks, P.M.H.; Chaddock, J.A.; Acharya, K.R.
Deposited on : 2008-10-29
Resolution : 2.59 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

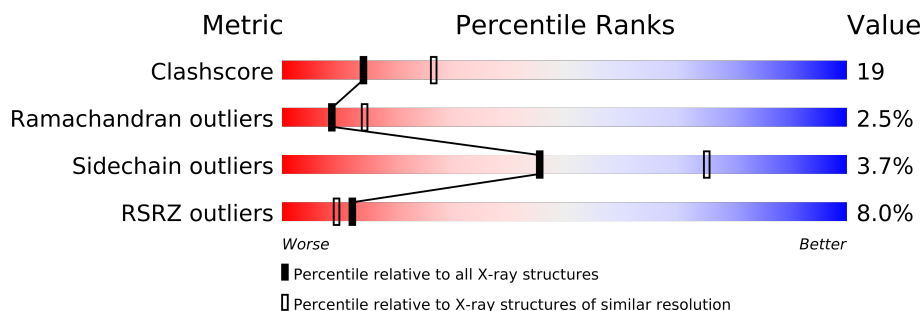
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.59 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	79885	2154 (2.60-2.60)
Ramachandran outliers	78287	2113 (2.60-2.60)
Sidechain outliers	78261	2113 (2.60-2.60)
RSRZ outliers	66119	1718 (2.60-2.60)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	450	
1	C	450	
2	B	431	
2	D	431	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
6	GOL	A	1436	X	-
6	GOL	A	1437	X	X
6	GOL	A	1438	X	X
6	GOL	B	1873	X	-
6	GOL	B	1874	X	X
7	ACT	C	1435	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 14160 atoms, of which 0 are hydrogen and 0 are deuterium.

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 BOTULINUM NEUROTOXIN A LIGHT CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	437	Total	C	N	O	S	0	0	1
			3534	2281	574	669	10			
1	C	436	Total	C	N	O	S	0	0	1
			3522	2273	573	666	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	GLU	PRO	CONFLICT	UNP A5HZZ9
A	432	ASP	ARG	CONFLICT	UNP A5HZZ9
C	2	GLU	PRO	CONFLICT	UNP A5HZZ9
C	432	ASP	ARG	CONFLICT	UNP A5HZZ9

- Molecule 2 is a protein called BOTULINUM NEUROTOXIN A HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	424	Total	C	N	O	S	0	0	1
			3429	2195	550	674	10			
2	D	408	Total	C	N	O	S	0	0	1
			3308	2119	529	650	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	452	LEU	ASP	CONFLICT	UNP A5HZZ9
B	453	GLN	LEU	CONFLICT	UNP A5HZZ9
D	452	LEU	ASP	CONFLICT	UNP A5HZZ9
D	453	GLN	LEU	CONFLICT	UNP A5HZZ9

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

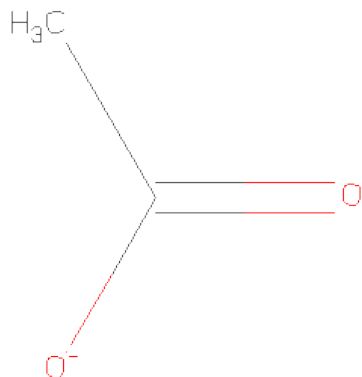
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Cl	0	0
			2	2		
5	D	1	Total	Cl	0	0
			1	1		

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 7 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

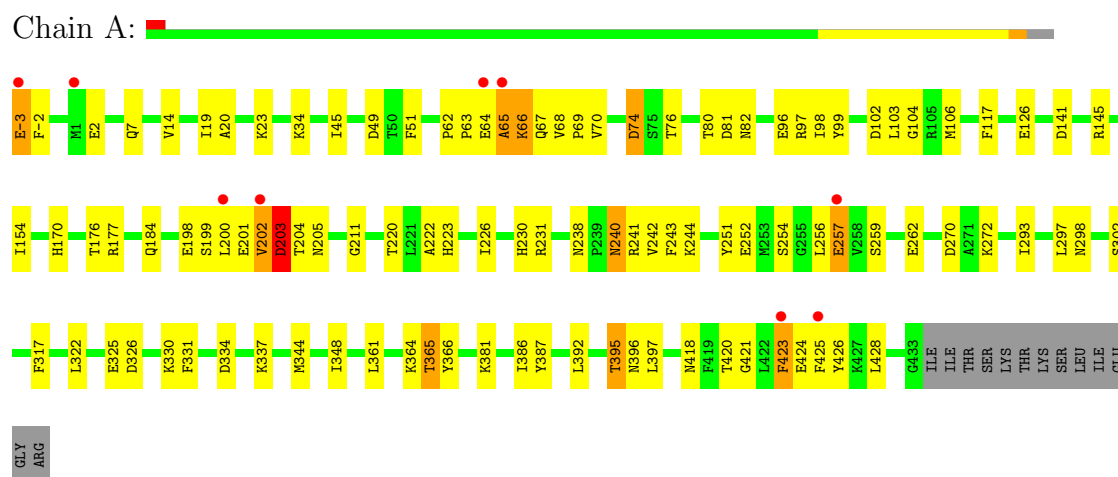
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	110	Total	O	0	0
			110	110		
8	B	49	Total	O	0	0
			49	49		
8	C	118	Total	O	0	0
			118	118		
8	D	31	Total	O	0	0
			31	31		

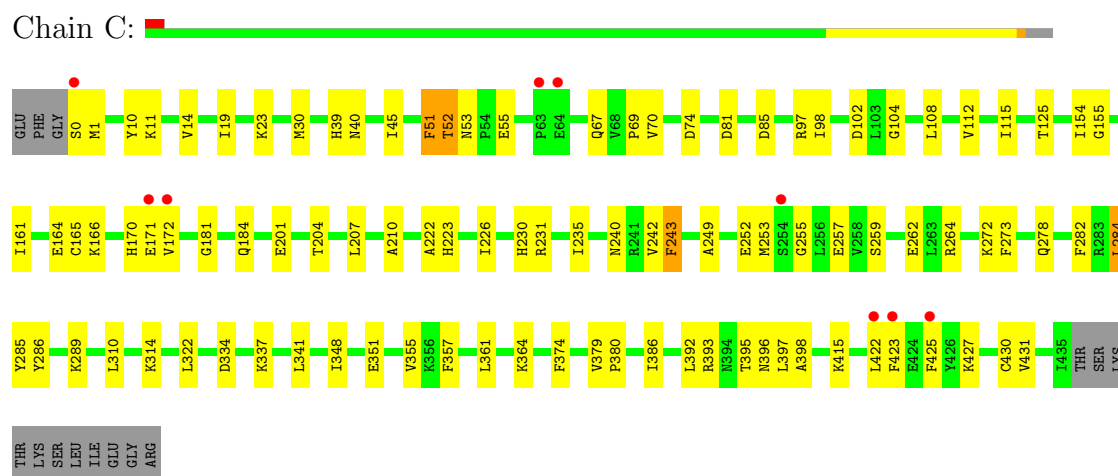
3 Residue-property plots

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

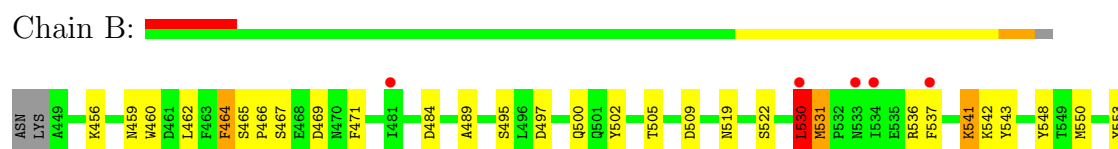
• Molecule 1: BOTULINUM NEUROTOXIN A LIGHT CHAIN

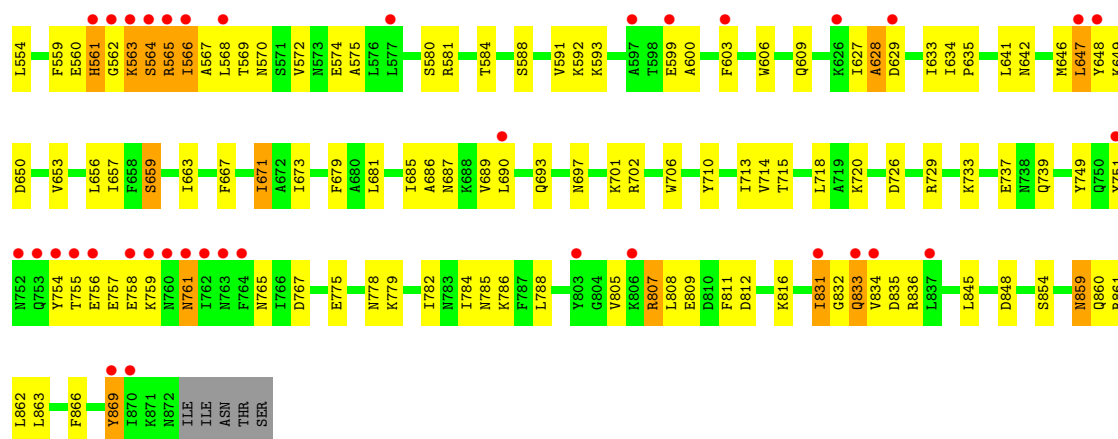


• Molecule 1: BOTULINUM NEUROTOXIN A LIGHT CHAIN



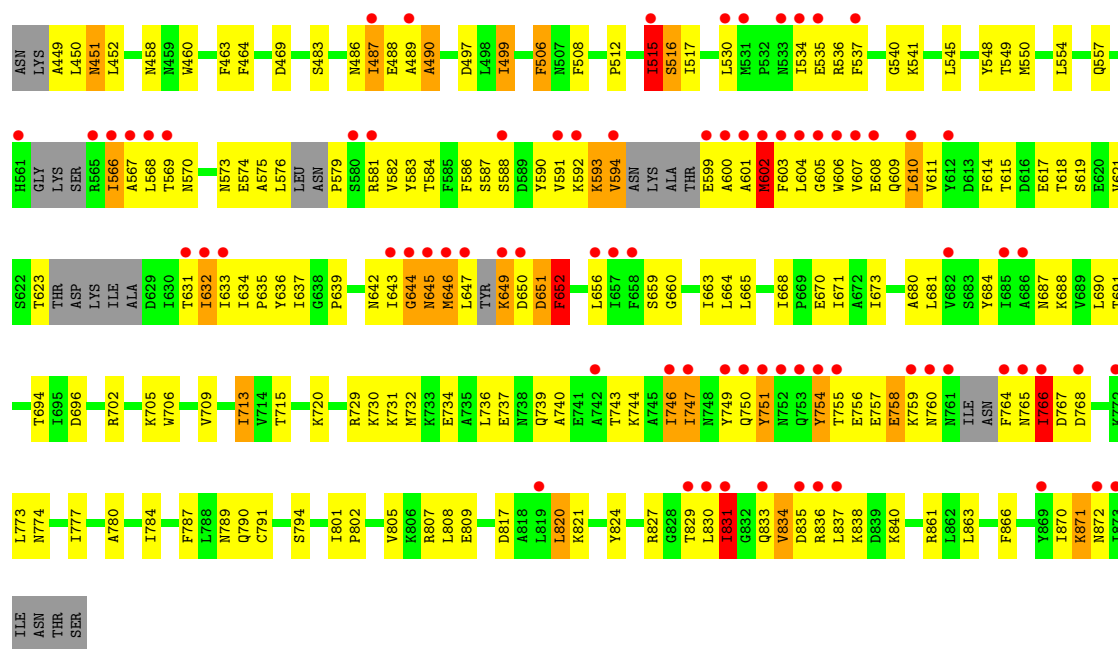
• Molecule 2: BOTULINUM NEUROTOXIN A HEAVY CHAIN





• Molecule 2: BOTULINUM NEUROTOXIN A HEAVY CHAIN

Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.36Å 156.91Å 211.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.11 – 2.59 49.11 – 2.59	Depositor EDS
% Data completeness (in resolution range)	91.2 (49.11-2.59) 91.1 (49.11-2.59)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.43 (at 2.58Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.212 , 0.253 0.212 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	34.9	Xtriage
Anisotropy	0.242	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 35.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 79768 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	14160	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.63% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, ZN, ACT, SO4, CL

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	0.49	0/3618	0.64	0/4896
1	C	0.51	0/3605	0.63	0/4879
2	B	0.48	0/3492	0.61	0/4735
2	D	0.45	0/3364	0.66	3/4552 (0.1%)
All	All	0.48	0/14079	0.64	3/19062 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	652	PHE	CB-CG-CD1	-9.74	113.98	120.80
2	D	652	PHE	CB-CG-CD2	6.02	125.02	120.80
2	D	831	ILE	CB-CA-C	-5.11	101.38	111.60

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3534	0	3471	108	0
1	C	3522	0	3467	79	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	3429	0	3392	136	0
2	D	3308	0	3258	249	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	C	10	0	0	0	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
5	A	2	0	0	1	0
5	D	1	0	0	0	0
6	A	18	0	9	1	0
6	B	12	0	6	0	0
7	C	4	0	3	0	0
8	A	110	0	0	4	0
8	B	49	0	0	3	0
8	C	118	0	0	1	0
8	D	31	0	0	4	0
All	All	14160	0	13606	534	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 19.

All (534) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:831:ILE:CG2	2:D:831:ILE:O	1.67	1.35
1:A:200:LEU:HG	2:B:861:ARG:NH1	1.46	1.30
2:D:649:LYS:O	2:D:649:LYS:CD	1.89	1.21
1:A:202:VAL:O	1:A:204:THR:N	1.74	1.19
2:D:659:SER:HB2	2:D:663:ILE:HD11	1.22	1.18
1:C:395:THR:HG22	1:C:397:LEU:H	1.13	1.11
2:D:450:LEU:HD22	2:D:451:ASN:H	1.13	1.10
1:A:395:THR:HG22	1:A:397:LEU:H	1.15	1.08
2:D:566:ILE:HG12	2:D:567:ALA:H	1.12	1.08
2:D:660:GLY:O	2:D:663:ILE:HD13	1.53	1.08
1:A:202:VAL:HG23	1:A:203:ASP:N	1.64	1.06
2:D:866:PHE:CZ	2:D:870:ILE:HD11	1.91	1.04
1:A:202:VAL:CG2	1:A:203:ASP:H	1.69	1.04
1:A:202:VAL:O	1:A:205:ASN:N	1.94	1.01
2:D:584:THR:HG22	2:D:586:PHE:H	1.21	1.01
2:D:649:LYS:O	2:D:649:LYS:HD3	1.61	0.99
2:D:602:MET:HG2	2:D:603:PHE:CE1	1.98	0.98
1:A:200:LEU:HG	2:B:861:ARG:HH11	1.07	0.97
2:D:831:ILE:HG23	2:D:831:ILE:O	1.14	0.96

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:202:VAL:HG23	1:A:203:ASP:H	0.82	0.96
1:A:201:GLU:OE2	1:A:364:LYS:NZ	2.02	0.93
2:B:530:LEU:HD12	2:B:530:LEU:H	1.34	0.93
2:D:709:VAL:O	2:D:713:ILE:HD13	1.71	0.91
2:D:743:THR:O	2:D:747:ILE:HD13	1.71	0.91
2:D:829:THR:O	2:D:830:LEU:HD23	1.69	0.91
2:D:487:ILE:HD13	2:D:487:ILE:H	1.36	0.90
1:A:200:LEU:CG	2:B:861:ARG:NH1	2.34	0.90
2:D:566:ILE:HG12	2:D:567:ALA:N	1.86	0.88
2:D:635:PRO:HA	2:D:652:PHE:HE2	1.37	0.88
1:A:74:ASP:OD1	1:A:76:THR:HG22	1.74	0.88
2:D:833:GLN:O	2:D:835:ASP:N	2.09	0.86
2:D:450:LEU:HD13	2:D:451:ASN:N	1.90	0.86
1:C:395:THR:HG22	1:C:397:LEU:N	1.90	0.86
2:D:487:ILE:HG12	2:D:488:GLU:H	1.39	0.85
2:D:637:ILE:HG12	2:D:732:MET:HE3	1.59	0.85
1:C:184:GLN:OE1	1:C:231:ARG:HD3	1.77	0.85
1:C:249:ALA:HB3	1:C:252:GLU:HG3	1.59	0.85
2:D:450:LEU:HD22	2:D:451:ASN:N	1.91	0.84
2:D:757:GLU:HG3	2:D:760:ASN:HB3	1.59	0.84
2:D:831:ILE:HG22	2:D:831:ILE:O	1.74	0.83
2:D:766:ILE:HG12	2:D:767:ASP:H	1.43	0.83
2:D:566:ILE:CG1	2:D:567:ALA:H	1.91	0.83
2:D:649:LYS:O	2:D:649:LYS:HD2	1.78	0.82
2:D:602:MET:HG2	2:D:603:PHE:CD1	2.14	0.82
1:A:395:THR:CG2	1:A:397:LEU:H	1.93	0.82
1:A:7:GLN:HG2	1:A:7:GLN:O	1.80	0.82
2:B:588:SER:O	2:B:591:VAL:HG22	1.81	0.81
2:D:601:ALA:O	2:D:602:MET:HB3	1.80	0.80
1:C:164:GLU:OE2	1:C:166:LYS:HE2	1.80	0.80
1:A:-2:PHE:HB2	1:A:387:TYR:HE2	1.47	0.80
1:A:14:VAL:HG13	1:A:20:ALA:HA	1.64	0.80
1:A:45:ILE:HB	1:A:154:ILE:HG22	1.62	0.80
1:C:210:ALA:HB2	2:D:774:ASN:OD1	1.82	0.80
2:D:593:LYS:O	2:D:599:GLU:HB2	1.82	0.79
1:A:200:LEU:HG	2:B:861:ARG:HH12	1.46	0.79
1:C:45:ILE:HB	1:C:154:ILE:HG12	1.65	0.79
2:B:572:VAL:HG13	2:B:574:GLU:H	1.47	0.78
2:D:599:GLU:HG2	2:D:600:ALA:N	1.98	0.78
1:A:200:LEU:CG	2:B:861:ARG:HH11	1.92	0.78
2:D:554:LEU:HD11	2:D:731:LYS:HG3	1.65	0.78
2:D:649:LYS:O	2:D:649:LYS:CG	2.29	0.77

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:565:ARG:HE	2:B:565:ARG:HA	1.49	0.77
2:D:515:ILE:O	2:D:516:SER:HB3	1.82	0.77
2:D:483:SER:HA	2:D:694:THR:HG22	1.66	0.77
2:D:633:ILE:O	2:D:635:PRO:HD3	1.85	0.77
1:A:201:GLU:HG2	1:A:361:LEU:HD11	1.67	0.76
1:A:428:LEU:HD23	2:B:542:LYS:HG3	1.68	0.76
2:D:588:SER:O	2:D:591:VAL:HG22	1.86	0.75
2:D:659:SER:CB	2:D:663:ILE:HD11	2.12	0.75
2:D:833:GLN:O	2:D:836:ARG:N	2.18	0.75
2:D:568:LEU:HD11	2:D:594:VAL:C	2.07	0.74
2:D:866:PHE:CE2	2:D:870:ILE:HD11	2.22	0.74
2:D:601:ALA:O	2:D:602:MET:CB	2.36	0.74
2:B:565:ARG:HG3	2:B:749:TYR:HB2	1.70	0.74
1:A:63:PRO:HG2	1:A:64:GLU:H	1.52	0.73
2:D:635:PRO:HA	2:D:652:PHE:CE2	2.22	0.73
2:D:488:GLU:HG3	2:D:489:ALA:H	1.53	0.73
1:A:395:THR:HG22	1:A:397:LEU:N	1.98	0.72
2:D:487:ILE:HG12	2:D:488:GLU:N	2.03	0.72
1:A:68:VAL:O	1:A:420:THR:HG21	1.89	0.72
2:D:599:GLU:HB3	2:D:610:LEU:HD11	1.70	0.72
1:A:66:LYS:O	1:A:68:VAL:N	2.22	0.72
2:D:599:GLU:HG2	2:D:600:ALA:H	1.52	0.72
2:D:833:GLN:O	2:D:834:VAL:C	2.27	0.72
2:D:673:ILE:HG22	2:D:807:ARG:HG3	1.70	0.72
2:D:690:LEU:O	2:D:694:THR:HG23	1.89	0.72
1:A:199:SER:O	1:A:200:LEU:CD1	2.37	0.71
2:D:450:LEU:CD2	2:D:451:ASN:H	1.97	0.71
2:D:584:THR:CG2	2:D:586:PHE:H	2.03	0.71
1:C:45:ILE:HD12	1:C:154:ILE:HD11	1.71	0.71
2:B:647:LEU:HD12	8:B:2031:HOH:O	1.90	0.70
2:D:829:THR:HG22	2:D:829:THR:O	1.90	0.70
2:D:573:ASN:O	2:D:576:LEU:HD12	1.90	0.69
2:D:664:LEU:HD11	2:D:791:CYS:SG	2.33	0.69
1:C:23:LYS:HA	1:C:30:MET:CE	2.22	0.69
2:D:632:ILE:HD13	2:D:632:ILE:H	1.58	0.68
1:A:201:GLU:CG	1:A:361:LEU:HD11	2.23	0.68
2:D:687:ASN:N	2:D:829:THR:HG21	2.08	0.68
2:B:785:ASN:HB3	2:B:861:ARG:NH2	2.09	0.68
2:B:575:ALA:HB1	2:B:581:ARG:O	1.93	0.68
2:D:570:ASN:HD22	2:D:581:ARG:HD3	1.57	0.68
2:D:866:PHE:O	2:D:870:ILE:HG12	1.93	0.68
2:D:487:ILE:CD1	2:D:487:ILE:H	2.05	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:515:ILE:HD11	2:D:517:ILE:HG23	1.76	0.68
1:A:202:VAL:CG2	1:A:203:ASP:N	2.39	0.68
2:B:563:LYS:HE2	2:B:563:LYS:HA	1.76	0.68
1:A:184:GLN:OE1	1:A:231:ARG:HD3	1.93	0.68
2:B:563:LYS:O	2:B:565:ARG:N	2.27	0.67
1:C:334:ASP:HB3	1:C:337:LYS:HB2	1.75	0.67
1:A:65:ALA:O	1:A:66:LYS:HB2	1.92	0.67
1:C:395:THR:CG2	1:C:397:LEU:H	1.99	0.67
1:A:66:LYS:O	1:A:68:VAL:HG22	1.95	0.67
2:B:627:ILE:HG12	2:B:657:ILE:HA	1.75	0.67
1:A:199:SER:O	1:A:200:LEU:HD12	1.95	0.67
2:D:870:ILE:O	2:D:871:LYS:HB2	1.94	0.66
1:A:201:GLU:HG3	1:A:361:LEU:CD1	2.25	0.66
1:A:202:VAL:O	1:A:203:ASP:C	2.34	0.66
2:B:785:ASN:HB3	2:B:861:ARG:HH22	1.59	0.66
2:D:566:ILE:HB	2:D:749:TYR:CE2	2.30	0.66
2:B:569:THR:HG23	2:B:584:THR:CG2	2.25	0.66
2:D:599:GLU:CG	2:D:600:ALA:N	2.57	0.66
1:A:201:GLU:O	1:A:202:VAL:C	2.34	0.66
2:D:619:SER:O	2:D:621:VAL:HG23	1.97	0.65
1:A:200:LEU:O	1:A:203:ASP:OD2	2.15	0.65
2:D:601:ALA:O	2:D:602:MET:SD	2.55	0.64
2:B:702:ARG:HD2	2:B:812:ASP:OD1	1.96	0.64
2:B:729:ARG:HB3	2:B:788:LEU:HD12	1.79	0.64
2:D:632:ILE:HD13	2:D:790:GLN:OE1	1.97	0.64
1:C:201:GLU:HG3	1:C:361:LEU:CD1	2.27	0.64
2:D:602:MET:CG	2:D:603:PHE:CE1	2.77	0.64
1:A:98:ILE:O	1:A:104:GLY:HA3	1.96	0.64
2:D:550:MET:CE	2:D:731:LYS:HG2	2.28	0.64
1:A:270:ASP:OD1	1:A:365:THR:HG23	1.97	0.64
2:D:780:ALA:O	2:D:784:ILE:HG12	1.98	0.64
2:B:553:TYR:OH	2:B:647:LEU:HB3	1.99	0.63
2:D:450:LEU:O	2:D:451:ASN:HB2	1.98	0.63
2:D:740:ALA:O	2:D:744:LYS:HG3	1.97	0.63
2:D:530:LEU:H	2:D:530:LEU:HD12	1.63	0.63
1:A:201:GLU:CG	1:A:361:LEU:CD1	2.76	0.63
2:B:697:ASN:O	2:B:701:LYS:HB2	1.98	0.63
2:B:584:THR:HG22	2:B:739:GLN:OE1	1.99	0.63
1:A:325:GLU:HA	1:A:330:LYS:O	1.99	0.62
2:D:530:LEU:H	2:D:530:LEU:CD1	2.12	0.62
2:D:644:GLY:O	2:D:645:ASN:C	2.37	0.62
2:B:634:ILE:HD12	2:B:784:ILE:HA	1.82	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:634:ILE:HB	2:D:637:ILE:HD12	1.82	0.62
2:D:734:GLU:HA	2:D:737:GLU:HG2	1.81	0.62
2:D:650:ASP:O	2:D:651:ASP:C	2.37	0.62
2:D:766:ILE:HG12	2:D:767:ASP:N	2.14	0.62
2:B:530:LEU:N	2:B:530:LEU:HD12	2.11	0.61
2:D:642:ASN:OD1	2:D:646:MET:HA	2.00	0.61
2:D:663:ILE:CD1	2:D:663:ILE:H	2.14	0.61
1:A:80:THR:HG22	1:A:82:ASN:H	1.65	0.61
2:B:464:PHE:CE2	2:B:466:PRO:HG3	2.35	0.61
1:A:34:LYS:HZ2	6:A:1436:GOL:H11	1.66	0.61
2:B:569:THR:HG23	2:B:584:THR:HG23	1.82	0.60
2:B:635:PRO:HB2	2:B:649:LYS:HB3	1.82	0.60
2:B:805:VAL:O	2:B:809:GLU:HG3	2.01	0.60
2:D:691:THR:CG2	2:D:830:LEU:HD11	2.32	0.60
2:D:647:LEU:O	2:D:650:ASP:HB3	2.01	0.60
2:D:688:LYS:NZ	2:D:831:ILE:HG22	2.16	0.60
2:D:730:LYS:HG2	8:D:2024:HOH:O	2.02	0.59
1:A:202:VAL:C	1:A:204:THR:N	2.54	0.59
2:D:483:SER:CA	2:D:694:THR:HG22	2.32	0.59
2:B:646:MET:C	2:B:647:LEU:HD23	2.22	0.59
2:B:714:VAL:O	2:B:718:LEU:HB2	2.03	0.59
2:D:515:ILE:HD13	2:D:516:SER:H	1.67	0.59
2:B:671:ILE:HG23	2:B:671:ILE:O	2.03	0.59
2:D:691:THR:HG21	2:D:830:LEU:HD11	1.83	0.59
2:D:805:VAL:O	2:D:809:GLU:HG3	2.03	0.58
2:D:736:LEU:HB3	2:D:777:ILE:HD12	1.85	0.58
2:D:663:ILE:HD12	2:D:663:ILE:H	1.69	0.58
2:D:668:ILE:H	2:D:668:ILE:HD12	1.67	0.58
2:D:550:MET:CE	2:D:731:LYS:HE2	2.32	0.58
2:B:706:TRP:CE3	2:B:808:LEU:HD13	2.39	0.58
1:C:249:ALA:HB3	1:C:252:GLU:CG	2.33	0.58
2:B:560:GLU:C	2:B:562:GLY:H	2.07	0.58
1:C:431:VAL:HG21	2:D:548:TYR:HE2	1.69	0.58
1:C:257:GLU:HG3	2:D:534:ILE:HG23	1.84	0.57
2:B:553:TYR:CE2	2:B:642:ASN:HB2	2.40	0.57
1:A:421:GLY:H	1:A:424:GLU:HG3	1.69	0.57
1:C:201:GLU:HG2	1:C:364:LYS:HZ3	1.70	0.57
2:B:689:VAL:HG12	2:B:693:GLN:OE1	2.05	0.57
1:A:200:LEU:HD12	2:B:785:ASN:ND2	2.19	0.57
2:B:860:GLN:CD	2:B:860:GLN:H	2.08	0.57
2:D:671:ILE:O	2:D:671:ILE:HG23	2.04	0.57
1:A:199:SER:O	1:A:200:LEU:HD13	2.04	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:200:LEU:HD13	2:B:729:ARG:HD3	1.87	0.57
1:C:23:LYS:HA	1:C:30:MET:HE2	1.87	0.57
2:D:489:ALA:O	2:D:490:ALA:HB3	2.05	0.56
2:B:559:PHE:HE1	2:B:565:ARG:HH22	1.52	0.56
2:D:663:ILE:HD12	2:D:663:ILE:N	2.20	0.56
2:D:599:GLU:N	2:D:599:GLU:OE1	2.38	0.56
2:D:550:MET:HE3	2:D:731:LYS:HG2	1.87	0.56
1:C:272:LYS:O	2:D:715:THR:HG23	2.05	0.56
2:D:610:LEU:HD12	2:D:610:LEU:H	1.70	0.56
2:D:530:LEU:N	2:D:530:LEU:HD12	2.20	0.56
1:A:272:LYS:O	2:B:715:THR:HG23	2.06	0.56
1:C:348:ILE:HG12	2:D:499:ILE:HD12	1.88	0.56
2:D:593:LYS:O	2:D:594:VAL:HG13	2.04	0.55
2:D:594:VAL:HG13	2:D:610:LEU:HD21	1.88	0.55
2:D:834:VAL:O	2:D:838:LYS:HG3	2.06	0.55
1:C:201:GLU:HG3	1:C:361:LEU:HD11	1.87	0.55
2:D:541:LYS:HE2	2:D:541:LYS:HA	1.88	0.55
1:A:177:ARG:HD3	1:A:238:ASN:HA	1.87	0.55
2:D:644:GLY:O	2:D:646:MET:N	2.39	0.55
2:B:754:TYR:HB2	2:B:758:GLU:HG2	1.88	0.55
1:A:74:ASP:CG	1:A:76:THR:HG22	2.26	0.55
2:D:584:THR:HG22	2:D:586:PHE:N	2.05	0.55
2:D:554:LEU:CD1	2:D:731:LYS:HG3	2.37	0.54
2:D:637:ILE:HG23	2:D:732:MET:CE	2.37	0.54
2:D:637:ILE:HG23	2:D:732:MET:HE1	1.89	0.54
2:B:569:THR:HG22	2:B:591:VAL:HG12	1.88	0.54
2:B:565:ARG:NE	2:B:565:ARG:HA	2.20	0.54
2:D:540:GLY:O	2:D:541:LYS:C	2.45	0.54
1:C:348:ILE:HG12	2:D:499:ILE:CD1	2.36	0.54
1:C:39:HIS:CD2	1:C:40:ASN:N	2.75	0.54
1:C:125:THR:HA	8:C:2039:HOH:O	2.08	0.54
2:B:570:ASN:ND2	2:B:572:VAL:HG12	2.23	0.54
2:D:515:ILE:HD13	2:D:516:SER:N	2.23	0.54
2:B:553:TYR:CZ	2:B:642:ASN:HB2	2.43	0.54
1:A:201:GLU:HG3	1:A:361:LEU:HD13	1.89	0.53
2:D:590:TYR:OH	2:D:743:THR:HG21	2.07	0.53
1:C:115:ILE:HD11	2:D:512:PRO:HB2	1.88	0.53
1:C:154:ILE:HG22	1:C:155:GLY:O	2.08	0.53
2:D:545:LEU:N	2:D:545:LEU:HD12	2.22	0.53
1:A:-3:GLU:HA	5:A:1435:CL:CL	2.45	0.53
2:D:696:ASP:OD2	2:D:840:LYS:HE2	2.08	0.53
1:C:67:GLN:HE22	2:D:537:PHE:H	1.56	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:729:ARG:HB3	2:B:788:LEU:CD1	2.38	0.53
2:B:755:THR:C	2:B:757:GLU:H	2.12	0.53
1:A:14:VAL:HG13	1:A:19:ILE:O	2.09	0.53
1:A:62:PRO:HG2	1:A:64:GLU:O	2.08	0.53
2:B:673:ILE:O	2:B:807:ARG:NH1	2.42	0.53
2:D:566:ILE:HB	2:D:749:TYR:CZ	2.44	0.53
2:B:530:LEU:O	2:B:531:MET:HB3	2.08	0.53
1:C:14:VAL:HA	1:C:19:ILE:HG22	1.91	0.53
2:B:497:ASP:O	2:B:500:GLN:HB3	2.08	0.53
2:B:778:ASN:O	2:B:782:ILE:HG13	2.09	0.53
2:B:687:ASN:ND2	2:B:690:LEU:HD22	2.23	0.53
1:C:255:GLY:HA3	2:D:537:PHE:CD2	2.43	0.53
2:D:450:LEU:C	2:D:450:LEU:HD13	2.29	0.53
2:B:679:PHE:CE2	2:B:701:LYS:HB3	2.45	0.53
1:A:381:LYS:HG2	8:A:2092:HOH:O	2.08	0.52
2:D:623:THR:O	2:D:623:THR:HG22	2.09	0.52
1:A:425:PHE:CZ	2:B:537:PHE:HB2	2.44	0.52
2:D:756:GLU:HG3	2:D:758:GLU:CD	2.30	0.52
2:D:634:ILE:HG21	2:D:784:ILE:CD1	2.40	0.52
2:B:471:PHE:CE2	2:B:720:LYS:HE3	2.45	0.52
2:D:643:ILE:HG22	2:D:663:ILE:HG22	1.91	0.52
2:D:583:TYR:CG	2:D:584:THR:N	2.77	0.52
2:D:687:ASN:H	2:D:829:THR:HG21	1.74	0.52
2:D:557:GLN:HA	2:D:582:VAL:HG13	1.90	0.52
1:C:422:LEU:O	1:C:423:PHE:CD1	2.63	0.52
1:C:249:ALA:O	1:C:427:LYS:NZ	2.42	0.52
1:C:230:HIS:CE1	1:C:264:ARG:HD3	2.44	0.52
2:B:567:ALA:O	2:B:568:LEU:HD22	2.11	0.51
2:D:668:ILE:N	2:D:668:ILE:HD12	2.25	0.51
2:D:535:GLU:CD	2:D:535:GLU:H	2.14	0.51
2:B:570:ASN:ND2	2:B:572:VAL:H	2.08	0.51
2:B:648:TYR:O	2:B:649:LYS:C	2.48	0.51
1:C:97:ARG:HA	1:C:386:ILE:HG23	1.91	0.51
1:C:23:LYS:HA	1:C:30:MET:HE1	1.90	0.51
2:D:569:THR:HG23	2:D:584:THR:OG1	2.10	0.51
2:B:702:ARG:HG3	2:B:811:PHE:HE1	1.75	0.51
1:A:421:GLY:N	1:A:424:GLU:HG3	2.25	0.51
2:B:593:LYS:NZ	2:B:609:GLN:HG3	2.27	0.51
1:C:115:ILE:HD11	2:D:512:PRO:CB	2.41	0.50
2:B:809:GLU:OE2	2:B:848:ASP:OD2	2.30	0.50
1:A:69:PRO:HG3	2:B:460:TRP:CD2	2.46	0.50
2:B:489:ALA:HA	2:D:872:ASN:O	2.12	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:684:TYR:CE1	2:D:694:THR:HG21	2.47	0.50
2:D:605:GLY:O	2:D:608:GLU:HB3	2.11	0.50
1:A:240:ASN:C	1:A:240:ASN:ND2	2.64	0.50
2:D:659:SER:HB2	2:D:663:ILE:CD1	2.16	0.50
2:D:450:LEU:CD1	2:D:451:ASN:N	2.69	0.50
2:D:610:LEU:HD13	2:D:747:ILE:HD11	1.94	0.50
2:B:600:ALA:HA	2:B:754:TYR:CZ	2.47	0.50
1:C:392:LEU:O	1:C:395:THR:HB	2.10	0.50
1:A:-2:PHE:HB2	1:A:387:TYR:CE2	2.37	0.50
1:C:181:GLY:HA2	1:C:231:ARG:O	2.11	0.50
2:D:739:GLN:NE2	2:D:739:GLN:HA	2.27	0.50
1:C:154:ILE:CG2	1:C:155:GLY:N	2.74	0.50
1:C:430:CYS:HB3	2:D:452:LEU:HB3	1.94	0.50
1:A:334:ASP:OD1	1:A:337:LYS:HG3	2.12	0.50
1:C:70:VAL:HG23	1:C:161:ILE:HD12	1.93	0.49
2:B:541:LYS:HG2	2:B:543:TYR:CE1	2.48	0.49
2:B:647:LEU:HB2	8:B:2031:HOH:O	2.12	0.49
1:A:102:ASP:O	1:A:106:MET:HG3	2.12	0.49
2:B:603:PHE:O	2:B:606:TRP:HB3	2.12	0.49
1:A:426:TYR:CE1	2:B:456:LYS:HG2	2.47	0.49
1:A:298:ASN:HD21	1:A:331:PHE:H	1.60	0.49
2:D:656:LEU:HD11	2:D:787:PHE:HZ	1.76	0.49
1:C:392:LEU:O	1:C:398:ALA:HB2	2.12	0.49
1:C:310:LEU:HG	1:C:314:LYS:HD2	1.94	0.49
2:D:602:MET:O	2:D:603:PHE:CB	2.61	0.49
2:D:592:LYS:C	2:D:594:VAL:H	2.15	0.49
1:A:200:LEU:HD12	2:B:785:ASN:CG	2.34	0.49
1:A:392:LEU:O	1:A:395:THR:HB	2.13	0.49
1:C:67:GLN:HG2	1:C:425:PHE:CE1	2.48	0.49
1:A:256:LEU:O	1:A:257:GLU:HB2	2.13	0.49
2:D:688:LYS:HZ1	2:D:831:ILE:HG22	1.77	0.48
2:B:706:TRP:CD2	2:B:808:LEU:HD13	2.48	0.48
2:D:636:TYR:C	2:D:639:PRO:HD2	2.32	0.48
2:D:757:GLU:CG	2:D:760:ASN:HB3	2.38	0.48
2:D:606:TRP:C	2:D:608:GLU:H	2.15	0.48
2:D:756:GLU:O	2:D:758:GLU:HG2	2.13	0.48
1:A:103:LEU:HD11	1:A:348:ILE:HG22	1.95	0.48
1:C:108:LEU:O	1:C:112:VAL:HG23	2.14	0.48
1:A:63:PRO:CG	1:A:64:GLU:H	2.24	0.48
2:B:751:TYR:CZ	2:B:758:GLU:HB3	2.49	0.48
1:C:67:GLN:OE1	2:D:536:ARG:HA	2.13	0.48
1:C:172:VAL:HG22	1:C:172:VAL:O	2.14	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:450:LEU:O	2:D:451:ASN:CB	2.61	0.48
1:A:14:VAL:CG1	1:A:20:ALA:HA	2.41	0.48
2:D:765:ASN:HD21	2:D:766:ILE:HD11	1.78	0.48
2:D:458:ASN:OD1	2:D:460:TRP:CD1	2.67	0.48
2:D:663:ILE:N	2:D:663:ILE:CD1	2.77	0.48
2:D:830:LEU:HD13	2:D:837:LEU:CD1	2.44	0.48
2:B:570:ASN:HD21	2:B:572:VAL:HG12	1.79	0.48
1:A:326:ASP:OD2	1:A:330:LYS:HB3	2.14	0.48
2:D:607:VAL:O	2:D:607:VAL:HG12	2.13	0.48
1:C:253:MET:HG3	2:D:463:PHE:O	2.14	0.48
1:A:344:MET:HE3	2:B:502:TYR:HD2	1.78	0.48
2:B:505:THR:HG22	1:C:204:THR:HG23	1.96	0.47
1:C:154:ILE:HG22	1:C:155:GLY:N	2.28	0.47
1:A:293:ILE:HD12	8:A:2044:HOH:O	2.13	0.47
2:B:765:ASN:OD1	2:B:767:ASP:HB2	2.14	0.47
2:B:754:TYR:HB2	2:B:758:GLU:CG	2.44	0.47
2:B:495:SER:HB2	2:D:497:ASP:OD2	2.14	0.47
2:D:576:LEU:C	2:D:579:PRO:HD2	2.35	0.47
2:B:627:ILE:CD1	2:B:633:ILE:HD11	2.45	0.47
2:B:633:ILE:HD13	2:B:656:LEU:HD23	1.97	0.47
1:C:285:TYR:CZ	1:C:289:LYS:HE3	2.50	0.47
1:C:395:THR:CG2	1:C:396:ASN:N	2.78	0.47
2:D:601:ALA:HB1	2:D:602:MET:HE1	1.96	0.47
2:D:834:VAL:HG12	2:D:838:LYS:HE3	1.97	0.47
1:C:0:SER:OG	1:C:1:MET:N	2.48	0.47
2:B:464:PHE:O	2:B:465:SER:HB3	2.15	0.47
2:B:681:LEU:HD22	2:B:681:LEU:N	2.29	0.47
2:D:817:ASP:O	2:D:821:LYS:HG3	2.14	0.47
2:B:685:ILE:O	2:B:686:ALA:HB3	2.14	0.47
2:D:550:MET:HE1	2:D:731:LYS:HG2	1.96	0.47
2:B:536:ARG:HG3	2:B:536:ARG:HH11	1.79	0.47
2:D:647:LEU:O	2:D:650:ASP:CB	2.63	0.47
2:B:537:PHE:CD2	2:B:537:PHE:N	2.81	0.47
1:A:262:GLU:HG2	1:A:366:TYR:CZ	2.50	0.47
1:C:11:LYS:HE2	1:C:81:ASP:HB3	1.96	0.47
2:D:610:LEU:HD12	2:D:610:LEU:N	2.29	0.46
2:D:709:VAL:HG21	2:D:808:LEU:HD21	1.98	0.46
2:B:755:THR:O	2:B:757:GLU:N	2.48	0.46
1:C:351:GLU:O	1:C:355:VAL:HG23	2.15	0.46
1:A:252:GLU:O	2:B:462:LEU:HD13	2.15	0.46
1:C:98:ILE:O	1:C:104:GLY:HA3	2.15	0.46
2:D:833:GLN:C	8:D:2027:HOH:O	2.53	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:176:THR:HG21	1:A:231:ARG:NH1	2.30	0.46
2:B:653:VAL:O	2:B:657:ILE:HG13	2.15	0.46
2:D:499:ILE:N	2:D:499:ILE:HD13	2.30	0.46
1:A:126:GLU:HA	1:A:302:SER:O	2.15	0.46
2:D:602:MET:O	2:D:603:PHE:CD1	2.68	0.46
2:B:553:TYR:OH	2:B:647:LEU:HD13	2.15	0.46
1:A:240:ASN:C	1:A:240:ASN:HD22	2.19	0.46
1:C:284:LEU:C	1:C:284:LEU:HD13	2.36	0.46
1:A:220:THR:O	1:A:223:HIS:HB3	2.16	0.46
2:D:584:THR:HG23	2:D:739:GLN:OE1	2.16	0.46
1:C:67:GLN:HA	1:C:425:PHE:CE2	2.51	0.46
1:A:222:ALA:O	1:A:226:ILE:HG13	2.16	0.46
2:B:646:MET:O	2:B:647:LEU:HG	2.15	0.46
2:D:645:ASN:O	2:D:646:MET:O	2.33	0.46
1:A:49:ASP:HB2	1:A:154:ILE:HD12	1.98	0.46
1:C:243:PHE:CZ	1:C:273:PHE:HB3	2.51	0.46
2:D:670:GLU:O	2:D:720:LYS:NZ	2.47	0.46
2:D:550:MET:HE3	2:D:731:LYS:HE2	1.98	0.46
2:D:773:LEU:O	2:D:777:ILE:HG12	2.16	0.46
1:A:141:ASP:OD2	1:A:141:ASP:C	2.54	0.45
1:C:53:ASN:OD1	1:C:55:GLU:HB3	2.15	0.45
2:B:567:ALA:C	2:B:568:LEU:HD22	2.37	0.45
2:B:710:TYR:O	2:B:714:VAL:HG23	2.16	0.45
2:D:450:LEU:HD13	2:D:451:ASN:CA	2.47	0.45
1:A:23:LYS:HE3	2:B:522:SER:O	2.17	0.45
2:B:785:ASN:CB	2:B:861:ARG:HH22	2.28	0.45
2:D:706:TRP:CE3	2:D:808:LEU:HD13	2.51	0.45
1:A:387:TYR:CD2	1:C:0:SER:HB2	2.51	0.45
2:B:816:LYS:HA	2:B:845:LEU:HD12	1.97	0.45
2:B:567:ALA:HA	2:B:580:SER:O	2.17	0.45
2:D:861:ARG:NH2	8:D:2031:HOH:O	2.49	0.45
2:B:566:ILE:HG13	2:B:566:ILE:O	2.16	0.45
2:D:587:SER:HB3	2:D:617:GLU:OE2	2.16	0.45
2:D:680:ALA:C	2:D:681:LEU:HD12	2.37	0.45
1:A:201:GLU:O	1:A:202:VAL:O	2.35	0.45
2:D:593:LYS:O	2:D:610:LEU:HD21	2.16	0.45
2:B:560:GLU:O	2:B:562:GLY:N	2.47	0.45
2:B:831:ILE:HG23	2:B:832:GLY:N	2.31	0.45
2:D:606:TRP:C	2:D:608:GLU:N	2.70	0.45
2:D:584:THR:CG2	2:D:586:PHE:HB2	2.47	0.45
2:D:870:ILE:HG22	2:D:872:ASN:OD1	2.17	0.45
1:A:244:LYS:HG2	2:B:467:SER:OG	2.16	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:602:MET:HB3	2:D:750:GLN:NE2	2.32	0.44
2:D:637:ILE:HG12	2:D:732:MET:CE	2.40	0.44
2:D:576:LEU:O	2:D:579:PRO:HD2	2.16	0.44
2:B:464:PHE:HB3	8:B:2039:HOH:O	2.18	0.44
1:A:251:TYR:OH	2:B:667:PHE:HB3	2.16	0.44
2:D:765:ASN:O	2:D:766:ILE:C	2.54	0.44
1:C:201:GLU:OE2	1:C:201:GLU:HA	2.16	0.44
1:A:262:GLU:CG	1:A:366:TYR:CZ	3.00	0.44
2:D:614:PHE:O	2:D:618:THR:HG23	2.17	0.44
2:D:758:GLU:O	2:D:760:ASN:N	2.50	0.44
1:A:254:SER:N	2:B:459:ASN:OD1	2.43	0.44
1:C:259:SER:OG	1:C:262:GLU:HB2	2.17	0.44
2:D:506:PHE:HB3	2:D:508:PHE:CE1	2.52	0.44
2:D:575:ALA:HB1	2:D:581:ARG:CD	2.48	0.44
1:A:242:VAL:HA	1:A:259:SER:HA	2.00	0.44
2:D:764:PHE:HB2	2:D:768:ASP:CB	2.48	0.44
2:B:786:LYS:HE3	2:B:786:LYS:HB2	1.78	0.44
1:C:51:PHE:O	1:C:52:THR:C	2.56	0.44
2:D:751:TYR:O	2:D:754:TYR:HB2	2.17	0.44
2:D:702:ARG:O	2:D:705:LYS:HB3	2.18	0.44
2:D:631:THR:HB	2:D:790:GLN:HE22	1.80	0.44
1:C:67:GLN:HA	1:C:425:PHE:CZ	2.53	0.44
2:D:764:PHE:HB2	2:D:768:ASP:HB3	1.99	0.44
2:D:594:VAL:CG1	2:D:746:ILE:HG21	2.48	0.43
2:D:566:ILE:HD12	2:D:749:TYR:CD2	2.52	0.43
2:B:702:ARG:NH1	2:B:812:ASP:OD2	2.43	0.43
1:A:70:VAL:HA	1:A:418:ASN:HD21	1.83	0.43
1:C:374:PHE:CD1	1:C:415:LYS:HA	2.53	0.43
1:A:202:VAL:O	1:A:204:THR:CA	2.61	0.43
2:D:584:THR:HG21	2:D:586:PHE:HB2	2.01	0.43
2:D:746:ILE:O	2:D:749:TYR:HB3	2.18	0.43
2:D:601:ALA:C	2:D:602:MET:SD	2.97	0.43
2:B:569:THR:CG2	2:B:591:VAL:HG12	2.48	0.43
2:D:535:GLU:OE2	2:D:535:GLU:N	2.41	0.43
1:C:284:LEU:HD22	1:C:284:LEU:O	2.18	0.43
2:D:450:LEU:CD2	2:D:451:ASN:N	2.70	0.43
2:D:747:ILE:H	2:D:747:ILE:CD1	2.32	0.43
2:D:729:ARG:NH2	2:D:789:ASN:OD1	2.51	0.43
1:A:241:ARG:HD3	8:A:2054:HOH:O	2.17	0.43
2:B:726:ASP:O	2:B:729:ARG:HG3	2.18	0.43
2:D:487:ILE:CG1	2:D:488:GLU:N	2.79	0.43
2:B:550:MET:HG3	2:B:641:LEU:HB3	2.00	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:702:ARG:HD3	2:B:845:LEU:HD22	2.01	0.43
1:C:322:LEU:HD12	1:C:341:LEU:HB2	2.00	0.43
1:A:96:GLU:HA	1:A:96:GLU:OE1	2.19	0.43
2:D:739:GLN:O	2:D:743:THR:HG23	2.19	0.43
2:D:765:ASN:ND2	2:D:766:ILE:CD1	2.82	0.43
1:C:278:GLN:O	1:C:282:PHE:CD2	2.72	0.43
1:C:222:ALA:O	1:C:226:ILE:HG13	2.19	0.43
1:A:395:THR:CG2	1:A:396:ASN:N	2.82	0.42
2:D:747:ILE:N	2:D:747:ILE:CD1	2.82	0.42
2:D:673:ILE:CG2	2:D:807:ARG:HG3	2.44	0.42
2:D:530:LEU:CD1	2:D:530:LEU:N	2.79	0.42
2:B:854:SER:HA	2:B:863:LEU:HD21	2.00	0.42
2:B:530:LEU:O	2:B:531:MET:CB	2.67	0.42
2:D:645:ASN:O	2:D:646:MET:C	2.57	0.42
2:B:833:GLN:NE2	2:B:836:ARG:HH21	2.17	0.42
2:D:801:ILE:N	2:D:802:PRO:CD	2.83	0.42
2:D:824:TYR:O	2:D:827:ARG:HG2	2.19	0.42
2:B:866:PHE:O	2:B:869:TYR:HB3	2.19	0.42
2:D:601:ALA:O	2:D:602:MET:CG	2.67	0.42
2:D:834:VAL:CG1	2:D:838:LYS:HE3	2.49	0.42
1:C:39:HIS:CD2	1:C:40:ASN:H	2.37	0.42
2:B:733:LYS:O	2:B:737:GLU:HG3	2.19	0.42
2:D:602:MET:O	2:D:603:PHE:HB2	2.19	0.42
2:D:599:GLU:OE1	2:D:599:GLU:CA	2.68	0.42
2:B:536:ARG:NH1	2:B:536:ARG:HG3	2.34	0.42
2:B:702:ARG:HD3	2:B:845:LEU:CD2	2.50	0.42
1:A:117:PHE:HA	1:A:317:PHE:CE1	2.55	0.42
1:A:198:GLU:HA	1:A:198:GLU:OE2	2.19	0.42
2:B:628:ALA:O	2:B:629:ASP:HB2	2.19	0.42
2:D:820:LEU:HA	2:D:820:LEU:HD12	1.81	0.42
2:D:634:ILE:HG21	2:D:784:ILE:HD13	2.02	0.42
2:D:758:GLU:C	2:D:760:ASN:H	2.23	0.42
2:B:627:ILE:HD12	2:B:633:ILE:HD11	2.00	0.42
2:D:574:GLU:OE2	2:D:574:GLU:N	2.40	0.42
1:A:2:GLU:HG3	1:A:99:TYR:HE2	1.85	0.42
1:C:393:ARG:HA	1:C:398:ALA:HB2	2.02	0.41
2:D:765:ASN:HD21	2:D:766:ILE:CD1	2.33	0.41
1:C:379:VAL:HB	1:C:380:PRO:HD3	2.02	0.41
2:D:833:GLN:C	2:D:835:ASP:N	2.73	0.41
2:D:681:LEU:HD12	2:D:681:LEU:N	2.35	0.41
1:A:423:PHE:CD1	1:A:423:PHE:N	2.81	0.41
1:C:201:GLU:HG2	1:C:364:LYS:NZ	2.35	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:659:SER:OG	2:B:663:ILE:HD11	2.20	0.41
2:D:515:ILE:O	2:D:516:SER:CB	2.56	0.41
2:B:679:PHE:HE2	2:B:701:LYS:HB3	1.83	0.41
2:D:601:ALA:HB1	2:D:602:MET:CE	2.51	0.41
2:D:573:ASN:C	2:D:575:ALA:N	2.74	0.41
1:A:184:GLN:NE2	1:A:231:ARG:HB3	2.35	0.41
1:A:97:ARG:HA	1:A:386:ILE:HG23	2.02	0.41
2:D:450:LEU:HD13	2:D:451:ASN:CG	2.40	0.41
2:D:713:ILE:N	2:D:713:ILE:CD1	2.83	0.41
2:B:671:ILE:HD11	2:B:713:ILE:HG12	2.02	0.41
2:D:545:LEU:HD22	2:D:549:THR:CG2	2.51	0.41
1:C:170:HIS:CD2	1:C:171:GLU:HG2	2.56	0.41
1:C:235:ILE:HA	1:C:286:TYR:CE1	2.56	0.41
2:D:691:THR:HG22	2:D:830:LEU:HD11	2.03	0.41
2:D:790:GLN:O	2:D:794:SER:HB2	2.21	0.41
2:D:569:THR:HG23	2:D:584:THR:CB	2.51	0.41
2:D:691:THR:HB	2:D:830:LEU:HD21	2.03	0.41
2:B:553:TYR:CD2	2:B:642:ASN:HB2	2.56	0.41
2:B:633:ILE:CD1	2:B:656:LEU:HD23	2.51	0.41
1:A:226:ILE:HG22	1:A:230:HIS:CE1	2.56	0.41
2:B:834:VAL:HG23	2:B:835:ASP:N	2.34	0.41
2:D:566:ILE:O	2:D:567:ALA:HB2	2.21	0.41
2:B:591:VAL:HG23	2:B:592:LYS:N	2.36	0.41
2:D:684:TYR:HE1	2:D:694:THR:HG21	1.84	0.41
1:A:80:THR:HG22	1:A:81:ASP:N	2.36	0.41
1:C:69:PRO:HG3	2:D:460:TRP:CD2	2.56	0.41
1:C:242:VAL:HG12	1:C:243:PHE:N	2.36	0.41
1:C:10:TYR:HB2	1:C:85:ASP:HA	2.03	0.41
2:B:554:LEU:HD12	2:B:554:LEU:HA	1.91	0.41
2:B:859:ASN:OD1	2:B:862:LEU:HG	2.21	0.41
2:D:756:GLU:O	2:D:758:GLU:N	2.51	0.41
2:B:563:LYS:O	2:B:564:SER:C	2.60	0.41
2:B:627:ILE:O	2:B:628:ALA:HB3	2.20	0.41
1:A:117:PHE:HE1	1:A:297:LEU:HD21	1.86	0.41
1:C:102:ASP:HB2	1:C:357:PHE:CE2	2.56	0.41
2:B:859:ASN:OD1	2:B:861:ARG:HG2	2.21	0.40
2:B:627:ILE:HD11	2:B:657:ILE:HG12	2.02	0.40
1:A:322:LEU:O	1:A:337:LYS:HE3	2.20	0.40
1:A:344:MET:HG3	1:A:348:ILE:HD12	2.03	0.40
2:D:575:ALA:HB1	2:D:581:ARG:HD2	2.04	0.40
1:A:293:ILE:CD1	8:A:2044:HOH:O	2.69	0.40
2:B:775:GLU:O	2:B:779:LYS:HG3	2.21	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:660:GLY:O	2:D:663:ILE:CD1	2.45	0.40
2:D:488:GLU:O	2:D:489:ALA:HB3	2.20	0.40
2:B:565:ARG:CG	2:B:749:TYR:HB2	2.46	0.40
1:A:103:LEU:HA	1:A:106:MET:HE2	2.02	0.40
1:C:223:HIS:ND1	1:C:351:GLU:OE1	2.54	0.40
2:B:599:GLU:OE1	2:B:599:GLU:N	2.35	0.40
1:A:145:ARG:HA	2:B:519:ASN:OD1	2.21	0.40
2:D:870:ILE:O	2:D:871:LYS:CB	2.67	0.40
2:D:634:ILE:HG13	2:D:784:ILE:HD13	2.04	0.40
2:D:634:ILE:HB	2:D:637:ILE:CD1	2.50	0.40
2:D:449:ALA:N	8:D:2003:HOH:O	2.54	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	435/450 (97%)	405 (93%)	23 (5%)	7 (2%)	14	28
1	C	434/450 (96%)	408 (94%)	24 (6%)	2 (0%)	38	67
2	B	422/431 (98%)	379 (90%)	29 (7%)	14 (3%)	6	9
2	D	394/431 (91%)	341 (86%)	34 (9%)	19 (5%)	4	4
All	All	1685/1762 (96%)	1533 (91%)	110 (6%)	42 (2%)	9	14

All (42) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	65	ALA
1	A	66	LYS
1	A	67	GLN
1	A	202	VAL
1	A	203	ASP
2	B	464	PHE
2	B	564	SER

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Mol	Chain	Res	Type
2	B	759	LYS
2	D	566	ILE
2	D	602	MET
2	D	645	ASN
2	D	646	MET
2	D	759	LYS
2	D	831	ILE
2	D	834	VAL
1	A	211	GLY
2	B	530	LEU
2	B	566	ILE
2	B	756	GLU
2	B	831	ILE
1	C	74	ASP
2	D	451	ASN
2	D	644	GLY
2	D	766	ILE
2	D	871	LYS
2	B	561	HIS
2	D	490	ALA
2	D	593	LYS
2	D	651	ASP
2	D	751	TYR
1	A	74	ASP
2	B	531	MET
2	B	628	ALA
2	B	833	GLN
2	B	859	ASN
1	C	52	THR
2	B	761	ASN
2	D	516	SER
2	D	464	PHE
2	D	611	VAL
2	B	671	ILE
2	D	515	ILE

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	390/402 (97%)	380 (97%)	10 (3%)	59	85
1	C	389/402 (97%)	383 (98%)	6 (2%)	76	93
2	B	385/393 (98%)	370 (96%)	15 (4%)	43	74
2	D	372/393 (95%)	346 (93%)	26 (7%)	21	41
All	All	1536/1590 (97%)	1479 (96%)	57 (4%)	45	75

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-3	GLU
1	A	51	PHE
1	A	170	HIS
1	A	203	ASP
1	A	240	ASN
1	A	243	PHE
1	A	257	GLU
1	A	365	THR
1	A	395	THR
1	A	423	PHE
2	B	469	ASP
2	B	484	ASP
2	B	509	ASP
2	B	530	LEU
2	B	541	LYS
2	B	548	TYR
2	B	561	HIS
2	B	563	LYS
2	B	565	ARG
2	B	647	LEU
2	B	650	ASP
2	B	659	SER
2	B	761	ASN
2	B	807	ARG
2	B	869	TYR
1	C	51	PHE
1	C	165	CYS
1	C	207	LEU
1	C	240	ASN
1	C	243	PHE
1	C	284	LEU
2	D	469	ASP
2	D	486	ASN

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Mol	Chain	Res	Type
2	D	487	ILE
2	D	499	ILE
2	D	506	PHE
2	D	515	ILE
2	D	594	VAL
2	D	602	MET
2	D	604	LEU
2	D	609	GLN
2	D	610	LEU
2	D	615	THR
2	D	632	ILE
2	D	649	LYS
2	D	652	PHE
2	D	665	LEU
2	D	713	ILE
2	D	746	ILE
2	D	747	ILE
2	D	754	TYR
2	D	755	THR
2	D	758	GLU
2	D	766	ILE
2	D	820	LEU
2	D	831	ILE
2	D	863	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	GLN
1	A	40	ASN
1	A	86	ASN
1	A	170	HIS
1	A	311	GLN
2	B	501	GLN
2	B	570	ASN
2	B	722	ASN
2	B	761	ASN
2	B	763	ASN
2	B	789	ASN
2	B	843	ASN
2	B	852	GLN
1	C	29	GLN

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Mol	Chain	Res	Type
1	C	82	ASN
1	C	170	HIS
1	C	174	ASN
1	C	269	HIS
1	C	394	ASN
2	D	557	GLN
2	D	693	GLN
2	D	765	ASN
2	D	843	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 15 ligands modelled in this entry, 5 are monoatomic - leaving 10 for Mogul analysis.

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	SO4	A	1433	-	4,4,4	0.35	0	6,6,6	0.09	0
6	GOL	A	1436	-	5,5,5	4.57	5 (100%)	5,5,5	5.77	3 (60%)
6	GOL	A	1437	-	5,5,5	4.48	5 (100%)	5,5,5	5.77	3 (60%)
6	GOL	A	1438	-	5,5,5	4.57	5 (100%)	5,5,5	5.71	3 (60%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	B	1872	-	4,4,4	0.32	0	6,6,6	0.09	0
6	GOL	B	1873	-	5,5,5	4.48	5 (100%)	5,5,5	5.74	3 (60%)
6	GOL	B	1874	-	5,5,5	4.55	5 (100%)	5,5,5	5.73	3 (60%)
7	ACT	C	1435	-	1,3,3	1.61	0	0,3,3	0.00	-
3	SO4	C	1436	-	4,4,4	0.32	0	6,6,6	0.08	0
3	SO4	C	1437	-	4,4,4	0.24	0	6,6,6	0.07	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	SO4	A	1433	-	-	0/0/0/0	0/0/0/0
6	GOL	A	1436	-	-	0/4/4/4	0/0/0/0
6	GOL	A	1437	-	-	0/4/4/4	0/0/0/0
6	GOL	A	1438	-	-	0/4/4/4	0/0/0/0
3	SO4	B	1872	-	-	0/0/0/0	0/0/0/0
6	GOL	B	1873	-	-	0/4/4/4	0/0/0/0
6	GOL	B	1874	-	-	0/4/4/4	0/0/0/0
7	ACT	C	1435	-	-	0/0/0/0	0/0/0/0
3	SO4	C	1436	-	-	0/0/0/0	0/0/0/0
3	SO4	C	1437	-	-	0/0/0/0	0/0/0/0

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	1438	GOL	C3-C2	-7.60	1.20	1.52
6	B	1873	GOL	C3-C2	-7.59	1.21	1.52
6	B	1874	GOL	C3-C2	-7.56	1.21	1.52
6	A	1436	GOL	C3-C2	-7.53	1.21	1.52
6	A	1437	GOL	C3-C2	-7.49	1.21	1.52
6	B	1874	GOL	O1-C1	-4.55	1.22	1.42
6	A	1438	GOL	O1-C1	-4.49	1.22	1.42
6	A	1436	GOL	O1-C1	-4.45	1.23	1.42
6	A	1437	GOL	O1-C1	-4.26	1.23	1.42
6	B	1873	GOL	O1-C1	-4.20	1.24	1.42
6	A	1436	GOL	O3-C3	3.33	1.56	1.42
6	B	1873	GOL	O3-C3	3.32	1.56	1.42
6	A	1437	GOL	O3-C3	3.24	1.56	1.42
6	B	1874	GOL	O3-C3	3.21	1.56	1.42
6	A	1438	GOL	O3-C3	3.18	1.56	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	1436	GOL	C1-C2	-3.01	1.39	1.52
6	A	1438	GOL	C1-C2	-2.98	1.40	1.52
6	B	1874	GOL	C1-C2	-2.91	1.40	1.52
6	A	1437	GOL	C1-C2	-2.88	1.40	1.52
6	A	1436	GOL	O2-C2	-2.78	1.34	1.43
6	A	1438	GOL	O2-C2	-2.75	1.35	1.43
6	B	1873	GOL	O2-C2	-2.73	1.35	1.43
6	A	1437	GOL	O2-C2	-2.71	1.35	1.43
6	B	1874	GOL	O2-C2	-2.63	1.35	1.43
6	B	1873	GOL	C1-C2	-2.60	1.41	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	1437	GOL	O3-C3-C2	10.52	161.04	109.71
6	A	1436	GOL	O3-C3-C2	10.51	160.97	109.71
6	B	1873	GOL	O3-C3-C2	10.40	160.43	109.71
6	A	1438	GOL	O3-C3-C2	10.31	160.00	109.71
6	B	1874	GOL	O3-C3-C2	10.29	159.91	109.71
6	B	1874	GOL	O2-C2-C3	6.80	139.21	108.22
6	A	1438	GOL	O2-C2-C3	6.75	138.96	108.22
6	A	1436	GOL	O2-C2-C3	6.70	138.76	108.22
6	B	1873	GOL	O2-C2-C3	6.60	138.28	108.22
6	A	1437	GOL	O2-C2-C3	6.58	138.19	108.22
6	B	1873	GOL	O1-C1-C2	3.53	126.93	109.71
6	A	1437	GOL	O1-C1-C2	3.43	126.45	109.71
6	B	1874	GOL	O1-C1-C2	3.31	125.89	109.71
6	A	1438	GOL	O1-C1-C2	3.29	125.77	109.71
6	A	1436	GOL	O1-C1-C2	3.27	125.66	109.71

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	437/450 (97%)	-0.01	9 (2%) 60 58	12, 27, 54, 98	0
1	C	436/450 (96%)	0.00	9 (2%) 60 58	12, 30, 59, 75	0
2	B	424/431 (98%)	0.59	42 (9%) 8 6	20, 42, 78, 93	0
2	D	408/431 (94%)	1.12	78 (19%) 2 1	21, 48, 101, 117	0
All	All	1705/1762 (96%)	0.41	138 (8%) 12 9	12, 36, 78, 117	0

All (138) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	873	ILE	14.0
2	D	764	PHE	9.3
2	D	872	ASN	8.4
2	D	603	PHE	8.3
2	D	606	TRP	8.2
2	D	602	MET	8.2
2	D	753	GLN	7.7
2	D	755	THR	7.7
2	D	566	ILE	7.1
2	B	760	ASN	6.9
2	B	565	ARG	6.6
2	D	649	LYS	6.4
2	D	831	ILE	6.4
2	D	601	ALA	6.3
2	B	761	ASN	5.9
2	D	754	TYR	5.9
2	D	751	TYR	5.8
2	D	644	GLY	5.8
2	D	604	LEU	5.8
2	D	761	ASN	5.7
2	B	566	ILE	5.5

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Mol	Chain	Res	Type	RSRZ
2	D	830	LEU	5.4
2	B	563	LYS	5.4
2	B	647	LEU	5.2
1	C	0	SER	5.1
2	D	600	ALA	4.9
2	D	605	GLY	4.8
2	B	568	LEU	4.6
2	D	645	ASN	4.5
2	D	569	THR	4.5
2	B	764	PHE	4.5
2	D	565	ARG	4.4
2	B	537	PHE	4.4
2	D	607	VAL	4.3
2	B	534	ILE	4.3
2	D	534	ILE	4.3
2	D	837	LEU	4.2
1	A	-3	GLU	4.2
2	D	599	GLU	4.1
1	C	64	GLU	4.1
2	B	763	ASN	4.1
2	D	581	ARG	4.1
2	D	658	PHE	4.0
2	D	765	ASN	4.0
2	D	750	GLN	4.0
2	D	568	LEU	3.9
2	D	747	ILE	3.9
1	A	65	ALA	3.8
2	B	755	THR	3.8
2	D	647	LEU	3.8
2	D	752	ASN	3.8
1	A	64	GLU	3.8
2	D	561	HIS	3.8
2	D	833	GLN	3.7
2	B	762	ILE	3.7
2	D	766	ILE	3.7
2	B	577	LEU	3.7
2	B	831	ILE	3.6
1	C	423	PHE	3.5
2	D	533	ASN	3.4
2	D	749	TYR	3.4
2	B	751	TYR	3.4
2	D	746	ILE	3.4

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Mol	Chain	Res	Type	RSRZ
2	B	530	LEU	3.4
2	D	631	THR	3.4
1	C	63	PRO	3.3
2	B	759	LYS	3.3
2	D	608	GLU	3.3
2	D	760	ASN	3.2
2	B	754	TYR	3.2
2	B	837	LEU	3.2
2	B	562	GLY	3.2
1	A	423	PHE	3.2
2	B	648	TYR	3.1
2	D	829	THR	3.1
2	D	836	ARG	3.1
2	D	580	SER	3.1
2	D	869	TYR	3.0
2	B	564	SER	3.0
2	D	646	MET	2.9
2	D	656	LEU	2.9
2	B	834	VAL	2.9
2	D	657	ILE	2.9
2	B	753	GLN	2.8
1	A	1	MET	2.8
2	D	685	ILE	2.8
2	B	561	HIS	2.8
2	D	835	ASP	2.7
2	B	752	ASN	2.7
2	D	567	ALA	2.7
1	C	422	LEU	2.7
2	B	756	GLU	2.7
1	A	425	PHE	2.7
2	D	537	PHE	2.7
2	B	533	ASN	2.7
2	D	591	VAL	2.7
2	B	803	TYR	2.7
2	D	612	TYR	2.7
2	D	592	LYS	2.6
2	D	759	LYS	2.6
2	B	806	LYS	2.5
2	D	594	VAL	2.5
1	A	257	GLU	2.5
2	B	870	ILE	2.5
2	B	603	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
2	D	819	LEU	2.5
2	D	632	ILE	2.4
2	D	530	LEU	2.4
2	B	597	ALA	2.4
2	D	487	ILE	2.4
1	C	425	PHE	2.3
2	B	690	LEU	2.3
1	A	200	LEU	2.3
2	D	489	ALA	2.2
2	D	742	ALA	2.2
2	D	633	ILE	2.2
2	D	643	ILE	2.2
2	D	682	VAL	2.2
2	B	758	GLU	2.2
2	D	535	GLU	2.2
2	B	626	LYS	2.2
2	B	629	ASP	2.2
1	A	202	VAL	2.2
2	D	772	LYS	2.2
2	B	599	GLU	2.2
2	D	610	LEU	2.2
1	C	254	SER	2.2
2	D	588	SER	2.2
2	D	515	ILE	2.2
2	D	531	MET	2.1
2	D	650	ASP	2.1
2	B	869	TYR	2.1
1	C	172	VAL	2.1
2	D	686	ALA	2.1
2	D	768	ASP	2.1
1	C	171	GLU	2.1
2	B	833	GLN	2.0
2	B	481	ILE	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
6	GOL	A	1438	6/6	0.17	2.88	54,55,56,58	0
7	ACT	C	1435	4/4	0.18	2.78	49,50,51,52	0
6	GOL	A	1437	6/6	0.20	2.72	44,46,47,49	0
6	GOL	B	1874	6/6	0.16	2.07	55,59,59,60	0
6	GOL	B	1873	6/6	0.20	1.96	51,53,55,56	0
5	CL	A	1439	1/1	0.15	1.61	44,44,44,44	0
3	SO4	A	1433	5/5	0.18	1.54	68,69,70,72	0
6	GOL	A	1436	6/6	0.16	1.21	54,55,56,58	0
4	ZN	C	1438	1/1	0.18	0.15	29,29,29,29	0
3	SO4	C	1436	5/5	0.16	0.09	82,83,85,85	0
5	CL	A	1435	1/1	0.10	-0.73	44,44,44,44	0
3	SO4	C	1437	5/5	0.11	-0.84	69,69,70,70	0
4	ZN	A	1434	1/1	0.14	-1.26	23,23,23,23	0
3	SO4	B	1872	5/5	0.12	-2.12	47,48,50,51	0
5	CL	D	1872	1/1	0.06	-2.35	47,47,47,47	0

6.5 Other polymers ⓘ

There are no such residues in this entry.