



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:07 PM UTC

PDB ID : 3NB2 / pdb_00003nb2
Title : Crystal structure of E. coli O157:H7 effector protein NleL
Authors : Lin, D.Y.; Chen, J.
Deposited on : 2010-06-02
Resolution : 2.10 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

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

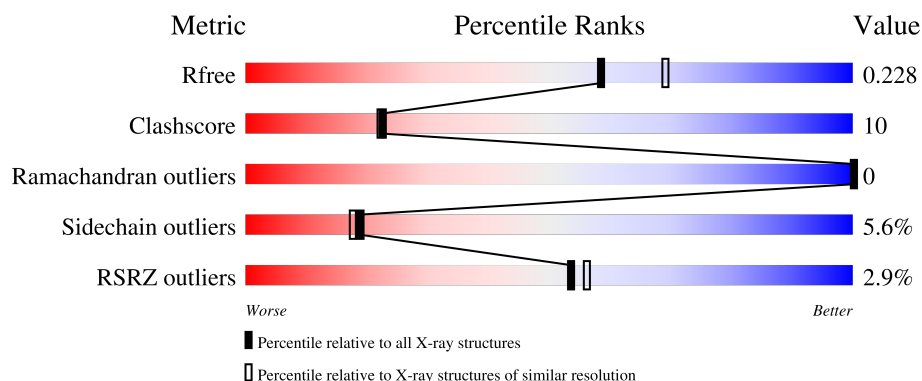
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	613	5% 77% 20% .
1	B	613	2% 79% 18% .
1	C	613	3% 81% 16% .
1	D	613	2% 78% 20% .

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GOL	C	11	-	-	X	-
2	GOL	C	21	-	-	X	-
5	DTT	A	785	-	-	X	-

2 Entry composition [i](#)

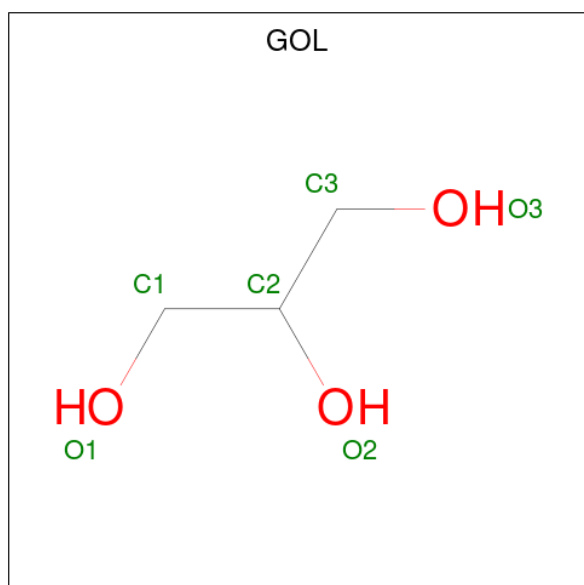
There are 6 unique types of molecules in this entry. The entry contains 20830 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 secreted effector protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	612	Total	C	N	O	S	0	0	0
			4842	3066	785	953	38			
1	B	612	Total	C	N	O	S	0	2	0
			4850	3072	788	952	38			
1	C	612	Total	C	N	O	S	0	0	0
			4846	3068	788	952	38			
1	D	612	Total	C	N	O	S	0	1	0
			4864	3082	789	955	38			

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

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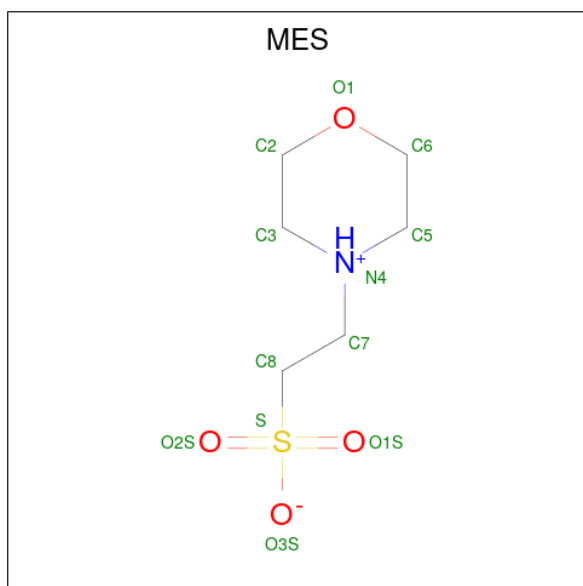
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total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

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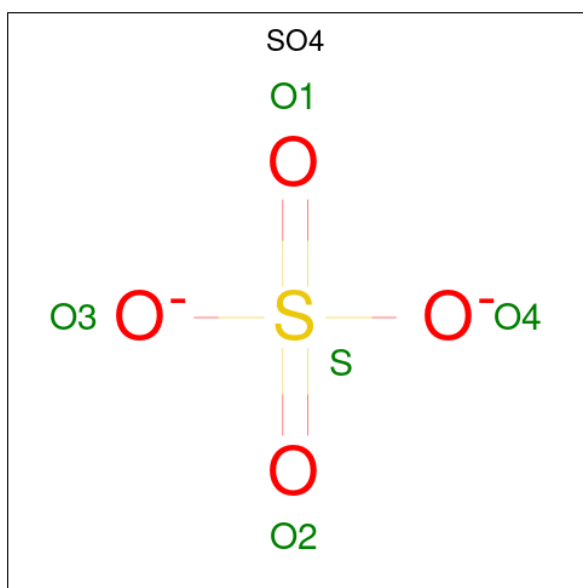
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	C	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	C	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	D	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	D	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



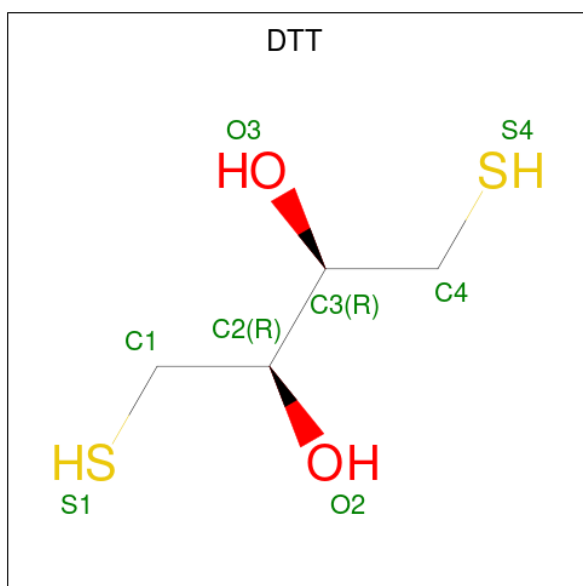
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is 2,3-DIHYDROXY-1,4-DITHIOBUTANE (CCD ID: DTT) (formula: $C_4H_{10}O_2S_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	O	S	0	0
			8	4	2	2		
5	C	1	Total	C	O	S	0	0
			8	4	2	2		

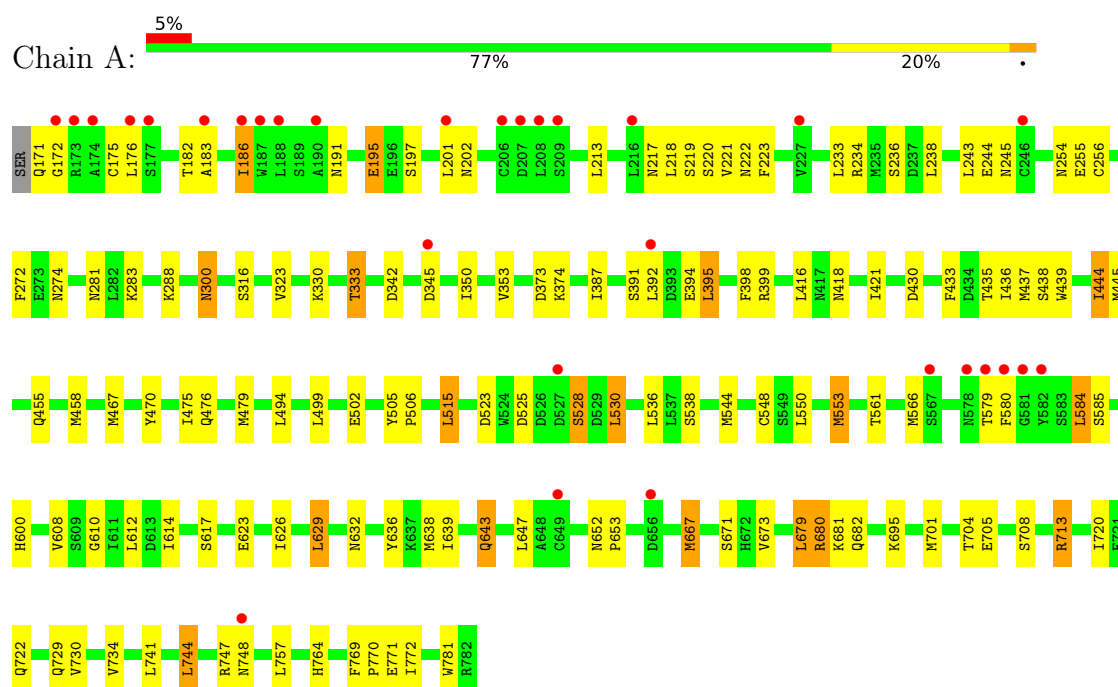
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	328	Total	O	0	0
			328	328		
6	B	273	Total	O	0	0
			273	273		
6	C	270	Total	O	0	0
			270	270		
6	D	227	Total	O	0	0
			227	227		

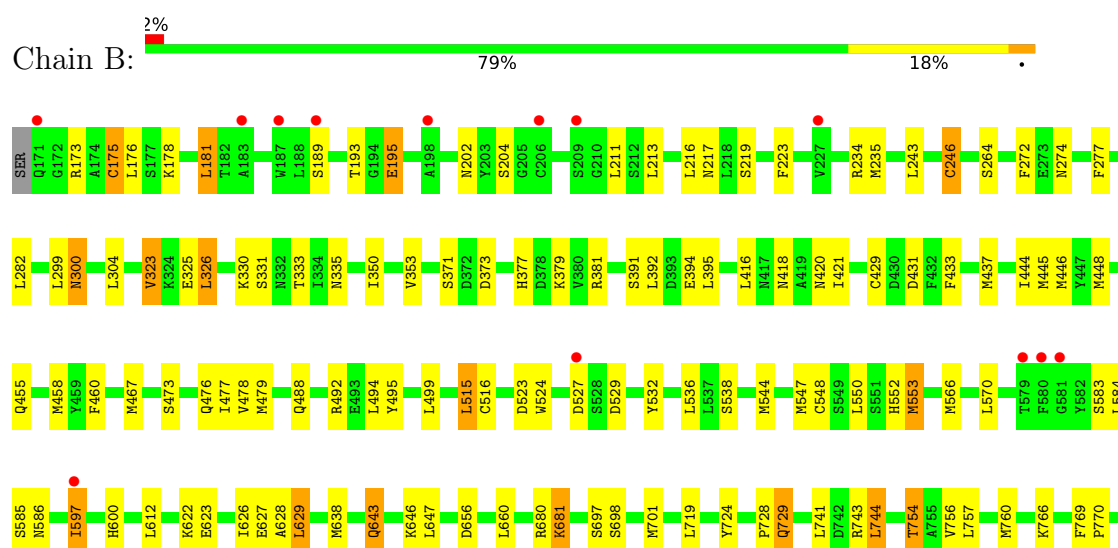
3 Residue-property plots

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

• Molecule 1: secreted effector protein

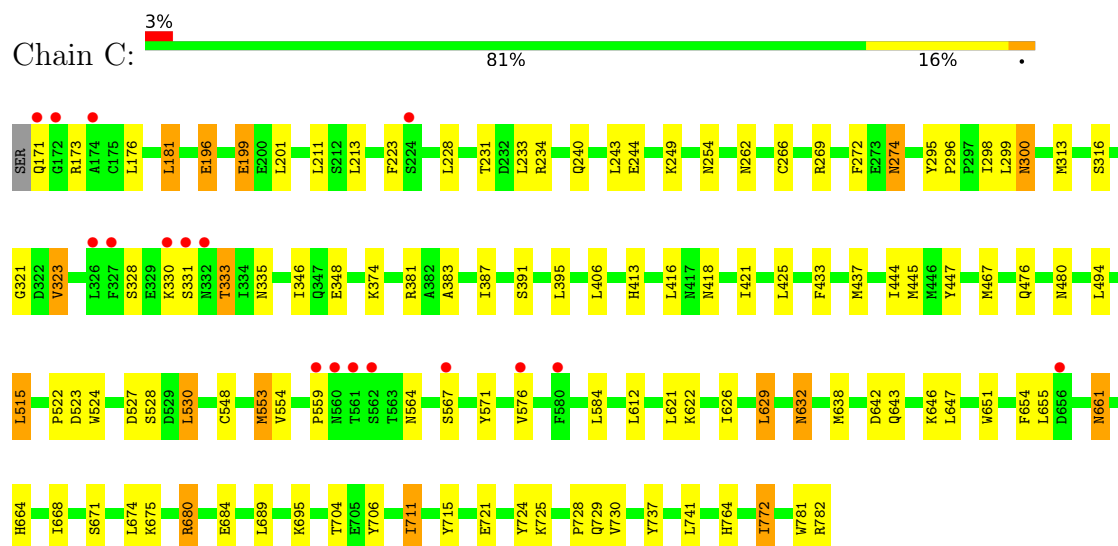


• Molecule 1: secreted effector protein

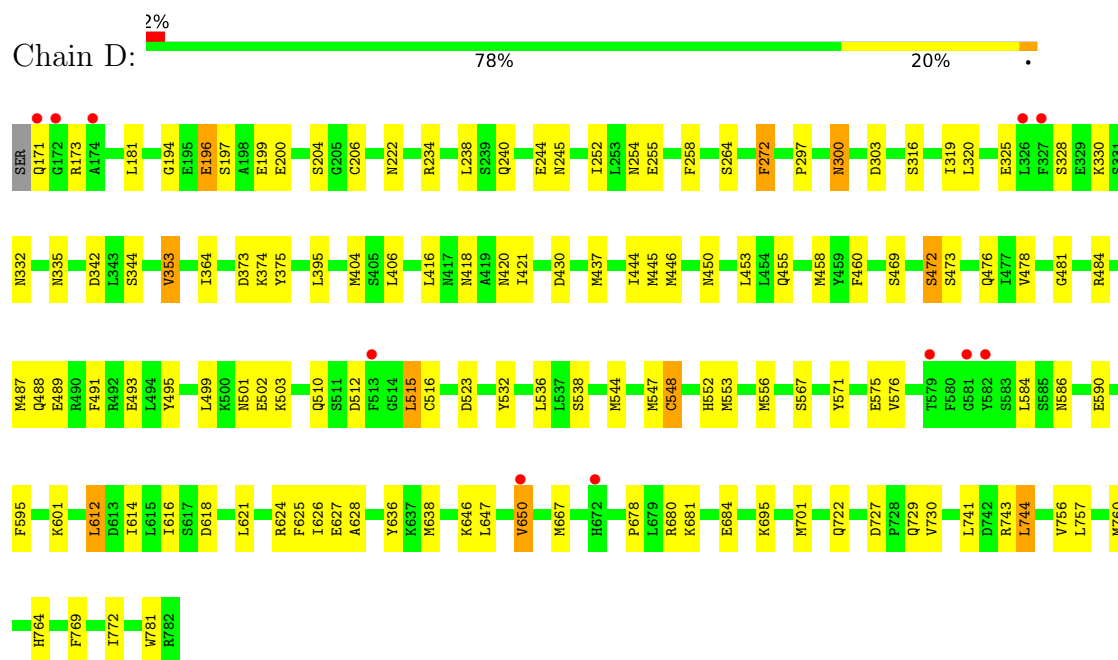


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- Molecule 1: secreted effector protein



- Molecule 1: secreted effector protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	177.01Å 124.20Å 154.90Å 90.00° 107.60° 90.00°	Depositor
Resolution (Å)	44.74 – 2.10 44.74 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.5 (44.74-2.10) 99.7 (44.74-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.48 (at 2.10Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, R_{free}	0.184 , 0.231 0.181 , 0.228	Depositor DCC
R_{free} test set	9326 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	35.4	Xtriage
Anisotropy	0.117	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 50.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	20830	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.25% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, GOL, MES, DTT

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	0.56	3/4946 (0.1%)	0.77	1/6695 (0.0%)
1	B	0.51	2/4960 (0.0%)	0.78	2/6713 (0.0%)
1	C	0.52	0/4950	0.75	0/6699
1	D	0.47	2/4972 (0.0%)	0.73	2/6727 (0.0%)
All	All	0.51	7/19828 (0.0%)	0.76	5/26834 (0.0%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	431	ASP	C-O	-5.89	1.17	1.24
1	A	720	ILE	C-O	-5.78	1.17	1.24
1	A	722	GLN	C-O	-5.55	1.17	1.24
1	D	430	ASP	C-O	-5.54	1.17	1.24
1	B	429	CYS	C-O	-5.28	1.17	1.24
1	A	430	ASP	C-O	-5.24	1.18	1.24
1	D	722	GLN	C-O	-5.03	1.18	1.24

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	580	PHE	N-CA-C	-5.69	108.13	114.62
1	D	616	ILE	N-CA-C	5.42	115.75	108.17
1	B	769	PHE	CA-C-N	5.37	125.01	119.05
1	B	769	PHE	C-N-CA	5.37	125.01	119.05
1	D	548	CYS	N-CA-C	5.33	115.58	108.38

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4842	0	4638	105	0
1	B	4850	0	4648	102	0
1	C	4846	0	4644	88	0
1	D	4864	0	4673	108	0
2	A	78	0	104	9	0
2	B	24	0	32	0	0
2	C	42	0	56	11	0
2	D	30	0	40	4	0
3	A	12	0	12	1	0
3	C	24	0	24	2	0
3	D	24	0	24	6	0
4	A	20	0	0	0	0
4	B	15	0	0	0	0
4	C	25	0	0	0	0
4	D	20	0	0	0	0
5	A	8	0	10	8	0
5	C	8	0	10	2	0
6	A	328	0	0	3	0
6	B	273	0	0	8	0
6	C	270	0	0	5	0
6	D	227	0	0	4	0
All	All	20830	0	18915	405	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 10.

All (405) 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:437:MET:HE3	1:B:473:SER:HB3	1.23	1.17
1:D:437:MET:HE3	1:D:473:SER:HB3	1.22	1.13
1:B:757:LEU:HA	1:B:760:MET:HE3	1.22	1.12
1:B:597:ILE:HD12	1:B:597:ILE:H	1.19	1.06
1:D:756:VAL:HG12	1:D:760:MET:HE2	1.33	1.06
1:B:756:VAL:HG12	1:B:760:MET:HE2	1.31	1.05

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:171:GLN:N	1:A:172:GLY:HA3	1.76	0.99
1:D:487:MET:HE3	1:D:491:PHE:CE2	1.98	0.98
1:D:374[B]:LYS:H	1:D:374[B]:LYS:HD2	1.31	0.96
1:A:771:GLU:HG2	2:A:16:GOL:O2	1.66	0.95
1:D:757:LEU:HA	1:D:760:MET:HE3	1.47	0.94
1:A:638:MET:HA	1:A:643:GLN:HG2	1.52	0.92
1:A:680:ARG:HD3	1:A:729:GLN:HE21	1.35	0.91
1:D:453:LEU:HG	1:D:487:MET:HE1	1.53	0.90
1:A:643:GLN:HA	1:A:643:GLN:HE21	1.37	0.89
1:D:437:MET:CE	1:D:473:SER:HB3	2.03	0.89
1:A:617:SER:HB3	2:A:14:GOL:H12	1.55	0.88
1:A:769:PHE:HA	2:A:16:GOL:H2	1.55	0.88
1:B:437:MET:CE	1:B:473:SER:HB3	2.04	0.88
1:B:638:MET:HA	1:B:643:GLN:HG2	1.56	0.87
1:D:330:LYS:HD3	1:D:332:ASN:H	1.40	0.85
1:C:548:CYS:SG	1:C:553:MET:HG3	2.17	0.84
1:C:299:LEU:H	2:C:11:GOL:H32	1.42	0.84
1:D:437:MET:HE3	1:D:473:SER:CB	2.07	0.83
1:B:418:ASN:HD22	1:B:421:ILE:H	1.27	0.83
1:D:437:MET:HE2	1:D:460:PHE:CZ	2.15	0.82
1:B:437:MET:HE2	1:B:460:PHE:CZ	2.13	0.82
1:D:437:MET:HE2	1:D:460:PHE:CE2	2.13	0.82
1:A:182:THR:O	1:A:186:ILE:HG22	1.77	0.82
1:A:636:TYR:HE2	1:A:638:MET:HE2	1.45	0.81
1:B:323:VAL:HG21	1:B:326:LEU:HD12	1.60	0.81
1:D:536:LEU:HD11	1:D:553:MET:HE1	1.60	0.81
1:C:480:ASN:OD1	2:C:21:GOL:H31	1.80	0.80
1:A:418:ASN:HD22	1:A:421:ILE:H	1.32	0.78
1:B:437:MET:HE2	1:B:460:PHE:CE2	2.19	0.78
1:B:330:LYS:O	1:B:331:SER:HB3	1.84	0.77
1:B:757:LEU:CA	1:B:760:MET:HE3	2.08	0.77
1:C:418:ASN:HD22	1:C:421:ILE:H	1.30	0.77
1:D:418:ASN:HD22	1:D:421:ILE:H	1.32	0.77
1:B:437:MET:HE3	1:B:473:SER:CB	2.12	0.76
1:C:782:ARG:HH22	2:C:23:GOL:H31	1.51	0.76
1:C:196:GLU:HG3	1:C:201:LEU:HD21	1.66	0.75
1:B:680:ARG:CZ	1:B:729:GLN:HG2	2.16	0.75
1:C:664:HIS:CE1	1:C:668:ILE:HD11	2.23	0.74
1:B:743:ARG:HD3	1:B:760:MET:HE1	1.70	0.74
1:A:502:GLU:HG2	6:A:826:HOH:O	1.87	0.73
1:B:467:MET:HE2	1:B:494:LEU:HD11	1.69	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:515:LEU:HD13	1:D:523:ASP:HB2	1.70	0.73
1:C:447:TYR:HD1	2:C:21:GOL:H2	1.54	0.72
1:A:614:ILE:O	2:A:14:GOL:H2	1.89	0.72
1:C:480:ASN:CG	2:C:21:GOL:H31	2.14	0.72
1:A:680:ARG:CD	1:A:729:GLN:HE21	2.02	0.72
1:D:330:LYS:HE2	1:D:332:ASN:HB2	1.69	0.72
1:C:299:LEU:H	2:C:11:GOL:C3	2.03	0.71
1:B:479:MET:HE3	1:B:550:LEU:HG	1.73	0.71
1:C:447:TYR:CD1	2:C:21:GOL:H2	2.25	0.70
5:A:785:DTT:H41	5:A:785:DTT:S1	2.31	0.70
1:B:643:GLN:HE21	1:B:643:GLN:HA	1.56	0.69
1:D:743:ARG:NH1	1:D:760:MET:HE1	2.07	0.69
1:D:453:LEU:CG	1:D:487:MET:HE1	2.23	0.69
1:D:300:ASN:H	1:D:300:ASN:HD22	1.40	0.69
1:C:559:PRO:HD2	6:C:858:HOH:O	1.93	0.69
5:C:785:DTT:S1	5:C:785:DTT:S4	2.91	0.68
1:A:244:GLU:HG2	1:A:245:ASN:OD1	1.95	0.67
1:A:548:CYS:SG	1:A:553:MET:HE2	2.35	0.67
1:A:747:ARG:O	1:A:748:ASN:HB2	1.95	0.67
1:B:467:MET:HE3	6:B:1182:HOH:O	1.96	0.66
1:A:373:ASP:CG	3:C:2:MES:H62	2.21	0.66
1:B:445:MET:H	1:B:476:GLN:NE2	1.94	0.66
1:D:374[B]:LYS:H	1:D:374[B]:LYS:CD	2.08	0.66
1:B:300:ASN:H	1:B:300:ASN:HD22	1.44	0.65
1:B:213:LEU:HB3	1:B:216:LEU:HD12	1.77	0.65
1:D:300:ASN:HD22	1:D:300:ASN:N	1.95	0.65
1:A:680:ARG:HD3	1:A:729:GLN:NE2	2.08	0.65
1:D:743:ARG:HD3	1:D:760:MET:HE1	1.80	0.64
1:A:515:LEU:HD13	1:A:523:ASP:HA	1.80	0.64
1:C:704:THR:HG23	1:C:706:TYR:H	1.61	0.64
1:A:300:ASN:HD22	1:A:300:ASN:H	1.46	0.64
1:C:300:ASN:HD22	1:C:300:ASN:H	1.44	0.63
1:C:643:GLN:HE21	1:C:643:GLN:HA	1.62	0.63
1:D:181:LEU:HD21	1:D:206:CYS:HB3	1.79	0.63
1:D:757:LEU:CA	1:D:760:MET:HE3	2.23	0.63
1:A:171:GLN:N	1:A:172:GLY:CA	2.59	0.63
1:B:698:SER:HA	1:B:754:THR:HG22	1.80	0.63
1:C:680:ARG:HD2	1:C:729:GLN:OE1	1.97	0.63
1:A:218:LEU:HB2	1:A:238:LEU:HD23	1.80	0.62
1:D:418:ASN:ND2	1:D:421:ILE:H	1.96	0.62
5:A:785:DTT:S1	5:A:785:DTT:C4	2.86	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:680:ARG:NH2	1:B:729:GLN:HG2	2.13	0.62
1:A:643:GLN:HA	1:A:643:GLN:NE2	2.13	0.62
1:D:501:ASN:OD1	1:D:503:LYS:HD2	1.99	0.62
1:C:655:LEU:HD12	1:C:711:ILE:HD11	1.82	0.62
1:D:238:LEU:HD22	1:D:258:PHE:CE1	2.35	0.61
1:D:636:TYR:HE2	1:D:638:MET:HE2	1.65	0.61
1:A:730:VAL:O	1:A:764:HIS:HE1	1.83	0.61
1:D:445:MET:H	1:D:476:GLN:NE2	1.98	0.61
1:C:571:TYR:CE1	1:C:576:VAL:HG22	2.36	0.61
1:A:467:MET:CE	1:A:494:LEU:HD11	2.30	0.61
1:D:244:GLU:HG3	1:D:264:SER:HB2	1.81	0.61
1:D:646:LYS:O	1:D:650:VAL:HG12	2.01	0.61
1:D:512:ASP:HB2	1:D:547:MET:HE1	1.82	0.61
1:B:492:ARG:HD2	1:B:524:TRP:CD2	2.36	0.61
1:C:211:LEU:O	1:C:231:THR:HG23	2.01	0.61
1:A:744:LEU:HD13	1:A:757:LEU:CD2	2.31	0.60
1:C:643:GLN:NE2	1:C:646:LYS:HE2	2.16	0.60
1:B:492:ARG:HD2	1:B:524:TRP:CG	2.35	0.60
1:C:298:ILE:HA	2:C:11:GOL:H32	1.84	0.60
1:D:374[B]:LYS:NZ	3:D:783:MES:H21	2.15	0.60
1:D:445:MET:H	1:D:476:GLN:HE22	1.48	0.60
1:B:444:ILE:HA	1:B:476:GLN:HE22	1.66	0.60
1:A:435:THR:HG22	5:A:785:DTT:S1	2.42	0.60
1:C:418:ASN:ND2	1:C:421:ILE:H	1.99	0.60
1:B:445:MET:H	1:B:476:GLN:HE22	1.50	0.60
1:D:234:ARG:HG2	1:D:254:ASN:HB2	1.84	0.60
1:D:374[B]:LYS:HZ3	3:D:783:MES:H21	1.67	0.60
1:D:437:MET:HE2	1:D:460:PHE:HE2	1.67	0.59
1:A:744:LEU:HD13	1:A:757:LEU:HD23	1.84	0.59
1:C:651:TRP:HB3	1:C:711:ILE:HD12	1.84	0.59
1:D:680:ARG:O	1:D:684:GLU:HG3	2.03	0.59
1:C:730:VAL:O	1:C:764:HIS:HE1	1.86	0.59
1:D:744:LEU:HD13	1:D:757:LEU:CD2	2.33	0.59
1:A:221:VAL:HG11	1:A:223:PHE:CZ	2.38	0.59
1:A:636:TYR:CE2	1:A:638:MET:HE2	2.32	0.58
1:B:437:MET:HE2	1:B:460:PHE:HZ	1.63	0.58
1:D:373:ASP:CG	3:D:783:MES:H62	2.29	0.58
1:A:479:MET:CE	1:A:553:MET:HE3	2.34	0.58
1:C:664:HIS:HE1	1:C:715:TYR:OH	1.85	0.58
1:C:300:ASN:HD22	1:C:300:ASN:N	2.02	0.58
1:D:330:LYS:CD	1:D:332:ASN:H	2.13	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:SER:O	1:A:220:SER:HB2	2.03	0.57
1:B:418:ASN:ND2	1:B:421:ILE:H	1.99	0.57
1:D:255:GLU:OE2	3:D:4:MES:H31	2.05	0.57
1:D:556:MET:HE2	1:D:595:PHE:CD1	2.39	0.57
1:D:234:ARG:HG3	1:D:252:ILE:HG22	1.87	0.57
1:D:538:SER:HB2	1:D:544:MET:HB2	1.87	0.57
1:C:621:LEU:HD11	1:C:638:MET:HE1	1.86	0.57
1:B:597:ILE:HD12	1:B:597:ILE:N	2.05	0.57
1:A:394:GLU:O	1:A:399:ARG:NE	2.38	0.57
1:C:467:MET:HE2	1:C:494:LEU:HD11	1.87	0.57
1:C:571:TYR:HE1	1:C:576:VAL:HG22	1.68	0.57
1:B:223:PHE:HB2	1:B:243:LEU:HD23	1.86	0.57
1:A:528:SER:OG	1:A:530:LEU:HB2	2.05	0.57
1:B:467:MET:CE	1:B:494:LEU:HD11	2.34	0.56
1:B:515:LEU:HD13	1:B:523:ASP:HB2	1.88	0.56
5:C:785:DTT:S4	5:C:785:DTT:C1	2.93	0.56
1:D:238:LEU:HD22	1:D:258:PHE:HE1	1.69	0.56
1:D:756:VAL:HG12	1:D:760:MET:CE	2.22	0.56
1:C:632:ASN:H	1:C:632:ASN:HD22	1.51	0.56
1:D:515:LEU:HD13	1:D:523:ASP:CB	2.34	0.56
1:C:445:MET:H	1:C:476:GLN:NE2	2.04	0.56
1:C:173:ARG:HD2	6:C:976:HOH:O	2.04	0.56
1:D:472:SER:HB3	6:D:903:HOH:O	2.06	0.56
1:C:528:SER:HB2	1:C:530:LEU:HD23	1.88	0.55
1:A:223:PHE:HB2	1:A:243:LEU:HD23	1.89	0.55
1:B:743:ARG:CD	1:B:760:MET:HE1	2.36	0.55
1:A:213:LEU:HB2	1:A:233:LEU:HD23	1.88	0.55
1:A:695:LYS:HG2	1:A:781:TRP:CE3	2.42	0.55
1:B:766:LYS:O	1:B:770:PRO:HG3	2.07	0.55
1:A:708:SER:OG	1:A:713:ARG:HD3	2.06	0.55
1:A:255:GLU:OE2	3:A:3:MES:H31	2.06	0.54
1:B:189:SER:HB2	1:B:217:ASN:H	1.72	0.54
1:C:684:GLU:OE1	1:C:772:ILE:HD13	2.07	0.54
1:B:681:LYS:HE2	6:B:896:HOH:O	2.08	0.54
1:B:274:ASN:HD21	1:D:316:SER:HB3	1.73	0.54
1:B:698:SER:CA	1:B:754:THR:HG22	2.38	0.54
1:D:437:MET:HE2	1:D:460:PHE:HZ	1.70	0.54
1:B:193:THR:HG22	1:B:195:GLU:H	1.73	0.53
1:A:439:TRP:CB	5:A:785:DTT:H11	2.39	0.53
1:C:171:GLN:N	1:C:199:GLU:O	2.42	0.53
1:A:730:VAL:HG22	1:A:769:PHE:CE2	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:743:ARG:NH1	1:B:760:MET:HE1	2.23	0.53
1:B:418:ASN:HD21	1:B:420:ASN:HB3	1.74	0.53
1:A:467:MET:CE	1:A:494:LEU:CD1	2.87	0.53
1:A:610:GLY:HA3	2:A:7:GOL:H12	1.91	0.53
1:A:330:LYS:HB2	1:A:333:THR:HG23	1.91	0.52
1:C:444:ILE:HA	1:C:476:GLN:HE22	1.74	0.52
1:D:244:GLU:HG2	1:D:245:ASN:ND2	2.24	0.52
1:D:320:LEU:HB2	1:D:375:TYR:CE1	2.44	0.52
1:D:727:ASP:OD1	1:D:729:GLN:HG2	2.09	0.52
1:A:479:MET:HE1	1:A:553:MET:HE3	1.92	0.52
1:D:744:LEU:HD13	1:D:757:LEU:HD23	1.92	0.52
1:C:413:HIS:HD2	6:C:841:HOH:O	1.92	0.52
1:B:628:ALA:HB1	1:B:701:MET:HE1	1.92	0.52
1:D:515:LEU:O	1:D:516:CYS:HB2	2.09	0.52
1:B:585:SER:OG	1:B:600:HIS:HE1	1.93	0.52
1:C:654:PHE:CD2	1:C:664:HIS:CD2	2.98	0.52
1:D:510:GLN:HG3	1:D:571:TYR:CE1	2.44	0.52
1:C:651:TRP:CB	1:C:711:ILE:HD12	2.40	0.51
1:C:173:ARG:HB2	1:C:176:LEU:HD13	1.92	0.51
1:B:529:ASP:HA	6:B:934:HOH:O	2.09	0.51
1:D:501:ASN:HD21	1:D:503:LYS:NZ	2.08	0.51
1:D:681:LYS:HD2	1:D:772:ILE:HD11	1.93	0.51
1:B:326:LEU:O	1:B:381[A]:ARG:NH1	2.44	0.51
1:A:391:SER:OG	1:A:394:GLU:HG3	2.11	0.51
1:A:467:MET:HE3	1:A:494:LEU:HD11	1.93	0.51
1:A:467:MET:HE3	6:A:924:HOH:O	2.11	0.51
1:D:501:ASN:HD21	1:D:503:LYS:HZ2	1.59	0.51
1:A:445:MET:H	1:A:476:GLN:NE2	2.09	0.51
1:C:622:LYS:HE2	6:C:1223:HOH:O	2.10	0.51
1:A:183:ALA:HA	1:A:186:ILE:CG2	2.41	0.51
1:A:438:SER:HB3	1:A:470:TYR:CZ	2.46	0.50
1:A:175:CYS:O	1:A:176:LEU:HD12	2.12	0.50
1:A:444:ILE:HA	1:A:476:GLN:HE22	1.76	0.50
1:C:330:LYS:HB2	1:C:333:THR:HG23	1.93	0.50
1:D:173:ARG:HD2	6:D:942:HOH:O	2.12	0.50
1:C:299:LEU:N	2:C:11:GOL:H32	2.18	0.50
1:D:614:ILE:HD13	1:D:667:MET:SD	2.50	0.50
1:B:455:GLN:HA	1:B:458:MET:CE	2.42	0.50
1:A:218:LEU:HB2	1:A:238:LEU:CD2	2.42	0.50
1:A:300:ASN:HD22	1:A:300:ASN:N	2.09	0.50
1:B:515:LEU:HD13	1:B:523:ASP:CB	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:481:GLY:HA2	1:D:487:MET:HE2	1.94	0.50
1:C:196:GLU:CG	1:C:201:LEU:HD21	2.37	0.50
1:D:197:SER:OG	1:D:200:GLU:HG3	2.11	0.50
1:C:299:LEU:HD11	1:C:313:MET:HE2	1.94	0.50
1:C:626:ILE:O	1:C:629:LEU:HB2	2.12	0.50
1:A:445:MET:H	1:A:476:GLN:HE22	1.58	0.50
1:C:695:LYS:HG2	1:C:781:TRP:CE3	2.47	0.50
1:C:661:ASN:C	1:C:661:ASN:HD22	2.20	0.49
1:A:734:VAL:CG2	2:A:24:GOL:H11	2.43	0.49
1:A:316:SER:HB3	1:C:274:ASN:HD21	1.77	0.49
1:A:632:ASN:HD21	1:A:695:LYS:NZ	2.11	0.49
1:B:660:LEU:HD11	1:B:719:LEU:HD23	1.95	0.49
1:B:446:MET:HE3	1:B:448:MET:HE2	1.94	0.49
1:C:522:PRO:HG2	1:C:524:TRP:NE1	2.28	0.49
1:C:721:GLU:O	1:C:725:LYS:HG3	2.12	0.49
1:D:743:ARG:HH11	1:D:760:MET:HE1	1.77	0.49
1:B:467:MET:CE	6:B:1182:HOH:O	2.58	0.48
1:D:618:ASP:OD1	1:D:618:ASP:C	2.56	0.48
1:A:505:TYR:HB3	1:A:506:PRO:HD3	1.95	0.48
1:B:178:LYS:HG2	1:B:211:LEU:HD21	1.95	0.48
1:D:444:ILE:HA	1:D:476:GLN:HE22	1.77	0.48
1:C:661:ASN:ND2	1:C:664:HIS:H	2.12	0.48
1:D:621:LEU:O	1:D:624:ARG:HB2	2.12	0.48
2:A:7:GOL:H11	6:A:909:HOH:O	2.13	0.48
1:B:243:LEU:O	1:B:246:CYS:HB2	2.12	0.48
1:D:730:VAL:O	1:D:764:HIS:HE1	1.97	0.48
1:A:281:ASN:OD1	1:A:283:LYS:HD3	2.14	0.48
1:B:626:ILE:O	1:B:629:LEU:HB2	2.13	0.48
1:A:538:SER:HB2	1:A:544:MET:HB2	1.96	0.47
1:B:213:LEU:HB3	1:B:216:LEU:CD1	2.43	0.47
1:C:632:ASN:HD22	1:C:632:ASN:N	2.13	0.47
1:C:643:GLN:HE22	1:C:646:LYS:HE2	1.78	0.47
1:A:667:MET:HE2	1:A:667:MET:HB2	1.82	0.47
1:B:743:ARG:CZ	1:B:760:MET:HE1	2.45	0.47
1:C:223:PHE:HB2	1:C:243:LEU:HD23	1.96	0.47
1:A:566:MET:HE1	1:A:579:THR:O	2.14	0.47
1:D:325:GLU:HG3	1:D:335:ASN:ND2	2.30	0.47
1:A:455:GLN:NE2	1:A:458:MET:HE1	2.29	0.47
1:C:447:TYR:HA	2:C:21:GOL:H32	1.96	0.47
1:D:450:ASN:OD1	1:D:484:ARG:HD2	2.14	0.47
1:D:532:TYR:CD1	1:D:548:CYS:HA	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:638:MET:CA	1:B:643:GLN:HG2	2.36	0.47
1:B:724:TYR:CE1	1:B:728:PRO:HB3	2.49	0.47
1:C:234:ARG:HG2	1:C:254:ASN:HB2	1.96	0.47
1:C:651:TRP:HB3	1:C:711:ILE:CD1	2.45	0.47
6:B:815:HOH:O	1:D:601:LYS:HE2	2.14	0.47
1:C:433:PHE:O	1:C:437:MET:HG2	2.15	0.47
1:C:671:SER:O	1:C:675:LYS:HB2	2.14	0.47
1:B:583:SER:HB3	1:B:586:ASN:HB2	1.96	0.46
1:D:624:ARG:HH21	2:D:27:GOL:H12	1.80	0.46
1:A:221:VAL:HG11	1:A:223:PHE:CE2	2.51	0.46
1:C:528:SER:O	1:C:530:LEU:HD22	2.15	0.46
1:A:350:ILE:O	1:A:353:VAL:HG22	2.15	0.46
1:C:330:LYS:O	1:C:331:SER:HB3	2.15	0.46
1:D:536:LEU:N	1:D:536:LEU:HD12	2.31	0.46
1:C:445:MET:H	1:C:476:GLN:HE22	1.64	0.46
1:B:455:GLN:HA	1:B:458:MET:HE2	1.97	0.46
1:C:724:TYR:CE1	1:C:728:PRO:HB3	2.51	0.46
1:A:433:PHE:O	1:A:437:MET:HG2	2.16	0.46
1:D:325:GLU:HG3	1:D:335:ASN:HD21	1.80	0.45
1:D:455:GLN:HA	1:D:458:MET:HE2	1.99	0.45
1:A:439:TRP:CG	5:A:785:DTT:H11	2.51	0.45
1:B:437:MET:HE1	1:B:477:ILE:HG13	1.98	0.45
1:B:488:GLN:O	1:B:492:ARG:HG3	2.16	0.45
1:A:175:CYS:C	1:A:176:LEU:HD12	2.41	0.45
1:B:478:VAL:HG11	1:B:495:TYR:HB2	1.98	0.45
1:B:597:ILE:H	1:B:597:ILE:CD1	1.97	0.45
1:C:515:LEU:CD1	1:C:523:ASP:HA	2.46	0.45
1:B:536:LEU:HD11	1:B:553:MET:HE1	1.98	0.45
1:B:643:GLN:NE2	1:B:646:LYS:HE2	2.31	0.45
1:D:194:GLY:N	1:D:196:GLU:OE1	2.50	0.45
1:C:243:LEU:O	1:C:266:CYS:SG	2.75	0.45
1:C:295:TYR:HA	1:C:296:PRO:HD3	1.83	0.45
1:A:639:ILE:HG12	1:A:704:THR:HG23	1.99	0.45
1:A:217:ASN:C	1:A:218:LEU:HD23	2.42	0.45
1:B:538:SER:HB2	1:B:544:MET:HB2	1.99	0.45
1:D:532:TYR:HD1	1:D:548:CYS:HA	1.80	0.45
2:A:4:GOL:H31	2:A:19:GOL:C1	2.47	0.45
1:B:515:LEU:O	1:B:516:CYS:HB2	2.16	0.45
1:C:244:GLU:HG3	1:C:262:ASN:ND2	2.32	0.45
1:D:204:SER:HB3	1:D:222:ASN:HD21	1.81	0.45
1:A:398:PHE:CZ	5:A:785:DTT:S1	3.08	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:638:MET:HA	1:A:643:GLN:CG	2.34	0.45
1:B:597:ILE:N	1:B:597:ILE:CD1	2.74	0.45
1:C:323:VAL:HG21	1:C:381:ARG:NH2	2.31	0.45
1:D:678:PRO:HD2	1:D:681:LYS:HB2	1.99	0.45
1:B:202:ASN:ND2	1:B:204:SER:H	2.15	0.44
1:B:766:LYS:HE2	2:D:26:GOL:O3	2.17	0.44
1:B:544:MET:HE3	1:B:570:LEU:HD11	1.98	0.44
1:D:420:ASN:HB2	6:D:841:HOH:O	2.16	0.44
1:D:453:LEU:HG	1:D:487:MET:CE	2.35	0.44
1:D:484:ARG:O	1:D:488:GLN:HG3	2.18	0.44
1:D:575:GLU:HB2	6:D:895:HOH:O	2.17	0.44
1:D:695:LYS:HG2	1:D:781:TRP:CE3	2.53	0.44
1:B:330:LYS:O	1:B:331:SER:CB	2.60	0.44
1:C:181:LEU:HD12	1:C:181:LEU:HA	1.86	0.44
1:C:711:ILE:HD13	1:C:711:ILE:O	2.18	0.44
1:C:321:GLY:CA	6:C:1244:HOH:O	2.66	0.44
1:D:272:PHE:CG	1:D:297:PRO:HG3	2.52	0.44
1:D:628:ALA:HB1	1:D:701:MET:CE	2.48	0.44
1:D:303:ASP:HA	1:D:319:ILE:HB	1.99	0.44
1:D:489:GLU:O	1:D:493:GLU:HG3	2.17	0.44
1:D:556:MET:HE2	1:D:595:PHE:CE1	2.53	0.44
1:A:455:GLN:CD	1:A:458:MET:CE	2.91	0.43
1:C:655:LEU:HD12	1:C:711:ILE:CD1	2.46	0.43
1:D:342:ASP:OD2	1:D:344:SER:OG	2.35	0.43
1:D:469:SER:HB3	2:D:26:GOL:H31	2.00	0.43
1:D:628:ALA:HB1	1:D:701:MET:HE1	1.99	0.43
1:A:221:VAL:HG12	1:A:222:ASN:N	2.32	0.43
1:C:643:GLN:HA	1:C:643:GLN:NE2	2.31	0.43
1:A:770:PRO:HD2	2:A:16:GOL:H2	1.99	0.43
1:B:350:ILE:O	1:B:353:VAL:HG22	2.18	0.43
1:B:566:MET:HE1	6:B:963:HOH:O	2.18	0.43
1:A:435:THR:CG2	5:A:785:DTT:S1	3.06	0.43
1:B:181:LEU:HD12	1:B:181:LEU:HA	1.83	0.43
1:D:300:ASN:N	1:D:300:ASN:ND2	2.64	0.43
1:B:282:LEU:HB2	1:B:304:LEU:HD12	2.00	0.43
1:C:330:LYS:O	1:C:331:SER:CB	2.67	0.43
1:A:730:VAL:CG2	1:A:769:PHE:CE2	3.01	0.43
6:B:1228:HOH:O	2:D:28:GOL:H2	2.19	0.43
1:C:328:SER:N	1:C:335:ASN:HD21	2.16	0.43
1:D:612:LEU:HD21	1:D:625:PHE:HB3	2.00	0.43
1:A:652:ASN:N	1:A:653:PRO:CD	2.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:346:ILE:O	1:C:348:GLU:N	2.45	0.42
1:C:383:ALA:O	1:C:387:ILE:HG23	2.19	0.42
1:A:374:LYS:HE3	1:A:374:LYS:HB2	1.91	0.42
1:B:330:LYS:HD2	1:B:333:THR:HG21	2.01	0.42
1:B:437:MET:HE1	1:B:477:ILE:CG1	2.49	0.42
1:A:530:LEU:HD12	1:A:530:LEU:HA	1.94	0.42
1:B:277:PHE:HB2	1:B:299:LEU:HD23	2.02	0.42
1:C:249:LYS:HA	1:C:269:ARG:HB3	2.02	0.42
1:D:730:VAL:HG13	1:D:769:PHE:CE2	2.54	0.42
1:A:221:VAL:CG1	1:A:223:PHE:CE2	3.03	0.42
1:B:193:THR:HG22	1:B:195:GLU:N	2.34	0.42
1:B:274:ASN:ND2	1:D:316:SER:HB3	2.34	0.42
1:D:552:HIS:CE1	1:D:567:SER:HB3	2.54	0.42
1:A:584:LEU:HD12	1:A:584:LEU:HA	1.73	0.42
1:B:515:LEU:HD13	1:B:523:ASP:HA	2.01	0.42
1:D:330:LYS:CE	1:D:332:ASN:HB2	2.46	0.42
1:D:404:MET:SD	1:D:446:MET:HE1	2.60	0.42
1:A:233:LEU:O	1:A:236:SER:HB2	2.20	0.42
1:A:626:ILE:O	1:A:629:LEU:HB2	2.20	0.42
1:A:681:LYS:HG3	1:A:772:ILE:HD11	2.02	0.42
1:C:196:GLU:HG3	1:C:201:LEU:CD2	2.43	0.42
1:C:564:ASN:OD1	1:C:567:SER:HB2	2.20	0.42
1:C:782:ARG:NH2	2:C:23:GOL:H31	2.28	0.42
1:A:436:ILE:HD13	5:A:785:DTT:H12	2.02	0.42
1:D:478:VAL:HG11	1:D:495:TYR:HB2	2.02	0.42
1:B:622:LYS:HG2	1:B:626:ILE:HD12	2.01	0.41
1:B:234:ARG:O	1:B:235:MET:HB2	2.19	0.41
1:A:373:ASP:OD2	3:C:2:MES:H62	2.21	0.41
1:A:632:ASN:C	1:A:701:MET:HE1	2.46	0.41
1:A:679:LEU:HD12	1:A:682:GLN:NE2	2.35	0.41
1:B:300:ASN:HD22	1:B:300:ASN:N	2.12	0.41
1:B:433:PHE:O	1:B:437:MET:HG2	2.20	0.41
1:A:233:LEU:O	1:A:256:CYS:SG	2.75	0.41
1:A:274:ASN:ND2	1:C:316:SER:HB3	2.36	0.41
1:A:455:GLN:CD	1:A:458:MET:HE3	2.46	0.41
1:A:585:SER:OG	1:A:600:HIS:HE1	2.02	0.41
1:B:552:HIS:HD2	6:B:1020:HOH:O	2.03	0.41
1:D:743:ARG:CZ	1:D:760:MET:HE1	2.51	0.41
1:C:213:LEU:HB2	1:C:233:LEU:HD23	2.01	0.41
1:D:171:GLN:N	1:D:199:GLU:O	2.54	0.41
1:A:202:ASN:OD1	1:A:202:ASN:C	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:ARG:HG2	1:A:254:ASN:HB2	2.02	0.41
1:A:274:ASN:HD21	1:C:316:SER:HB3	1.84	0.41
1:A:395:LEU:HD12	1:A:395:LEU:HA	1.77	0.41
1:A:475:ILE:HG13	1:A:536:LEU:HD23	2.02	0.41
1:B:373:ASP:OD1	3:D:4:MES:H22	2.20	0.41
1:B:701:MET:HE3	1:B:701:MET:HB3	1.88	0.41
1:D:586:ASN:O	1:D:590:GLU:HB2	2.20	0.41
1:A:195:GLU:H	1:A:195:GLU:HG3	1.70	0.41
1:B:173:ARG:C	1:B:175:CYS:H	2.28	0.41
1:B:325:GLU:HG3	1:B:335:ASN:ND2	2.36	0.41
1:B:532:TYR:HB3	1:B:547:MET:HG2	2.03	0.41
1:A:288:LYS:HE2	1:A:342:ASP:CG	2.46	0.41
1:A:525:ASP:O	1:A:525:ASP:CG	2.64	0.41
1:B:373:ASP:CG	3:D:4:MES:H22	2.46	0.41
1:C:228:LEU:HD22	1:C:233:LEU:HD11	2.03	0.41
1:B:371:SER:OG	1:B:379:LYS:HE2	2.22	0.40
1:C:737:TYR:CZ	1:C:741:LEU:HD11	2.56	0.40
1:D:353:VAL:HG13	1:D:364:ILE:HG22	2.03	0.40
1:A:679:LEU:HD12	1:A:679:LEU:HA	1.84	0.40
1:B:173:ARG:H	1:B:176:LEU:HD13	1.86	0.40
1:B:377:HIS:O	1:B:381[A]:ARG:HG3	2.20	0.40
1:B:532:TYR:CD1	1:B:548:CYS:HA	2.56	0.40
1:B:391:SER:OG	1:B:394[A]:GLU:HG3	2.21	0.40
1:B:548:CYS:SG	1:B:553:MET:HE2	2.61	0.40
1:C:680:ARG:HD2	1:C:729:GLN:CD	2.46	0.40
1:A:479:MET:HE2	1:A:550:LEU:HD13	2.03	0.40
1:B:744:LEU:HD13	1:B:757:LEU:CD2	2.51	0.40
1:D:328:SER:HB2	1:D:335:ASN:OD1	2.22	0.40
1:D:612:LEU:HD23	1:D:626:ILE:HD13	2.04	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	610/613 (100%)	583 (96%)	27 (4%)	0	100	100
1	B	612/613 (100%)	580 (95%)	32 (5%)	0	100	100
1	C	610/613 (100%)	585 (96%)	25 (4%)	0	100	100
1	D	611/613 (100%)	584 (96%)	27 (4%)	0	100	100
All	All	2443/2452 (100%)	2332 (96%)	111 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	541/546 (99%)	504 (93%)	37 (7%)	14	12
1	B	541/546 (99%)	509 (94%)	32 (6%)	18	16
1	C	541/546 (99%)	509 (94%)	32 (6%)	18	16
1	D	545/546 (100%)	525 (96%)	20 (4%)	30	33
All	All	2168/2184 (99%)	2047 (94%)	121 (6%)	19	18

All (121) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	186	ILE
1	A	191	ASN
1	A	195	GLU
1	A	197	SER
1	A	201	LEU
1	A	272	PHE
1	A	300	ASN
1	A	323	VAL
1	A	333	THR
1	A	345	ASP
1	A	387	ILE
1	A	392	LEU
1	A	395	LEU

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Mol	Chain	Res	Type
1	A	416	LEU
1	A	444	ILE
1	A	499	LEU
1	A	515	LEU
1	A	528	SER
1	A	530	LEU
1	A	553	MET
1	A	561	THR
1	A	584	LEU
1	A	608	VAL
1	A	612	LEU
1	A	623	GLU
1	A	629	LEU
1	A	643	GLN
1	A	647	LEU
1	A	667	MET
1	A	671	SER
1	A	673	VAL
1	A	679	LEU
1	A	680	ARG
1	A	705	GLU
1	A	713	ARG
1	A	741	LEU
1	A	744	LEU
1	B	175	CYS
1	B	181	LEU
1	B	195	GLU
1	B	219	SER
1	B	246	CYS
1	B	264	SER
1	B	272	PHE
1	B	300	ASN
1	B	323	VAL
1	B	326	LEU
1	B	392	LEU
1	B	395	LEU
1	B	416	LEU
1	B	499	LEU
1	B	515	LEU
1	B	527	ASP
1	B	553	MET
1	B	584	LEU

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Mol	Chain	Res	Type
1	B	597	ILE
1	B	612	LEU
1	B	623	GLU
1	B	627	GLU
1	B	629	LEU
1	B	643	GLN
1	B	647	LEU
1	B	656	ASP
1	B	681	LYS
1	B	697	SER
1	B	729	GLN
1	B	741	LEU
1	B	744	LEU
1	B	754	THR
1	C	181	LEU
1	C	196	GLU
1	C	199	GLU
1	C	240	GLN
1	C	272	PHE
1	C	274	ASN
1	C	300	ASN
1	C	323	VAL
1	C	333	THR
1	C	374	LYS
1	C	391	SER
1	C	395	LEU
1	C	406	LEU
1	C	416	LEU
1	C	425	LEU
1	C	515	LEU
1	C	527	ASP
1	C	530	LEU
1	C	553	MET
1	C	554	VAL
1	C	584	LEU
1	C	612	LEU
1	C	629	LEU
1	C	632	ASN
1	C	642	ASP
1	C	647	LEU
1	C	661	ASN
1	C	674	LEU

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Mol	Chain	Res	Type
1	C	680	ARG
1	C	689	LEU
1	C	711	ILE
1	C	772	ILE
1	D	196	GLU
1	D	240	GLN
1	D	272	PHE
1	D	300	ASN
1	D	353	VAL
1	D	395	LEU
1	D	406	LEU
1	D	416	LEU
1	D	472	SER
1	D	499	LEU
1	D	502	GLU
1	D	515	LEU
1	D	576	VAL
1	D	584	LEU
1	D	612	LEU
1	D	627	GLU
1	D	647	LEU
1	D	650	VAL
1	D	741	LEU
1	D	744	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (67) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	284	ASN
1	A	294	GLN
1	A	300	ASN
1	A	418	ASN
1	A	441	ASN
1	A	476	GLN
1	A	480	ASN
1	A	509	GLN
1	A	577	GLN
1	A	600	HIS
1	A	632	ASN
1	A	643	GLN
1	A	682	GLN
1	A	729	GLN

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Mol	Chain	Res	Type
1	A	764	HIS
1	B	202	ASN
1	B	245	ASN
1	B	300	ASN
1	B	418	ASN
1	B	420	ASN
1	B	441	ASN
1	B	476	GLN
1	B	480	ASN
1	B	509	GLN
1	B	540	GLN
1	B	552	HIS
1	B	600	HIS
1	B	632	ASN
1	B	643	GLN
1	B	663	GLN
1	B	672	HIS
1	B	717	ASN
1	B	745	GLN
1	C	217	ASN
1	C	262	ASN
1	C	284	ASN
1	C	300	ASN
1	C	335	ASN
1	C	418	ASN
1	C	420	ASN
1	C	476	GLN
1	C	480	ASN
1	C	488	GLN
1	C	509	GLN
1	C	577	GLN
1	C	632	ASN
1	C	643	GLN
1	C	661	ASN
1	C	664	HIS
1	C	717	ASN
1	C	745	GLN
1	C	764	HIS
1	D	245	ASN
1	D	284	ASN
1	D	300	ASN
1	D	332	ASN

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Mol	Chain	Res	Type
1	D	418	ASN
1	D	441	ASN
1	D	476	GLN
1	D	501	ASN
1	D	600	HIS
1	D	632	ASN
1	D	661	ASN
1	D	717	ASN
1	D	745	GLN
1	D	749	ASN
1	D	764	HIS

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

52 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	GOL	A	1	-	5,5,5	0.46	0	5,5,5	0.61	0
4	SO4	A	783	-	4,4,4	0.24	0	6,6,6	0.05	0
3	MES	D	783	-	12,12,12	2.22	1 (8%)	15,16,16	2.43	6 (40%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	D	784	-	4,4,4	0.24	0	6,6,6	0.20	0
4	SO4	D	785	-	4,4,4	0.25	0	6,6,6	0.17	0
2	GOL	C	17	-	5,5,5	0.35	0	5,5,5	0.39	0
2	GOL	C	21	-	5,5,5	0.32	0	5,5,5	0.64	0
3	MES	C	1	-	12,12,12	2.25	1 (8%)	15,16,16	2.34	6 (40%)
2	GOL	A	4	-	5,5,5	0.44	0	5,5,5	0.29	0
2	GOL	B	3	-	5,5,5	0.37	0	5,5,5	0.33	0
4	SO4	C	784	-	4,4,4	0.21	0	6,6,6	0.25	0
2	GOL	A	13	-	5,5,5	0.38	0	5,5,5	0.43	0
2	GOL	D	26	-	5,5,5	0.32	0	5,5,5	0.45	0
2	GOL	A	7	-	5,5,5	0.40	0	5,5,5	0.52	0
4	SO4	B	784	-	4,4,4	0.23	0	6,6,6	0.17	0
2	GOL	D	5	-	5,5,5	0.41	0	5,5,5	0.40	0
4	SO4	B	783	-	4,4,4	0.23	0	6,6,6	0.27	0
4	SO4	C	783	-	4,4,4	0.30	0	6,6,6	0.08	0
2	GOL	B	8	-	5,5,5	0.39	0	5,5,5	0.53	0
2	GOL	A	18	-	5,5,5	0.39	0	5,5,5	0.40	0
5	DTT	C	785	-	7,7,7	0.93	0	4,8,8	0.85	0
2	GOL	A	24	-	5,5,5	0.34	0	5,5,5	0.29	0
4	SO4	A	5	-	4,4,4	0.22	0	6,6,6	0.19	0
2	GOL	D	28	-	5,5,5	0.36	0	5,5,5	0.33	0
2	GOL	D	9	-	5,5,5	0.35	0	5,5,5	0.29	0
4	SO4	C	10	-	4,4,4	0.23	0	6,6,6	0.10	0
2	GOL	A	15	-	5,5,5	0.39	0	5,5,5	0.51	0
4	SO4	C	7	-	4,4,4	0.23	0	6,6,6	0.22	0
4	SO4	D	16	-	4,4,4	0.23	0	6,6,6	0.13	0
4	SO4	A	784	-	4,4,4	0.25	0	6,6,6	0.13	0
4	SO4	C	3	-	4,4,4	0.24	0	6,6,6	0.12	0
3	MES	C	2	-	12,12,12	2.36	1 (8%)	15,16,16	2.41	6 (40%)
2	GOL	A	16	-	5,5,5	0.40	0	5,5,5	0.39	0
2	GOL	A	19	-	5,5,5	0.41	0	5,5,5	0.28	0
4	SO4	B	12	-	4,4,4	0.25	0	6,6,6	0.08	0
2	GOL	B	10	-	5,5,5	0.39	0	5,5,5	0.33	0
2	GOL	A	2	-	5,5,5	0.31	0	5,5,5	0.42	0
2	GOL	C	29	-	5,5,5	0.43	0	5,5,5	0.48	0
2	GOL	D	27	-	5,5,5	0.38	0	5,5,5	0.33	0
2	GOL	A	12	-	5,5,5	0.34	0	5,5,5	0.29	0
2	GOL	C	11	-	5,5,5	0.29	0	5,5,5	0.82	0
2	GOL	C	22	-	5,5,5	0.39	0	5,5,5	0.50	0
5	DTT	A	785	-	7,7,7	0.74	0	4,8,8	1.10	0
2	GOL	C	23	-	5,5,5	0.36	0	5,5,5	0.33	0
2	GOL	A	14	-	5,5,5	0.36	0	5,5,5	0.48	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	MES	A	3	-	12,12,12	2.23	1 (8%)	15,16,16	2.38	6 (40%)
4	SO4	A	17	-	4,4,4	0.31	0	6,6,6	0.25	0
2	GOL	A	25	-	5,5,5	0.40	0	5,5,5	0.27	0
2	GOL	B	6	-	5,5,5	0.39	0	5,5,5	0.21	0
3	MES	D	4	-	12,12,12	2.09	1 (8%)	15,16,16	2.35	5 (33%)
4	SO4	D	15	-	4,4,4	0.23	0	6,6,6	0.12	0
2	GOL	C	20	-	5,5,5	0.40	0	5,5,5	0.28	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
2	GOL	A	1	-	-	2/4/4/4	-
3	MES	D	783	-	-	1/6/14/14	0/1/1/1
2	GOL	C	17	-	-	2/4/4/4	-
2	GOL	C	21	-	-	0/4/4/4	-
3	MES	C	1	-	-	1/6/14/14	0/1/1/1
2	GOL	A	4	-	-	2/4/4/4	-
2	GOL	B	3	-	-	0/4/4/4	-
2	GOL	A	13	-	-	4/4/4/4	-
2	GOL	D	26	-	-	2/4/4/4	-
2	GOL	A	7	-	-	0/4/4/4	-
2	GOL	D	5	-	-	2/4/4/4	-
2	GOL	B	8	-	-	2/4/4/4	-
2	GOL	A	18	-	-	2/4/4/4	-
5	DTT	C	785	-	-	4/8/8/8	-
2	GOL	A	24	-	-	3/4/4/4	-
2	GOL	D	28	-	-	2/4/4/4	-
2	GOL	D	9	-	-	4/4/4/4	-
2	GOL	A	15	-	-	2/4/4/4	-
3	MES	C	2	-	-	1/6/14/14	0/1/1/1
2	GOL	A	16	-	-	2/4/4/4	-
2	GOL	A	19	-	-	2/4/4/4	-
2	GOL	B	10	-	-	2/4/4/4	-
2	GOL	A	2	-	-	4/4/4/4	-
2	GOL	C	29	-	-	4/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	D	27	-	-	3/4/4/4	-
2	GOL	A	12	-	-	2/4/4/4	-
2	GOL	C	11	-	-	0/4/4/4	-
2	GOL	C	22	-	-	3/4/4/4	-
5	DTT	A	785	-	-	5/8/8/8	-
2	GOL	C	23	-	-	2/4/4/4	-
2	GOL	A	14	-	-	4/4/4/4	-
3	MES	A	3	-	-	4/6/14/14	0/1/1/1
2	GOL	A	25	-	-	4/4/4/4	-
2	GOL	B	6	-	-	2/4/4/4	-
3	MES	D	4	-	-	4/6/14/14	0/1/1/1
2	GOL	C	20	-	-	2/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	2	MES	C8-S	-7.89	1.66	1.77
3	C	1	MES	C8-S	-7.43	1.67	1.77
3	A	3	MES	C8-S	-7.39	1.67	1.77
3	D	783	MES	C8-S	-7.38	1.67	1.77
3	D	4	MES	C8-S	-6.95	1.67	1.77

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1	MES	C5-N4-C3	5.51	120.72	108.84
3	C	2	MES	C5-N4-C3	5.45	120.57	108.84
3	A	3	MES	C5-N4-C3	5.00	119.61	108.84
3	D	4	MES	C5-N4-C3	4.97	119.55	108.84
3	D	783	MES	C5-N4-C3	4.61	118.78	108.84
3	D	783	MES	C2-C3-N4	-4.07	103.94	110.12
3	A	3	MES	C7-N4-C3	3.61	120.87	111.24
3	C	2	MES	C7-N4-C3	3.58	120.77	111.24
3	D	4	MES	C7-N4-C3	3.56	120.74	111.24
3	D	783	MES	C7-N4-C3	3.56	120.72	111.24
3	D	783	MES	C7-N4-C5	3.45	120.43	111.24
3	C	2	MES	C2-C3-N4	-3.37	104.99	110.12
3	C	1	MES	C7-N4-C3	3.32	120.09	111.24
3	D	4	MES	C6-C5-N4	-3.24	105.19	110.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	4	MES	O1S-S-C8	3.24	111.62	106.73
3	A	3	MES	C6-C5-N4	-3.24	105.20	110.12
3	D	4	MES	C7-N4-C5	3.18	119.70	111.24
3	A	3	MES	C7-N4-C5	3.10	119.51	111.24
3	C	1	MES	C7-N4-C5	2.98	119.18	111.24
3	C	1	MES	C6-C5-N4	-2.96	105.62	110.12
3	D	783	MES	C6-C5-N4	-2.91	105.70	110.12
3	C	1	MES	C2-C3-N4	-2.82	105.84	110.12
3	C	2	MES	C7-N4-C5	2.77	118.61	111.24
3	C	2	MES	C6-C5-N4	-2.76	105.92	110.12
3	D	783	MES	O1S-S-C8	2.48	110.47	106.73
3	A	3	MES	C2-C3-N4	-2.34	106.56	110.12
3	C	2	MES	O2S-S-C8	2.21	110.08	106.73
3	C	1	MES	C8-C7-N4	-2.18	104.12	112.36
3	A	3	MES	O1S-S-C8	2.13	109.95	106.73

There are no chirality outliers.

All (85) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1	GOL	O1-C1-C2-C3
2	A	2	GOL	C1-C2-C3-O3
2	A	2	GOL	O2-C2-C3-O3
2	A	13	GOL	O1-C1-C2-C3
2	A	14	GOL	O1-C1-C2-O2
2	A	14	GOL	O1-C1-C2-C3
2	A	18	GOL	O1-C1-C2-C3
2	A	19	GOL	O1-C1-C2-C3
2	A	24	GOL	O1-C1-C2-C3
2	A	25	GOL	O1-C1-C2-O2
2	A	25	GOL	O1-C1-C2-C3
2	B	6	GOL	C1-C2-C3-O3
2	C	17	GOL	O1-C1-C2-C3
2	C	20	GOL	O1-C1-C2-C3
2	C	22	GOL	O1-C1-C2-C3
2	C	23	GOL	O1-C1-C2-C3
2	C	29	GOL	O1-C1-C2-C3
2	C	29	GOL	C1-C2-C3-O3
2	C	29	GOL	O2-C2-C3-O3
2	D	9	GOL	O1-C1-C2-O2
2	D	9	GOL	O1-C1-C2-C3
2	D	9	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
2	D	26	GOL	O1-C1-C2-C3
2	D	27	GOL	O1-C1-C2-C3
2	D	28	GOL	O1-C1-C2-C3
3	A	3	MES	C8-C7-N4-C5
3	A	3	MES	C7-C8-S-O3S
3	D	4	MES	C8-C7-N4-C5
3	D	4	MES	C7-C8-S-O2S
3	D	4	MES	C7-C8-S-O3S
5	A	785	DTT	S1-C1-C2-O2
5	A	785	DTT	S1-C1-C2-C3
5	A	785	DTT	C1-C2-C3-C4
5	C	785	DTT	S1-C1-C2-O2
5	C	785	DTT	S1-C1-C2-C3
5	C	785	DTT	C2-C3-C4-S4
2	A	24	GOL	O1-C1-C2-O2
2	B	8	GOL	O1-C1-C2-O2
2	A	2	GOL	O1-C1-C2-C3
2	A	4	GOL	C1-C2-C3-O3
2	A	12	GOL	C1-C2-C3-O3
2	A	13	GOL	C1-C2-C3-O3
2	A	14	GOL	C1-C2-C3-O3
2	A	15	GOL	O1-C1-C2-C3
2	A	25	GOL	C1-C2-C3-O3
2	B	8	GOL	O1-C1-C2-C3
2	D	5	GOL	O1-C1-C2-C3
2	A	1	GOL	O1-C1-C2-O2
2	A	13	GOL	O1-C1-C2-O2
2	A	15	GOL	O1-C1-C2-O2
2	A	19	GOL	O1-C1-C2-O2
2	B	6	GOL	O2-C2-C3-O3
2	C	17	GOL	O1-C1-C2-O2
2	C	22	GOL	O1-C1-C2-O2
2	C	29	GOL	O1-C1-C2-O2
2	D	5	GOL	O1-C1-C2-O2
2	D	9	GOL	O2-C2-C3-O3
2	D	27	GOL	O1-C1-C2-O2
2	A	12	GOL	O2-C2-C3-O3
2	A	13	GOL	O2-C2-C3-O3
2	C	20	GOL	O1-C1-C2-O2
2	C	23	GOL	O1-C1-C2-O2
2	D	26	GOL	O1-C1-C2-O2
2	D	28	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
3	C	1	MES	C8-C7-N4-C3
2	A	4	GOL	O2-C2-C3-O3
2	A	18	GOL	O1-C1-C2-O2
2	A	25	GOL	O2-C2-C3-O3
2	A	16	GOL	O1-C1-C2-O2
2	B	10	GOL	O2-C2-C3-O3
3	A	3	MES	C7-C8-S-O1S
3	A	3	MES	C7-C8-S-O2S
3	D	4	MES	C7-C8-S-O1S
2	D	27	GOL	O2-C2-C3-O3
5	A	785	DTT	O3-C3-C4-S4
5	C	785	DTT	O3-C3-C4-S4
2	A	2	GOL	O1-C1-C2-O2
2	A	14	GOL	O2-C2-C3-O3
3	C	2	MES	C8-C7-N4-C5
3	D	783	MES	C8-C7-N4-C3
5	A	785	DTT	C2-C3-C4-S4
2	A	24	GOL	C1-C2-C3-O3
2	A	16	GOL	O1-C1-C2-C3
2	B	10	GOL	C1-C2-C3-O3
2	C	22	GOL	O2-C2-C3-O3

There are no ring outliers.

18 monomers are involved in 43 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	783	MES	3	0
2	C	21	GOL	5	0
2	A	4	GOL	1	0
2	D	26	GOL	2	0
2	A	7	GOL	2	0
5	C	785	DTT	2	0
2	A	24	GOL	1	0
2	D	28	GOL	1	0
3	C	2	MES	2	0
2	A	16	GOL	3	0
2	A	19	GOL	1	0
2	D	27	GOL	1	0
2	C	11	GOL	4	0
5	A	785	DTT	8	0
2	C	23	GOL	2	0
2	A	14	GOL	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	3	MES	1	0
3	D	4	MES	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	612/613 (99%)	0.02	30 (4%) 35 37	21, 36, 78, 128	0
1	B	612/613 (99%)	0.09	13 (2%) 63 66	24, 40, 80, 139	2 (0%)
1	C	612/613 (99%)	0.05	17 (2%) 55 57	22, 40, 74, 145	0
1	D	612/613 (99%)	0.23	11 (1%) 67 70	26, 48, 80, 122	1 (0%)
All	All	2448/2452 (99%)	0.10	71 (2%) 53 56	21, 41, 79, 145	3 (0%)

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	580	PHE	6.8
1	B	579	THR	6.7
1	B	580	PHE	4.8
1	C	580	PHE	4.6
1	A	172	GLY	4.4
1	B	171	GLN	4.1
1	C	326	LEU	4.0
1	C	327	PHE	3.9
1	A	581	GLY	3.9
1	D	582	TYR	3.8
1	C	561	THR	3.8
1	D	327	PHE	3.7
1	A	183	ALA	3.7
1	A	206	CYS	3.5
1	A	173	ARG	3.5
1	A	177	SER	3.5
1	C	171	GLN	3.4
1	A	187	TRP	3.4
1	A	392	LEU	3.4
1	B	227	VAL	3.3
1	D	171	GLN	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	560	ASN	3.1
1	C	331	SER	3.1
1	A	174	ALA	3.0
1	D	672	HIS	3.0
1	A	176	LEU	3.0
1	C	172	GLY	3.0
1	A	527	ASP	3.0
1	D	581	GLY	2.9
1	D	174	ALA	2.9
1	A	186	ILE	2.8
1	A	579	THR	2.6
1	A	748	ASN	2.6
1	A	649	CYS	2.6
1	A	216	LEU	2.6
1	A	582	TYR	2.5
1	D	172	GLY	2.5
1	A	227	VAL	2.5
1	A	578	ASN	2.5
1	C	330	LYS	2.5
1	D	326	LEU	2.4
1	B	597	ILE	2.4
1	C	576	VAL	2.4
1	A	246	CYS	2.4
1	C	559	PRO	2.4
1	D	579	THR	2.3
1	A	208	LEU	2.3
1	B	527	ASP	2.3
1	D	513	PHE	2.3
1	A	567	SER	2.3
1	A	207	ASP	2.3
1	C	567	SER	2.2
1	B	206	CYS	2.2
1	B	187	TRP	2.2
1	C	332	ASN	2.2
1	A	345	ASP	2.2
1	A	188	LEU	2.2
1	B	183	ALA	2.1
1	B	198	ALA	2.1
1	A	209	SER	2.1
1	B	189	SER	2.1
1	C	656	ASP	2.1
1	A	190	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	224	SER	2.1
1	A	656	ASP	2.1
1	B	209	SER	2.1
1	A	201	LEU	2.0
1	B	581	GLY	2.0
1	D	650	VAL	2.0
1	C	562	SER	2.0
1	C	174	ALA	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	SO4	A	784	5/5	0.64	0.12	116,120,121,121	0
4	SO4	D	15	5/5	0.68	0.17	160,161,162,162	0
2	GOL	A	14	6/6	0.70	0.18	43,65,73,78	0
4	SO4	C	783	5/5	0.71	0.13	91,99,102,105	0
2	GOL	D	27	6/6	0.71	0.13	72,79,85,89	0
4	SO4	B	12	5/5	0.73	0.14	138,140,141,142	0
4	SO4	A	783	5/5	0.74	0.12	143,146,148,148	0
4	SO4	C	10	5/5	0.79	0.11	107,110,111,114	0
4	SO4	B	784	5/5	0.80	0.14	77,83,88,100	0
4	SO4	D	16	5/5	0.80	0.07	112,115,118,121	0
2	GOL	C	20	6/6	0.81	0.25	82,89,91,93	0
2	GOL	D	5	6/6	0.81	0.18	50,58,60,69	0
4	SO4	C	3	5/5	0.81	0.09	109,109,113,114	0
2	GOL	A	18	6/6	0.82	0.16	49,71,77,79	0
2	GOL	C	11	6/6	0.82	0.18	41,51,59,67	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	SO4	D	785	5/5	0.83	0.15	88,94,96,107	0
2	GOL	D	26	6/6	0.84	0.15	76,79,83,85	0
2	GOL	C	17	6/6	0.84	0.15	46,59,76,81	0
2	GOL	C	21	6/6	0.85	0.15	25,54,61,64	0
2	GOL	A	13	6/6	0.86	0.14	42,51,56,58	0
2	GOL	A	4	6/6	0.86	0.12	41,54,59,64	0
2	GOL	A	7	6/6	0.86	0.17	43,54,62,67	0
2	GOL	A	25	6/6	0.86	0.16	76,81,84,87	0
2	GOL	C	29	6/6	0.86	0.18	61,75,83,87	0
2	GOL	B	8	6/6	0.86	0.16	49,57,65,70	0
2	GOL	A	24	6/6	0.87	0.10	56,59,62,62	0
4	SO4	C	784	5/5	0.87	0.15	27,76,88,95	0
2	GOL	C	23	6/6	0.87	0.12	61,83,85,91	0
2	GOL	A	16	6/6	0.87	0.17	47,58,68,71	0
2	GOL	A	1	6/6	0.87	0.14	27,35,44,44	0
2	GOL	D	28	6/6	0.88	0.18	78,81,87,89	0
2	GOL	C	22	6/6	0.88	0.11	51,68,74,79	0
2	GOL	A	19	6/6	0.88	0.11	47,51,56,62	0
2	GOL	B	10	6/6	0.88	0.12	55,60,62,64	0
3	MES	D	783	12/12	0.89	0.13	59,62,90,94	0
2	GOL	B	6	6/6	0.90	0.12	39,56,60,62	0
2	GOL	A	2	6/6	0.90	0.15	34,40,47,50	0
5	DTT	C	785	8/8	0.90	0.16	42,53,75,85	0
2	GOL	A	15	6/6	0.91	0.14	52,62,67,74	0
2	GOL	D	9	6/6	0.91	0.12	55,66,68,68	0
2	GOL	B	3	6/6	0.91	0.15	45,70,78,81	0
2	GOL	A	12	6/6	0.92	0.09	50,55,69,72	0
5	DTT	A	785	8/8	0.92	0.14	26,50,85,114	0
4	SO4	D	784	5/5	0.92	0.09	62,63,67,72	0
3	MES	C	2	12/12	0.93	0.10	48,50,68,74	0
3	MES	A	3	12/12	0.94	0.09	50,57,62,64	0
4	SO4	C	7	5/5	0.94	0.08	47,54,59,63	0
4	SO4	A	17	5/5	0.95	0.09	51,51,58,63	0
3	MES	D	4	12/12	0.95	0.09	50,62,67,69	0
3	MES	C	1	12/12	0.96	0.11	35,56,61,63	0
4	SO4	B	783	5/5	0.97	0.08	45,47,58,65	0
4	SO4	A	5	5/5	0.99	0.04	28,38,44,44	0

6.5 Other polymers ⓘ

There are no such residues in this entry.