



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 30, 2021 – 07:18 PM EST

PDB ID : 3L0Q
Title : The crystal structure of xylulose kinase from Yersinia pseudotuberculosis
Authors : Zhang, Z.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center
for Structural Genomics (NYSGXRC)
Deposited on : 2009-12-10
Resolution : 1.61 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.16
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.16

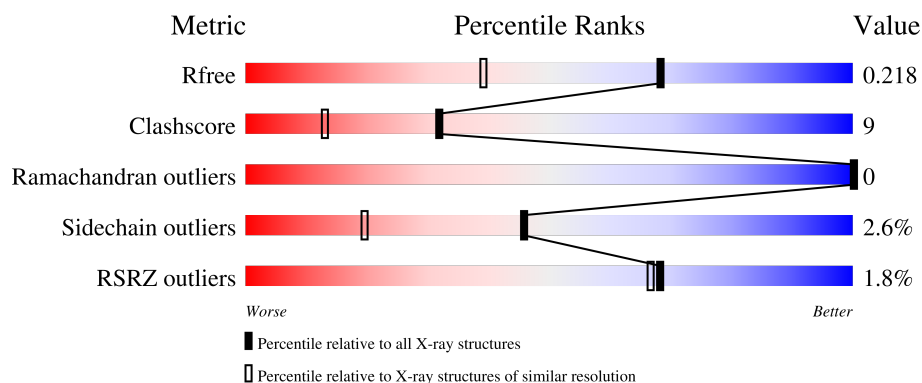
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.61 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4693 (1.64-1.60)
Clashscore	141614	5002 (1.64-1.60)
Ramachandran outliers	138981	4888 (1.64-1.60)
Sidechain outliers	138945	4887 (1.64-1.60)
RSRZ outliers	127900	4609 (1.64-1.60)

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 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	554	<div> <div>%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div>..</div> </div> </div>
1	B	554	<div> <div>2%</div> <div> <div></div> <div>84%</div> <div>13%</div> <div>..</div> </div> </div>

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	XUL	A	553	X	-	-	-
4	GOL	B	554	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9257 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.

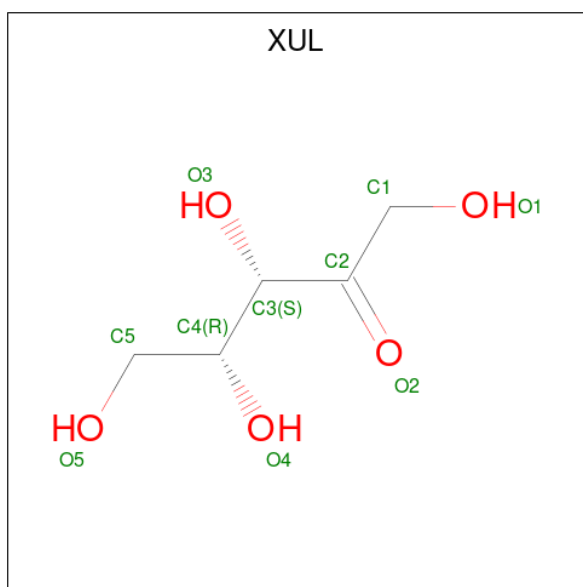
- Molecule 1 is a protein called xylulose kinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	541	Total	C	N	O	S	Se	0	5	0
			4160	2624	725	780	6	25			
1	B	542	Total	C	N	O	S	Se	0	3	0
			4166	2628	727	781	6	24			

There are 22 discrepancies between the modelled and reference sequences:

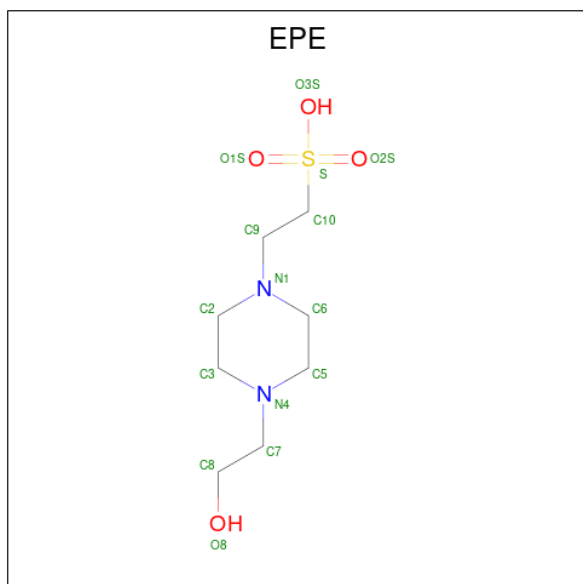
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MSE	-	expression tag	UNP Q665C6
A	0	SER	-	expression tag	UNP Q665C6
A	1	LEU	-	expression tag	UNP Q665C6
A	545	GLU	-	expression tag	UNP Q665C6
A	546	GLY	-	expression tag	UNP Q665C6
A	547	HIS	-	expression tag	UNP Q665C6
A	548	HIS	-	expression tag	UNP Q665C6
A	549	HIS	-	expression tag	UNP Q665C6
A	550	HIS	-	expression tag	UNP Q665C6
A	551	HIS	-	expression tag	UNP Q665C6
A	552	HIS	-	expression tag	UNP Q665C6
B	-1	MSE	-	expression tag	UNP Q665C6
B	0	SER	-	expression tag	UNP Q665C6
B	1	LEU	-	expression tag	UNP Q665C6
B	545	GLU	-	expression tag	UNP Q665C6
B	546	GLY	-	expression tag	UNP Q665C6
B	547	HIS	-	expression tag	UNP Q665C6
B	548	HIS	-	expression tag	UNP Q665C6
B	549	HIS	-	expression tag	UNP Q665C6
B	550	HIS	-	expression tag	UNP Q665C6
B	551	HIS	-	expression tag	UNP Q665C6
B	552	HIS	-	expression tag	UNP Q665C6

- Molecule 2 is D-XYLULOSE (three-letter code: XUL) (formula: C₅H₁₀O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			10	5	5		

- Molecule 3 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: $C_8H_{18}N_2O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

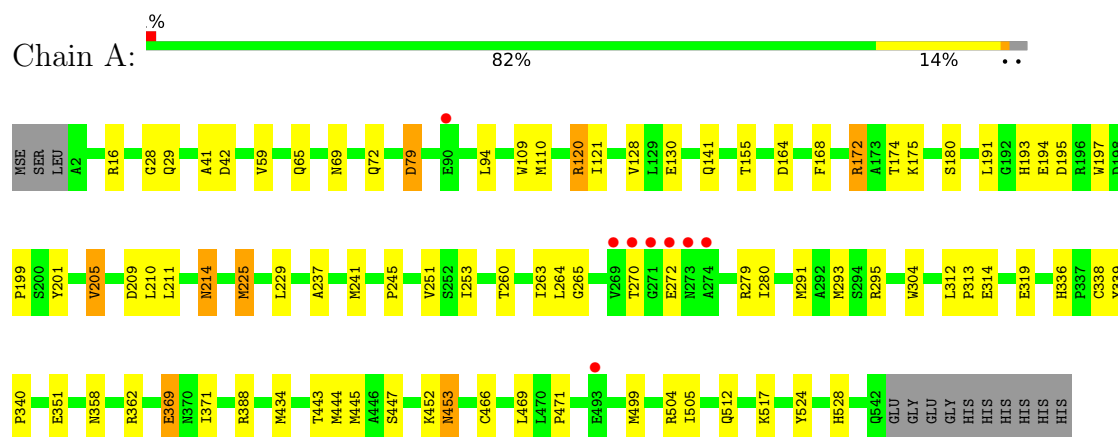
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	430	Total	O	0	0
			430	430		
5	B	470	Total	O	0	0
			470	470		

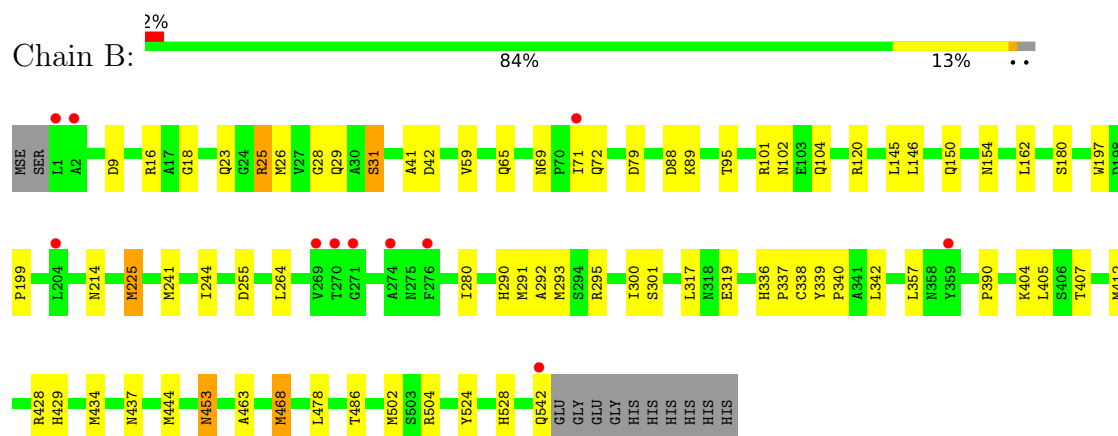
3 Residue-property plots

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: xylulose kinase



• Molecule 1: xylulose kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.88Å 110.00Å 80.53Å 90.00° 99.71° 90.00°	Depositor
Resolution (Å)	40.85 – 1.61 40.85 – 1.61	Depositor EDS
% Data completeness (in resolution range)	92.3 (40.85-1.61) 98.9 (40.85-1.61)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 1.61Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.192 , 0.222 0.190 , 0.218	Depositor DCC
R_{free} test set	6830 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	22.2	Xtriage
Anisotropy	0.362	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9257	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.25% 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: XUL, GOL, EPE

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	0.73	0/4245	0.82	2/5725 (0.0%)
1	B	0.77	0/4244	0.82	4/5722 (0.1%)
All	All	0.75	0/8489	0.82	6/11447 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	468	MSE	CG-SE-CE	-6.82	83.89	98.90
1	A	209	ASP	CB-CG-OD1	5.91	123.62	118.30
1	B	405	LEU	CA-CB-CG	5.83	128.72	115.30
1	B	295	ARG	NE-CZ-NH1	-5.75	117.42	120.30
1	A	504	ARG	NE-CZ-NH1	5.75	123.17	120.30
1	B	88	ASP	CB-CG-OD1	5.54	123.29	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4160	0	4104	64	0
1	B	4166	0	4114	83	0
2	A	10	0	10	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	15	0	18	2	0
4	B	6	0	8	11	0
5	A	430	0	0	9	0
5	B	470	0	0	9	0
All	All	9257	0	8254	148	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 9.

All (148) 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:291:MSE:HE2	1:B:317:LEU:HD11	1.28	1.14
1:B:291:MSE:HE1	1:B:317:LEU:HD21	1.31	1.06
1:B:154:ASN:HB2	4:B:554:GOL:H12	1.31	1.05
1:B:101:ARG:HH21	1:B:104:GLN:HE21	1.13	0.95
1:B:154:ASN:HB2	4:B:554:GOL:C3	1.97	0.93
1:B:18:GLY:HA3	1:B:26:MSE:HE3	1.49	0.93
1:B:154:ASN:HB2	4:B:554:GOL:C1	1.99	0.91
1:B:154:ASN:CB	4:B:554:GOL:H12	2.06	0.85
1:A:120:ARG:HH21	1:A:120:ARG:HG2	1.40	0.83
1:A:336:HIS:HD2	1:A:338:CYS:H	1.26	0.83
1:B:336:HIS:HD2	1:B:338:CYS:H	1.28	0.81
1:B:154:ASN:HB2	4:B:554:GOL:H31	1.63	0.78
1:A:120:ARG:HH21	1:A:120:ARG:CG	1.96	0.78
1:B:180:SER:OG	1:B:225:MSE:HE2	1.84	0.77
1:A:29:GLN:H	1:A:65:GLN:HE21	1.33	0.76
1:B:291:MSE:CE	1:B:317:LEU:HD21	2.15	0.75
1:B:357:LEU:HD21	1:B:412:MSE:HE3	1.70	0.73
1:B:154:ASN:CB	4:B:554:GOL:C3	2.67	0.72
1:B:59:VAL:HG21	1:B:241:MSE:HE2	1.71	0.72
1:B:18:GLY:CA	1:B:26:MSE:HE3	2.20	0.71
1:B:291:MSE:CE	1:B:317:LEU:HD11	2.16	0.71
1:A:237:ALA:O	1:A:241:MSE:HG3	1.92	0.70
1:A:16:ARG:NE	5:A:580:HOH:O	2.24	0.69
1:B:154:ASN:HB2	4:B:554:GOL:C2	2.22	0.69
1:B:154:ASN:CB	4:B:554:GOL:H31	2.22	0.68
1:B:26:MSE:HE2	1:B:29:GLN:HB2	1.76	0.68
1:A:120:ARG:NH2	1:A:120:ARG:HG2	2.06	0.67
1:B:120:ARG:HH12	1:B:146:LEU:HD11	1.60	0.66
1:B:342:LEU:HD22	1:B:412:MSE:HE1	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:ASN:HB2	1:B:72:GLN:HG3	1.78	0.65
1:B:336:HIS:CD2	1:B:338:CYS:H	2.12	0.65
1:A:180:SER:OG	1:A:225:MSE:HE2	1.95	0.65
1:A:471:PRO:HD2	5:A:593:HOH:O	1.97	0.64
1:A:445[B]:MSE:HE1	1:A:505:ILE:HG12	1.79	0.63
1:A:336:HIS:CD2	1:A:338:CYS:H	2.12	0.62
1:B:444:MSE:HE3	1:B:468:MSE:HE3	1.81	0.62
1:A:444:MSE:HE2	1:A:466:CYS:HB2	1.82	0.62
1:B:453:ASN:C	1:B:453:ASN:HD22	2.02	0.61
1:B:25:ARG:HD2	5:B:1021:HOH:O	1.99	0.61
1:B:71:ILE:HD12	1:B:244:ILE:HG13	1.83	0.61
1:B:26:MSE:HE2	1:B:29:GLN:CB	2.31	0.60
1:B:180:SER:HG	1:B:225:MSE:HE2	1.66	0.60
1:B:291:MSE:HE1	1:B:317:LEU:CD2	2.21	0.60
1:B:528:HIS:HD2	5:B:607:HOH:O	1.86	0.59
1:A:517:LYS:HE2	5:A:685:HOH:O	2.02	0.59
1:B:524:TYR:CZ	1:B:528:HIS:HE1	2.20	0.59
1:B:89:LYS:HE3	5:B:948:HOH:O	2.02	0.59
1:A:263:ILE:HD12	1:A:445[B]:MSE:HG3	1.84	0.58
1:B:342:LEU:HD22	1:B:412:MSE:CE	2.34	0.57
1:B:292:ALA:HB3	1:B:434:MSE:HE1	1.85	0.57
1:A:264:LEU:HD12	1:A:445[A]:MSE:HG3	1.87	0.56
1:B:337:PRO:HG2	1:B:407:THR:HG21	1.87	0.56
1:B:95:THR:H	1:B:102:ASN:HD22	1.53	0.56
1:A:453:ASN:C	1:A:453:ASN:HD22	2.10	0.55
1:A:293:MSE:HE1	1:A:312:LEU:HD12	1.88	0.55
1:B:154:ASN:CA	4:B:554:GOL:H12	2.37	0.55
1:B:29:GLN:H	1:B:65:GLN:NE2	2.05	0.55
1:A:41:ALA:O	1:A:42:ASP:HB2	2.07	0.54
1:B:291:MSE:HE2	1:B:317:LEU:CD1	2.20	0.54
1:A:443:THR:HG23	1:A:469:LEU:HD13	1.89	0.54
1:A:291:MSE:SE	1:A:319:GLU:HG3	2.57	0.54
1:B:429:HIS:HD2	5:B:580:HOH:O	1.90	0.54
1:A:229:LEU:HD11	1:A:251:VAL:HA	1.90	0.53
1:A:263:ILE:HD13	1:A:447:SER:HB3	1.90	0.53
1:A:362:ARG:HG2	1:A:371:ILE:HD11	1.90	0.52
1:B:337:PRO:HG2	1:B:407:THR:CG2	2.39	0.52
1:A:369:GLU:HB2	5:A:638:HOH:O	2.09	0.52
1:B:26:MSE:HE1	1:B:478:LEU:HD13	1.91	0.52
1:B:486:THR:HG21	1:B:502:MSE:SE	2.60	0.51
1:B:291:MSE:HE3	1:B:319:GLU:HB2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:PHE:CZ	1:A:172:ARG:HG3	2.46	0.50
1:B:197:TRP:O	1:B:199:PRO:HD3	2.11	0.50
1:A:524:TYR:CZ	1:A:528:HIS:HE1	2.30	0.50
1:B:301:SER:HB3	5:B:904:HOH:O	2.11	0.50
1:B:255:ASP:OD1	3:B:553:EPE:H32	2.12	0.50
1:B:428:ARG:HD2	1:B:463:ALA:O	2.12	0.49
1:A:351:GLU:OE2	5:A:793:HOH:O	2.20	0.49
1:B:26:MSE:HE1	1:B:478:LEU:CD1	2.42	0.49
1:B:357:LEU:CD2	1:B:412:MSE:HE3	2.40	0.48
1:A:279:ARG:HD3	1:A:445[A]:MSE:SE	2.64	0.48
1:B:154:ASN:CB	4:B:554:GOL:H32	2.42	0.48
1:A:445[B]:MSE:HE1	1:A:505:ILE:CG1	2.44	0.48
1:A:245:PRO:HG3	5:A:726:HOH:O	2.14	0.48
1:B:120:ARG:NH1	1:B:146:LEU:HD21	2.28	0.48
3:B:553:EPE:H61	3:B:553:EPE:H101	1.50	0.48
1:B:101:ARG:NH2	1:B:104:GLN:HE21	1.96	0.48
1:A:358:ASN:HD21	1:A:453:ASN:ND2	2.12	0.47
1:A:265:GLY:HA3	1:A:499:MSE:HE2	1.97	0.47
1:B:120:ARG:NH1	1:B:150:GLN:NE2	2.63	0.47
1:A:339:TYR:HB3	1:A:340:PRO:HD3	1.97	0.47
1:B:339:TYR:HB3	1:B:340:PRO:HD3	1.97	0.47
1:A:197:TRP:O	1:A:199:PRO:HD3	2.15	0.47
1:A:279:ARG:CD	1:A:445[A]:MSE:SE	3.13	0.46
1:A:28:GLY:HA2	1:A:65:GLN:HG2	1.96	0.46
1:B:28:GLY:HA2	1:B:65:GLN:HE21	1.80	0.46
1:A:94:LEU:HG	1:A:155:THR:HG23	1.96	0.46
1:B:154:ASN:CG	4:B:554:GOL:H31	2.36	0.46
1:B:291:MSE:HE3	1:B:319:GLU:HG3	1.97	0.46
1:B:26:MSE:HE2	1:B:29:GLN:CG	2.46	0.46
1:A:121:ILE:HG12	1:A:205:VAL:HG22	1.97	0.46
1:B:69:ASN:HB3	5:B:981:HOH:O	2.15	0.45
1:A:109:TRP:CZ2	1:A:110:MSE:HE3	2.50	0.45
1:A:141:GLN:HA	1:A:141:GLN:OE1	2.16	0.45
1:B:225:MSE:N	1:B:225:MSE:HE3	2.30	0.45
1:A:174:THR:O	1:A:175:LYS:HB2	2.15	0.45
1:B:71:ILE:HD12	1:B:71:ILE:HA	1.88	0.45
1:B:280:ILE:HD11	1:B:290:HIS:HB3	1.97	0.45
1:B:120:ARG:NH1	1:B:146:LEU:HD11	2.30	0.45
1:B:59:VAL:HG21	1:B:241:MSE:CE	2.44	0.45
1:A:59:VAL:HG21	1:A:241:MSE:HE2	1.98	0.45
1:B:9:ASP:OD1	1:B:79:ASP:OD2	2.34	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:280:ILE:HD12	1:A:434:MSE:SE	2.67	0.45
1:A:65:GLN:HG3	1:A:65:GLN:O	2.16	0.45
1:A:362:ARG:CG	1:A:371:ILE:HD11	2.47	0.44
1:B:300:ILE:HG12	1:B:437:ASN:HB3	2.00	0.44
1:B:524:TYR:CZ	1:B:528:HIS:CE1	3.01	0.44
1:B:342:LEU:CD2	1:B:412:MSE:HE1	2.47	0.44
1:B:528:HIS:CD2	5:B:607:HOH:O	2.67	0.44
1:A:445[A]:MSE:HB3	1:A:445[A]:MSE:HE3	1.78	0.44
1:A:260:THR:OG1	1:A:291:MSE:HG3	2.17	0.44
1:B:41:ALA:O	1:B:42:ASP:HB2	2.17	0.44
1:B:291:MSE:HE3	1:B:319:GLU:CG	2.48	0.43
1:B:95:THR:N	1:B:102:ASN:HD22	2.16	0.43
1:B:145:LEU:HD22	1:B:162:LEU:HD13	2.01	0.43
1:A:270:THR:HG23	5:A:910:HOH:O	2.18	0.43
1:B:25:ARG:HE	1:B:25:ARG:HB2	1.55	0.43
1:A:109:TRP:CZ2	1:A:110:MSE:CE	3.02	0.42
1:A:369:GLU:OE1	1:A:512:GLN:NE2	2.52	0.42
1:A:128:VAL:HG12	1:A:191:LEU:HD21	2.02	0.42
1:B:16:ARG:HG2	1:B:31:SER:OG	2.19	0.42
1:A:304:TRP:CZ3	1:A:388:ARG:HD3	2.54	0.42
1:B:404:LYS:HE3	5:B:616:HOH:O	2.20	0.42
1:A:29:GLN:N	1:A:65:GLN:HE21	2.10	0.42
1:B:89:LYS:HD2	5:B:647:HOH:O	2.21	0.41
1:A:201:TYR:O	1:A:205:VAL:HB	2.19	0.41
1:A:280:ILE:CD1	1:A:434:MSE:SE	3.18	0.41
1:A:79:ASP:HA	1:A:253:ILE:O	2.19	0.41
1:A:193:HIS:CD2	1:A:194:GLU:HG3	2.55	0.41
1:A:69:ASN:O	1:A:72:GLN:HB2	2.20	0.41
1:A:130:GLU:OE2	5:A:651:HOH:O	2.21	0.41
1:A:214:ASN:ND2	5:A:696:HOH:O	2.53	0.41
1:B:95:THR:H	1:B:102:ASN:ND2	2.18	0.41
1:A:524:TYR:CZ	1:A:528:HIS:CE1	3.09	0.41
1:B:264:LEU:HD23	1:B:293:MSE:CE	2.49	0.41
1:A:313:PRO:O	1:A:314:GLU:HB2	2.21	0.40
1:A:279:ARG:HD3	1:A:445[A]:MSE:HG3	2.03	0.40
1:A:295:ARG:HH11	1:A:295:ARG:HD2	1.72	0.40
1:B:468:MSE:HE3	1:B:468:MSE:HB2	1.74	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	544/554 (98%)	526 (97%)	18 (3%)	0	100	100
1	B	543/554 (98%)	528 (97%)	15 (3%)	0	100	100
All	All	1087/1108 (98%)	1054 (97%)	33 (3%)	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	441/423 (104%)	427 (97%)	14 (3%)	39	14
1	B	441/423 (104%)	432 (98%)	9 (2%)	55	29
All	All	882/846 (104%)	859 (97%)	23 (3%)	46	19

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	79	ASP
1	A	120	ARG
1	A	164	ASP
1	A	172	ARG
1	A	195	ASP
1	A	205	VAL
1	A	210	LEU
1	A	211	LEU

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Mol	Chain	Res	Type
1	A	214	ASN
1	A	225	MSE
1	A	272	GLU
1	A	369	GLU
1	A	452	LYS
1	A	453	ASN
1	B	23	GLN
1	B	25	ARG
1	B	31	SER
1	B	214	ASN
1	B	225	MSE
1	B	390	PRO
1	B	453	ASN
1	B	504	ARG
1	B	542	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
1	A	65	GLN
1	A	151	HIS
1	A	154	ASN
1	A	336	HIS
1	A	345	GLN
1	A	453	ASN
1	A	514	ASN
1	A	528	HIS
1	B	23	GLN
1	B	65	GLN
1	B	69	ASN
1	B	102	ASN
1	B	104	GLN
1	B	235	GLN
1	B	336	HIS
1	B	345	GLN
1	B	429	HIS
1	B	453	ASN
1	B	514	ASN
1	B	528	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
4	GOL	B	554	-	5,5,5	0.46	0	5,5,5	0.51	0
3	EPE	B	553	-	15,15,15	0.83	1 (6%)	18,20,20	2.10	6 (33%)
2	XUL	A	553	-	7,9,9	0.51	0	5,11,11	0.99	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
4	GOL	B	554	-	-	4/4/4/4	-
3	EPE	B	553	-	-	6/9/19/19	0/1/1/1
2	XUL	A	553	-	1/1/3/3	8/12/12/12	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	553	EPE	C10-S	2.90	1.81	1.77

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	553	EPE	C5-N4-C3	5.50	121.21	108.83
3	B	553	EPE	C7-N4-C3	3.46	120.07	111.23
3	B	553	EPE	C2-C3-N4	2.97	116.73	110.64
3	B	553	EPE	C7-N4-C5	2.61	117.91	111.23
3	B	553	EPE	O2S-S-C10	2.29	109.67	106.92
3	B	553	EPE	O3S-S-C10	2.05	109.08	105.77

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	553	XUL	C3

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	554	GOL	C1-C2-C3-O3
3	B	553	EPE	C10-C9-N1-C6
3	B	553	EPE	C8-C7-N4-C3
3	B	553	EPE	S-C10-C9-N1
3	B	553	EPE	C9-C10-S-O1S
2	A	553	XUL	O1-C1-C2-C3
2	A	553	XUL	O1-C1-C2-O2
2	A	553	XUL	C1-C2-C3-C4
2	A	553	XUL	O2-C2-C3-O3
2	A	553	XUL	C2-C3-C4-C5
2	A	553	XUL	C2-C3-C4-O4
2	A	553	XUL	O3-C3-C4-O4
3	B	553	EPE	C9-C10-S-O3S
4	B	554	GOL	O2-C2-C3-O3
4	B	554	GOL	O1-C1-C2-O2
2	A	553	XUL	O3-C3-C4-C5
3	B	553	EPE	C9-C10-S-O2S
4	B	554	GOL	O1-C1-C2-C3

There are no ring outliers.

2 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	554	GOL	11	0
3	B	553	EPE	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	517/554 (93%)	-0.02	8 (1%) 73 72	12, 22, 32, 44	0
1	B	518/554 (93%)	-0.02	11 (2%) 63 62	13, 23, 34, 50	0
All	All	1035/1108 (93%)	-0.02	19 (1%) 68 67	12, 22, 34, 50	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	270	THR	5.9
1	A	273	ASN	5.0
1	B	1	LEU	4.9
1	B	269	VAL	4.7
1	A	271	GLY	3.8
1	B	204	LEU	3.4
1	A	272	GLU	3.1
1	B	2	ALA	3.0
1	A	269	VAL	3.0
1	B	271	GLY	3.0
1	B	542	GLN	3.0
1	A	274	ALA	2.5
1	A	90	GLU	2.5
1	B	276	PHE	2.3
1	B	71	ILE	2.3
1	B	274	ALA	2.2
1	B	270	THR	2.2
1	B	359	TYR	2.2
1	A	493	GLU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	GOL	B	554	6/6	0.84	0.26	27,30,31,31	0
3	EPE	B	553	15/15	0.86	0.33	19,24,26,27	15
2	XUL	A	553	10/10	0.87	0.12	23,26,29,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.