



Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 03:42 AM GMT

PDB ID : 1ISW
Title : Crystal structure of xylanase from Streptomyces olivaceoviridis E-86 complexed with xylobiose
Authors : Fujimoto, Z.; Kuno, A.; Kaneko, S.; Kobayashi, H.; Kusakabe, I.; Mizuno, H.
Deposited on : 2001-12-27
Resolution : 2.10 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>

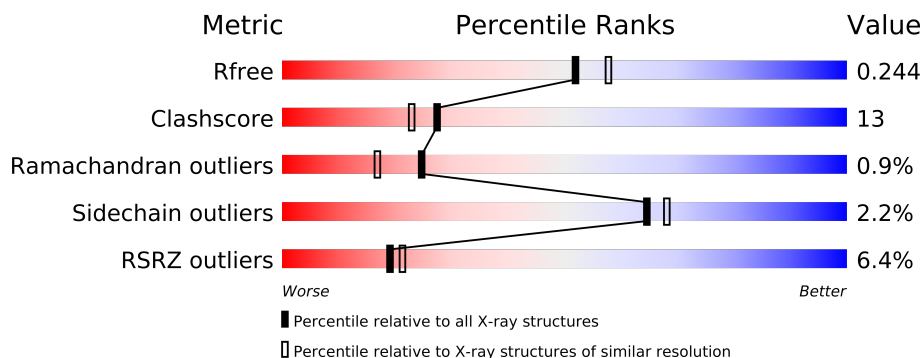
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.10 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}	66092	3012 (2.10-2.10)
Clashscore	79885	3649 (2.10-2.10)
Ramachandran outliers	78287	3610 (2.10-2.10)
Sidechain outliers	78261	3611 (2.10-2.10)
RSRZ outliers	66119	3013 (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 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	436	
1	B	436	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7466 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.

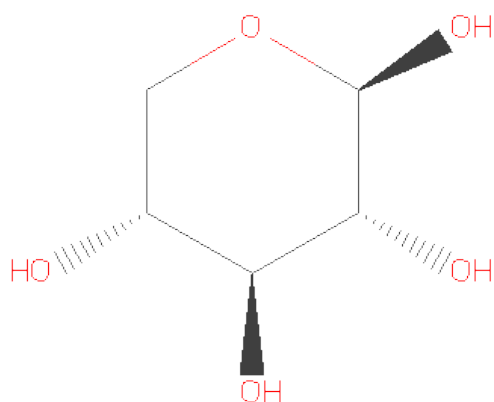
- Molecule 1 is a protein called endo-1,4-beta-D-xylanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	436	Total	C	N	O	S	0	4	0
			3297	2024	600	656	17			
1	B	436	Total	C	N	O	S	0	5	0
			3294	2024	597	656	17			

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

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	2	Total	C	O	0	0
			19	10	9		
2	A	2	Total	C	O	0	0
			19	10	9		
2	A	2	Total	C	O	0	0
			19	10	9		
2	B	2	Total	C	O	0	0
			19	10	9		
2	B	2	Total	C	O	0	0
			19	10	9		

- Molecule 3 is SUGAR (XYLOPYRANOSE) (three-letter code: XYP) (formula: C₅H₁₀O₅).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	357	Total	O	0	0
			357	357		
4	B	383	Total	O	0	0
			383	383		

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.

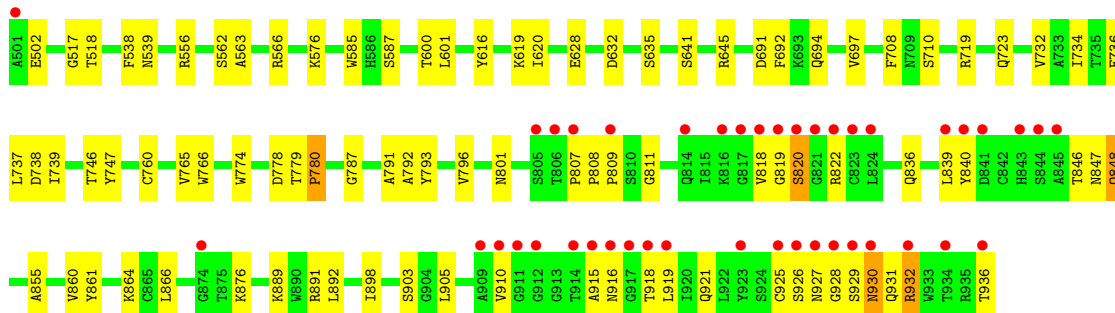
- Molecule 1: endo-1,4-beta-D-xylanase

Chain A: 



- Molecule 1: endo-1,4-beta-D-xylanase

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	74.96Å 94.26Å 137.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.74 – 2.10 44.70 – 2.00	Depositor EDS
% Data completeness (in resolution range)	87.8 (29.74-2.10) 86.5 (44.70-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.69 (at 2.00Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.194 , 0.242 0.198 , 0.244	Depositor DCC
R_{free} test set	5881 reflections (11.34%)	DCC
Wilson B-factor (Å ²)	24.1	Xtriage
Anisotropy	0.151	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 48.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	1 of 57749 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7466	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.16% 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: XYP

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	0.32	0/3390	0.60	0/4600
1	B	0.32	0/3395	0.60	0/4607
All	All	0.32	0/6785	0.60	0/9207

There are no bond length outliers.

There are no bond angle outliers.

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	3297	0	3098	81	0
1	B	3294	0	3084	87	0
2	A	57	0	51	3	0
2	B	38	0	34	3	0
3	A	20	0	20	3	0
3	B	20	0	20	2	0
4	A	357	0	0	9	0
4	B	383	0	0	7	0
All	All	7466	0	6307	168	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 13.

All (168) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:50:ASP:HA	1:A:90:PRO:HG3	1.43	1.00
1:A:24[B]:LYS:HE2	1:A:24[B]:LYS:NZ	1.33	0.98
1:A:24[B]:LYS:HE3	1:A:24[B]:LYS:NZ	1.33	0.97
1:B:925:CYS:HA	1:B:931:GLN:HE22	1.36	0.88
1:A:144:GLN:HE22	1:A:148:ASN:HA	1.37	0.88
1:B:616:TYR:HB3	1:B:620[A]:ILE:HD13	1.56	0.85
1:B:616:TYR:HB3	1:B:620[A]:ILE:CD1	2.09	0.81
1:A:144:GLN:NE2	1:A:148:ASN:HA	2.04	0.72
1:A:234:ILE:HD12	1:A:263:ILE:HG12	1.72	0.72
1:A:2:GLU:HB2	1:A:301:ASN:OD1	1.90	0.72
1:B:708:PHE:HB2	1:B:739[A]:ILE:HD12	1.71	0.72
1:A:49:ILE:HG22	1:A:90:PRO:HD3	1.71	0.72
1:A:174:ILE:H	1:A:174:ILE:HD13	1.55	0.70
1:A:55:GLN:HB3	1:A:58:GLN:HB2	1.74	0.70
1:A:66:ARG:NH1	4:A:1818:HOH:O	2.25	0.69
1:B:929:SER:HA	1:B:932:ARG:HE	1.57	0.68
1:A:236:GLU:OE1	2:A:452:XYP:H1B	1.93	0.68
1:B:839:LEU:HD12	1:B:918:THR:OG1	1.94	0.67
1:B:818:VAL:HG23	1:B:819:GLY:N	2.09	0.67
1:A:236:GLU:HG2	1:A:266:TRP:CE3	2.30	0.67
1:B:819:GLY:HA3	1:B:929:SER:HB2	1.77	0.67
1:B:903:SER:HB2	1:B:905:LEU:HD13	1.76	0.66
1:A:1:ALA:HB1	1:A:7:ALA:HB1	1.78	0.65
1:B:818:VAL:HG23	1:B:819:GLY:H	1.62	0.64
1:B:836:GLN:HB2	1:B:876:LYS:HE3	1.80	0.64
1:B:840:TYR:CE2	3:B:961:XYP:H1B	2.33	0.63
1:A:102:ARG:O	1:A:106:ILE:HG12	1.99	0.63
1:B:719:ARG:O	1:B:723:GLN:HG3	2.00	0.62
1:B:736:GLU:OE1	2:B:952:XYP:H1B	2.01	0.61
1:B:915:ALA:O	1:B:918:THR:HG23	2.01	0.59
1:A:238:ASP:HB2	1:A:280:PRO:HB2	1.85	0.59
1:B:502:GLU:HB2	1:B:801:ASN:OD1	2.03	0.58
1:A:435:ARG:HD3	4:A:1726:HOH:O	2.03	0.57
1:B:556:ARG:HH12	1:B:600:THR:HG22	1.69	0.57
1:A:426:SER:O	1:A:427:ASN:HB2	2.05	0.57
1:B:635:SER:HB3	4:B:1335:HOH:O	2.03	0.57
1:B:778:ASP:O	1:B:779:THR:C	2.43	0.56
1:A:340:TYR:CZ	3:A:1461:XYP:H5B2	2.40	0.56
1:A:231:ASP:OD1	1:A:259:ARG:HG3	2.04	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:925:CYS:HA	1:B:931:GLN:NE2	2.12	0.56
1:B:910:VAL:HG23	1:B:919:LEU:O	2.05	0.56
1:A:274:TRP:CZ3	2:A:452:XYP:H5B1	2.41	0.56
1:A:49:ILE:HG22	1:A:90:PRO:CD	2.36	0.55
1:B:891:ARG:NH1	4:B:1330:HOH:O	2.28	0.55
1:A:190:ARG:O	1:A:194:GLN:HG3	2.06	0.55
1:A:1:ALA:CB	1:A:7:ALA:HB1	2.37	0.55
1:B:738:ASP:HB2	1:B:780:PRO:HB2	1.88	0.54
1:B:846:THR:C	1:B:848:GLN:H	2.10	0.54
1:B:822:ARG:HG2	1:B:822:ARG:HH11	1.72	0.54
1:B:820:SER:OG	1:B:839:LEU:HD13	2.08	0.54
1:A:311:GLY:HA2	1:A:351:THR:HG23	1.89	0.54
1:A:172:TYR:HB3	1:A:205:GLN:OE1	2.07	0.54
1:A:105:MET:HE3	1:A:143:LEU:HD22	1.90	0.53
1:B:930:ASN:N	1:B:930:ASN:HD22	2.06	0.53
1:B:860:VAL:O	1:B:861:TYR:HB2	2.09	0.53
1:B:563:ALA:HA	1:B:566:ARG:NH1	2.24	0.53
1:A:24[A]:LYS:HE2	1:A:272:ASP:OD1	2.09	0.52
1:B:739[B]:ILE:HD13	1:B:746:THR:HG22	1.90	0.52
1:A:278:ASP:O	1:A:279:THR:C	2.48	0.52
1:A:279:THR:N	1:A:280:PRO:HD3	2.25	0.51
1:A:199:ILE:HD13	1:A:200:ASP:N	2.25	0.51
1:A:128:GLU:HA	1:A:171:ASP:CG	2.31	0.51
1:B:619:LYS:C	1:B:620[A]:ILE:HD12	2.31	0.51
1:B:840:TYR:CD1	1:B:916:ASN:HB3	2.46	0.51
1:A:24[B]:LYS:HE3	1:A:24[B]:LYS:HZ1	1.24	0.50
1:A:409:ALA:HB3	1:A:430:ASN:HB2	1.92	0.50
1:A:425:CYS:HA	1:A:431:GLN:OE1	2.12	0.50
1:A:101:LEU:HD21	1:A:142:ASN:OD1	2.11	0.50
1:A:294:THR:OG1	1:A:391:ARG:HD2	2.12	0.50
1:A:340:TYR:CE2	3:A:1461:XYP:H1B	2.47	0.50
1:B:820:SER:C	1:B:822:ARG:H	2.15	0.49
1:A:105:MET:CE	1:A:143:LEU:HD22	2.42	0.49
1:A:100:THR:HG23	4:A:1661:HOH:O	2.12	0.49
1:B:926:SER:C	1:B:928:GLY:H	2.15	0.49
1:A:53:GLU:OE2	1:A:56:ARG:HA	2.13	0.49
1:B:619:LYS:HE2	4:B:1281:HOH:O	2.11	0.49
1:A:1:ALA:HB3	4:A:1637:HOH:O	2.11	0.49
1:B:860:VAL:O	1:B:864:LYS:HB2	2.12	0.49
1:B:925:CYS:CA	1:B:931:GLN:HE22	2.18	0.49
1:B:616:TYR:HB3	1:B:620[A]:ILE:HD11	1.90	0.49
1:A:336:GLN:HB2	1:A:376:LYS:HE2	1.95	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:708:PHE:HB2	1:B:739[A]:ILE:CD1	2.42	0.49
1:A:24[B]:LYS:HE2	1:A:24[B]:LYS:HZ3	1.23	0.49
1:A:142:ASN:O	1:A:146:THR:HG23	2.13	0.49
1:B:779:THR:N	1:B:780:PRO:HD3	2.28	0.48
1:A:330:SER:OG	1:A:332:THR:HG23	2.13	0.48
1:B:585:TRP:CE2	1:B:587:SER:HB3	2.48	0.48
1:B:628:GLU:OE2	2:B:952:XYP:O4A	2.31	0.48
1:A:199:ILE:C	1:A:199:ILE:HD13	2.34	0.48
1:B:818:VAL:CG2	1:B:819:GLY:N	2.76	0.48
1:A:85:TRP:CE2	1:A:87:SER:HB3	2.49	0.48
1:B:556:ARG:HH12	1:B:600:THR:CG2	2.26	0.48
1:A:127:ASN:ND2	1:A:128:GLU:HG3	2.29	0.48
1:A:149:ASP:O	1:A:153:VAL:HG23	2.13	0.48
1:A:368:ALA:HB1	1:A:422:LEU:HD11	1.96	0.48
1:A:374:GLY:HA2	1:A:421:GLN:OE1	2.13	0.47
1:A:2:GLU:CB	1:A:301:ASN:OD1	2.62	0.47
1:B:855:ALA:O	1:B:889:LYS:HD2	2.13	0.47
1:A:236:GLU:HG2	1:A:266:TRP:CZ3	2.49	0.47
1:A:387:ASN:HB3	1:A:402:GLN:OE1	2.15	0.47
1:B:738:ASP:O	1:B:739[A]:ILE:HD13	2.13	0.47
1:A:391:ARG:NH2	4:A:1822:HOH:O	2.48	0.47
1:A:189:VAL:O	1:A:193:LYS:HG2	2.15	0.47
1:B:926:SER:C	1:B:928:GLY:N	2.69	0.46
1:B:930:ASN:HD22	1:B:930:ASN:H	1.63	0.46
1:B:708:PHE:HD2	1:B:739[A]:ILE:CD1	2.28	0.46
1:B:819:GLY:CA	1:B:929:SER:HB2	2.45	0.46
1:A:308:PRO:HA	1:A:309:PRO:HD3	1.83	0.46
1:B:926:SER:O	1:B:928:GLY:N	2.49	0.46
1:A:291:ALA:HA	1:A:391:ARG:HG2	1.98	0.45
1:B:539:ASN:HA	1:B:576:LYS:HG3	1.98	0.45
1:B:562:SER:O	1:B:566:ARG:HG3	2.17	0.45
1:B:791:ALA:HA	1:B:891:ARG:HG2	1.99	0.44
1:B:818:VAL:CG2	1:B:819:GLY:H	2.28	0.44
1:B:517:GLY:HA2	1:B:538:PHE:HB3	1.99	0.44
1:B:921:GLN:CD	4:B:1324:HOH:O	2.55	0.44
1:B:840:TYR:HA	1:B:916:ASN:OD1	2.17	0.44
1:A:174:ILE:N	1:A:174:ILE:HD13	2.28	0.44
1:A:66:ARG:NH1	4:A:1623:HOH:O	2.51	0.44
1:A:171:ASP:O	1:A:205:GLN:HG3	2.18	0.44
1:A:307:PRO:HA	1:A:308:PRO:HD3	1.92	0.44
1:B:787:GLY:HA3	4:B:1152:HOH:O	2.18	0.44
1:B:932:ARG:HH11	1:B:932:ARG:HG2	1.82	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:384:GLY:O	1:A:389:LYS:HE2	2.18	0.43
1:B:866:LEU:HA	1:B:866:LEU:HD12	1.94	0.43
1:B:691:ASP:HA	1:B:694:GLN:HG2	2.01	0.43
1:B:732:VAL:HG23	1:B:760:CYS:HA	2.00	0.43
1:B:692:PHE:HB3	1:B:697:VAL:HB	2.00	0.43
1:B:518:THR:HA	1:B:765:VAL:O	2.19	0.43
1:A:79:ARG:HB3	1:A:79:ARG:NH1	2.33	0.43
1:B:739[B]:ILE:HD13	1:B:746:THR:CG2	2.49	0.43
1:B:566:ARG:NH1	4:B:1118:HOH:O	2.52	0.42
1:B:774:TRP:HH2	2:B:956:XYP:H4B	1.85	0.42
1:B:807:PRO:HA	1:B:808:PRO:HD3	1.84	0.42
1:B:819:GLY:HA3	1:B:929:SER:CB	2.48	0.42
1:A:340:TYR:CE1	3:A:1461:XYP:H5B2	2.54	0.42
1:B:892:LEU:CD2	1:B:898:ILE:HG12	2.49	0.42
1:B:641:SER:O	1:B:645:ARG:HG3	2.19	0.42
1:B:738:ASP:C	1:B:739[A]:ILE:HD13	2.40	0.42
1:B:929:SER:HA	1:B:932:ARG:NE	2.30	0.42
1:A:309:PRO:HB3	1:A:436:THR:O	2.19	0.42
1:B:892:LEU:HD23	1:B:898:ILE:HG12	2.01	0.42
1:B:793:TYR:C	1:B:793:TYR:CD1	2.93	0.42
1:A:49:ILE:CG2	1:A:90:PRO:HD3	2.47	0.42
1:A:378:GLN:NE2	1:B:710:SER:HB2	2.35	0.42
1:A:259:ARG:HB2	4:A:1556:HOH:O	2.19	0.42
1:B:734:ILE:CG2	1:B:737:LEU:HB2	2.50	0.41
1:B:736:GLU:HG2	1:B:766:TRP:CE3	2.55	0.41
1:B:910:VAL:HA	4:B:1324:HOH:O	2.20	0.41
1:A:419:LEU:C	1:A:420[A]:ILE:HD12	2.41	0.41
1:A:232:VAL:HG23	1:A:260:CYS:HA	2.03	0.41
1:B:822:ARG:HB2	1:B:839:LEU:HB3	2.03	0.41
1:B:936:THR:HG22	1:B:936:THR:OXT	2.20	0.41
1:A:68:TYR:CZ	1:A:72:VAL:HG21	2.55	0.41
1:B:792:ALA:O	1:B:796:VAL:HG23	2.21	0.41
1:A:218:PHE:HB2	4:A:1691:HOH:O	2.21	0.41
2:A:472:XYP:H2B	4:A:1633:HOH:O	2.20	0.41
1:B:840:TYR:CZ	3:B:961:XYP:H5B2	2.56	0.41
1:B:929:SER:CA	1:B:932:ARG:HE	2.28	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	438/436 (100%)	417 (95%)	18 (4%)	3 (1%)	30	23
1	B	439/436 (101%)	415 (94%)	19 (4%)	5 (1%)	21	13
All	All	877/872 (101%)	832 (95%)	37 (4%)	8 (1%)	25	17

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	311	GLY
1	B	820	SER
1	B	927	ASN
1	A	2	GLU
1	B	811	GLY
1	B	809	PRO
1	B	847	ASN
1	A	312	GLY

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	345/341 (101%)	337 (98%)	8 (2%)	63	66
1	B	346/341 (102%)	339 (98%)	7 (2%)	68	72
All	All	691/682 (101%)	676 (98%)	15 (2%)	64	68

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	174	ILE

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Mol	Chain	Res	Type
1	A	199	ILE
1	A	243	SER
1	A	247	TYR
1	A	259	ARG
1	A	274	TRP
1	A	280	PRO
1	A	332	THR
1	B	601	LEU
1	B	632	ASP
1	B	747	TYR
1	B	780	PRO
1	B	848	GLN
1	B	930	ASN
1	B	932	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	GLN
1	A	69	ASN
1	A	173	ASN
1	A	284	ASN
1	A	314	GLN
1	A	343	HIS
1	B	511	GLN
1	B	603	GLN
1	B	673	ASN
1	B	784	ASN
1	B	814	GLN
1	B	927	ASN
1	B	930	ASN
1	B	931	GLN

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 ⓘ

10 carbohydrates are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	XYP	A	451	2	8,9,10	0.56	0	8,12,14	0.75	0
2	XYP	A	452	2	10,10,10	0.55	0	14,14,14	0.61	0
2	XYP	A	456	2	8,9,10	0.65	0	8,12,14	0.72	0
2	XYP	A	457	2	10,10,10	0.57	0	14,14,14	0.59	0
2	XYP	A	471	2	8,9,10	0.54	0	8,12,14	0.74	0
2	XYP	A	472	2	10,10,10	0.67	0	14,14,14	0.59	0
2	XYP	B	951	2	8,9,10	0.55	0	8,12,14	0.72	0
2	XYP	B	952	2	10,10,10	0.60	0	14,14,14	0.66	0
2	XYP	B	956	2	8,9,10	0.67	0	8,12,14	0.72	0
2	XYP	B	957	2	10,10,10	0.56	0	14,14,14	0.61	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	XYP	A	451	2	-	0/0/14/17	0/1/1/1
2	XYP	A	452	2	-	0/0/17/17	0/1/1/1
2	XYP	A	456	2	-	0/0/14/17	0/1/1/1
2	XYP	A	457	2	-	0/0/17/17	0/1/1/1
2	XYP	A	471	2	-	0/0/14/17	0/1/1/1
2	XYP	A	472	2	-	0/0/17/17	0/1/1/1
2	XYP	B	951	2	-	0/0/14/17	0/1/1/1
2	XYP	B	952	2	-	0/0/17/17	0/1/1/1
2	XYP	B	956	2	-	0/0/14/17	0/1/1/1
2	XYP	B	957	2	-	0/0/17/17	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.6 Ligand geometry

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	XYP	A	1461	-	10,10,10	0.55	0	14,14,14	0.60	0
3	XYP	A	1491	-	10,10,10	0.64	0	14,14,14	0.60	0
3	XYP	B	961	-	10,10,10	0.66	0	14,14,14	0.60	0
3	XYP	B	971	-	10,10,10	0.67	0	14,14,14	0.60	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	XYP	A	1461	-	-	0/0/17/17	0/1/1/1
3	XYP	A	1491	-	-	0/0/17/17	0/1/1/1
3	XYP	B	961	-	-	0/0/17/17	0/1/1/1
3	XYP	B	971	-	-	0/0/17/17	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	436/436 (100%)	0.06	15 (3%) 43 47	14, 23, 43, 83	0
1	B	436/436 (100%)	0.16	42 (9%) 8 9	14, 22, 65, 77	0
All	All	872/872 (100%)	0.11	57 (6%) 19 20	14, 23, 57, 83	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	306	THR	8.6
1	A	305	SER	7.7
1	B	915	ALA	6.2
1	B	818	VAL	5.8
1	A	304	SER	5.7
1	B	822	ARG	5.4
1	B	936	THR	5.3
1	B	820	SER	5.2
1	A	436	THR	5.0
1	A	303	GLY	4.5
1	B	819	GLY	4.2
1	B	923	TYR	4.0
1	A	309	PRO	3.9
1	B	843	HIS	3.9
1	B	911	GLY	3.9
1	B	823	CYS	3.8
1	B	839	LEU	3.8
1	B	840	TYR	3.7
1	B	844	SER	3.7
1	B	928	GLY	3.6
1	A	307	PRO	3.5
1	A	97	SER	3.5
1	B	845	ALA	3.3
1	A	308	PRO	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	932	ARG	3.3
1	B	805	SER	3.2
1	B	909	ALA	3.2
1	B	929	SER	3.2
1	B	816	LYS	3.1
1	B	917	GLY	3.1
1	B	918	THR	3.1
1	B	910	VAL	3.0
1	B	874	GLY	2.9
1	A	100	THR	2.9
1	B	807	PRO	2.9
1	B	930	ASN	2.8
1	A	311	GLY	2.8
1	B	914	THR	2.8
1	B	824	LEU	2.7
1	B	926	SER	2.7
1	B	809	PRO	2.6
1	B	912	GLY	2.6
1	A	133	ASP	2.6
1	B	919	LEU	2.6
1	B	501	ALA	2.5
1	B	817	GLY	2.5
1	B	916	ASN	2.5
1	B	841	ASP	2.4
1	B	821	GLY	2.4
1	B	814	GLN	2.4
1	B	927	ASN	2.3
1	B	806	THR	2.3
1	A	344	SER	2.2
1	B	925	CYS	2.2
1	A	95	SER	2.2
1	B	934	THR	2.2
1	A	198	PRO	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 ⓘ

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
2	XYP	A	456	9/10	0.23	13.71	60,60,62,62	0
2	XYP	A	472	10/10	0.22	6.81	37,51,52,55	0
2	XYP	A	457	10/10	0.23	4.87	57,59,60,61	0
2	XYP	A	452	10/10	0.15	4.80	26,33,36,41	0
2	XYP	B	952	10/10	0.12	3.03	16,22,27,32	0
2	XYP	B	956	9/10	0.21	2.27	48,50,51,52	0
2	XYP	B	957	10/10	0.20	2.13	42,45,46,46	0
2	XYP	A	471	9/10	0.10	0.99	23,27,29,30	0
2	XYP	B	951	9/10	0.07	-1.42	16,17,19,19	0
2	XYP	A	451	9/10	0.08	-2.03	22,24,26,28	0

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
3	XYP	A	1491	10/10	0.20	0.61	57,58,59,60	0
3	XYP	B	971	10/10	0.29	0.33	54,55,55,55	0
3	XYP	A	1461	10/10	0.11	-0.15	28,30,31,33	0
3	XYP	B	961	10/10	0.21	-0.21	64,65,66,66	0

6.5 Other polymers ⓘ

There are no such residues in this entry.