



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 12, 2026 – 01:13 pm BST

PDB ID : 30CW / pdb_000030cw
Title : Class 3 L-asparaginase from Colletotrichum fructicola with active site Ser51 modified with boric acid - glycerol ester
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Deposited on : 2026-04-21
Resolution : 1.95 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 Phenix2.2
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 2.2
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.51

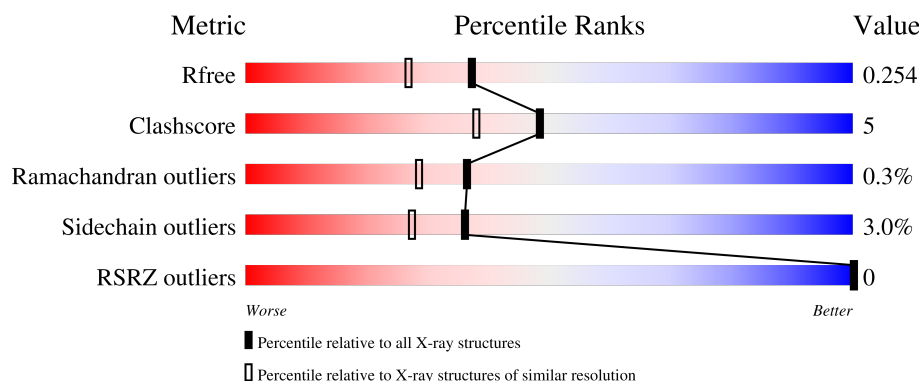
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	363	 85% 10% . .
1	D	363	 82% 14% . .
2	B	363	 81% 15% . .
2	C	363	 82% 14% .

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 10855 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 L-asparaginase ii.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	349	Total	C	N	O	S	0	1	0
			2625	1633	473	501	18			
1	D	350	Total	C	N	O	S	0	0	0
			2626	1634	473	501	18			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	GLY	-	expression tag	UNP L2FY35
A	-4	ILE	-	expression tag	UNP L2FY35
A	-3	ASP	-	expression tag	UNP L2FY35
A	-2	PRO	-	expression tag	UNP L2FY35
A	-1	PHE	-	expression tag	UNP L2FY35
A	0	THR	-	expression tag	UNP L2FY35
D	-5	GLY	-	expression tag	UNP L2FY35
D	-4	ILE	-	expression tag	UNP L2FY35
D	-3	ASP	-	expression tag	UNP L2FY35
D	-2	PRO	-	expression tag	UNP L2FY35
D	-1	PHE	-	expression tag	UNP L2FY35
D	0	THR	-	expression tag	UNP L2FY35

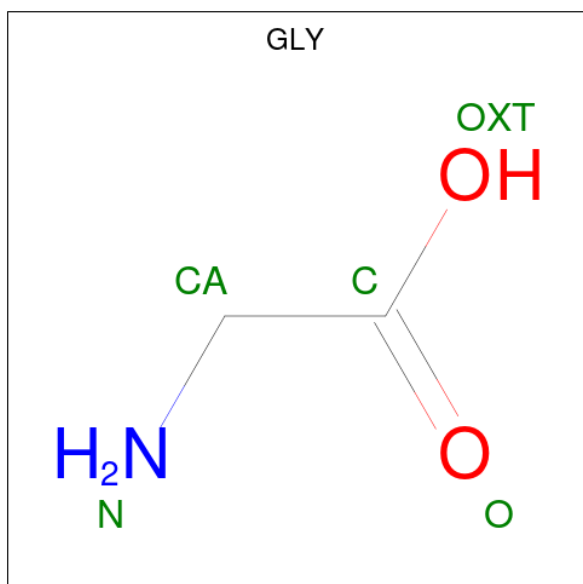
- Molecule 2 is a protein called L-asparaginase ii.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	349	Total 2630	B 1	C 1634	N 473	O 504	S 18	0	1	0
2	C	349	Total 2630	B 1	C 1634	N 473	O 504	S 18	0	1	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-5	GLY	-	expression tag	UNP L2FY35
B	-4	ILE	-	expression tag	UNP L2FY35
B	-3	ASP	-	expression tag	UNP L2FY35
B	-2	PRO	-	expression tag	UNP L2FY35
B	-1	PHE	-	expression tag	UNP L2FY35
B	0	THR	-	expression tag	UNP L2FY35
C	-5	GLY	-	expression tag	UNP L2FY35
C	-4	ILE	-	expression tag	UNP L2FY35
C	-3	ASP	-	expression tag	UNP L2FY35
C	-2	PRO	-	expression tag	UNP L2FY35
C	-1	PHE	-	expression tag	UNP L2FY35
C	0	THR	-	expression tag	UNP L2FY35

- Molecule 3 is GLYCINE (CCD ID: GLY) (formula: $C_2H_5NO_2$).

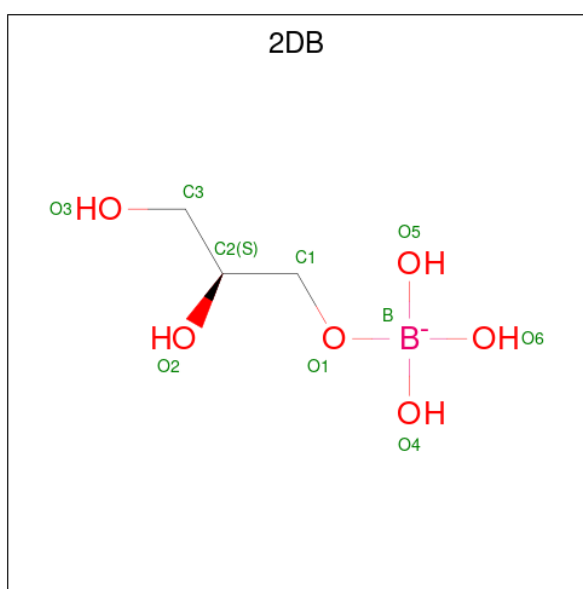


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			5	2	1	2		
3	C	1	Total	C	N	O	0	0
			5	2	1	2		
3	D	1	Total	C	N	O	0	0
			5	2	1	2		

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

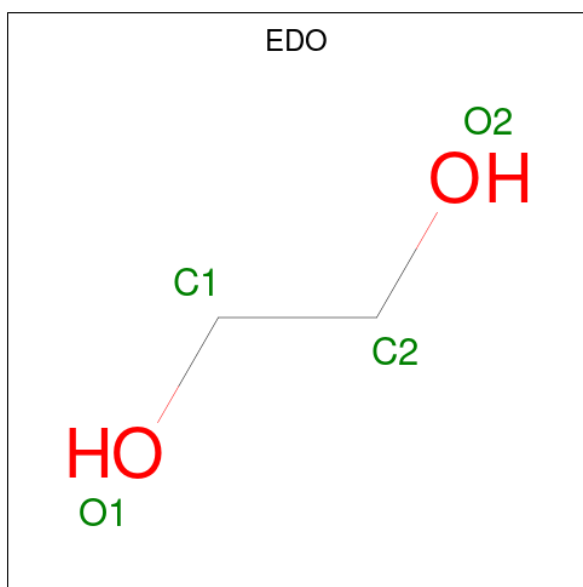
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		
4	B	1	Total	Zn	0	0
			1	1		
4	C	1	Total	Zn	0	0
			1	1		
4	D	1	Total	Zn	0	0
			1	1		

- Molecule 5 is (S)-(2,3-DIHYDROXYPROPOXY)TRIHYDROXYBORATE (CCD ID: 2DB) (formula: $C_3H_{10}BO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	B	C	O	0	0
			9	1	3	5		
5	D	1	Total	B	C	O	0	0
			9	1	3	5		

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

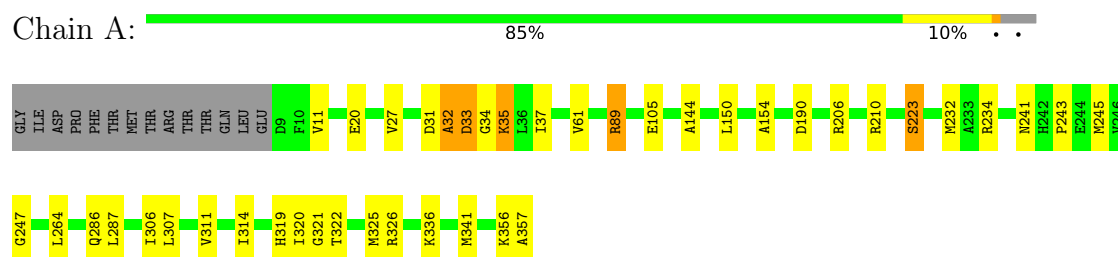
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	77	Total	O	0	0
			77	77		
7	B	79	Total	O	0	0
			79	79		
7	C	64	Total	O	0	0
			64	64		
7	D	79	Total	O	0	0
			79	79		

3 Residue-property plots [i](#)

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.

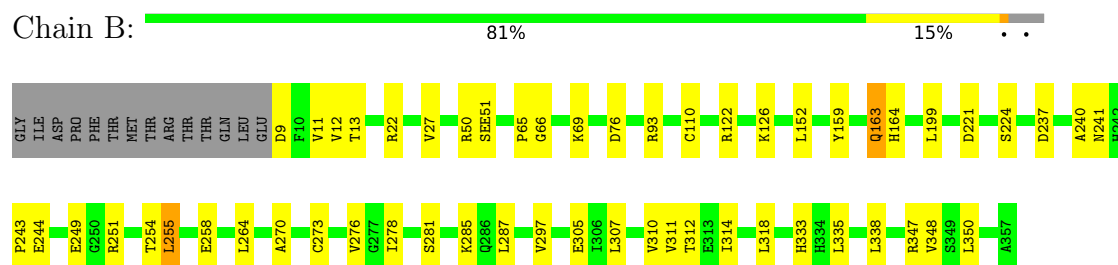
• Molecule 1: L-asparaginase ii



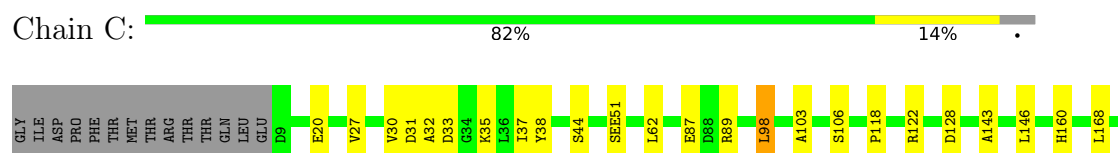
• Molecule 1: L-asparaginase ii



• Molecule 2: L-asparaginase ii



• Molecule 2: L-asparaginase ii





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	88.64Å 71.35Å 212.35Å 90.00° 97.14° 90.00°	Depositor
Resolution (Å)	70.23 – 1.95 70.23 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.3 (70.23-1.95) 99.5 (70.23-1.95)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 1.95Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.200 , 0.252 0.204 , 0.254	Depositor DCC
R_{free} test set	1000 reflections (1.04%)	wwPDB-VP
Wilson B-factor (Å ²)	32.3	Xtriage
Anisotropy	0.377	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 38.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10855	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.38% 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: ZN, SEE, EDO, 2DB

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	0.55	0/2672	0.70	0/3605
1	D	0.49	0/2673	0.65	0/3606
2	B	0.49	0/2666	0.66	0/3595
2	C	0.50	0/2666	0.66	0/3595
All	All	0.51	0/10677	0.67	0/14401

There are no bond length outliers.

There are no bond angle outliers.

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	2625	0	2573	25	0
1	D	2626	0	2576	34	0
2	B	2630	0	2578	29	0
2	C	2630	0	2578	29	0
3	A	5	0	2	0	0
3	C	5	0	2	0	0
3	D	5	0	2	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	9	0	7	0	0
5	D	9	0	7	2	0
6	C	8	0	12	0	0
7	A	77	0	0	5	0
7	B	79	0	0	2	0
7	C	64	0	0	1	0
7	D	79	0	0	3	0
All	All	10855	0	10337	114	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 5.

All (114) 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:A:322:THR:H	1:A:325:MET:HE3	1.37	0.89
2:B:347:ARG:NH1	7:B:501:HOH:O	2.21	0.74
1:A:326:ARG:NH2	7:A:503:HOH:O	2.21	0.67
1:A:341:MET:HE3	2:B:110:CYS:C	2.23	0.64
1:D:171:LYS:NZ	1:D:175:GLU:OE2	2.28	0.63
2:B:13:THR:HG22	2:B:348:VAL:HG22	1.81	0.62
1:D:230:ARG:NH2	7:D:503:HOH:O	2.32	0.61
1:D:270:ALA:N	5:D:403:2DB:H12	2.15	0.61
2:B:312:THR:HG21	2:B:333:HIS:HB2	1.83	0.60
1:D:76:ASP:OD1	1:D:93:ARG:NH1	2.35	0.59
2:C:321:GLY:O	2:C:326:ARG:NH1	2.36	0.59
2:B:27:VAL:HG21	2:B:314:ILE:HD11	1.86	0.58
2:B:264:LEU:C	2:B:264:LEU:HD12	2.29	0.58
2:C:322:THR:H	2:C:325:MET:CE	2.16	0.57
1:D:199:LEU:N	1:D:199:LEU:HD23	2.19	0.57
2:B:12:VAL:HG22	2:B:22:ARG:HG2	1.86	0.57
2:B:65:PRO:HD2	2:B:152:LEU:HD23	1.88	0.56
1:D:80:MET:HE3	1:D:144:ALA:HB1	1.88	0.56
1:A:321:GLY:O	1:A:326:ARG:NH1	2.39	0.56
1:A:31:ASP:HB3	1:A:35:LYS:HG2	1.88	0.55
2:B:9:ASP:N	7:B:504:HOH:O	2.39	0.55
1:D:119:ASN:OD1	1:D:120:VAL:N	2.39	0.55
2:C:322:THR:H	2:C:325:MET:HE3	1.72	0.54
1:D:13:THR:HG22	1:D:348:VAL:HG22	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:37:ILE:HG22	2:C:38:TYR:CD2	2.43	0.54
2:B:335:LEU:HD12	2:B:348:VAL:HG21	1.88	0.54
1:A:287:LEU:HD11	1:A:320:ILE:HD11	1.89	0.53
2:C:287:LEU:HD11	2:C:320:ILE:HD11	1.91	0.53
1:A:241:ASN:C	1:A:243:PRO:HD3	2.34	0.53
2:C:338:LEU:HD23	2:C:344:GLU:HA	1.90	0.53
2:C:98:LEU:HD13	2:C:143:ALA:HB2	1.92	0.52
1:A:37:ILE:O	1:A:356:LYS:HD2	2.10	0.51
2:B:199:LEU:HD23	2:B:199:LEU:N	2.26	0.51
2:C:31:ASP:O	2:C:33:ASP:N	2.44	0.50
1:A:34:GLY:N	1:A:206:ARG:HD3	2.26	0.50
1:A:307:LEU:O	1:A:311:VAL:HG23	2.12	0.49
2:C:87:GLU:OE1	2:C:89:ARG:HD3	2.11	0.49
2:C:160:HIS:HB2	2:C:194:LEU:HD23	1.94	0.49
2:C:27:VAL:HG21	2:C:314:ILE:HD11	1.93	0.49
2:C:264:LEU:C	2:C:264:LEU:HD12	2.37	0.48
2:B:76:ASP:OD1	2:B:93:ARG:NH1	2.42	0.48
1:A:31:ASP:O	1:A:33:ASP:N	2.47	0.48
1:A:264:LEU:HD12	1:A:264:LEU:C	2.39	0.48
2:B:251:ARG:O	2:B:255:LEU:HG	2.14	0.48
2:B:281:SER:O	2:B:285:LYS:HG3	2.14	0.48
2:C:273:CYS:HA	2:C:297:VAL:O	2.14	0.48
1:A:105:GLU:OE1	7:A:501:HOH:O	2.20	0.47
2:B:50:ARG:CZ	2:B:270:ALA:HB3	2.44	0.47
1:A:357:ALA:O	7:A:502:HOH:O	2.20	0.46
2:B:159:TYR:HA	2:B:164:HIS:CD2	2.50	0.46
2:C:313:GLU:HB2	2:C:350:LEU:HD13	1.98	0.46
1:D:74:GLU:HG2	1:D:238:ALA:HA	1.98	0.46
2:B:163:GLN:O	2:B:163:GLN:HG3	2.16	0.46
1:A:32:ALA:HB3	1:A:35:LYS:NZ	2.31	0.45
1:D:282:GLU:O	1:D:286:GLN:HG3	2.16	0.45
1:A:31:ASP:HB3	1:A:35:LYS:CG	2.46	0.45
1:A:89:ARG:HG3	1:A:245:MET:HE1	1.98	0.45
2:C:62:LEU:HD12	2:C:235:ILE:HD11	1.99	0.45
2:B:273:CYS:HA	2:B:297:VAL:O	2.17	0.45
2:C:98:LEU:HD13	2:C:143:ALA:CB	2.46	0.45
2:C:254:THR:O	2:C:258:GLU:HG3	2.17	0.45
2:C:234:ARG:NH2	7:C:508:HOH:O	2.41	0.45
1:D:80:MET:HE3	1:D:144:ALA:CB	2.47	0.44
1:D:11:VAL:HG11	1:D:306:ILE:HD12	1.99	0.44
1:D:180:LEU:HB3	1:D:184:ASP:HB2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:83:SER:HB3	5:D:403:2DB:H31	1.99	0.44
2:B:11:VAL:HG21	2:B:310:VAL:CG2	2.48	0.44
2:B:237:ASP:O	2:B:241:ASN:HB2	2.18	0.44
1:D:12:VAL:HG22	1:D:22:ARG:HG2	2.00	0.44
1:D:201:LEU:HD22	1:D:296:ALA:HB1	1.99	0.44
2:C:37:ILE:HG22	2:C:38:TYR:CE2	2.53	0.43
1:D:85:ASN:CB	1:D:87:GLU:HG3	2.48	0.43
2:C:264:LEU:HD12	2:C:264:LEU:O	2.19	0.43
1:A:11:VAL:HG11	1:A:306:ILE:HD12	2.00	0.43
1:A:61:VAL:HG11	1:A:144:ALA:HB1	2.00	0.43
1:D:89:ARG:CG	7:D:548:HOH:O	2.66	0.43
2:C:128:ASP:OD1	1:D:126:LYS:NZ	2.48	0.43
2:B:66:GLY:O	2:B:69:LYS:HG2	2.19	0.42
2:C:118:PRO:O	2:C:122:ARG:HB2	2.19	0.42
1:D:236:PHE:C	1:D:236:PHE:CD1	2.97	0.42
1:D:38:TYR:CE2	1:D:356:LYS:HG3	2.55	0.42
1:A:210:ARG:HG2	1:A:232:MET:HE1	2.01	0.42
2:C:199:LEU:N	2:C:199:LEU:HD23	2.35	0.42
2:B:244:GLU:OE1	2:B:244:GLU:N	2.40	0.42
2:B:122:ARG:O	2:B:126:LYS:HG3	2.19	0.42
1:D:46:LEU:HD12	1:D:199:LEU:C	2.45	0.42
1:D:183:GLN:CD	1:D:183:GLN:H	2.28	0.42
2:B:240:ALA:O	2:B:243:PRO:HD3	2.20	0.42
1:D:199:LEU:HD23	1:D:199:LEU:H	1.83	0.42
1:D:287:LEU:HD11	1:D:320:ILE:HD11	2.01	0.41
1:D:31:ASP:OD1	1:D:31:ASP:C	2.62	0.41
1:D:164:HIS:ND1	1:D:166:ILE:HG22	2.35	0.41
2:B:254:THR:O	2:B:258:GLU:HG3	2.20	0.41
2:C:30:VAL:HG21	2:C:206:ARG:HA	2.02	0.41
1:A:319:HIS:CD2	7:A:511:HOH:O	2.73	0.41
1:D:61:VAL:HG11	1:D:144:ALA:HB1	2.03	0.41
2:B:287:LEU:HD21	2:B:318:LEU:O	2.21	0.41
2:B:307:LEU:O	2:B:311:VAL:HG23	2.21	0.41
1:D:89:ARG:HG3	7:D:548:HOH:O	2.20	0.41
1:D:150:LEU:HD22	1:D:156:PHE:CE1	2.55	0.41
2:C:280:GLU:O	2:C:281:SER:HB3	2.20	0.41
2:C:300:GLU:OE2	1:D:22:ARG:HD2	2.20	0.41
1:A:27:VAL:HG21	1:A:314:ILE:HD11	2.02	0.41
1:A:150:LEU:HA	1:A:154:ALA:O	2.21	0.41
1:A:223:SER:O	1:A:223:SER:OG	2.37	0.41
1:D:27:VAL:HB	1:D:40:VAL:HB	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:119:ASN:OD1	1:D:119:ASN:C	2.63	0.41
2:B:276:VAL:HG12	2:B:278:ILE:HG13	2.02	0.40
2:C:35:LYS:HE2	2:C:37:ILE:HD11	2.03	0.40
1:D:281:SER:O	1:D:285:LYS:HD3	2.20	0.40
1:A:286:GLN:NE2	7:A:524:HOH:O	2.54	0.40
2:C:103:ALA:CB	2:C:146:LEU:HD13	2.51	0.40
2:B:11:VAL:HG22	2:B:350:LEU:HD23	2.04	0.40
2:C:305:GLU:HG3	2:C:306:ILE:N	2.36	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	348/363 (96%)	331 (95%)	15 (4%)	2 (1%)	21	12
1	D	348/363 (96%)	339 (97%)	8 (2%)	1 (0%)	36	28
2	B	347/363 (96%)	338 (97%)	9 (3%)	0	100	100
2	C	347/363 (96%)	334 (96%)	12 (4%)	1 (0%)	36	28
All	All	1390/1452 (96%)	1342 (96%)	44 (3%)	4 (0%)	36	28

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	32	ALA
2	C	32	ALA
1	A	247	GLY
1	D	247	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	269/281 (96%)	261 (97%)	8 (3%)	36	27
1	D	269/281 (96%)	259 (96%)	10 (4%)	30	20
2	B	268/280 (96%)	261 (97%)	7 (3%)	40	33
2	C	268/280 (96%)	261 (97%)	7 (3%)	40	33
All	All	1074/1122 (96%)	1042 (97%)	32 (3%)	36	27

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	GLU
1	A	33	ASP
1	A	35	LYS
1	A	89	ARG
1	A	190	ASP
1	A	223	SER
1	A	234	ARG
1	A	336	LYS
2	B	163	GLN
2	B	221	ASP
2	B	224	SER
2	B	249	GLU
2	B	255	LEU
2	B	305	GLU
2	B	338	LEU
2	C	20	GLU
2	C	44	SER
2	C	98	LEU
2	C	106	SER
2	C	168	LEU
2	C	190	ASP
2	C	282	GLU
1	D	8	GLU
1	D	89	ARG

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Mol	Chain	Res	Type
1	D	128	ASP
1	D	150	LEU
1	D	285	LYS
1	D	316	GLU
1	D	347	ARG
1	D	350	LEU
1	D	353	GLN
1	D	356	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	328	GLN
2	C	163	GLN
2	C	334	HIS
1	D	319	HIS
1	D	328	GLN
1	D	334	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	SEE	B	51	2	8,9,10	1.26	1 (12%)	6,12,14	0.99	0
2	SEE	C	51	2	8,9,10	1.26	1 (12%)	6,12,14	0.46	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	SEE	B	51	2	-	0/0/8/10	-
2	SEE	C	51	2	-	0/0/8/10	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	51	SEE	B-OG	-3.40	1.40	1.46
2	B	51	SEE	B-OG	-3.31	1.40	1.46

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 4 are monoatomic - leaving 7 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
6	EDO	C	403	-	3,3,3	0.44	0	2,2,2	0.50	0
5	2DB	A	403	1	7,8,9	3.61	3 (42%)	6,9,12	1.21	1 (16%)
3	GLY	A	401	4	4,4,4	1.27	1 (25%)	3,4,4	1.44	0
6	EDO	C	402	-	3,3,3	0.33	0	2,2,2	1.16	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	2DB	D	403	1	7,8,9	3.67	3 (42%)	6,9,12	1.03	1 (16%)
3	GLY	D	401	4	4,4,4	1.15	1 (25%)	3,4,4	1.64	1 (33%)
3	GLY	C	401	4	4,4,4	1.26	1 (25%)	3,4,4	1.49	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
6	EDO	C	403	-	-	1/1/1/1	-
5	2DB	A	403	1	-	1/4/7/8	-
3	GLY	A	401	4	-	0/2/2/2	-
6	EDO	C	402	-	-	0/1/1/1	-
5	2DB	D	403	1	-	4/4/7/8	-
3	GLY	D	401	4	-	0/2/2/2	-
3	GLY	C	401	4	-	0/2/2/2	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	403	2DB	B-O1	7.62	1.47	1.37
5	A	403	2DB	B-O1	7.58	1.47	1.37
5	D	403	2DB	B-O5	4.35	1.47	1.37
5	A	403	2DB	B-O4	4.30	1.47	1.37
5	D	403	2DB	B-O4	4.15	1.47	1.37
5	A	403	2DB	B-O5	3.91	1.46	1.37
3	A	401	GLY	OXT-C	-2.49	1.22	1.30
3	C	401	GLY	OXT-C	-2.29	1.23	1.30
3	D	401	GLY	OXT-C	-2.12	1.23	1.30

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	403	2DB	O5-B-O4	-2.59	110.51	119.67
5	D	403	2DB	O5-B-O4	-2.43	111.09	119.67
3	D	401	GLY	OXT-C-O	-2.05	118.19	123.30

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	D	403	2DB	O1-C1-C2-C3
5	D	403	2DB	O1-C1-C2-O2
5	D	403	2DB	C1-C2-C3-O3
5	D	403	2DB	O2-C2-C3-O3
5	A	403	2DB	C1-C2-C3-O3
6	C	403	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	403	2DB	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	349/363 (96%)	-1.14	0 100 100	18, 35, 60, 133	1 (0%)
1	D	350/363 (96%)	-1.14	0 100 100	28, 38, 56, 97	0
2	B	348/363 (95%)	-1.11	0 100 100	18, 39, 55, 106	1 (0%)
2	C	348/363 (95%)	-1.11	0 100 100	20, 37, 59, 117	1 (0%)
All	All	1395/1452 (96%)	-1.12	0 100 100	18, 37, 58, 133	3 (0%)

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [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
2	SEE	B	51	10/11	0.99	0.02	30,34,35,36	0
2	SEE	C	51	10/11	0.99	0.03	27,30,31,37	0

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($\text{\AA}^2$)	Q<0.9
6	EDO	C	403	4/4	0.97	0.06	49,50,51,60	0
3	GLY	C	401	5/5	0.99	0.04	33,37,52,54	0
3	GLY	D	401	5/5	0.99	0.04	42,46,92,93	0
5	2DB	A	403	9/10	0.99	0.04	27,35,50,51	0
5	2DB	D	403	9/10	0.99	0.05	35,42,52,63	0
6	EDO	C	402	4/4	0.99	0.03	33,35,41,41	0
3	GLY	A	401	5/5	0.99	0.03	30,31,41,43	0
4	ZN	A	402	1/1	1.00	0.01	31,31,31,31	0
4	ZN	B	401	1/1	1.00	0.02	31,31,31,31	0
4	ZN	C	404	1/1	1.00	0.01	33,33,33,33	0
4	ZN	D	402	1/1	1.00	0.01	30,30,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.