



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:43 PM UTC

PDB ID : 4PCR / pdb_00004pcr
Title : Crystal structure of Canavalia brasiliensis seed lectin (ConBr) complexed with Gamma-Aminobutyric Acid (GABA)
Authors : Delatorre, P.; Rocha, B.A.M.; Silva-Filho, J.C.; Teixeira, C.S.; Cavada, B.S.; Nascimento, K.S.; Nbreaga, R.B.; Nagano, C.S.; Sampaio, A.H.; Leal, R.B.; Neto, I.L.B.
Deposited on : 2014-04-16
Resolution : 2.15 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

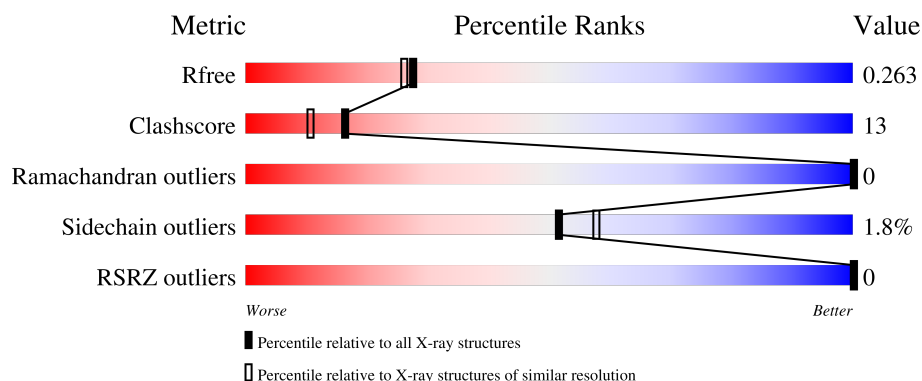
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	237	 78% 17% ..
1	D	237	 76% 22% ..

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
4	IPA	A	303	-	-	X	-
6	PEG	D	308	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 3706 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 Concanavalin-Br.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	234	Total	C	N	O	S	0	0	0
			1781	1125	300	354	2			
1	D	232	Total	C	N	O	S	0	0	0
			1767	1118	297	350	2			

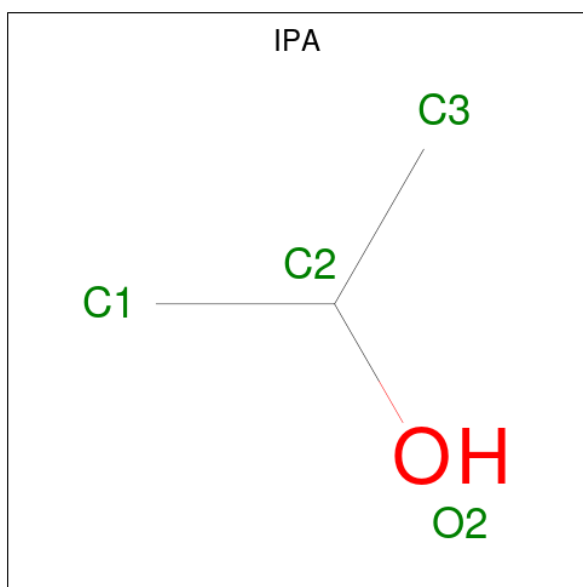
- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

- Molecule 3 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

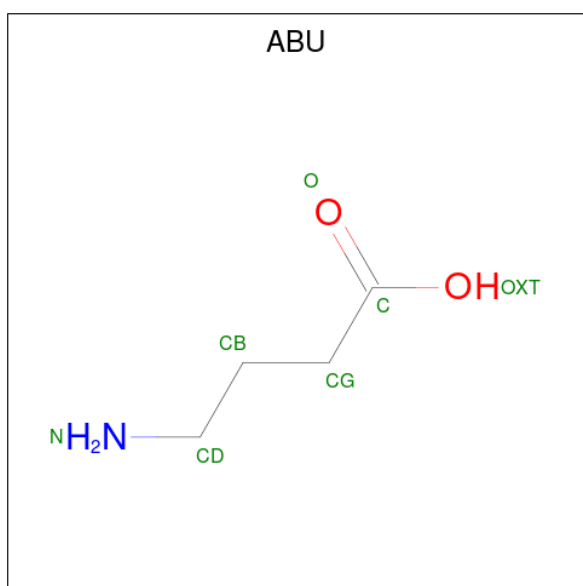
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mn	0	0
			1	1		
3	D	1	Total	Mn	0	0
			1	1		

- Molecule 4 is ISOPROPYL ALCOHOL (CCD ID: IPA) (formula: C₃H₈O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	3	1		
4	D	1	Total	C	O	0	0
			4	3	1		

- Molecule 5 is GAMMA-AMINO-BUTANOIC ACID (CCD ID: ABU) (formula: $C_4H_9NO_2$).



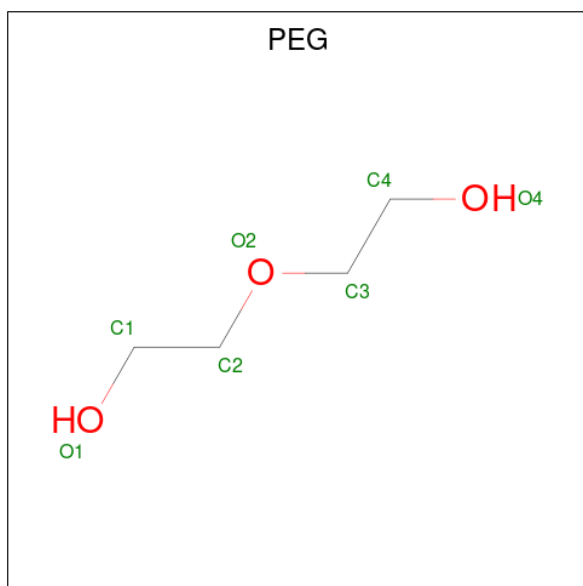
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			7	4	1	2		
5	A	1	Total	C	N	O	0	0
			7	4	1	2		

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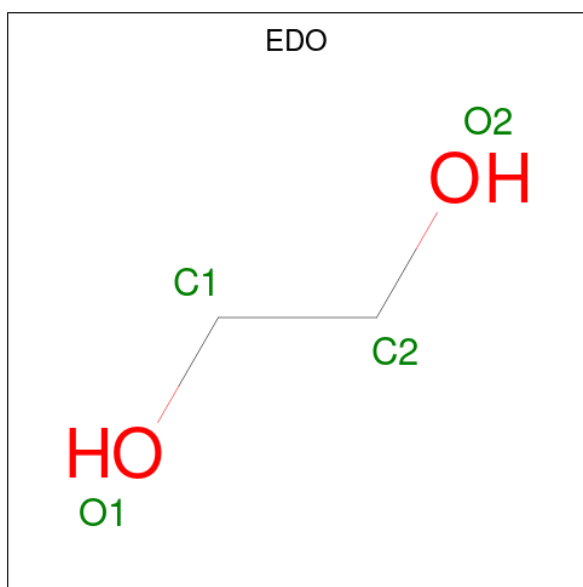
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	D	1	Total	C	N	O	0	0
			7	4	1	2		

- Molecule 6 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			7	4	3		
6	D	1	Total	C	O	0	0
			7	4	3		
6	D	1	Total	C	O	0	0
			7	4	3		

- Molecule 7 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

- Molecule 8 is water.

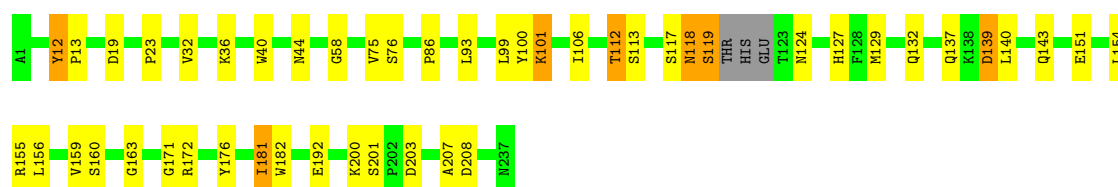
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	46	Total	O	0	0
			46	46		
8	D	50	Total	O	0	0
			50	50		

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.

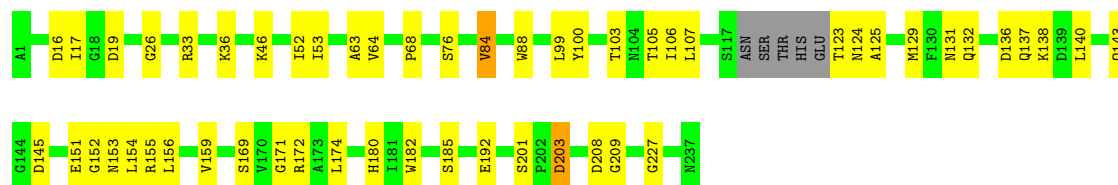
• Molecule 1: Concanavalin-Br

Chain A: 



• Molecule 1: Concanavalin-Br

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	120.41Å 72.13Å 68.39Å 90.00° 124.47° 90.00°	Depositor
Resolution (Å)	20.64 – 2.15 20.64 – 2.15	Depositor EDS
% Data completeness (in resolution range)	89.2 (20.64-2.15) 89.4 (20.64-2.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.63 (at 2.15Å)	Xtriage
Refinement program	REFMAC 5.8.0069	Depositor
R, R_{free}	0.200 , 0.236 0.223 , 0.263	Depositor DCC
R_{free} test set	2411 reflections (9.14%)	wwPDB-VP
Wilson B-factor (Å ²)	40.3	Xtriage
Anisotropy	0.102	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 31.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.455 for -h-2*1,-k,l	Xtriage
Reported twinning fraction	0.579 for H, K, L 0.421 for H+4/2L, -K, -L	Depositor
Outliers	1 of 23715 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	3706	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 7.17% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PEG, EDO, ABU, MN, CA, IPA

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	1.18	4/1821 (0.2%)	1.18	7/2477 (0.3%)
1	D	1.26	4/1807 (0.2%)	1.18	8/2458 (0.3%)
All	All	1.22	8/3628 (0.2%)	1.18	15/4935 (0.3%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	182	TRP	C-O	-7.57	1.14	1.23
1	D	63	ALA	CA-C	-6.11	1.44	1.52
1	D	209	GLY	N-CA	6.07	1.53	1.44
1	A	93	LEU	CA-C	6.06	1.60	1.52
1	A	58	GLY	C-O	-5.59	1.17	1.24
1	D	64	VAL	C-O	5.47	1.29	1.24
1	A	112	THR	C-O	-5.11	1.17	1.23
1	A	12	TYR	CA-C	5.01	1.56	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	118	ASN	N-CA-C	11.07	123.35	111.28
1	D	84	VAL	CB-CA-C	-9.18	100.71	110.62
1	A	19	ASP	CA-C-N	-7.01	113.48	120.21
1	A	19	ASP	C-N-CA	-7.01	113.48	120.21
1	D	19	ASP	CA-C-N	-6.28	111.99	119.84
1	D	19	ASP	C-N-CA	-6.28	111.99	119.84
1	D	84	VAL	N-CA-C	5.88	119.41	113.47
1	D	185	SER	N-CA-C	5.78	120.02	113.15
1	D	103	THR	N-CA-C	-5.62	101.97	110.28
1	A	118	ASN	CA-C-N	5.54	131.67	121.70
1	A	118	ASN	C-N-CA	5.54	131.67	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	26	GLY	N-CA-C	5.48	119.28	110.90
1	A	117	SER	CA-C-N	5.42	127.55	120.28
1	A	117	SER	C-N-CA	5.42	127.55	120.28
1	D	153	ASN	N-CA-C	5.29	118.62	110.42

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	1781	0	1740	53	0
1	D	1767	0	1729	48	0
2	A	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	D	1	0	0	0	0
4	A	4	0	8	6	0
4	D	4	0	8	2	0
5	A	14	0	0	2	0
5	D	7	0	0	1	0
6	A	7	0	10	0	0
6	D	14	0	20	7	0
7	D	8	0	12	1	0
8	A	46	0	0	0	0
8	D	50	0	0	2	0
All	All	3706	0	3527	93	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 13.

All (93) 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:D:129:MET:HE2	1:D:131:ASN:HD21	1.08	1.10

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:118:ASN:CB	1:A:119:SER:HB2	1.88	1.03
1:A:118:ASN:HB2	1:A:119:SER:HB2	1.39	1.03
1:D:129:MET:HE2	1:D:131:ASN:ND2	1.76	0.99
1:A:129:MET:CE	1:D:123:THR:HB	2.03	0.88
1:A:129:MET:HE2	1:D:123:THR:HB	1.54	0.87
1:D:46:LYS:HB3	6:D:307:PEG:H41	1.59	0.85
1:A:86:PRO:O	1:A:181:ILE:HD11	1.78	0.84
1:A:129:MET:HG2	1:D:125:ALA:HB3	1.60	0.83
1:A:181:ILE:CD1	1:A:182:TRP:CD1	2.62	0.83
1:A:155:ARG:HD3	1:A:159:VAL:HG11	1.61	0.82
1:A:118:ASN:CA	1:A:119:SER:HB2	2.11	0.80
1:A:181:ILE:HD12	1:A:182:TRP:CD1	2.17	0.79
1:D:155:ARG:HD3	1:D:159:VAL:HG11	1.65	0.77
1:A:100:TYR:H	4:A:303:IPA:H12	1.48	0.76
1:A:113:SER:HG	5:A:304:ABU:N	1.84	0.76
1:A:112:THR:OG1	1:A:127:HIS:HD2	1.70	0.74
1:A:155:ARG:HD3	1:A:159:VAL:CG1	2.19	0.72
1:D:68:PRO:HG2	6:D:307:PEG:H42	1.72	0.72
1:A:129:MET:HE2	1:D:123:THR:O	1.90	0.72
1:D:151:GLU:OE2	1:D:155:ARG:NH2	2.22	0.71
1:A:112:THR:OG1	1:A:127:HIS:CD2	2.44	0.70
1:A:36:LYS:HD3	1:A:76:SER:O	1.92	0.68
1:A:143:GLN:HE22	1:A:172:ARG:HH21	1.45	0.64
1:A:151:GLU:OE2	1:A:155:ARG:NH2	2.30	0.63
1:D:107:LEU:O	6:D:308:PEG:C4	2.47	0.62
1:D:156:LEU:O	1:D:171:GLY:HA3	1.98	0.61
1:A:99:LEU:H	4:A:303:IPA:H12	1.65	0.61
1:A:137:GLN:HG3	1:A:139:ASP:OD1	2.01	0.60
1:A:156:LEU:O	1:A:171:GLY:HA3	2.02	0.59
1:D:100:TYR:H	4:D:303:IPA:H12	1.67	0.59
1:A:99:LEU:HB2	4:A:303:IPA:H33	1.85	0.59
1:D:84:VAL:HG12	1:D:84:VAL:O	2.01	0.59
1:A:112:THR:HB	1:A:192:GLU:OE1	2.03	0.58
1:D:143:GLN:HE22	1:D:172:ARG:HH21	1.50	0.58
1:D:129:MET:CE	1:D:131:ASN:HD21	1.99	0.58
1:D:155:ARG:HD3	1:D:159:VAL:CG1	2.35	0.57
1:A:106:ILE:HB	1:A:154:LEU:HB3	1.86	0.57
1:A:181:ILE:CD1	1:A:182:TRP:HD1	2.17	0.57
1:A:129:MET:HE1	1:D:123:THR:HB	1.85	0.57
1:A:139:ASP:O	1:A:176:TYR:HB2	2.04	0.57
1:D:107:LEU:O	6:D:308:PEG:H42	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:36:LYS:HD3	1:D:76:SER:O	2.06	0.55
1:D:137:GLN:HG2	1:D:140:LEU:HD12	1.88	0.54
1:D:155:ARG:NH1	8:D:436:HOH:O	2.40	0.54
1:A:118:ASN:CA	1:A:119:SER:CB	2.85	0.54
1:A:118:ASN:N	1:A:119:SER:HB2	2.23	0.53
1:D:33:ARG:HA	8:D:444:HOH:O	2.08	0.53
1:A:118:ASN:N	1:A:119:SER:CB	2.72	0.53
1:D:52:ILE:O	1:D:192:GLU:HB3	2.10	0.52
1:D:107:LEU:O	6:D:308:PEG:H41	2.10	0.52
1:D:106:ILE:HB	1:D:154:LEU:HB3	1.92	0.51
1:A:99:LEU:H	4:A:303:IPA:C1	2.23	0.51
1:A:129:MET:O	1:A:129:MET:HG3	2.10	0.51
1:A:137:GLN:HG2	1:A:140:LEU:HD12	1.92	0.50
1:D:152:GLY:O	6:D:308:PEG:O4	2.29	0.50
1:A:155:ARG:CD	1:A:159:VAL:HG11	2.38	0.50
1:A:99:LEU:CB	4:A:303:IPA:H33	2.41	0.50
1:A:129:MET:O	1:D:124:ASN:HA	2.11	0.50
1:D:131:ASN:C	1:D:132:GLN:HG2	2.36	0.50
1:A:155:ARG:CD	1:A:159:VAL:CG1	2.88	0.49
1:A:207:ALA:CB	4:A:303:IPA:H11	2.43	0.49
1:D:36:LYS:HA	7:D:306:EDO:H22	1.95	0.49
1:D:46:LYS:CB	6:D:307:PEG:H41	2.38	0.48
1:D:155:ARG:CD	1:D:159:VAL:HG11	2.39	0.48
1:D:52:ILE:C	1:D:53:ILE:HG13	2.40	0.47
1:A:101:LYS:N	1:A:101:LYS:HD3	2.30	0.47
1:D:105:THR:OG1	1:D:155:ARG:NH1	2.48	0.47
1:A:160:SER:O	1:A:163:GLY:N	2.49	0.45
1:D:88:TRP:CH2	1:D:180:HIS:HB2	2.52	0.45
1:A:23:PRO:HB2	1:A:40:TRP:O	2.17	0.45
1:D:84:VAL:O	1:D:84:VAL:CG1	2.64	0.45
1:D:88:TRP:CZ3	1:D:180:HIS:HB2	2.51	0.44
1:D:145:ASP:OD2	1:D:169:SER:OG	2.25	0.44
1:A:207:ALA:HB1	1:A:208:ASP:CG	2.43	0.44
1:D:99:LEU:HB2	4:D:303:IPA:H33	2.00	0.43
1:D:136:ASP:OD2	1:D:138:LYS:NZ	2.44	0.43
1:A:129:MET:CG	1:D:125:ALA:HB3	2.38	0.43
1:A:12:TYR:HA	1:A:13:PRO:HD3	1.86	0.43
1:A:124:ASN:HA	1:D:129:MET:O	2.19	0.43
1:D:201:SER:OG	1:D:203:ASP:HB2	2.19	0.42
1:A:75:VAL:O	1:A:75:VAL:HG13	2.19	0.42
5:A:304:ABU:OXT	5:A:304:ABU:CD	2.68	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:SER:OG	1:A:203:ASP:HB2	2.20	0.41
1:A:129:MET:HE2	1:D:123:THR:C	2.46	0.41
1:A:44:ASN:OD1	1:A:200:LYS:HA	2.21	0.41
1:A:181:ILE:HD13	1:A:182:TRP:HD1	1.84	0.41
1:A:132:GLN:HB3	5:D:304:ABU:CB	2.51	0.41
1:A:181:ILE:HD12	1:A:182:TRP:N	2.36	0.40
1:D:17:ILE:O	1:D:33:ARG:HG2	2.21	0.40
1:D:99:LEU:HD12	1:D:99:LEU:HA	1.91	0.40
1:D:174:LEU:HD12	1:D:174:LEU:N	2.37	0.40
1:D:208:ASP:OD1	1:D:227:GLY:HA2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	230/237 (97%)	222 (96%)	8 (4%)	0	100	100
1	D	228/237 (96%)	222 (97%)	6 (3%)	0	100	100
All	All	458/474 (97%)	444 (97%)	14 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	199/202 (98%)	194 (98%)	5 (2%)	42	45
1	D	197/202 (98%)	195 (99%)	2 (1%)	68	75
All	All	396/404 (98%)	389 (98%)	7 (2%)	51	58

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	VAL
1	A	101	LYS
1	A	119	SER
1	A	139	ASP
1	A	181	ILE
1	D	16	ASP
1	D	203	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	ASN
1	A	43	GLN
1	A	127	HIS
1	A	143	GLN
1	D	43	GLN
1	D	51	HIS
1	D	127	HIS
1	D	131	ASN
1	D	143	GLN
1	D	237	ASN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 14 ligands modelled in this entry, 4 are monoatomic - leaving 10 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
7	EDO	D	306	-	3,3,3	0.68	0	2,2,2	0.28	0
7	EDO	D	305	-	3,3,3	0.73	0	2,2,2	0.44	0
5	ABU	A	304	-	6,6,6	1.70	1 (16%)	6,6,6	1.66	2 (33%)
4	IPA	A	303	-	3,3,3	0.56	0	3,3,3	0.33	0
6	PEG	D	308	-	6,6,6	0.38	0	5,5,5	0.25	0
4	IPA	D	303	-	3,3,3	0.59	0	3,3,3	0.28	0
5	ABU	D	304	-	6,6,6	0.81	0	6,6,6	1.18	0
6	PEG	D	307	-	6,6,6	0.36	0	5,5,5	0.77	0
6	PEG	A	306	-	6,6,6	0.53	0	5,5,5	0.80	0
5	ABU	A	305	-	6,6,6	0.81	0	6,6,6	1.19	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
7	EDO	D	305	-	-	1/1/1/1	-
5	ABU	A	304	-	-	3/4/4/4	-
6	PEG	D	308	-	-	0/4/4/4	-
7	EDO	D	306	-	-	1/1/1/1	-
5	ABU	D	304	-	-	2/4/4/4	-
6	PEG	D	307	-	-	1/4/4/4	-
6	PEG	A	306	-	-	3/4/4/4	-
5	ABU	A	305	-	-	4/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	304	ABU	O-C	4.07	1.35	1.22

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	304	ABU	OXT-C-O	2.73	130.35	123.33
5	A	304	ABU	CB-CG-C	-2.17	108.86	114.51

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	305	ABU	CD-CB-CG-C
7	D	305	EDO	O1-C1-C2-O2
7	D	306	EDO	O1-C1-C2-O2
6	D	307	PEG	O1-C1-C2-O2
6	A	306	PEG	O2-C3-C4-O4
6	A	306	PEG	C1-C2-O2-C3
6	A	306	PEG	O1-C1-C2-O2
5	A	305	ABU	O-C-CG-CB
5	D	304	ABU	OXT-C-CG-CB
5	A	304	ABU	OXT-C-CG-CB
5	A	304	ABU	O-C-CG-CB
5	D	304	ABU	O-C-CG-CB
5	A	305	ABU	OXT-C-CG-CB
5	A	305	ABU	CG-CB-CD-N
5	A	304	ABU	CD-CB-CG-C

There are no ring outliers.

7 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	D	306	EDO	1	0
5	A	304	ABU	2	0
4	A	303	IPA	6	0
6	D	308	PEG	4	0
4	D	303	IPA	2	0
5	D	304	ABU	1	0
6	D	307	PEG	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 [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	234/237 (98%)	-1.33	0 100 100	17, 27, 42, 65	0
1	D	232/237 (97%)	-1.35	0 100 100	16, 27, 38, 69	0
All	All	466/474 (98%)	-1.34	0 100 100	16, 27, 41, 69	0

There are no RSRZ outliers to report.

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
5	ABU	A	305	7/7	0.97	0.04	18,20,27,28	0
5	ABU	D	304	7/7	0.97	0.06	19,21,28,28	0
6	PEG	D	307	7/7	0.97	0.06	12,13,16,22	0
4	IPA	D	303	4/4	0.98	0.03	15,15,17,18	0
6	PEG	A	306	7/7	0.98	0.03	20,21,24,24	0
4	IPA	A	303	4/4	0.98	0.03	13,16,17,18	0
7	EDO	D	306	4/4	0.98	0.03	17,20,23,28	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	PEG	D	308	7/7	0.99	0.04	13,17,21,23	0
7	EDO	D	305	4/4	0.99	0.02	14,15,15,15	0
5	ABU	A	304	7/7	0.99	0.03	16,17,19,20	0
3	MN	A	302	1/1	1.00	0.01	23,23,23,23	0
3	MN	D	302	1/1	1.00	0.01	24,24,24,24	0
2	CA	A	301	1/1	1.00	0.02	27,27,27,27	0
2	CA	D	301	1/1	1.00	0.03	27,27,27,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.