



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:11 AM UTC

PDB ID : 1NF5 / pdb_00001nf5
Title : Crystal Structure of Lactose Synthase, Complex with Glucose
Authors : Ramakrishnan, B.; Qasba, P.K.
Deposited on : 2002-12-13
Resolution : 2.00 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

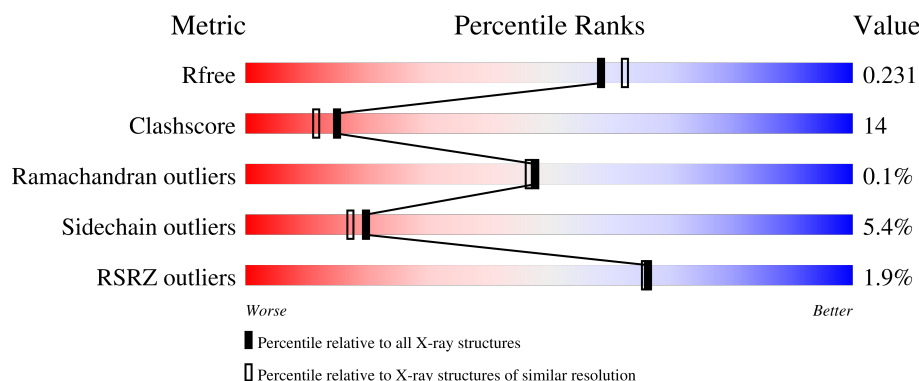
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	123	2% 72% 24% ..
1	C	123	2% 76% 23% .
2	B	286	3% 66% 24% 5% 5%
2	D	286	% 67% 26% . 5%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7037 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 Alpha-lactalbumin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	123	Total	C	N	O	S	0	0	0
			980	620	156	195	9			
1	C	123	Total	C	N	O	S	0	0	0
			980	620	156	195	9			

- Molecule 2 is a protein called beta-1,4-galactosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	272	Total	C	N	O	S	0	0	0
			2218	1424	382	398	14			
2	D	272	Total	C	N	O	S	0	0	0
			2218	1424	382	398	14			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	117	ALA	-	see remark 999	UNP P08037
B	118	SER	-	see remark 999	UNP P08037
B	119	MET	-	see remark 999	UNP P08037
B	120	THR	-	see remark 999	UNP P08037
B	121	GLY	-	see remark 999	UNP P08037
B	122	GLY	-	see remark 999	UNP P08037
B	123	GLN	-	see remark 999	UNP P08037
B	124	GLN	-	see remark 999	UNP P08037
B	125	MET	-	see remark 999	UNP P08037
B	126	GLY	-	see remark 999	UNP P08037
B	127	ARG	-	see remark 999	UNP P08037
B	128	GLY	-	see remark 999	UNP P08037
B	129	SER	-	see remark 999	UNP P08037
D	117	ALA	-	see remark 999	UNP P08037
D	118	SER	-	see remark 999	UNP P08037
D	119	MET	-	see remark 999	UNP P08037

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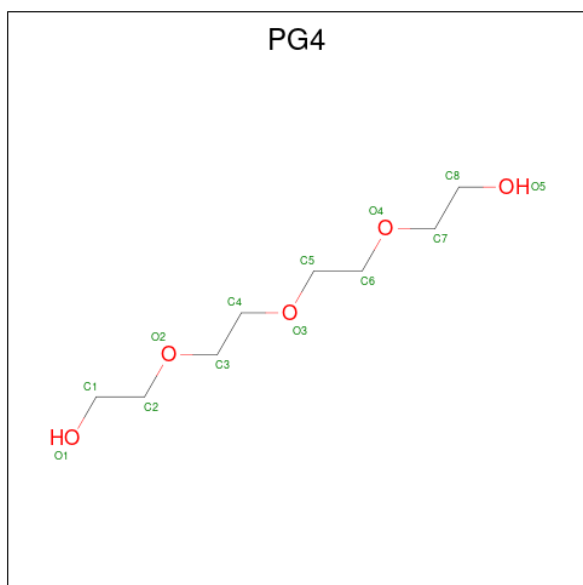
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Chain	Residue	Modelled	Actual	Comment	Reference
D	120	THR	-	see remark 999	UNP P08037
D	121	GLY	-	see remark 999	UNP P08037
D	122	GLY	-	see remark 999	UNP P08037
D	123	GLN	-	see remark 999	UNP P08037
D	124	GLN	-	see remark 999	UNP P08037
D	125	MET	-	see remark 999	UNP P08037
D	126	GLY	-	see remark 999	UNP P08037
D	127	ARG	-	see remark 999	UNP P08037
D	128	GLY	-	see remark 999	UNP P08037
D	129	SER	-	see remark 999	UNP P08037

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

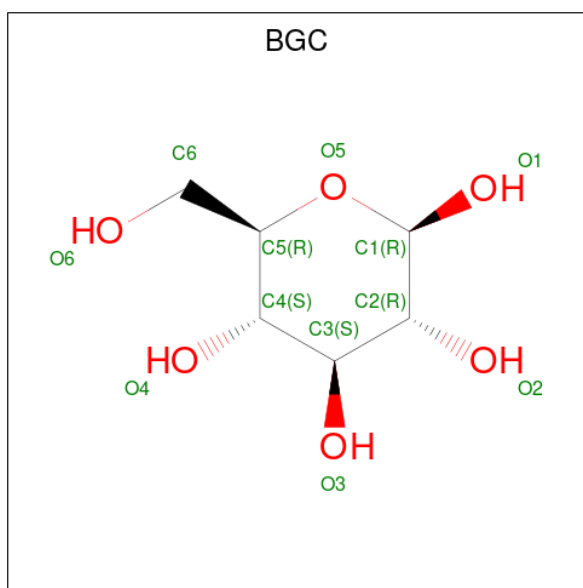
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Ca 1 1	0	0
3	C	1	Total Ca 1 1	0	0

- Molecule 4 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



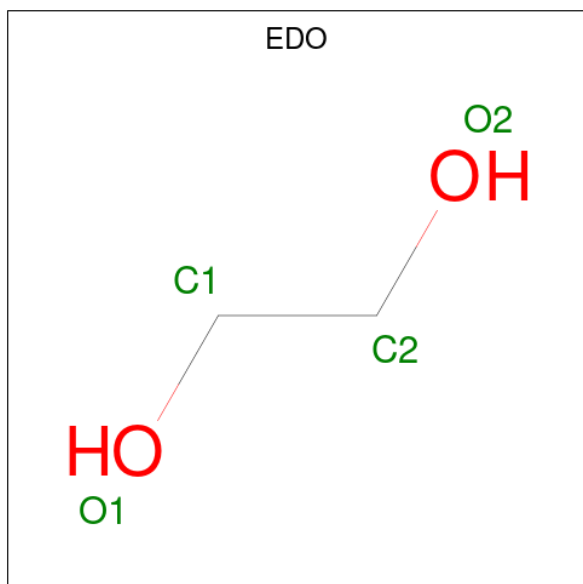
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 10 6 4	0	0
4	B	1	Total C O 10 6 4	0	0

- Molecule 5 is beta-D-glucopyranose (CCD ID: BGC) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			12	6	6		
5	D	1	Total	C	O	0	0
			12	6	6		

- 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		

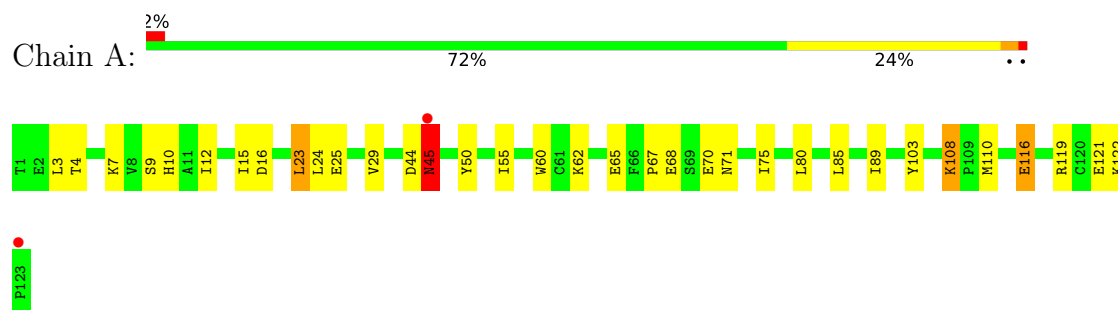
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	106	Total 106	O 106	0	0
7	B	181	Total 181	O 181	0	0
7	C	119	Total 119	O 119	0	0
7	D	185	Total 185	O 185	0	0

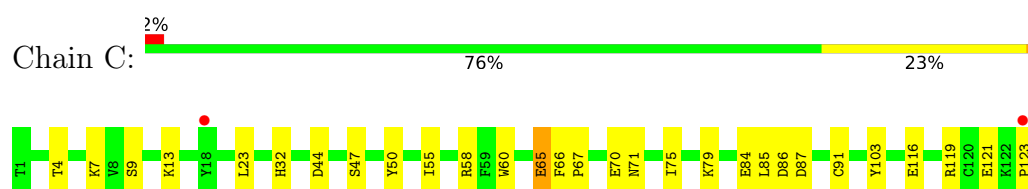
3 Residue-property plots [i](#)

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.

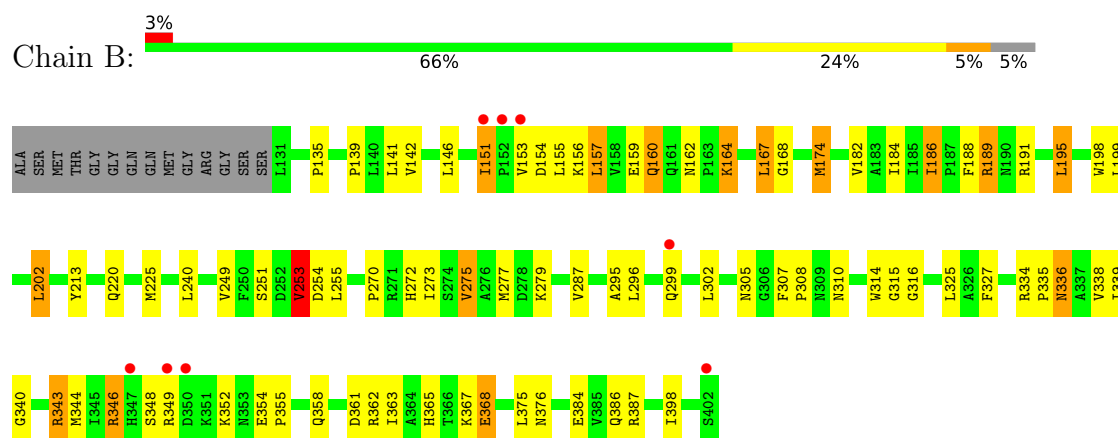
- Molecule 1: Alpha-lactalbumin



- Molecule 1: Alpha-lactalbumin

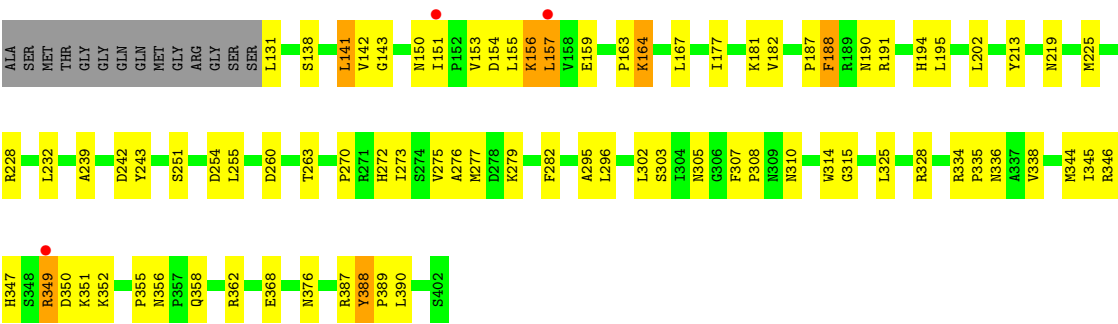


- Molecule 2: beta-1,4-galactosyltransferase



- Molecule 2: beta-1,4-galactosyltransferase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.22Å 93.90Å 99.97Å 90.00° 101.50° 90.00°	Depositor
Resolution (Å)	24.73 – 2.00 24.73 – 2.00	Depositor EDS
% Data completeness (in resolution range)	81.7 (24.73-2.00) 81.7 (24.73-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 1.99Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.190 , 0.241 0.182 , 0.231	Depositor DCC
R_{free} test set	2893 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	25.8	Xtriage
Anisotropy	0.236	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 49.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7037	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.55% 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: EDO, PG4, BGC, CA

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	0.66	0/1001	1.08	4/1350 (0.3%)
1	C	0.67	0/1001	1.19	6/1350 (0.4%)
2	B	0.61	0/2278	1.08	16/3085 (0.5%)
2	D	0.64	0/2278	1.12	19/3085 (0.6%)
All	All	0.64	0/6558	1.11	45/8870 (0.5%)

There are no bond length outliers.

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	315	GLY	N-CA-C	10.67	125.39	112.48
1	C	66	PHE	CA-C-N	10.32	129.97	119.24
1	C	66	PHE	C-N-CA	10.32	129.97	119.24
1	C	103	TYR	N-CA-C	-9.06	101.40	111.28
2	D	182	VAL	N-CA-C	8.62	120.24	108.17
1	A	50	TYR	N-CA-C	8.24	122.83	109.40
2	B	273	ILE	N-CA-C	8.20	120.27	111.58
2	D	315	GLY	N-CA-C	8.20	122.40	112.48
1	A	60	TRP	N-CA-C	7.91	123.45	112.45
2	D	273	ILE	N-CA-C	7.88	119.93	111.58
1	C	65	GLU	N-CA-C	7.79	120.86	111.82
2	D	213	TYR	N-CA-C	7.63	121.01	109.41
1	A	103	TYR	N-CA-C	-7.60	102.99	111.28
1	C	50	TYR	N-CA-C	7.43	121.52	109.40
2	B	182	VAL	N-CA-C	7.36	118.48	108.17
2	B	310	ASN	N-CA-C	6.88	120.63	112.93
2	B	202	LEU	N-CA-C	6.75	118.72	111.36
2	D	202	LEU	N-CA-C	6.62	118.57	111.36
2	D	356	ASN	CA-C-N	6.38	126.59	119.32
2	D	356	ASN	C-N-CA	6.38	126.59	119.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	314	TRP	N-CA-C	6.26	119.54	110.28
2	D	335	PRO	N-CA-C	-6.25	102.07	111.57
2	D	143	GLY	CA-C-N	6.24	126.45	119.90
2	D	143	GLY	C-N-CA	6.24	126.45	119.90
2	B	213	TYR	N-CA-C	6.08	118.65	109.41
1	A	45	ASN	O-C-N	-6.06	114.53	122.59
1	C	60	TRP	N-CA-C	5.83	120.55	112.45
2	D	156	LYS	N-CA-C	-5.74	105.03	111.28
2	B	186	ILE	N-CA-C	5.64	113.63	107.60
2	B	314	TRP	N-CA-C	5.62	118.60	110.28
2	B	251	SER	N-CA-C	5.61	118.24	108.76
2	D	251	SER	N-CA-C	5.56	118.16	108.76
2	B	346	ARG	N-CA-C	5.54	117.69	110.43
2	D	388	TYR	CA-C-N	-5.53	112.91	119.05
2	D	388	TYR	C-N-CA	-5.53	112.91	119.05
2	B	335	PRO	N-CA-C	-5.52	101.21	111.03
2	D	263	THR	N-CA-C	5.49	118.19	109.96
2	D	275	VAL	N-CA-C	5.45	116.94	111.00
2	D	190	ASN	N-CA-C	5.35	118.78	111.39
2	B	162	ASN	CA-C-N	5.19	124.85	119.56
2	B	162	ASN	C-N-CA	5.19	124.85	119.56
2	B	188	PHE	N-CA-C	5.13	116.86	108.76
2	D	188	PHE	N-CA-C	5.12	117.50	108.75
2	B	275	VAL	N-CA-C	5.05	116.51	111.00
2	B	253	VAL	CB-CA-C	-5.04	100.32	111.77

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	980	0	936	30	0
1	C	980	0	936	28	0
2	B	2218	0	2185	67	0
2	D	2218	0	2185	53	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	1	0	0	0	0
3	C	1	0	0	0	0
4	A	10	0	13	0	0
4	B	10	0	13	0	0
5	B	12	0	12	0	0
5	D	12	0	12	2	0
6	C	4	0	6	0	0
7	A	106	0	0	4	0
7	B	181	0	0	1	0
7	C	119	0	0	2	0
7	D	185	0	0	2	0
All	All	7037	0	6298	175	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 14.

All (175) 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:44:ASP:C	1:A:45:ASN:HD22	1.52	1.18
2:B:343:ARG:HG2	2:B:343:ARG:HH11	1.16	1.09
2:B:151:ILE:HD12	2:B:151:ILE:H	1.28	0.99
2:B:336:ASN:HD22	2:B:338:VAL:H	1.20	0.89
2:B:343:ARG:HH11	2:B:343:ARG:CG	1.87	0.88
1:A:71:ASN:HD21	1:A:75:ILE:H	1.16	0.86
1:C:13:LYS:HD3	1:C:23:LEU:HD11	1.57	0.85
1:A:44:ASP:C	1:A:45:ASN:ND2	2.36	0.82
1:A:4:THR:H	1:A:7:LYS:HD2	1.45	0.80
2:B:151:ILE:HD12	2:B:151:ILE:N	1.97	0.79
2:B:157:LEU:HD13	2:B:160:GLN:NE2	1.99	0.78
2:D:305:ASN:HD21	2:D:376:ASN:H	1.30	0.78
2:D:167:LEU:HD11	2:D:387:ARG:HB3	1.67	0.76
2:B:305:ASN:HD21	2:B:376:ASN:H	1.35	0.75
1:A:45:ASN:HD22	1:A:45:ASN:N	1.84	0.74
1:C:47:SER:HB3	1:C:58:ARG:NH1	2.03	0.74
2:B:336:ASN:ND2	2:B:338:VAL:H	1.87	0.71
1:C:32:HIS:HD2	5:D:527:BGC:O1	1.73	0.71
2:D:155:LEU:O	2:D:159:GLU:HG3	1.91	0.71
2:D:349:ARG:HE	2:D:350:ASP:H	1.38	0.70
2:B:155:LEU:O	2:B:159:GLU:HG3	1.92	0.69
2:B:361:ASP:HB2	7:B:1307:HOH:O	1.91	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:71:ASN:HD21	1:C:75:ILE:H	1.40	0.69
2:B:343:ARG:HG2	2:B:343:ARG:NH1	1.96	0.68
2:B:327:PHE:CZ	2:B:367:LYS:HB2	2.29	0.68
2:B:151:ILE:H	2:B:151:ILE:CD1	1.96	0.67
2:B:191:ARG:HG2	2:B:191:ARG:HH11	1.58	0.67
2:B:189:ARG:NH2	2:B:352:LYS:HB3	2.10	0.66
2:D:336:ASN:ND2	2:D:338:VAL:HB	2.10	0.66
1:C:9:SER:O	1:C:13:LYS:HG2	1.96	0.66
1:A:44:ASP:O	1:A:45:ASN:ND2	2.29	0.65
2:B:186:ILE:CG2	2:B:253:VAL:HG13	2.26	0.65
2:D:279:LYS:HD2	2:D:346:ARG:HH12	1.64	0.63
2:B:336:ASN:HD22	2:B:338:VAL:N	1.93	0.63
2:D:164:LYS:HE3	7:D:825:HOH:O	1.99	0.62
1:A:65:GLU:HG3	7:A:1020:HOH:O	2.01	0.61
2:B:255:LEU:HD21	2:B:277:MET:HE3	1.82	0.61
2:D:295:ALA:C	2:D:296:LEU:HD12	2.26	0.61
1:A:16:ASP:HB2	1:A:23:LEU:HD23	1.83	0.61
2:B:275:VAL:HG22	2:B:340:GLY:HA3	1.81	0.61
1:A:12:ILE:O	1:A:15:ILE:HG22	2.01	0.61
2:B:186:ILE:HG21	2:B:253:VAL:HG13	1.82	0.60
2:B:354:GLU:CD	2:B:355:PRO:HD2	2.26	0.59
1:A:68:GLU:HG2	1:A:68:GLU:O	2.03	0.59
2:B:164:LYS:HE2	2:B:174:MET:HE3	1.83	0.59
2:D:167:LEU:CD1	2:D:387:ARG:HB3	2.33	0.59
2:B:167:LEU:HD13	2:B:168:GLY:N	2.19	0.58
2:D:167:LEU:C	2:D:167:LEU:HD13	2.29	0.58
2:B:346:ARG:HH11	2:B:346:ARG:HG2	1.68	0.58
2:B:384:GLU:OE2	2:B:386:GLN:HG2	2.03	0.57
2:B:336:ASN:ND2	2:B:339:ILE:H	2.03	0.57
1:C:116:GLU:OE2	1:C:119:ARG:NH2	2.38	0.57
1:C:13:LYS:HD3	1:C:23:LEU:CD1	2.31	0.56
2:B:164:LYS:CE	2:B:174:MET:HE3	2.36	0.56
2:D:167:LEU:HD11	2:D:387:ARG:CB	2.35	0.56
2:D:154:ASP:HB3	2:D:157:LEU:HB2	1.86	0.56
2:D:336:ASN:HD22	2:D:338:VAL:H	1.53	0.56
2:B:272:HIS:HB3	2:B:334:ARG:HG2	1.87	0.56
2:B:275:VAL:CG2	2:B:340:GLY:HA3	2.35	0.56
2:D:255:LEU:HD21	2:D:277:MET:HE3	1.87	0.56
1:A:4:THR:N	1:A:7:LYS:HD2	2.18	0.55
2:B:358:GLN:O	2:B:362:ARG:HG3	2.07	0.55
2:D:191:ARG:HH11	2:D:194:HIS:HD2	1.54	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:164:LYS:N	2:D:164:LYS:HD3	2.23	0.54
2:D:336:ASN:HD21	2:D:338:VAL:HB	1.72	0.54
2:B:336:ASN:HD22	2:B:336:ASN:C	2.15	0.54
2:D:351:LYS:HB2	2:D:352:LYS:HD2	1.90	0.54
2:D:163:PRO:HG2	2:D:164:LYS:HD3	1.90	0.53
1:C:47:SER:CB	1:C:58:ARG:NH1	2.72	0.53
2:D:187:PRO:HD3	2:D:232:LEU:HD21	1.90	0.53
1:A:110:MET:HE3	7:A:1080:HOH:O	2.09	0.52
2:D:225:MET:HA	2:D:352:LYS:O	2.08	0.52
2:D:159:GLU:OE1	2:D:389:PRO:HG2	2.10	0.52
2:B:186:ILE:HG21	2:B:253:VAL:CG1	2.40	0.52
2:D:272:HIS:HB3	2:D:334:ARG:HG2	1.93	0.51
1:A:55:ILE:HB	1:A:80:LEU:HD13	1.93	0.51
2:D:138:SER:HB3	2:D:141:LEU:HD13	1.91	0.51
1:A:71:ASN:HD21	1:A:75:ILE:N	1.97	0.51
2:B:307:PHE:HB3	2:B:308:PRO:HD2	1.92	0.51
1:A:108:LYS:HE3	7:A:1233:HOH:O	2.10	0.51
2:B:167:LEU:CD1	2:B:387:ARG:HB3	2.40	0.51
2:B:154:ASP:OD1	2:B:156:LYS:HB2	2.10	0.51
1:A:4:THR:HG23	1:A:7:LYS:HE3	1.92	0.51
2:B:327:PHE:CE2	2:B:367:LYS:HD3	2.46	0.50
1:A:116:GLU:OE1	1:A:119:ARG:NH2	2.45	0.50
2:B:336:ASN:HD21	2:B:338:VAL:HB	1.76	0.50
2:B:327:PHE:CD2	2:B:367:LYS:HD3	2.47	0.50
1:C:47:SER:HB3	1:C:58:ARG:HH12	1.73	0.50
1:A:62:LYS:HE3	1:A:67:PRO:O	2.12	0.49
1:A:10:HIS:NE2	1:A:122:LYS:O	2.44	0.49
2:D:345:ILE:O	2:D:347:HIS:HD2	1.94	0.49
2:B:365:HIS:O	2:B:368:GLU:HG2	2.13	0.49
2:B:195:LEU:HD22	2:B:199:LEU:HG	1.95	0.48
2:D:279:LYS:HD2	2:D:346:ARG:NH1	2.27	0.48
1:A:4:THR:HG23	1:A:7:LYS:CE	2.44	0.48
2:B:336:ASN:ND2	2:B:338:VAL:N	2.59	0.48
2:D:279:LYS:NZ	7:D:1285:HOH:O	2.46	0.48
1:A:9:SER:HA	1:A:23:LEU:HD11	1.96	0.48
1:C:65:GLU:O	1:C:67:PRO:HD3	2.14	0.47
1:C:4:THR:H	1:C:7:LYS:HE2	1.79	0.47
1:C:4:THR:H	1:C:7:LYS:CE	2.28	0.47
2:B:249:VAL:HG22	2:B:295:ALA:HB2	1.97	0.47
1:A:25:GLU:O	1:A:29:VAL:HG23	2.15	0.47
2:B:343:ARG:CG	2:B:343:ARG:NH1	2.56	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:4:THR:HG23	1:C:7:LYS:HE2	1.96	0.47
2:B:191:ARG:HG2	2:B:191:ARG:NH1	2.25	0.47
2:B:198:TRP:CZ2	2:B:202:LEU:HG	2.50	0.47
2:D:159:GLU:HG2	2:D:390:LEU:HG	1.97	0.47
2:D:295:ALA:O	2:D:296:LEU:HD12	2.14	0.47
2:D:131:LEU:HB2	2:D:177:ILE:CD1	2.46	0.46
2:B:277:MET:CE	2:B:344:MET:HE3	2.46	0.46
2:B:336:ASN:HD21	2:B:339:ILE:H	1.64	0.46
2:B:225:MET:HA	2:B:352:LYS:O	2.15	0.46
1:C:119:ARG:HG2	1:C:121:GLU:OE2	2.16	0.46
1:A:122:LYS:HG3	1:C:86:ASP:OD1	2.16	0.45
2:B:316:GLY:HA2	2:B:363:ILE:HD11	1.98	0.45
2:B:254:ASP:O	2:B:344:MET:HA	2.16	0.45
1:C:70:GLU:O	1:C:71:ASN:C	2.59	0.45
2:B:384:GLU:OE2	2:B:386:GLN:CG	2.65	0.45
2:B:336:ASN:ND2	2:B:336:ASN:C	2.75	0.44
2:B:157:LEU:HD13	2:B:160:GLN:HE21	1.82	0.44
1:C:32:HIS:HE1	7:C:877:HOH:O	2.00	0.44
2:D:349:ARG:NE	2:D:350:ASP:H	2.10	0.44
2:B:349:ARG:NH1	2:B:354:GLU:OE2	2.51	0.44
2:D:270:PRO:HG2	2:D:325:LEU:HD22	1.99	0.44
7:A:1005:HOH:O	1:C:79:LYS:HE3	2.18	0.44
2:D:239:ALA:O	2:D:242:ASP:HB2	2.17	0.44
1:C:55:ILE:HD13	1:C:91:CYS:SG	2.58	0.44
1:C:65:GLU:C	1:C:67:PRO:HD3	2.42	0.44
1:A:3:LEU:HD22	1:A:7:LYS:HD3	2.01	0.43
2:D:279:LYS:CD	2:D:346:ARG:HH12	2.29	0.43
2:B:220:GLN:CD	2:B:398:ILE:HD11	2.43	0.43
2:B:270:PRO:HG2	2:B:325:LEU:HD22	2.00	0.43
2:D:187:PRO:HD3	2:D:232:LEU:CD2	2.48	0.43
2:D:191:ARG:HH11	2:D:194:HIS:CD2	2.34	0.43
1:A:121:GLU:HG2	1:C:87:ASP:HA	2.01	0.43
2:B:135:PRO:HG3	2:B:139:PRO:HD3	2.00	0.43
2:D:181:LYS:HE2	2:D:243:TYR:CZ	2.53	0.43
2:D:150:ASN:O	2:D:151:ILE:C	2.61	0.43
2:D:307:PHE:HB3	2:D:308:PRO:HD2	2.01	0.43
2:B:157:LEU:CD1	2:B:160:GLN:NE2	2.77	0.42
1:A:122:LYS:NZ	1:C:84:GLU:OE1	2.43	0.42
2:D:167:LEU:HD22	2:D:388:TYR:O	2.19	0.42
2:D:310:ASN:HB3	2:D:362:ARG:NH2	2.35	0.42
2:D:188:PHE:O	2:D:219:ASN:HA	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:181:LYS:HE2	2:D:243:TYR:CE2	2.55	0.42
2:D:358:GLN:OE1	2:D:362:ARG:NH1	2.53	0.42
1:C:121:GLU:O	1:C:123:PRO:N	2.52	0.42
2:B:157:LEU:O	2:B:160:GLN:HB3	2.20	0.41
2:B:346:ARG:HG2	2:B:346:ARG:NH1	2.35	0.41
2:D:228:ARG:O	2:D:232:LEU:HG	2.21	0.41
2:D:358:GLN:O	2:D:362:ARG:HG3	2.20	0.41
1:C:32:HIS:CD2	5:D:527:BGC:O1	2.63	0.41
2:D:131:LEU:CB	2:D:177:ILE:CD1	2.99	0.41
2:D:155:LEU:HD23	2:D:155:LEU:HA	1.83	0.41
1:A:45:ASN:ND2	1:A:45:ASN:N	2.57	0.41
2:B:295:ALA:C	2:B:296:LEU:HD12	2.45	0.41
2:B:348:SER:O	2:B:349:ARG:C	2.63	0.41
2:B:160:GLN:HE21	2:B:160:GLN:HB3	1.59	0.41
2:D:303:SER:O	2:D:328:ARG:HG3	2.20	0.41
1:A:24:LEU:CD2	1:A:119:ARG:HA	2.50	0.41
2:D:254:ASP:O	2:D:344:MET:HA	2.21	0.41
1:A:12:ILE:HG13	1:A:23:LEU:HD21	2.03	0.41
2:B:154:ASP:O	2:B:157:LEU:N	2.54	0.41
2:B:279:LYS:O	2:B:279:LYS:HG2	2.20	0.41
1:C:67:PRO:HG3	7:C:1117:HOH:O	2.20	0.41
2:D:142:VAL:HG22	2:D:260:ASP:OD1	2.20	0.41
2:B:184:ILE:HD12	2:B:184:ILE:N	2.36	0.41
2:B:336:ASN:ND2	2:B:338:VAL:HB	2.36	0.41
2:D:151:ILE:HG13	2:D:151:ILE:O	2.21	0.41
1:A:12:ILE:HG22	1:A:89:ILE:HD11	2.04	0.40
1:C:47:SER:CB	1:C:58:ARG:HH12	2.32	0.40
2:D:276:ALA:HB1	2:D:282:PHE:C	2.47	0.40
1:C:44:ASP:OD1	1:C:44:ASP:C	2.64	0.40
1:C:121:GLU:O	1:C:123:PRO:HD3	2.21	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	121/123 (98%)	119 (98%)	1 (1%)	1 (1%)	16	11
1	C	121/123 (98%)	117 (97%)	4 (3%)	0	100	100
2	B	270/286 (94%)	259 (96%)	11 (4%)	0	100	100
2	D	270/286 (94%)	263 (97%)	7 (3%)	0	100	100
All	All	782/818 (96%)	758 (97%)	23 (3%)	1 (0%)	48	46

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	45	ASN

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	109/109 (100%)	103 (94%)	6 (6%)	19	17
1	C	109/109 (100%)	108 (99%)	1 (1%)	70	78
2	B	245/254 (96%)	224 (91%)	21 (9%)	10	6
2	D	245/254 (96%)	235 (96%)	10 (4%)	27	26
All	All	708/726 (98%)	670 (95%)	38 (5%)	20	17

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	LEU
1	A	45	ASN
1	A	70	GLU
1	A	85	LEU
1	A	108	LYS
1	A	116	GLU
2	B	141	LEU

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Mol	Chain	Res	Type
2	B	142	VAL
2	B	146	LEU
2	B	151	ILE
2	B	153	VAL
2	B	157	LEU
2	B	160	GLN
2	B	164	LYS
2	B	167	LEU
2	B	174	MET
2	B	189	ARG
2	B	195	LEU
2	B	240	LEU
2	B	253	VAL
2	B	287	VAL
2	B	299	GLN
2	B	302	LEU
2	B	336	ASN
2	B	343	ARG
2	B	368	GLU
2	B	375	LEU
1	C	85	LEU
2	D	141	LEU
2	D	153	VAL
2	D	156	LYS
2	D	157	LEU
2	D	164	LYS
2	D	195	LEU
2	D	302	LEU
2	D	349	ARG
2	D	355	PRO
2	D	368	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	ASN
1	A	71	ASN
2	B	160	GLN
2	B	219	ASN
2	B	305	ASN
2	B	310	ASN
2	B	336	ASN

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Mol	Chain	Res	Type
1	C	32	HIS
1	C	71	ASN
2	D	161	GLN
2	D	180	HIS
2	D	194	HIS
2	D	219	ASN
2	D	305	ASN
2	D	310	ASN
2	D	336	ASN
2	D	347	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 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	$\# Z > 2$
4	PG4	A	126	-	9,9,12	0.49	0	8,8,11	0.26	0
5	BGC	D	527	-	12,12,12	1.15	1 (8%)	17,17,17	2.32	9 (52%)
6	EDO	C	128	-	3,3,3	0.39	0	2,2,2	0.14	0
4	PG4	B	404	-	9,9,12	0.44	0	8,8,11	0.30	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	BGC	B	403	-	12,12,12	1.04	1 (8%)	17,17,17	1.84	3 (17%)

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	PG4	A	126	-	-	3/7/7/10	-
5	BGC	D	527	-	-	0/2/22/22	0/1/1/1
6	EDO	C	128	-	-	1/1/1/1	-
4	PG4	B	404	-	-	4/7/7/10	-
5	BGC	B	403	-	-	1/2/22/22	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	527	BGC	O5-C5	2.42	1.50	1.44
5	B	403	BGC	O5-C5	2.35	1.50	1.44

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	403	BGC	O5-C5-C6	-4.77	94.62	106.44
5	D	527	BGC	O5-C5-C6	-4.30	95.79	106.44
5	D	527	BGC	O5-C1-C2	4.24	117.75	110.30
5	D	527	BGC	O3-C3-C4	-3.28	102.65	110.38
5	D	527	BGC	O1-C1-C2	3.27	118.48	108.98
5	B	403	BGC	C6-C5-C4	2.71	119.68	113.02
5	D	527	BGC	C6-C5-C4	2.53	119.24	113.02
5	B	403	BGC	O3-C3-C4	-2.48	104.53	110.38
5	D	527	BGC	O1-C1-O5	2.37	117.44	110.41
5	D	527	BGC	O2-C2-C1	2.24	114.42	109.25
5	D	527	BGC	O4-C4-C5	-2.07	104.22	109.32
5	D	527	BGC	O5-C5-C4	-2.04	106.03	109.70

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	404	PG4	O3-C5-C6-O4

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Mol	Chain	Res	Type	Atoms
4	A	126	PG4	O1-C1-C2-O2
4	B	404	PG4	O1-C1-C2-O2
5	B	403	BGC	C4-C5-C6-O6
4	B	404	PG4	C6-C5-O3-C4
4	B	404	PG4	C1-C2-O2-C3
4	A	126	PG4	O2-C3-C4-O3
6	C	128	EDO	O1-C1-C2-O2
4	A	126	PG4	C1-C2-O2-C3

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	527	BGC	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	123/123 (100%)	-0.26	2 (1%) 70 70	18, 26, 51, 63	0
1	C	123/123 (100%)	-0.34	2 (1%) 70 70	15, 23, 43, 70	0
2	B	272/286 (95%)	-0.08	8 (2%) 53 52	18, 28, 52, 73	0
2	D	272/286 (95%)	-0.26	3 (1%) 78 77	17, 27, 51, 61	0
All	All	790/818 (96%)	-0.21	15 (1%) 66 66	15, 27, 51, 73	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	123	PRO	3.8
2	D	151	ILE	3.5
1	A	123	PRO	2.8
2	B	349	ARG	2.7
2	B	350	ASP	2.6
2	B	299	GLN	2.6
1	C	18	TYR	2.4
2	D	157	LEU	2.3
2	B	152	PRO	2.3
2	B	151	ILE	2.2
1	A	45	ASN	2.1
2	B	153	VAL	2.1
2	D	349	ARG	2.0
2	B	347	HIS	2.0
2	B	402	SER	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	EDO	C	128	4/4	0.83	0.11	50,50,51,52	0
4	PG4	B	404	10/13	0.86	0.13	50,52,53,54	0
4	PG4	A	126	10/13	0.93	0.11	44,47,50,50	0
5	BGC	B	403	12/12	0.96	0.06	19,21,24,24	0
5	BGC	D	527	12/12	0.97	0.05	17,19,20,21	0
3	CA	A	124	1/1	0.99	0.03	24,24,24,24	0
3	CA	C	526	1/1	1.00	0.02	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.