



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:26 AM UTC

PDB ID : 1OQM / pdb_00001oqm
Title : A 1:1 complex between alpha-lactalbumin and beta1,4-galactosyltransferase
in the presence of UDP-N-acetyl-galactosamine
Authors : Ramakrishnan, B.; Qasba, P.K.
Deposited on : 2003-03-10
Resolution : 2.10 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

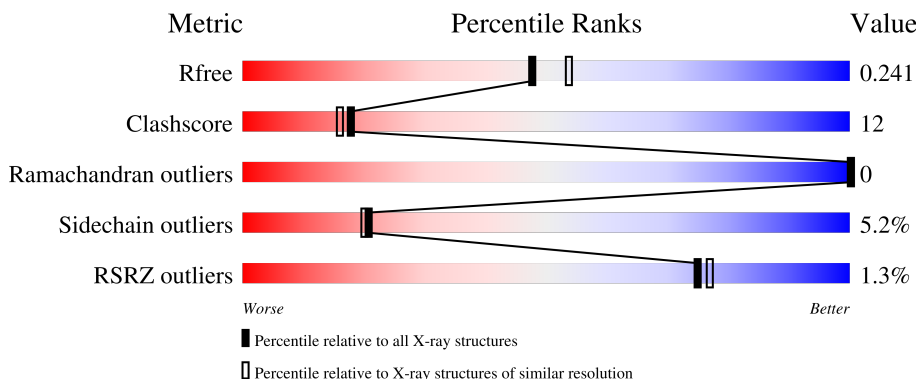
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

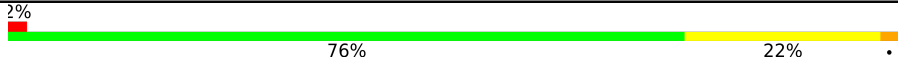
The reported resolution of this entry is 2.10 Å.

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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (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 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	
1	C	123	
2	B	286	
2	D	286	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7121 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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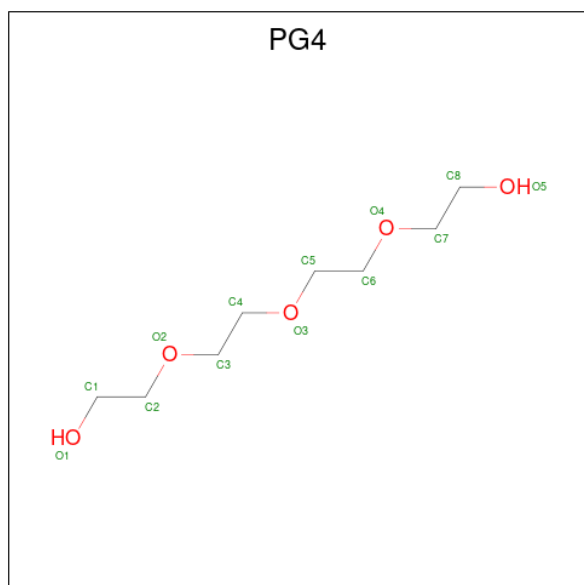
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Chain	Residue	Modelled	Actual	Comment	Reference
D	120	THR	-	SEE REMARK 999	UNP P08037
D	121	GLY	-	SEE REMARK 999	UNP P08037
D	122	GLY	-	SEE REMARK 999	UNP P08037
D	123	GLN	-	SEE REMARK 999	UNP P08037
D	124	GLN	-	SEE REMARK 999	UNP P08037
D	125	MET	-	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).

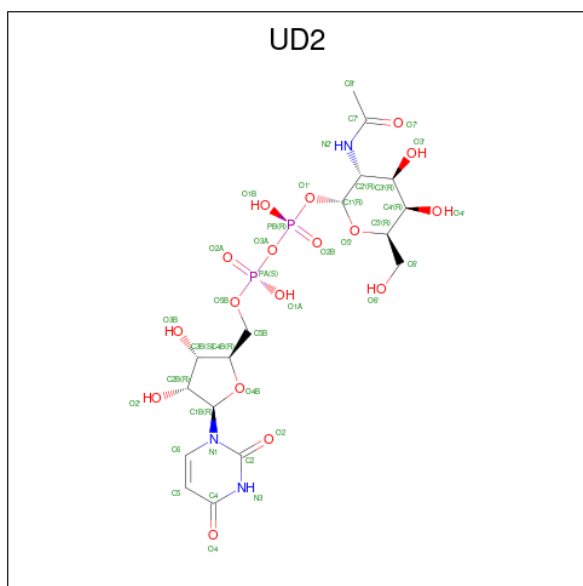


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 7 4 3	0	0
4	B	1	Total C O 7 4 3	0	0

- Molecule 5 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Mn	0	0
			1	1		
5	D	1	Total	Mn	0	0
			1	1		

- Molecule 6 is URIDINE-DIPHOSPHATE-N-ACETYLGLACTOSAMINE (CCD ID: UD2) (formula: C₁₇H₂₇N₃O₁₇P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	B	1	Total	C	N	O	P	0	0
			39	17	3	17	2		
6	D	1	Total	C	N	O	P	0	0
			39	17	3	17	2		

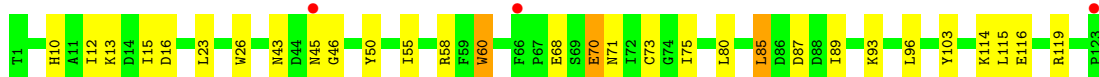
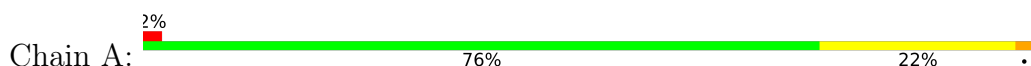
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	114	Total	O	0	0
			114	114		
7	B	191	Total	O	0	0
			191	191		
7	C	134	Total	O	0	0
			134	134		
7	D	190	Total	O	0	0
			190	190		

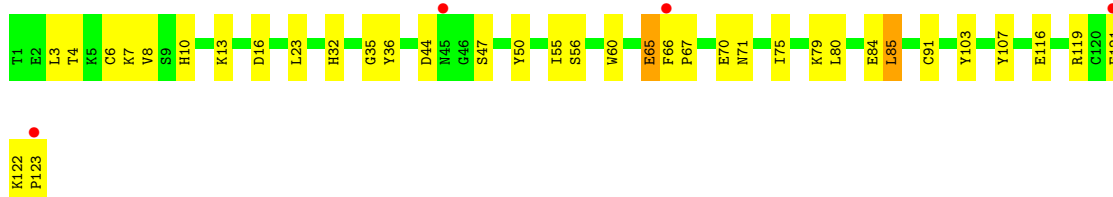
3 Residue-property plots [i](#)

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.

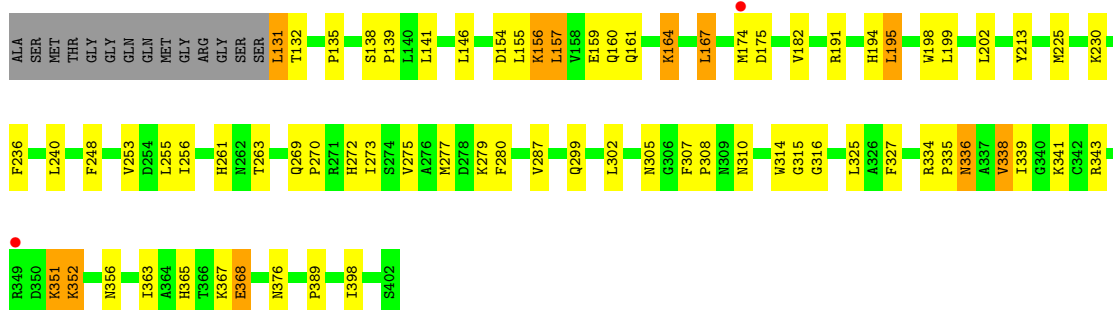
- Molecule 1: Alpha-lactalbumin



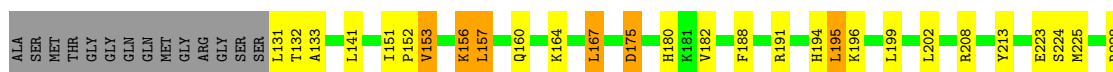
- Molecule 1: Alpha-lactalbumin

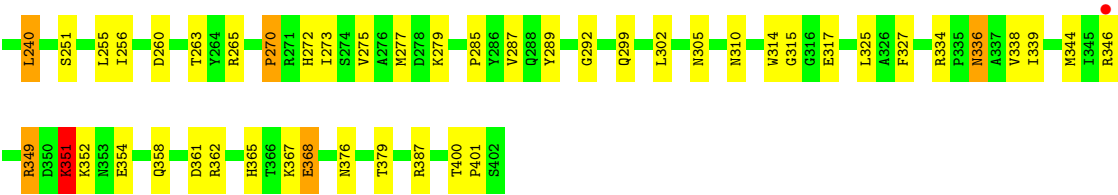


- Molecule 2: beta-1,4-galactosyltransferase



- 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	57.47Å 95.68Å 100.58Å 90.00° 101.40° 90.00°	Depositor
Resolution (Å)	19.56 – 2.10 19.56 – 2.10	Depositor EDS
% Data completeness (in resolution range)	90.1 (19.56-2.10) 90.0 (19.56-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 2.09Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.194 , 0.244 0.191 , 0.241	Depositor DCC
R_{free} test set	5711 reflections (10.17%)	wwPDB-VP
Wilson B-factor (Å ²)	29.3	Xtriage
Anisotropy	0.526	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 49.4	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	7121	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.37% 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: MN, CA, UD2, PG4

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	0/1001	0.98	5/1350 (0.4%)
1	C	0.61	0/1001	1.09	8/1350 (0.6%)
2	B	0.55	0/2278	1.05	15/3085 (0.5%)
2	D	0.56	0/2278	1.08	15/3085 (0.5%)
All	All	0.56	0/6558	1.06	43/8870 (0.5%)

There are no bond length outliers.

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	44	ASP	N-CA-C	-10.31	98.85	108.75
2	D	182	VAL	N-CA-C	8.81	120.51	108.17
2	B	273	ILE	N-CA-C	8.35	119.21	111.45
2	B	202	LEU	N-CA-C	8.18	120.28	111.36
2	B	182	VAL	N-CA-C	8.16	119.64	107.80
1	C	50	TYR	N-CA-C	7.89	122.27	109.40
2	D	213	TYR	N-CA-C	7.84	121.33	109.41
1	C	60	TRP	N-CA-C	7.78	123.26	112.45
2	D	315	GLY	N-CA-C	7.74	122.44	112.68
1	A	60	TRP	N-CA-C	7.49	122.86	112.45
2	D	188	PHE	N-CA-C	7.27	120.40	108.99
2	D	202	LEU	N-CA-C	6.92	118.91	111.36
1	C	103	TYR	N-CA-C	-6.86	103.81	111.28
2	D	273	ILE	N-CA-C	6.83	118.82	111.58
1	C	107	TYR	N-CA-C	6.80	119.31	111.02
2	D	310	ASN	N-CA-C	6.55	120.88	113.02
2	B	175	ASP	N-CA-C	6.52	118.39	111.28
1	A	50	TYR	N-CA-C	6.44	119.89	109.40
2	B	213	TYR	N-CA-C	6.30	119.45	109.81
2	D	351	LYS	N-CA-C	-6.29	101.17	110.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	335	PRO	N-CA-C	-6.25	102.07	111.57
2	B	314	TRP	N-CA-C	6.14	119.37	110.28
2	B	310	ASN	N-CA-C	6.10	120.41	113.15
2	B	351	LYS	N-CA-C	-6.06	101.95	110.50
2	D	151	ILE	N-CA-C	6.01	114.44	107.89
1	C	85	LEU	N-CA-C	5.98	120.19	113.01
2	B	356	ASN	CA-C-N	5.81	125.48	119.56
2	B	356	ASN	C-N-CA	5.81	125.48	119.56
1	C	65	GLU	N-CA-C	5.79	120.49	112.90
2	D	175	ASP	N-CA-C	5.71	118.72	111.69
1	A	85	LEU	N-CA-C	5.67	120.06	113.20
2	B	338	VAL	N-CA-C	-5.60	106.24	111.45
2	D	275	VAL	N-CA-C	5.54	115.73	110.53
2	D	263	THR	N-CA-C	5.53	118.34	110.10
2	D	314	TRP	N-CA-C	5.48	118.39	110.28
2	B	261	HIS	N-CA-C	-5.34	106.27	112.89
1	A	46	GLY	N-CA-C	-5.26	107.85	115.27
2	B	132	THR	N-CA-C	5.20	117.44	109.95
1	A	103	TYR	N-CA-C	-5.20	105.61	111.28
2	D	251	SER	N-CA-C	5.20	117.92	109.24
2	B	315	GLY	N-CA-C	5.18	125.46	113.18
2	D	256	ILE	CB-CA-C	-5.05	105.23	110.13
1	C	56	SER	N-CA-C	5.00	118.37	109.96

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	20	0
1	C	980	0	936	24	0
2	B	2218	0	2185	53	0
2	D	2218	0	2185	61	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	7	0	9	3	0
4	B	7	0	9	0	0
5	B	1	0	0	0	0
5	D	1	0	0	0	0
6	B	39	0	25	5	0
6	D	39	0	25	6	0
7	A	114	0	0	4	0
7	B	191	0	0	4	0
7	C	134	0	0	2	0
7	D	190	0	0	8	0
All	All	7121	0	6310	158	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 12.

All (158) 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:114:LYS:HA	4:A:814:PG4:H22	1.23	1.11
2:D:277:MET:SD	6:D:528:UD2:H8'1	1.93	1.08
2:D:255:LEU:HD11	6:D:528:UD2:H8'3	1.36	1.04
1:A:10:HIS:HA	1:A:13:LYS:HE2	1.49	0.94
2:D:349:ARG:HH11	2:D:349:ARG:HG2	1.33	0.92
1:A:45:ASN:HA	7:A:1167:HOH:O	1.77	0.84
1:A:71:ASN:HD21	1:A:75:ILE:H	1.24	0.84
1:A:89:ILE:HG22	7:A:1418:HOH:O	1.76	0.83
1:C:71:ASN:HD21	1:C:75:ILE:H	1.23	0.82
2:B:154:ASP:HB3	2:B:157:LEU:HB2	1.62	0.82
2:D:336:ASN:HD22	2:D:338:VAL:H	1.26	0.82
2:B:336:ASN:HD22	2:B:338:VAL:H	1.28	0.80
2:B:327:PHE:CZ	2:B:367:LYS:HB2	2.16	0.80
2:D:132:THR:HG22	2:D:133:ALA:H	1.46	0.78
2:D:225:MET:HE2	2:D:352:LYS:HE3	1.66	0.78
2:D:305:ASN:HD21	2:D:376:ASN:H	1.32	0.78
2:B:240:LEU:HD11	2:B:248:PHE:CZ	2.21	0.75
2:B:351:LYS:O	2:B:352:LYS:HB2	1.86	0.75
1:C:65:GLU:O	1:C:67:PRO:HD3	1.87	0.74
2:B:191:ARG:HH11	2:B:194:HIS:HD2	1.36	0.73
2:D:336:ASN:HD21	2:D:338:VAL:HB	1.53	0.72
1:C:65:GLU:C	1:C:67:PRO:HD3	2.14	0.72
2:B:240:LEU:HD11	2:B:248:PHE:HZ	1.53	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:132:THR:HG22	2:D:133:ALA:N	2.06	0.70
2:B:156:LYS:O	2:B:156:LYS:HE3	1.92	0.69
2:B:365:HIS:O	2:B:368:GLU:HG2	1.92	0.69
2:B:277:MET:CE	6:B:404:UD2:H8'1	2.24	0.68
1:A:26:TRP:HZ2	1:A:96:LEU:HD11	1.60	0.67
2:B:255:LEU:HD11	6:B:404:UD2:H8'3	1.77	0.67
2:B:305:ASN:HD21	2:B:376:ASN:H	1.41	0.66
1:A:12:ILE:O	1:A:15:ILE:HG22	1.96	0.66
1:C:122:LYS:HD3	1:C:123:PRO:HD2	1.78	0.65
1:C:4:THR:H	1:C:7:LYS:HE2	1.61	0.65
1:C:10:HIS:HB2	7:C:1432:HOH:O	1.97	0.64
2:B:191:ARG:NH1	2:B:194:HIS:HD2	1.95	0.63
2:D:344:MET:CE	6:D:528:UD2:H8'2	2.27	0.63
2:D:336:ASN:ND2	2:D:338:VAL:H	1.95	0.62
2:B:138:SER:HB3	2:B:141:LEU:HD13	1.83	0.61
2:D:349:ARG:HH11	2:D:349:ARG:CG	2.12	0.60
2:B:272:HIS:HB3	2:B:334:ARG:HG2	1.84	0.60
2:B:299:GLN:HG3	7:B:1129:HOH:O	2.02	0.59
2:B:338:VAL:O	2:B:341:LYS:HG3	2.01	0.59
2:B:336:ASN:ND2	2:B:338:VAL:H	1.98	0.59
2:B:155:LEU:O	2:B:159:GLU:HG3	2.03	0.59
2:B:157:LEU:HD12	2:B:161:GLN:OE1	2.03	0.59
2:B:255:LEU:HD11	6:B:404:UD2:C8'	2.31	0.59
2:D:327:PHE:CE2	2:D:367:LYS:HD2	2.38	0.59
2:D:224:SER:HB2	7:D:980:HOH:O	2.01	0.58
2:D:191:ARG:HH11	2:D:194:HIS:HD2	1.51	0.58
2:D:225:MET:HE1	2:D:354:GLU:OE2	2.06	0.56
2:D:344:MET:HE2	6:D:528:UD2:H8'2	1.88	0.55
2:D:361:ASP:HB3	7:D:1433:HOH:O	2.07	0.55
2:D:379:THR:HA	7:D:1295:HOH:O	2.06	0.55
2:B:236:PHE:O	2:B:240:LEU:HD13	2.07	0.54
1:C:121:GLU:O	1:C:122:LYS:C	2.50	0.54
2:D:255:LEU:HD11	6:D:528:UD2:C8'	2.23	0.54
2:B:277:MET:HE1	6:B:404:UD2:H8'1	1.91	0.53
1:A:116:GLU:HB3	4:A:814:PG4:H11	1.91	0.53
2:B:287:VAL:HB	7:C:1417:HOH:O	2.07	0.53
1:C:32:HIS:HD2	7:D:1156:HOH:O	1.92	0.52
2:D:270:PRO:HG2	2:D:325:LEU:HD22	1.92	0.52
1:A:58:ARG:HB3	7:A:1296:HOH:O	2.10	0.52
1:A:68:GLU:HG2	1:A:68:GLU:O	2.08	0.52
1:C:4:THR:OG1	1:C:7:LYS:HG3	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:131:LEU:HD22	2:D:175:ASP:O	2.10	0.52
2:D:191:ARG:HD2	2:D:194:HIS:CD2	2.44	0.52
2:D:152:PRO:HA	7:D:1458:HOH:O	2.09	0.51
2:D:228:ARG:HD3	2:D:317:GLU:OE1	2.10	0.51
2:B:352:LYS:HD2	2:B:352:LYS:N	2.24	0.51
2:D:225:MET:HE2	2:D:352:LYS:CE	2.36	0.51
1:A:93:LYS:HD3	7:A:1418:HOH:O	2.11	0.50
1:C:10:HIS:O	1:C:13:LYS:HG3	2.11	0.50
2:D:351:LYS:O	2:D:352:LYS:HB2	2.12	0.50
2:D:336:ASN:HD22	2:D:338:VAL:N	2.03	0.50
2:D:339:ILE:HD11	7:D:1077:HOH:O	2.11	0.50
1:C:70:GLU:OE1	1:C:70:GLU:N	2.45	0.49
1:C:116:GLU:OE2	1:C:119:ARG:NH2	2.45	0.49
1:A:119:ARG:HG3	1:A:119:ARG:HH21	1.77	0.49
2:D:153:VAL:HG22	2:D:196:LYS:HB3	1.93	0.49
2:D:132:THR:CG2	2:D:133:ALA:H	2.22	0.49
1:C:16:ASP:HB2	1:C:23:LEU:HD21	1.95	0.49
2:D:365:HIS:O	2:D:368:GLU:HG3	2.13	0.49
1:A:55:ILE:HB	1:A:80:LEU:HD13	1.94	0.48
2:D:153:VAL:CG2	2:D:196:LYS:HB3	2.43	0.48
2:B:154:ASP:N	7:B:1408:HOH:O	2.45	0.48
2:D:349:ARG:HG2	2:D:349:ARG:NH1	2.13	0.48
2:B:279:LYS:HE2	7:B:1297:HOH:O	2.13	0.47
2:B:198:TRP:HA	2:B:256:ILE:HD11	1.97	0.47
1:C:4:THR:N	1:C:7:LYS:HE2	2.30	0.47
2:B:191:ARG:NH1	2:B:194:HIS:CD2	2.79	0.46
1:A:115:LEU:H	4:A:814:PG4:H32	1.80	0.46
2:B:327:PHE:CE1	2:B:367:LYS:HB2	2.50	0.46
2:D:167:LEU:HD22	2:D:387:ARG:HB3	1.98	0.46
2:B:191:ARG:HD2	2:B:194:HIS:CD2	2.51	0.46
2:D:156:LYS:O	2:D:160:GLN:HG2	2.16	0.46
2:B:164:LYS:HD2	2:B:174:MET:HE1	1.97	0.46
2:D:327:PHE:CZ	2:D:367:LYS:HB2	2.50	0.46
2:D:358:GLN:OE1	2:D:362:ARG:NH1	2.49	0.46
2:D:180:HIS:CE1	2:D:265:ARG:HD2	2.51	0.45
2:D:164:LYS:HE2	7:D:1342:HOH:O	2.16	0.45
2:B:174:MET:HE3	7:B:1163:HOH:O	2.17	0.45
1:C:4:THR:H	1:C:7:LYS:CE	2.26	0.45
2:D:358:GLN:O	2:D:362:ARG:HG3	2.16	0.45
2:B:269:GLN:O	2:B:270:PRO:C	2.60	0.45
2:D:336:ASN:ND2	2:D:336:ASN:C	2.75	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:270:PRO:CG	2:B:325:LEU:HD22	2.46	0.45
2:D:164:LYS:H	2:D:164:LYS:HD3	1.81	0.45
2:B:195:LEU:HD22	2:B:199:LEU:CD1	2.47	0.45
1:A:60:TRP:O	1:A:73:CYS:HB2	2.17	0.44
2:B:157:LEU:O	2:B:160:GLN:HB3	2.16	0.44
2:B:157:LEU:CD1	2:B:161:GLN:OE1	2.66	0.44
2:B:270:PRO:HG2	2:B:325:LEU:HD22	1.99	0.44
1:C:116:GLU:OE1	1:C:119:ARG:NH2	2.51	0.44
2:B:253:VAL:HG12	6:B:404:UD2:H2B	1.98	0.44
1:C:35:GLY:O	1:C:36:TYR:HB2	2.17	0.44
1:C:122:LYS:HD3	1:C:123:PRO:CD	2.44	0.44
2:D:191:ARG:NH1	2:D:194:HIS:HD2	2.14	0.44
2:D:272:HIS:HB3	2:D:334:ARG:HG2	2.00	0.43
1:A:16:ASP:CG	1:A:23:LEU:HG	2.43	0.43
1:A:70:GLU:O	1:A:71:ASN:C	2.60	0.43
2:D:336:ASN:ND2	2:D:338:VAL:N	2.62	0.43
1:C:55:ILE:HD13	1:C:91:CYS:SG	2.57	0.43
2:B:336:ASN:ND2	2:B:339:ILE:H	2.16	0.43
2:B:307:PHE:HB3	2:B:308:PRO:HD2	2.01	0.43
2:B:336:ASN:HD22	2:B:338:VAL:N	2.06	0.43
2:D:195:LEU:HD22	2:D:199:LEU:HG	2.00	0.43
2:D:400:THR:HB	2:D:401:PRO:HD2	2.00	0.43
2:B:336:ASN:HD22	2:B:336:ASN:C	2.26	0.43
2:B:343:ARG:HG2	2:B:343:ARG:HH11	1.84	0.43
2:B:316:GLY:HA2	2:B:363:ILE:HD11	2.00	0.42
2:D:336:ASN:ND2	2:D:338:VAL:HB	2.29	0.42
1:C:70:GLU:O	1:C:71:ASN:C	2.61	0.42
1:C:55:ILE:HB	1:C:80:LEU:HD13	2.01	0.42
2:D:362:ARG:HB3	7:D:1239:HOH:O	2.19	0.42
1:C:3:LEU:HA	1:C:7:LYS:HE3	2.02	0.42
1:C:6:CYS:SG	1:C:122:LYS:HB2	2.59	0.42
2:D:365:HIS:HD2	2:D:368:GLU:OE1	2.02	0.42
1:A:116:GLU:OE2	1:A:119:ARG:NE	2.42	0.42
2:B:167:LEU:HA	2:B:389:PRO:HA	2.01	0.42
2:D:292:GLY:HA2	6:D:528:UD2:H8'3	2.02	0.42
2:B:230:LYS:HD3	2:B:398:ILE:HB	2.02	0.42
2:D:157:LEU:HD22	2:D:157:LEU:HA	1.86	0.42
1:A:75:ILE:CD1	1:A:87:ASP:HB2	2.50	0.42
1:C:8:VAL:HG21	1:C:36:TYR:CD1	2.55	0.41
2:B:131:LEU:CD2	2:B:131:LEU:N	2.83	0.41
2:D:279:LYS:CD	2:D:346:ARG:HH12	2.33	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:285:PRO:HG2	2:D:289:TYR:CG	2.55	0.41
1:A:119:ARG:HG3	1:A:119:ARG:NH2	2.35	0.41
2:B:277:MET:HE2	2:B:280:PHE:CD1	2.55	0.41
2:D:228:ARG:NH1	2:D:317:GLU:OE1	2.52	0.41
2:B:195:LEU:HD22	2:B:199:LEU:HG	2.02	0.41
2:D:208:ARG:NH2	2:D:260:ASP:OD1	2.50	0.41
2:D:240:LEU:HD12	2:D:240:LEU:HA	1.83	0.41
2:B:135:PRO:HG3	2:B:139:PRO:HD3	2.03	0.40
2:D:132:THR:CG2	2:D:133:ALA:N	2.77	0.40
2:D:164:LYS:HD3	2:D:164:LYS:N	2.36	0.40
2:D:400:THR:HB	2:D:401:PRO:CD	2.51	0.40
2:B:225:MET:HE3	2:B:352:LYS:O	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%)	114 (94%)	7 (6%)	0	100	100
1	C	121/123 (98%)	117 (97%)	4 (3%)	0	100	100
2	B	270/286 (94%)	262 (97%)	8 (3%)	0	100	100
2	D	270/286 (94%)	261 (97%)	9 (3%)	0	100	100
All	All	782/818 (96%)	754 (96%)	28 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar

resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	109/109 (100%)	106 (97%)	3 (3%)	38	43
1	C	109/109 (100%)	104 (95%)	5 (5%)	24	25
2	B	245/254 (96%)	232 (95%)	13 (5%)	20	19
2	D	245/254 (96%)	229 (94%)	16 (6%)	15	13
All	All	708/726 (98%)	671 (95%)	37 (5%)	21	20

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	ASN
1	A	70	GLU
1	A	85	LEU
2	B	131	LEU
2	B	146	LEU
2	B	156	LYS
2	B	157	LEU
2	B	164	LYS
2	B	167	LEU
2	B	195	LEU
2	B	263	THR
2	B	275	VAL
2	B	302	LEU
2	B	336	ASN
2	B	352	LYS
2	B	368	GLU
1	C	47	SER
1	C	66	PHE
1	C	79	LYS
1	C	84	GLU
1	C	85	LEU
2	D	141	LEU
2	D	153	VAL
2	D	156	LYS
2	D	157	LEU
2	D	167	LEU
2	D	195	LEU
2	D	223	GLU

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Mol	Chain	Res	Type
2	D	240	LEU
2	D	270	PRO
2	D	287	VAL
2	D	299	GLN
2	D	302	LEU
2	D	336	ASN
2	D	349	ARG
2	D	351	LYS
2	D	368	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	ASN
1	A	71	ASN
2	B	150	ASN
2	B	160	GLN
2	B	194	HIS
2	B	219	ASN
2	B	305	ASN
2	B	310	ASN
2	B	336	ASN
1	C	32	HIS
1	C	71	ASN
2	D	160	GLN
2	D	180	HIS
2	D	190	ASN
2	D	194	HIS
2	D	261	HIS
2	D	305	ASN
2	D	310	ASN
2	D	336	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 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	UD2	D	528	5	40,41,41	1.58	7 (17%)	59,62,62	1.66	14 (23%)
4	PG4	B	813	-	6,6,12	0.65	0	5,5,11	0.14	0
6	UD2	B	404	5	40,41,41	1.59	6 (15%)	59,62,62	1.68	15 (25%)
4	PG4	A	814	-	6,6,12	0.46	0	5,5,11	0.33	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	UD2	D	528	5	-	8/26/63/63	0/3/3/3
4	PG4	B	813	-	-	3/4/4/10	-
6	UD2	B	404	5	-	8/26/63/63	0/3/3/3
4	PG4	A	814	-	-	2/4/4/10	-

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	D	528	UD2	PA-O3A	5.44	1.65	1.59
6	B	404	UD2	PA-O3A	5.30	1.65	1.59
6	D	528	UD2	PB-O1'	-4.30	1.46	1.59
6	B	404	UD2	PB-O3A	-3.39	1.55	1.59
6	B	404	UD2	PB-O1'	-3.33	1.49	1.59
6	D	528	UD2	O5'-C1'	2.53	1.48	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	404	UD2	C2-N1	2.37	1.42	1.38
6	D	528	UD2	C4-N3	2.31	1.42	1.38
6	D	528	UD2	C1'-C2'	2.21	1.56	1.53
6	B	404	UD2	O5'-C5'	2.18	1.49	1.44
6	D	528	UD2	C2-N1	2.11	1.41	1.38
6	D	528	UD2	O5'-C5'	2.08	1.49	1.44
6	B	404	UD2	C4'-C5'	2.04	1.57	1.53

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	404	UD2	O5'-C5'-C6'	-4.65	94.92	106.44
6	D	528	UD2	O1B-PB-O1'	4.33	124.32	106.70
6	B	404	UD2	O1B-PB-O1'	4.23	123.93	106.70
6	D	528	UD2	O5'-C5'-C6'	-3.92	96.72	106.44
6	D	528	UD2	O5'-C1'-O1'	3.86	116.41	111.36
6	D	528	UD2	O5B-C5B-C4B	3.23	120.00	108.99
6	D	528	UD2	C8'-C7'-N2'	-3.23	110.76	116.12
6	B	404	UD2	C8'-C7'-N2'	-3.17	110.85	116.12
6	D	528	UD2	O7'-C7'-N2'	3.07	127.41	121.98
6	B	404	UD2	O4B-C4B-C3B	-3.07	99.07	105.15
6	B	404	UD2	O5'-C1'-O1'	2.97	115.25	111.36
6	B	404	UD2	O2-C2-N1	-2.95	118.95	122.80
6	B	404	UD2	C3'-C2'-N2'	-2.88	105.30	110.62
6	B	404	UD2	O1B-PB-O3A	-2.86	99.55	107.27
6	D	528	UD2	O2-C2-N1	-2.79	119.17	122.80
6	B	404	UD2	O5B-C5B-C4B	2.76	118.40	108.99
6	D	528	UD2	O4B-C4B-C3B	-2.75	99.70	105.15
6	D	528	UD2	O1B-PB-O3A	-2.74	99.87	107.27
6	B	404	UD2	O7'-C7'-N2'	2.71	126.78	121.98
6	D	528	UD2	O2-C2-N3	2.43	125.97	121.49
6	D	528	UD2	O3'-C3'-C4'	-2.38	104.76	110.38
6	B	404	UD2	O2-C2-N3	2.37	125.86	121.49
6	B	404	UD2	C6'-C5'-C4'	2.18	118.38	113.02
6	B	404	UD2	C2B-C1B-N1	2.18	119.31	113.25
6	D	528	UD2	O1A-PA-O3A	2.09	112.91	107.27
6	B	404	UD2	O1A-PA-O3A	2.06	112.85	107.27
6	B	404	UD2	O3'-C3'-C2'	2.06	113.68	109.58
6	D	528	UD2	C4'-C3'-C2'	-2.03	107.44	110.40
6	D	528	UD2	O3'-C3'-C2'	2.03	113.62	109.58

There are no chirality outliers.

All (21) torsion outliers are listed below:

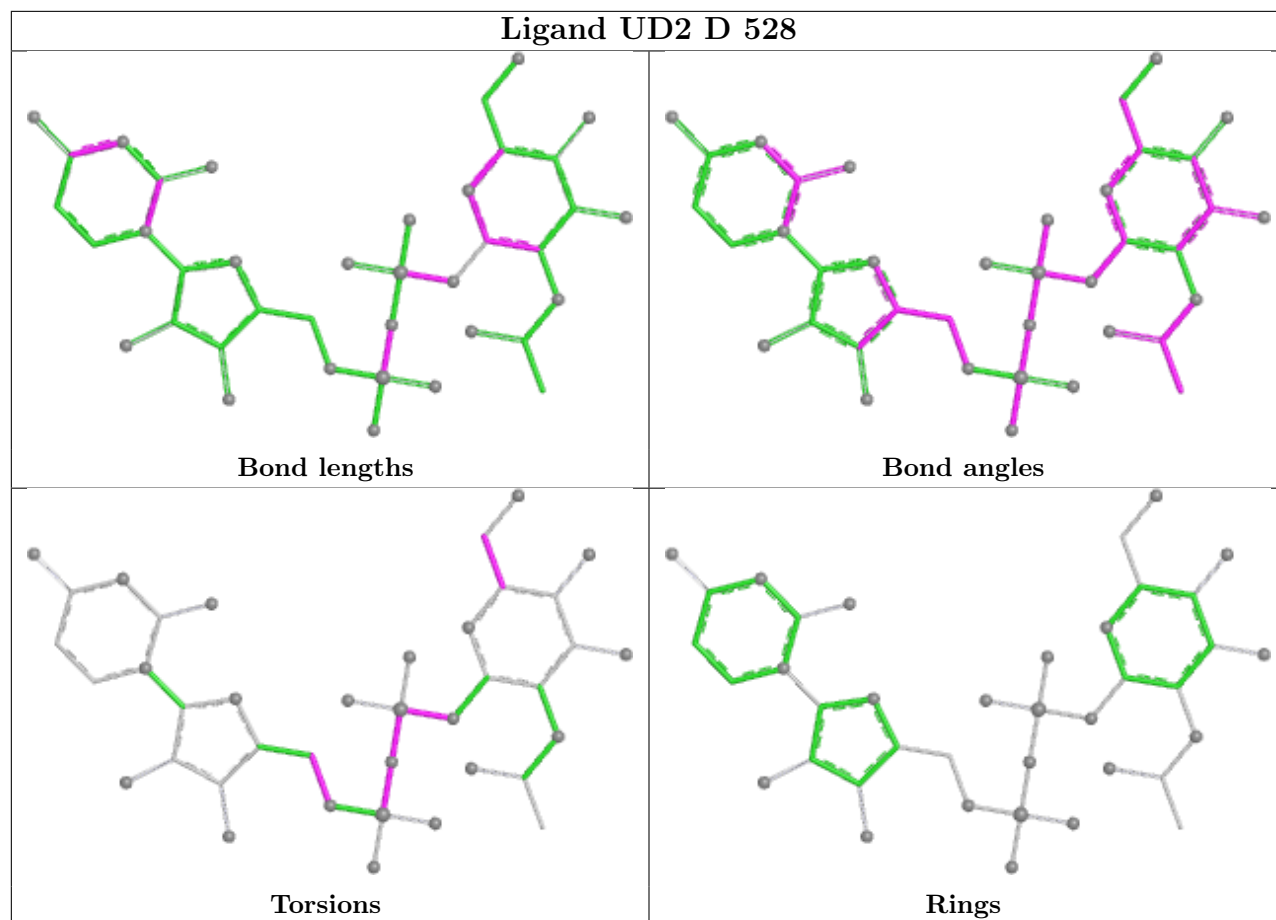
Mol	Chain	Res	Type	Atoms
6	B	404	UD2	C1'-C2'-N2'-C7'
6	B	404	UD2	C1'-O1'-PB-O3A
6	B	404	UD2	C1'-O1'-PB-O1B
6	D	528	UD2	C1'-O1'-PB-O3A
6	D	528	UD2	C1'-O1'-PB-O1B
4	B	813	PG4	O1-C1-C2-O2
4	A	814	PG4	O1-C1-C2-O2
6	D	528	UD2	O5'-C5'-C6'-O6'
6	B	404	UD2	O5'-C5'-C6'-O6'
6	D	528	UD2	C1'-O1'-PB-O2B
4	A	814	PG4	C1-C2-O2-C3
6	B	404	UD2	C4B-C5B-O5B-PA
6	B	404	UD2	C5B-O5B-PA-O3A
4	B	813	PG4	C4-C3-O2-C2
6	D	528	UD2	C4B-C5B-O5B-PA
6	B	404	UD2	PB-O3A-PA-O1A
6	D	528	UD2	PB-O3A-PA-O1A
6	D	528	UD2	PA-O3A-PB-O1B
4	B	813	PG4	O2-C3-C4-O3
6	B	404	UD2	PB-O3A-PA-O2A
6	D	528	UD2	PB-O3A-PA-O2A

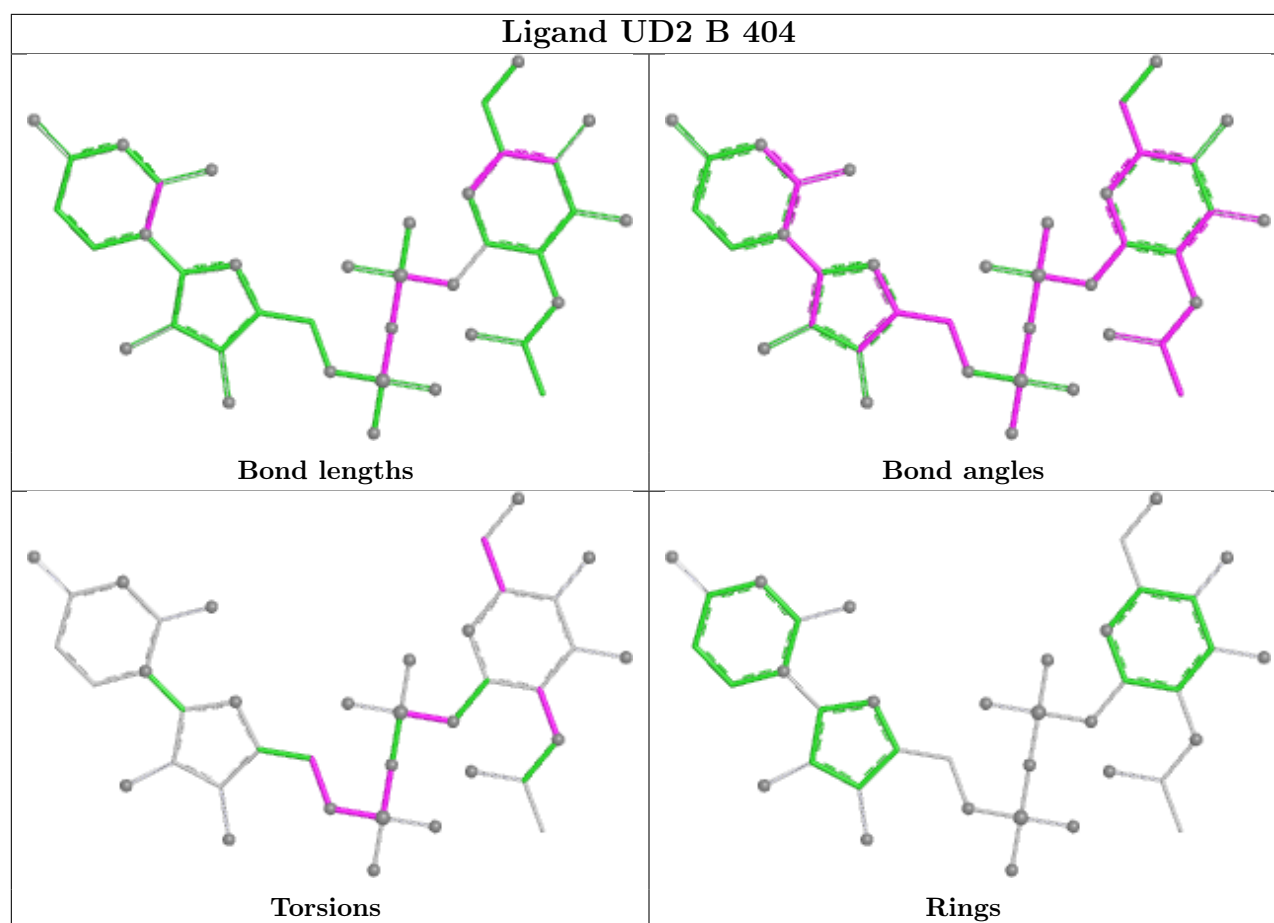
There are no ring outliers.

3 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	528	UD2	6	0
6	B	404	UD2	5	0
4	A	814	PG4	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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%)	-0.09	3 (2%) 59 62	21, 30, 58, 69	0
1	C	123/123 (100%)	-0.26	4 (3%) 49 51	19, 28, 51, 82	0
2	B	272/286 (95%)	-0.15	2 (0%) 84 86	20, 32, 56, 67	0
2	D	272/286 (95%)	-0.15	1 (0%) 88 90	20, 33, 51, 62	0
All	All	790/818 (96%)	-0.16	10 (1%) 75 77	19, 31, 55, 82	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	123	PRO	4.2
2	B	349	ARG	2.8
1	C	66	PHE	2.7
1	A	66	PHE	2.6
1	C	121	GLU	2.6
2	B	174	MET	2.6
1	A	45	ASN	2.3
2	D	346	ARG	2.2
1	A	123	PRO	2.2
1	C	45	ASN	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](#)

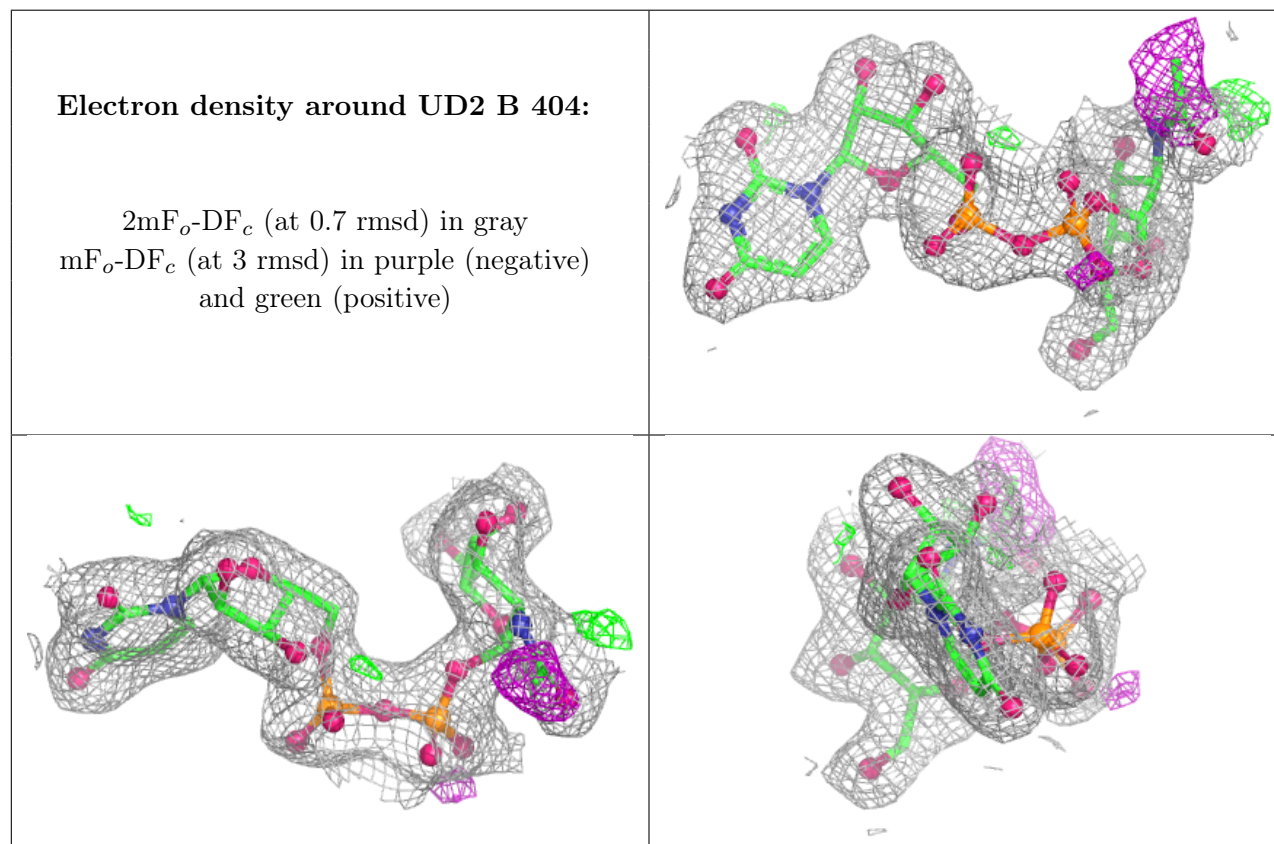
There are no oligosaccharides in this entry.

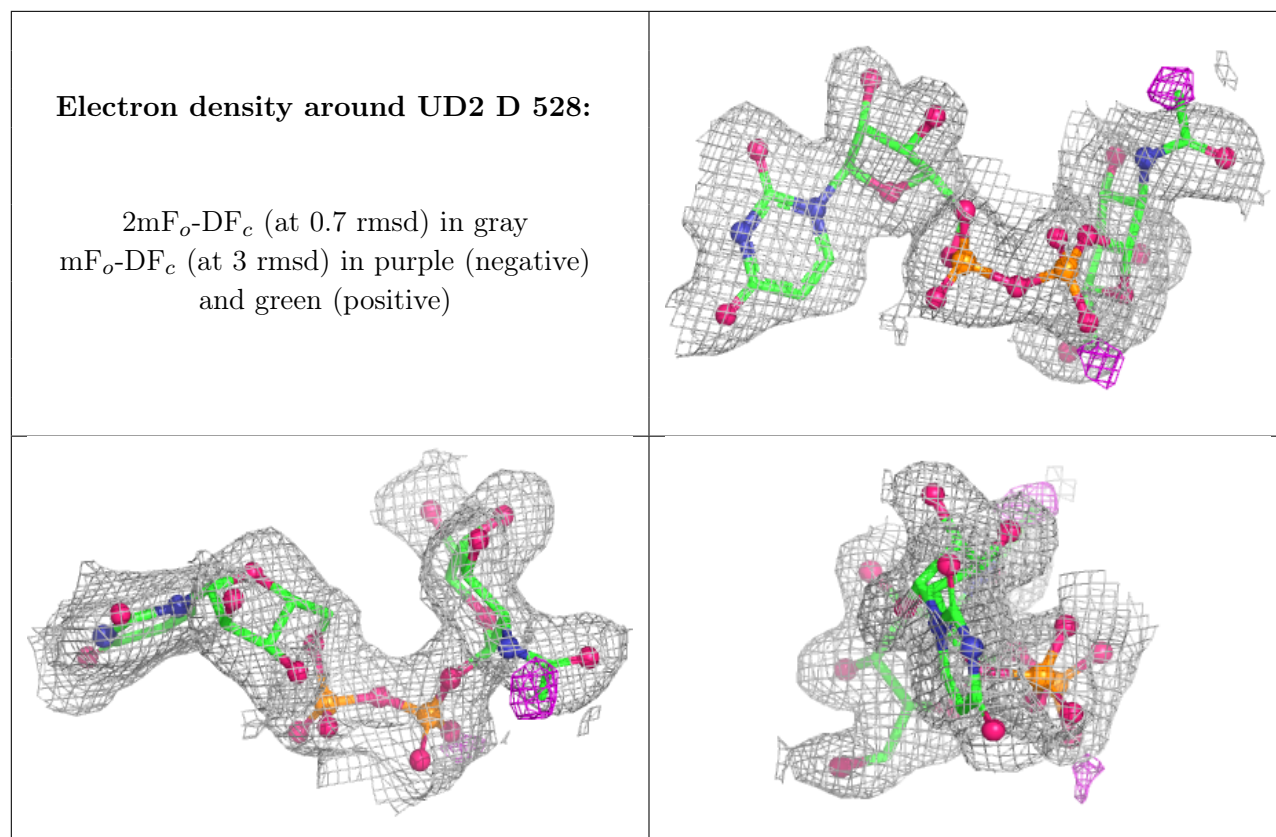
6.4 Ligands ⓘ

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
4	PG4	A	814	7/13	0.81	0.14	47,53,55,58	0
4	PG4	B	813	7/13	0.82	0.12	48,50,52,53	0
6	UD2	B	404	39/39	0.96	0.07	20,30,38,41	0
6	UD2	D	528	39/39	0.97	0.06	20,24,30,36	0
5	MN	D	403	1/1	0.99	0.02	29,29,29,29	0
3	CA	A	124	1/1	0.99	0.02	24,24,24,24	0
3	CA	C	124	1/1	0.99	0.01	26,26,26,26	0
5	MN	B	403	1/1	1.00	0.01	31,31,31,31	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.





6.5 Other polymers [i](#)

There are no such residues in this entry.