



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:32 AM UTC

PDB ID : 3MJT / pdb_00003mjt
Title : Structure of A-type Ketoreductases from Modular Polyketide Synthase
Authors : Zheng, J.; Taylor, C.A.; Piasecki, S.K.; Keatinge-Clay, A.T.
Deposited on : 2010-04-13
Resolution : 1.60 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

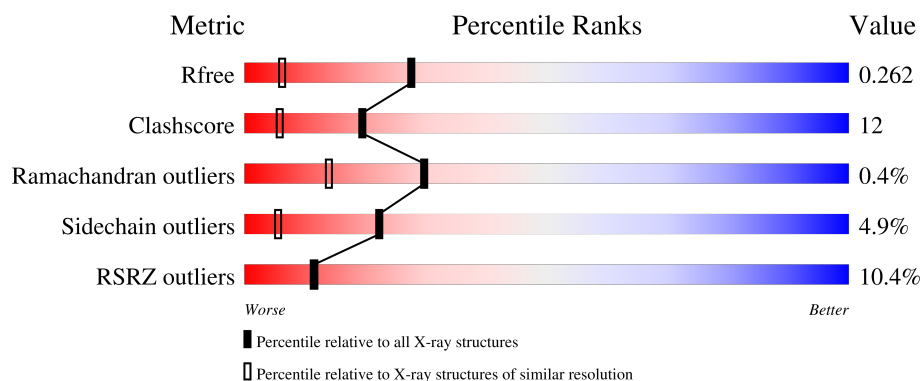
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (1.60-1.60)

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	496	11% 75% 19% • •
1	B	496	9% 73% 19% • • •

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
2	NDP	A	476	X	-	-	-
2	NDP	B	476	X	-	-	-
3	GOL	A	477	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7581 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 AmphB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	476	Total	C	N	O	S	0	0	0
			3466	2164	636	657	9			
1	B	475	Total	C	N	O	S	0	0	0
			3460	2161	635	655	9			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	expression tag	UNP Q93NW7
A	-19	GLY	-	expression tag	UNP Q93NW7
A	-18	SER	-	expression tag	UNP Q93NW7
A	-17	SER	-	expression tag	UNP Q93NW7
A	-16	HIS	-	expression tag	UNP Q93NW7
A	-15	HIS	-	expression tag	UNP Q93NW7
A	-14	HIS	-	expression tag	UNP Q93NW7
A	-13	HIS	-	expression tag	UNP Q93NW7
A	-12	HIS	-	expression tag	UNP Q93NW7
A	-11	HIS	-	expression tag	UNP Q93NW7
A	-10	SER	-	expression tag	UNP Q93NW7
A	-9	SER	-	expression tag	UNP Q93NW7
A	-8	GLY	-	expression tag	UNP Q93NW7
A	-7	LEU	-	expression tag	UNP Q93NW7
A	-6	VAL	-	expression tag	UNP Q93NW7
A	-5	PRO	-	expression tag	UNP Q93NW7
A	-4	ARG	-	expression tag	UNP Q93NW7
A	-3	GLY	-	expression tag	UNP Q93NW7
A	-2	SER	-	expression tag	UNP Q93NW7
A	-1	HIS	-	expression tag	UNP Q93NW7
A	0	MET	-	expression tag	UNP Q93NW7
A	364	HIS	GLN	engineered mutation	UNP Q93NW7
B	-20	MET	-	expression tag	UNP Q93NW7
B	-19	GLY	-	expression tag	UNP Q93NW7
B	-18	SER	-	expression tag	UNP Q93NW7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-17	SER	-	expression tag	UNP Q93NW7
B	-16	HIS	-	expression tag	UNP Q93NW7
B	-15	HIS	-	expression tag	UNP Q93NW7
B	-14	HIS	-	expression tag	UNP Q93NW7
B	-13	HIS	-	expression tag	UNP Q93NW7
B	-12	HIS	-	expression tag	UNP Q93NW7
B	-11	HIS	-	expression tag	UNP Q93NW7
B	-10	SER	-	expression tag	UNP Q93NW7
B	-9	SER	-	expression tag	UNP Q93NW7
B	-8	GLY	-	expression tag	UNP Q93NW7
B	-7	LEU	-	expression tag	UNP Q93NW7
B	-6	VAL	-	expression tag	UNP Q93NW7
B	-5	PRO	-	expression tag	UNP Q93NW7
B	-4	ARG	-	expression tag	UNP Q93NW7
B	-3	GLY	-	expression tag	UNP Q93NW7
B	-2	SER	-	expression tag	UNP Q93NW7
B	-1	HIS	-	expression tag	UNP Q93NW7
B	0	MET	-	expression tag	UNP Q93NW7
B	364	HIS	GLN	engineered mutation	UNP Q93NW7

- # NDP

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 48	C 21	N 7	O 17	P 3	0	0



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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

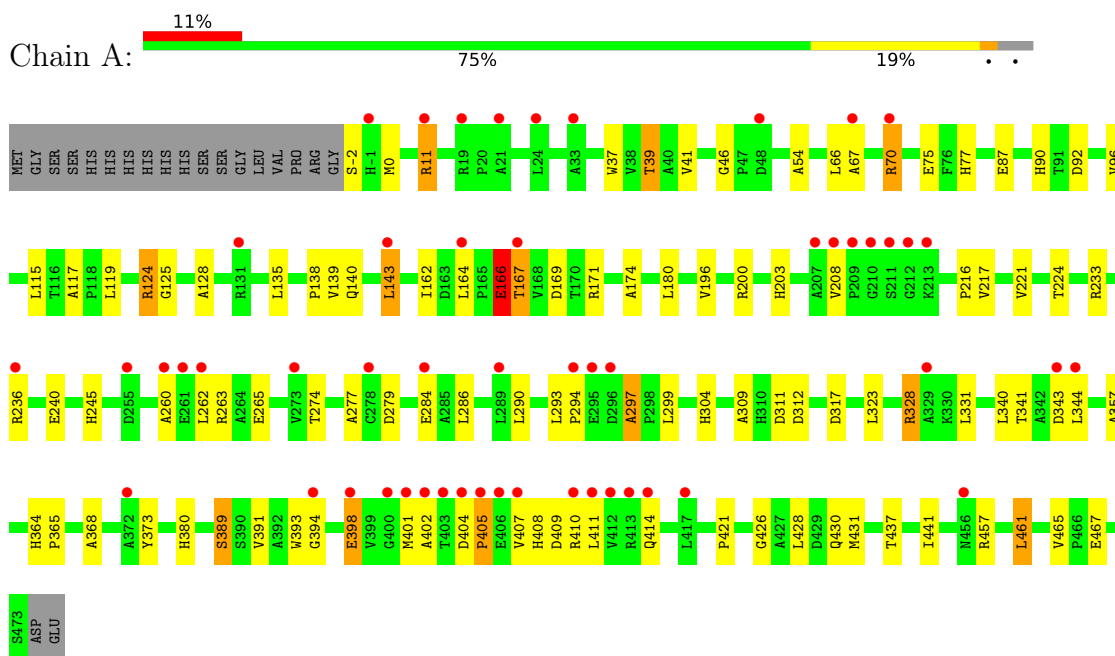
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	268	Total	O	0	0
			268	268		
4	B	279	Total	O	0	0
			279	279		

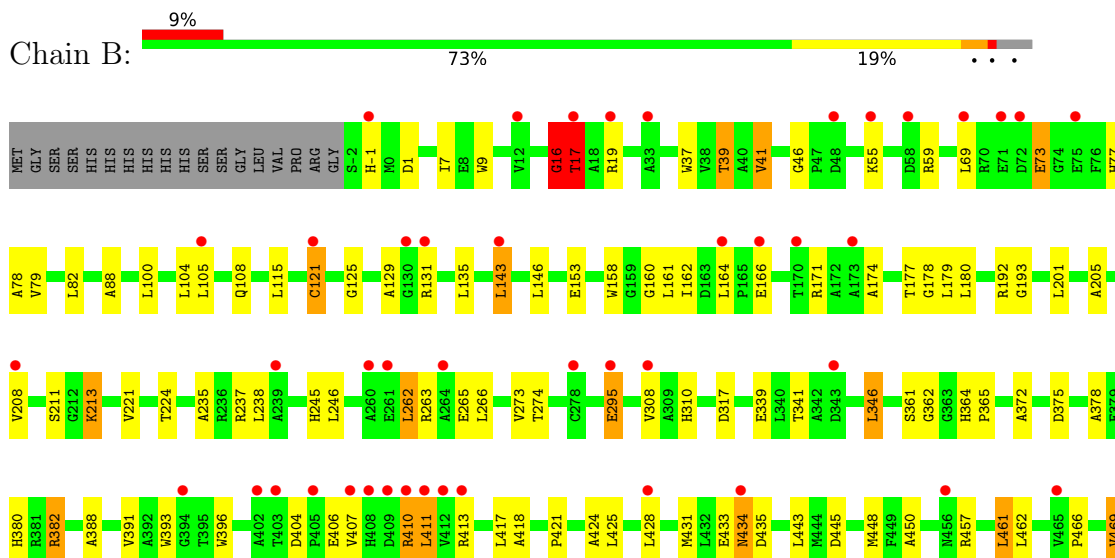
3 Residue-property plots [i](#)

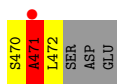
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: AmphB



• Molecule 1: AmphB





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	61.40Å 63.61Å 71.85Å 72.84° 67.20° 89.77°	Depositor
Resolution (Å)	62.79 – 1.60 62.79 – 1.60	Depositor EDS
% Data completeness (in resolution range)	94.7 (62.79-1.60) 94.7 (62.79-1.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.32 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.223 , 0.265 0.222 , 0.262	Depositor DCC
R_{free} test set	5935 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å ²)	16.5	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 40.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.109 for h,-k,h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7581	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.74% 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: GOL, NDP

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	1.60	23/3537 (0.7%)	1.37	18/4832 (0.4%)
1	B	1.63	25/3531 (0.7%)	1.43	22/4824 (0.5%)
All	All	1.61	48/7068 (0.7%)	1.40	40/9656 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	3

All (48) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	17	THR	N-CA	9.90	1.58	1.46
1	B	158	TRP	CA-C	-8.37	1.42	1.52
1	A	96	VAL	C-O	-8.24	1.15	1.24
1	A	357	ALA	CA-CB	7.78	1.66	1.53
1	B	17	THR	CA-C	7.73	1.63	1.52
1	A	54	ALA	CA-CB	7.18	1.61	1.52
1	A	39	THR	CB-CG2	-7.06	1.29	1.52
1	A	277	ALA	CA-CB	7.01	1.61	1.53
1	A	465	VAL	CA-CB	6.88	1.60	1.53
1	B	471	ALA	CA-CB	6.76	1.64	1.53
1	A	357	ALA	C-O	-6.73	1.15	1.24
1	A	162	ILE	CA-CB	6.39	1.63	1.54
1	B	317	ASP	N-CA	6.38	1.54	1.45
1	A	430	GLN	N-CA	6.31	1.53	1.46
1	B	457	ARG	CB-CG	-6.25	1.33	1.52
1	B	424	ALA	C-O	6.23	1.31	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	297	ALA	N-CA	6.12	1.53	1.46
1	B	88	ALA	CA-CB	-6.12	1.43	1.53
1	A	39	THR	CA-CB	6.07	1.62	1.53
1	B	179	LEU	N-CA	6.03	1.53	1.46
1	B	153	GLU	C-O	-6.01	1.16	1.24
1	B	129	ALA	C-O	-5.89	1.17	1.24
1	A	312	ASP	C-O	5.86	1.31	1.23
1	B	469	VAL	CA-CB	-5.82	1.48	1.54
1	A	441	ILE	C-O	5.75	1.30	1.24
1	A	117	ALA	CA-C	5.62	1.59	1.53
1	A	389	SER	C-O	5.56	1.30	1.23
1	A	323	LEU	CA-C	5.55	1.59	1.52
1	B	418	ALA	CA-CB	5.50	1.61	1.53
1	A	317	ASP	N-CA	5.37	1.53	1.46
1	B	79	VAL	CA-CB	5.35	1.60	1.54
1	B	78	ALA	CA-CB	-5.34	1.44	1.53
1	B	224	THR	N-CA	-5.30	1.39	1.45
1	B	372	ALA	C-O	5.30	1.30	1.24
1	B	9	TRP	C-O	5.23	1.29	1.24
1	A	465	VAL	CA-C	-5.21	1.48	1.53
1	A	119	LEU	CA-CB	5.21	1.60	1.53
1	B	7	ILE	CB-CG2	5.19	1.69	1.52
1	B	104	LEU	N-CA	5.19	1.52	1.46
1	B	162	ILE	C-O	5.16	1.29	1.24
1	A	208	VAL	C-O	5.16	1.30	1.24
1	B	450	ALA	CA-C	-5.14	1.46	1.52
1	A	143	LEU	N-CA	-5.14	1.39	1.46
1	A	309	ALA	N-CA	5.11	1.53	1.45
1	B	362	GLY	N-CA	5.10	1.51	1.45
1	B	361	SER	C-O	5.09	1.29	1.23
1	A	221	VAL	N-CA	5.02	1.52	1.46
1	B	1	ASP	C-O	5.00	1.30	1.24

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	382	ARG	NE-CZ-NH1	11.43	132.93	121.50
1	B	121	CYS	CA-CB-SG	-10.62	89.99	114.40
1	B	382	ARG	NE-CZ-NH2	-10.51	109.74	119.20
1	B	17	THR	N-CA-C	8.59	129.09	110.80
1	A	174	ALA	N-CA-C	-8.06	102.07	112.23
1	A	293	LEU	N-CA-C	7.92	118.70	109.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	46	GLY	CA-C-N	-7.86	111.73	120.45
1	B	46	GLY	C-N-CA	-7.86	111.73	120.45
1	A	328	ARG	N-CA-C	7.63	120.27	111.11
1	B	382	ARG	CD-NE-CZ	7.50	134.90	124.40
1	B	16	GLY	CA-C-N	-6.89	108.38	121.54
1	B	16	GLY	C-N-CA	-6.89	108.38	121.54
1	B	382	ARG	CB-CG-CD	6.71	126.73	111.30
1	B	434	ASN	N-CA-C	6.63	121.20	113.18
1	A	368	ALA	CA-C-O	-6.62	113.88	120.70
1	A	46	GLY	CA-C-N	-6.29	113.47	120.45
1	A	46	GLY	C-N-CA	-6.29	113.47	120.45
1	B	433	GLU	N-CA-C	-6.05	104.80	111.82
1	B	146	LEU	N-CA-C	5.89	117.70	111.28
1	B	205	ALA	CA-C-N	5.83	125.76	119.76
1	B	205	ALA	C-N-CA	5.83	125.76	119.76
1	B	39	THR	N-CA-C	5.79	117.59	111.28
1	B	16	GLY	O-C-N	-5.74	116.19	122.68
1	A	70	ARG	NE-CZ-NH1	-5.64	115.86	121.50
1	A	196	VAL	CB-CA-C	-5.62	103.19	110.84
1	A	260	ALA	CA-C-N	5.54	128.02	120.54
1	A	260	ALA	C-N-CA	5.54	128.02	120.54
1	A	297	ALA	CA-C-N	5.46	125.41	119.78
1	A	297	ALA	C-N-CA	5.46	125.41	119.78
1	A	169	ASP	N-CA-CB	5.46	117.91	110.38
1	B	375	ASP	O-C-N	5.34	127.78	122.12
1	A	200	ARG	NE-CZ-NH2	-5.33	114.41	119.20
1	B	365	PRO	CA-C-N	-5.18	113.64	120.22
1	B	365	PRO	C-N-CA	-5.18	113.64	120.22
1	A	166	GLU	CA-C-N	-5.17	115.72	123.00
1	A	166	GLU	C-N-CA	-5.17	115.72	123.00
1	B	41	VAL	N-CA-C	5.16	115.37	110.42
1	B	174	ALA	N-CA-C	-5.13	105.87	111.82
1	A	461	LEU	CD1-CG-CD2	5.06	121.93	110.80
1	A	167	THR	CB-CA-C	5.01	118.47	110.16

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	16	GLY	Peptide
1	B	17	THR	Peptide
1	B	471	ALA	Peptide

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	3466	0	3425	66	0
1	B	3460	0	3420	95	0
2	A	48	0	25	1	0
2	B	48	0	25	1	0
3	A	6	0	8	4	0
3	B	6	0	8	0	0
4	A	268	0	0	12	0
4	B	279	0	0	8	0
All	All	7581	0	6911	161	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 (161) 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:B:121:CYS:HB3	4:B:756:HOH:O	1.11	1.27
1:B:121:CYS:SG	1:B:121:CYS:O	2.04	1.14
1:B:461:LEU:O	1:B:461:LEU:HD13	1.54	1.08
1:B:470:SER:O	1:B:471:ALA:HB3	1.55	1.03
1:B:461:LEU:HD13	1:B:461:LEU:C	1.88	0.98
1:B:121:CYS:SG	1:B:161:LEU:HD23	2.02	0.98
1:B:208:VAL:HG23	1:B:211:SER:OG	1.65	0.96
1:A:263:ARG:HD3	4:A:666:HOH:O	1.67	0.94
1:B:407:VAL:O	1:B:411:LEU:HD13	1.67	0.93
1:B:262:LEU:O	1:B:266:LEU:HD13	1.69	0.92
1:B:470:SER:O	1:B:471:ALA:CB	2.16	0.92
1:B:17:THR:HG21	1:B:177:THR:OG1	1.68	0.92
1:A:90:HIS:HD2	1:A:92:ASP:H	1.13	0.91
1:B:461:LEU:C	1:B:461:LEU:CD1	2.47	0.87
1:B:391:VAL:HG21	1:B:428:LEU:HD13	1.57	0.87
1:B:201:LEU:HB2	1:B:461:LEU:HD11	1.56	0.85
1:A:404:ASP:HB3	1:A:407:VAL:HG22	1.59	0.84
1:B:59:ARG:HB2	1:B:105:LEU:HD12	1.58	0.84
1:B:121:CYS:SG	1:B:161:LEU:HA	2.18	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:391:VAL:HG21	1:A:428:LEU:HD13	1.58	0.83
1:A:341:THR:HA	1:A:344:LEU:HD13	1.59	0.82
1:A:216:PRO:HA	3:A:477:GOL:H31	1.64	0.79
1:B:295:GLU:H	1:B:295:GLU:CD	1.89	0.79
1:A:11:ARG:HH21	1:A:11:ARG:HG3	1.48	0.78
1:B:411:LEU:CD1	1:B:411:LEU:H	1.97	0.78
1:B:41:VAL:HG21	1:B:164:LEU:HD11	1.65	0.77
1:A:240:GLU:OE1	4:A:511:HOH:O	2.03	0.76
1:B:59:ARG:HB2	1:B:105:LEU:CD1	2.17	0.75
1:B:471:ALA:O	1:B:472:LEU:C	2.31	0.74
1:A:224:THR:OG1	1:A:304:HIS:HD2	1.73	0.71
1:A:39:THR:HB	4:A:731:HOH:O	1.90	0.71
1:B:41:VAL:CG2	1:B:164:LEU:HD11	2.19	0.71
1:B:407:VAL:O	1:B:411:LEU:CD1	2.39	0.71
1:A:217:VAL:H	3:A:477:GOL:H11	1.57	0.69
1:B:39:THR:HB	4:B:500:HOH:O	1.92	0.68
1:B:201:LEU:CB	1:B:461:LEU:HD11	2.21	0.68
1:A:263:ARG:NH2	4:A:666:HOH:O	2.06	0.67
1:A:166:GLU:H	1:A:166:GLU:CD	2.03	0.67
1:A:203:HIS:HE1	1:A:467:GLU:OE2	1.76	0.67
1:A:263:ARG:CD	4:A:666:HOH:O	2.35	0.65
1:A:411:LEU:O	1:A:414:GLN:HG2	1.95	0.65
1:B:37:TRP:CE3	1:B:164:LEU:HD12	2.31	0.65
1:B:461:LEU:HD12	1:B:462:LEU:HD23	1.78	0.64
1:A:11:ARG:HH21	1:A:11:ARG:CG	2.09	0.64
1:B:466:PRO:O	1:B:469:VAL:CG2	2.45	0.64
1:A:398:GLU:HG3	1:A:421:PRO:CD	2.28	0.64
1:B:121:CYS:SG	1:B:161:LEU:CD2	2.81	0.63
1:A:77:HIS:HE1	1:A:180:LEU:O	1.81	0.63
1:A:203:HIS:HD2	4:A:512:HOH:O	1.81	0.63
1:A:284:GLU:HB2	4:A:671:HOH:O	1.99	0.63
1:A:286:LEU:HD23	1:A:340:LEU:HD12	1.80	0.62
1:B:100:LEU:HD12	1:B:143:LEU:HD13	1.81	0.62
1:A:135:LEU:O	1:A:380:HIS:HD2	1.83	0.61
1:B:411:LEU:HD13	1:B:411:LEU:H	1.65	0.61
1:A:402:ALA:O	1:A:408:HIS:HB2	2.01	0.61
1:B:393:TRP:HB2	2:B:476:NDP:C5N	2.31	0.61
1:A:245:HIS:HE1	1:A:274:THR:OG1	1.83	0.60
1:A:224:THR:OG1	1:A:304:HIS:CD2	2.53	0.60
1:B:466:PRO:HA	1:B:469:VAL:CG2	2.31	0.60
1:B:208:VAL:HG23	1:B:211:SER:HG	1.66	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:17:THR:CG2	1:B:177:THR:OG1	2.47	0.59
1:A:414:GLN:NE2	4:A:591:HOH:O	2.35	0.59
1:B:411:LEU:CD1	1:B:411:LEU:N	2.64	0.59
1:B:364:HIS:HD2	4:B:550:HOH:O	1.86	0.58
1:A:391:VAL:HG22	1:A:431:MET:SD	2.44	0.58
1:B:245:HIS:HE1	1:B:274:THR:OG1	1.85	0.58
1:A:67:ALA:O	1:A:70:ARG:HG3	2.03	0.58
1:B:16:GLY:O	1:B:17:THR:C	2.47	0.57
1:A:90:HIS:CD2	1:A:92:ASP:H	2.06	0.56
1:B:411:LEU:N	1:B:411:LEU:HD12	2.20	0.56
1:B:466:PRO:O	1:B:469:VAL:HG22	2.05	0.56
1:B:17:THR:HG23	1:B:178:GLY:N	2.21	0.56
1:B:266:LEU:HD12	1:B:266:LEU:N	2.20	0.55
1:B:391:VAL:HG22	1:B:431:MET:SD	2.47	0.55
1:B:77:HIS:HE1	1:B:180:LEU:O	1.90	0.55
1:B:105:LEU:HD13	1:B:108:GLN:OE1	2.06	0.55
1:A:391:VAL:HG21	1:A:428:LEU:CD1	2.34	0.55
1:B:171:ARG:NH1	4:B:484:HOH:O	2.39	0.54
1:B:266:LEU:N	1:B:266:LEU:CD1	2.70	0.54
1:B:471:ALA:O	1:B:472:LEU:O	2.25	0.54
1:A:279:ASP:OD1	1:A:328:ARG:NH2	2.41	0.54
1:B:308:VAL:HG13	1:B:310:HIS:ND1	2.23	0.54
1:A:37:TRP:CG	1:A:124:ARG:HG3	2.43	0.53
1:B:466:PRO:O	1:B:469:VAL:HG23	2.07	0.53
1:B:466:PRO:HA	1:B:469:VAL:HG22	1.89	0.53
1:A:405:PRO:HA	1:A:409:ASP:OD1	2.08	0.53
1:A:70:ARG:HG2	1:A:115:LEU:HD11	1.91	0.52
1:A:401:MET:O	1:A:407:VAL:HG21	2.10	0.52
1:B:262:LEU:HD22	1:B:266:LEU:HD11	1.92	0.52
1:B:263:ARG:HG3	1:B:273:VAL:HG11	1.92	0.52
1:B:382:ARG:HD3	1:B:388:ALA:O	2.09	0.52
1:B:19:ARG:CZ	1:B:19:ARG:HB3	2.41	0.51
1:B:391:VAL:HG21	1:B:428:LEU:CD1	2.35	0.51
1:B:341:THR:HB	1:B:346:LEU:HD22	1.92	0.51
1:A:410:ARG:HG3	1:A:411:LEU:H	1.75	0.51
1:B:221:VAL:HG11	1:B:238:LEU:HD13	1.91	0.51
1:A:236:ARG:CD	4:A:523:HOH:O	2.58	0.50
1:A:398:GLU:HG3	1:A:421:PRO:HD3	1.94	0.50
1:A:217:VAL:H	3:A:477:GOL:C1	2.22	0.50
1:B:39:THR:CB	4:B:500:HOH:O	2.56	0.50
1:A:66:LEU:O	1:A:70:ARG:HG2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:308:VAL:CG1	1:B:310:HIS:CE1	2.95	0.50
1:B:413:ARG:O	1:B:448:MET:HE2	2.12	0.49
1:B:417:LEU:HB2	1:B:443:LEU:HD23	1.95	0.49
1:B:69:LEU:O	1:B:73:GLU:HG2	2.12	0.49
1:B:235:ALA:HB1	1:B:246:LEU:HD13	1.93	0.49
1:A:124:ARG:H	1:A:140:GLN:NE2	2.10	0.49
1:A:404:ASP:CB	1:A:407:VAL:HG22	2.36	0.49
1:B:135:LEU:O	1:B:380:HIS:HD2	1.95	0.49
1:A:236:ARG:HD2	1:A:265:GLU:OE1	2.12	0.49
1:B:378:ALA:O	1:B:382:ARG:HG2	2.12	0.49
1:B:41:VAL:CG2	1:B:164:LEU:CD1	2.90	0.48
1:A:216:PRO:HA	3:A:477:GOL:H11	1.97	0.47
1:B:100:LEU:CD1	1:B:143:LEU:HD13	2.45	0.47
1:B:213:LYS:HD2	1:B:435:ASP:HB2	1.96	0.47
1:A:286:LEU:CD2	1:A:340:LEU:HD12	2.45	0.46
1:B:341:THR:HB	1:B:346:LEU:CD2	2.45	0.46
1:A:404:ASP:HB3	1:A:407:VAL:CG2	2.39	0.46
1:B:237:ARG:CD	1:B:425:LEU:HD13	2.46	0.46
1:A:39:THR:CB	4:A:731:HOH:O	2.55	0.46
1:A:138:PRO:HB2	1:A:373:TYR:CE1	2.50	0.46
1:B:37:TRP:HE3	1:B:164:LEU:HD12	1.81	0.46
1:B:262:LEU:HD22	1:B:266:LEU:CD1	2.46	0.46
1:B:308:VAL:HG13	1:B:310:HIS:CE1	2.50	0.46
1:A:236:ARG:HD3	4:A:523:HOH:O	2.16	0.46
1:B:404:ASP:C	1:B:406:GLU:H	2.23	0.45
1:A:294:PRO:HG2	1:A:297:ALA:HB3	1.98	0.45
1:A:393:TRP:HB2	2:A:476:NDP:C5N	2.46	0.45
1:A:389:SER:HA	1:A:437:THR:O	2.15	0.45
1:A:125:GLY:HA2	4:A:552:HOH:O	2.16	0.45
1:B:121:CYS:SG	1:B:160:GLY:O	2.75	0.45
1:A:124:ARG:H	1:A:140:GLN:HE22	1.63	0.44
1:B:41:VAL:HG21	1:B:164:LEU:CD1	2.41	0.44
1:B:396:TRP:HH2	1:B:425:LEU:HD21	1.83	0.44
1:B:143:LEU:HD13	1:B:143:LEU:HA	1.89	0.43
1:B:466:PRO:C	1:B:469:VAL:HG22	2.43	0.43
1:B:339:GLU:OE1	4:B:497:HOH:O	2.22	0.43
1:A:41:VAL:HG21	1:A:164:LEU:HD11	2.01	0.43
1:A:135:LEU:O	1:A:380:HIS:CD2	2.68	0.42
1:A:233:ARG:HD2	1:A:233:ARG:HA	1.84	0.42
1:B:125:GLY:HA2	4:B:551:HOH:O	2.20	0.42
1:B:201:LEU:CB	1:B:461:LEU:CD1	2.95	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:398:GLU:HG3	1:A:421:PRO:HD2	2.02	0.42
1:A:290:LEU:HD22	1:A:299:LEU:HD22	2.01	0.42
1:B:466:PRO:CA	1:B:469:VAL:HG22	2.50	0.42
1:A:128:ALA:HB2	1:A:135:LEU:HB2	2.02	0.41
1:B:308:VAL:HG11	1:B:310:HIS:CE1	2.54	0.41
1:B:105:LEU:HD13	1:B:105:LEU:HA	1.82	0.41
1:B:406:GLU:O	1:B:410:ARG:HG2	2.19	0.41
1:B:59:ARG:CB	1:B:105:LEU:HD12	2.40	0.41
1:B:213:LYS:HD2	4:B:542:HOH:O	2.21	0.41
1:A:341:THR:CA	1:A:344:LEU:HD13	2.40	0.41
1:B:266:LEU:CD1	1:B:266:LEU:H	2.33	0.41
1:B:192:ARG:O	1:B:193:GLY:C	2.63	0.41
1:B:469:VAL:HG23	1:B:470:SER:N	2.36	0.41
1:A:87:GLU:HB3	1:A:139:VAL:HG21	2.03	0.40
1:A:364:HIS:N	1:A:365:PRO:CD	2.85	0.40
1:B:208:VAL:CG2	1:B:211:SER:OG	2.53	0.40
1:A:0:MET:SD	1:A:426:GLY:HA3	2.61	0.40
1:A:331:LEU:HA	1:A:373:TYR:CD2	2.55	0.40
1:B:417:LEU:HD11	1:B:445:ASP:HB2	2.03	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	474/496 (96%)	459 (97%)	13 (3%)	2 (0%)	30	14
1	B	473/496 (95%)	459 (97%)	12 (2%)	2 (0%)	30	14
All	All	947/992 (96%)	918 (97%)	25 (3%)	4 (0%)	30	14

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	17	THR
1	B	471	ALA
1	A	394	GLY
1	A	405	PRO

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	334/351 (95%)	320 (96%)	14 (4%)	26	7
1	B	333/351 (95%)	314 (94%)	19 (6%)	18	4
All	All	667/702 (95%)	634 (95%)	33 (5%)	22	5

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-2	SER
1	A	11	ARG
1	A	75	GLU
1	A	124	ARG
1	A	143	LEU
1	A	166	GLU
1	A	167	THR
1	A	171	ARG
1	A	262	LEU
1	A	311	ASP
1	A	343	ASP
1	A	398	GLU
1	A	457	ARG
1	A	461	LEU
1	B	-1	HIS
1	B	17	THR
1	B	55	LYS
1	B	73	GLU
1	B	82	LEU
1	B	115	LEU

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Mol	Chain	Res	Type
1	B	131	ARG
1	B	143	LEU
1	B	166	GLU
1	B	213	LYS
1	B	262	LEU
1	B	265	GLU
1	B	295	GLU
1	B	346	LEU
1	B	410	ARG
1	B	411	LEU
1	B	421	PRO
1	B	434	ASN
1	B	461	LEU

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	68	GLN
1	A	77	HIS
1	A	90	HIS
1	A	140	GLN
1	A	203	HIS
1	A	245	HIS
1	A	304	HIS
1	A	380	HIS
1	A	414	GLN
1	A	434	ASN
1	A	456	ASN
1	B	32	HIS
1	B	77	HIS
1	B	241	GLN
1	B	245	HIS
1	B	336	HIS
1	B	364	HIS
1	B	380	HIS
1	B	408	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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](#)

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	GOL	B	477	-	5,5,5	0.57	0	5,5,5	0.87	0
3	GOL	A	477	-	5,5,5	0.51	0	5,5,5	0.49	0
2	NDP	B	476	-	51,52,52	1.99	9 (17%)	71,80,80	2.19	23 (32%)
2	NDP	A	476	-	51,52,52	1.74	9 (17%)	71,80,80	2.29	21 (29%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	477	-	-	0/4/4/4	-
3	GOL	A	477	-	-	4/4/4/4	-
2	NDP	B	476	-	1/1/14/17	9/34/77/77	0/5/5/5
2	NDP	A	476	-	1/1/14/17	7/34/77/77	0/5/5/5

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	476	NDP	PA-O3	7.55	1.67	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	476	NDP	C4N-C3N	-5.23	1.40	1.50
2	A	476	NDP	C4N-C3N	-5.22	1.40	1.50
2	A	476	NDP	PN-O3	-5.18	1.53	1.59
2	A	476	NDP	PA-O3	4.65	1.64	1.59
2	B	476	NDP	C8A-N7A	4.17	1.39	1.31
2	B	476	NDP	C5A-N7A	-3.99	1.31	1.39
2	B	476	NDP	C4N-C5N	-3.66	1.39	1.49
2	B	476	NDP	PN-O3	-3.65	1.55	1.59
2	B	476	NDP	O4D-C1D	3.57	1.50	1.42
2	A	476	NDP	C4N-C5N	-3.18	1.40	1.49
2	A	476	NDP	C8A-N7A	3.05	1.37	1.31
2	A	476	NDP	C5A-C4A	-2.65	1.34	1.39
2	B	476	NDP	C1D-N1N	2.47	1.53	1.46
2	A	476	NDP	C4A-N3A	2.26	1.38	1.34
2	A	476	NDP	C2D-C3D	-2.07	1.47	1.53
2	B	476	NDP	O5B-C5B	-2.03	1.36	1.44
2	A	476	NDP	O3B-C3B	2.00	1.47	1.43

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	476	NDP	N3A-C2A-N1A	-7.00	117.98	128.58
2	A	476	NDP	O4B-C4B-C3B	6.25	117.55	105.15
2	B	476	NDP	N9A-C8A-N7A	-6.24	105.08	113.94
2	B	476	NDP	N3A-C2A-N1A	-6.04	119.44	128.58
2	A	476	NDP	N9A-C8A-N7A	-5.76	105.77	113.94
2	A	476	NDP	C2B-C3B-C4B	-5.43	90.31	101.99
2	A	476	NDP	C4B-O4B-C1B	-5.00	98.43	109.47
2	B	476	NDP	C2B-C3B-C4B	-4.94	91.38	101.99
2	B	476	NDP	O4B-C4B-C3B	4.88	114.83	105.15
2	B	476	NDP	C2A-N1A-C6A	4.80	126.62	118.73
2	B	476	NDP	C4B-O4B-C1B	-3.85	100.95	109.47
2	B	476	NDP	C5A-N7A-C8A	3.76	109.36	103.45
2	A	476	NDP	C4A-N9A-C8A	3.56	109.48	105.74
2	B	476	NDP	C5B-C4B-C3B	3.42	127.52	115.21
2	B	476	NDP	O2B-P2B-O1X	-3.41	97.18	109.33
2	A	476	NDP	C5A-N7A-C8A	3.39	108.78	103.45
2	A	476	NDP	C5A-C4A-N3A	-3.29	122.18	126.72
2	A	476	NDP	O4B-C4B-C5B	3.27	119.81	109.33
2	B	476	NDP	O7N-C7N-C3N	-3.22	114.83	120.90
2	A	476	NDP	C6N-N1N-C2N	3.19	122.74	119.32
2	A	476	NDP	C2A-N1A-C6A	3.18	123.95	118.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	476	NDP	O2B-P2B-O1X	-3.11	98.25	109.33
2	B	476	NDP	C4A-N9A-C8A	3.04	108.93	105.74
2	A	476	NDP	C2A-N3A-C4A	2.98	119.12	111.83
2	A	476	NDP	O7N-C7N-C3N	-2.98	115.28	120.90
2	A	476	NDP	O2A-PA-O3	2.98	115.33	107.27
2	A	476	NDP	C3N-C2N-N1N	-2.77	119.14	123.20
2	B	476	NDP	O2B-C2B-C1B	2.72	119.60	110.05
2	B	476	NDP	C3N-C2N-N1N	-2.63	119.35	123.20
2	B	476	NDP	O4B-C4B-C5B	2.58	117.61	109.33
2	B	476	NDP	C2D-C3D-C4D	2.53	107.50	102.61
2	B	476	NDP	O3-PN-O1N	-2.49	103.21	110.70
2	B	476	NDP	C5A-C4A-N3A	-2.42	123.38	126.72
2	B	476	NDP	O2N-PN-O3	2.30	113.50	107.27
2	B	476	NDP	C4A-C5A-N7A	-2.22	108.04	110.58
2	A	476	NDP	O4D-C1D-C2D	-2.21	101.88	106.62
2	A	476	NDP	O3X-P2B-O2X	2.21	116.09	107.80
2	B	476	NDP	O3D-C3D-C2D	-2.21	104.75	111.82
2	A	476	NDP	O2D-C2D-C3D	2.19	118.84	111.82
2	B	476	NDP	C2A-N3A-C4A	2.10	116.95	111.83
2	A	476	NDP	C2D-C3D-C4D	2.07	106.60	102.61
2	A	476	NDP	C4D-O4D-C1D	2.05	113.99	109.47
2	B	476	NDP	O2D-C2D-C3D	2.05	118.38	111.82
2	B	476	NDP	O3X-P2B-O2X	2.03	115.42	107.80

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	476	NDP	C4B
2	B	476	NDP	C4B

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	476	NDP	C5D-O5D-PN-O3
2	A	476	NDP	C5D-O5D-PN-O1N
2	A	476	NDP	C5D-O5D-PN-O2N
2	B	476	NDP	C5D-O5D-PN-O3
2	B	476	NDP	C5D-O5D-PN-O1N
2	B	476	NDP	C5D-O5D-PN-O2N
3	A	477	GOL	O1-C1-C2-C3
3	A	477	GOL	C1-C2-C3-O3
3	A	477	GOL	O2-C2-C3-O3

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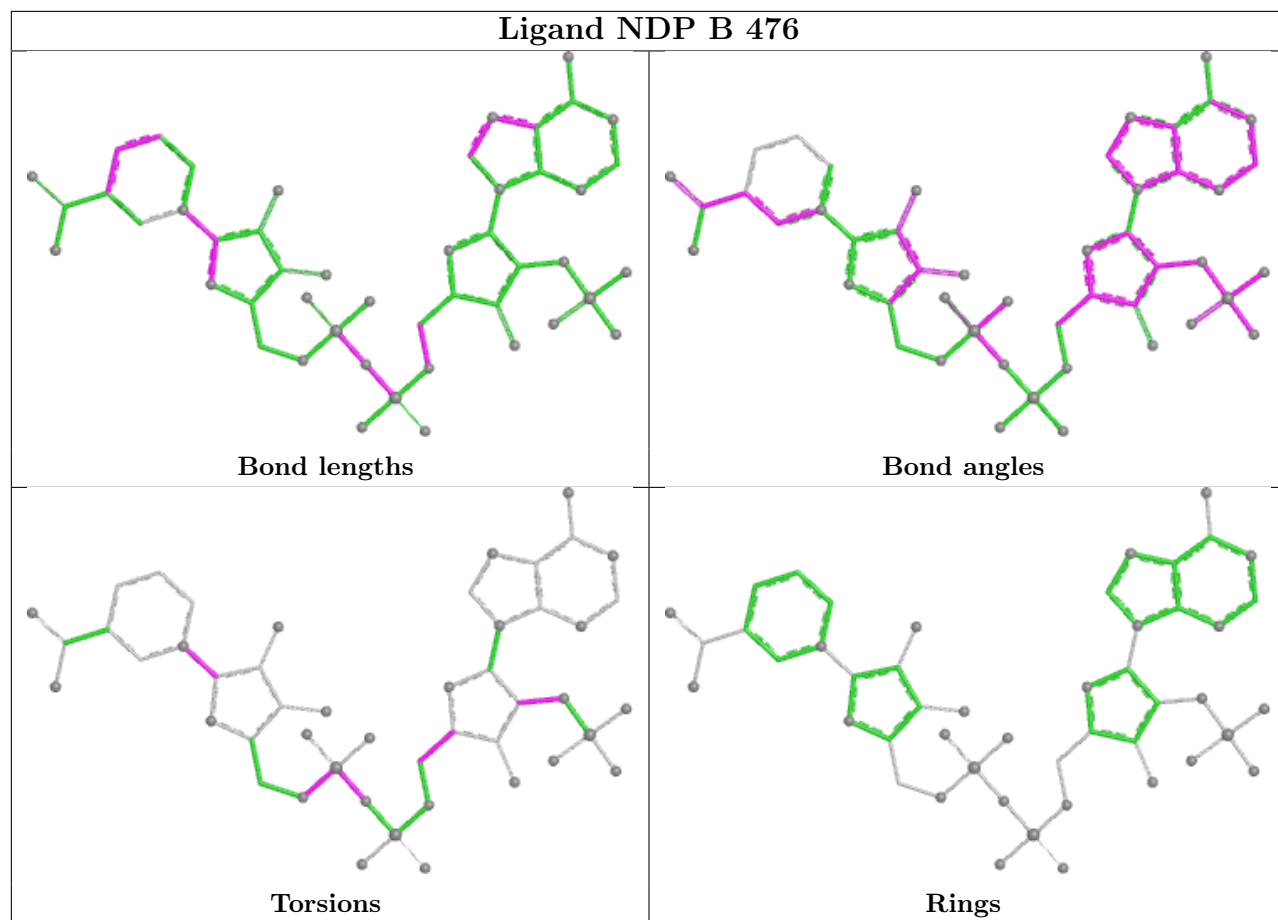
Mol	Chain	Res	Type	Atoms
3	A	477	GOL	O1-C1-C2-O2
2	A	476	NDP	O4B-C4B-C5B-O5B
2	B	476	NDP	O4B-C4B-C5B-O5B
2	B	476	NDP	C3B-C4B-C5B-O5B
2	A	476	NDP	O4D-C1D-N1N-C6N
2	B	476	NDP	O4D-C1D-N1N-C6N
2	B	476	NDP	C1B-C2B-O2B-P2B
2	A	476	NDP	PA-O3-PN-O1N
2	B	476	NDP	PA-O3-PN-O2N
2	A	476	NDP	PA-O3-PN-O2N
2	B	476	NDP	PA-O3-PN-O1N

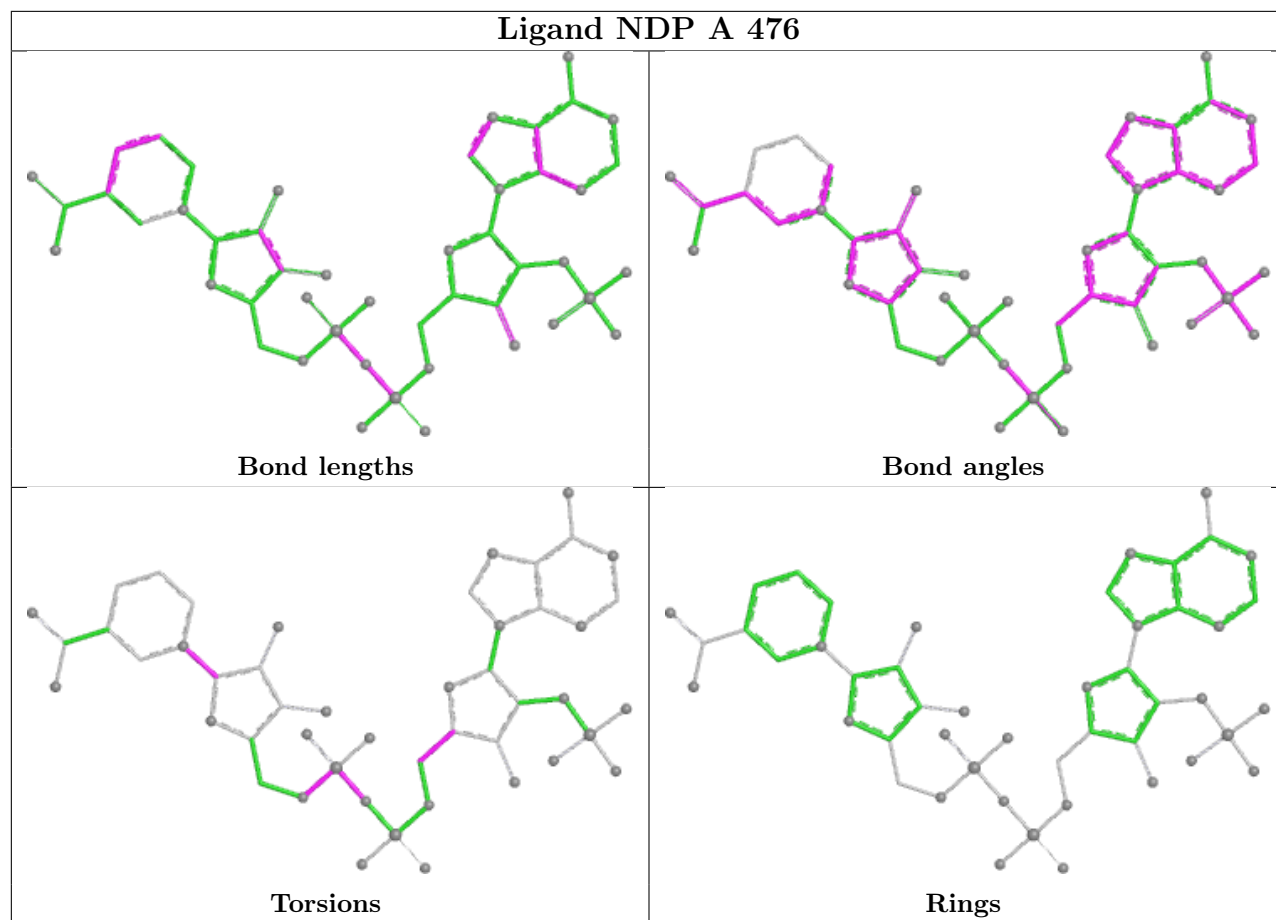
There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	477	GOL	4	0
2	B	476	NDP	1	0
2	A	476	NDP	1	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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	476/496 (95%)	0.88	53 (11%)	10 10	9, 17, 34, 60	0
1	B	475/496 (95%)	0.80	46 (9%)	13 13	8, 16, 34, 61	0
All	All	951/992 (95%)	0.84	99 (10%)	11 11	8, 17, 34, 61	0

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	402	ALA	9.2
1	B	471	ALA	8.9
1	A	411	LEU	7.8
1	B	412	VAL	6.7
1	A	402	ALA	6.5
1	A	412	VAL	6.1
1	B	17	THR	5.9
1	B	411	LEU	5.7
1	A	209	PRO	5.2
1	A	407	VAL	5.1
1	A	405	PRO	5.1
1	B	405	PRO	5.1
1	B	403	THR	4.6
1	A	-1	HIS	4.5
1	A	208	VAL	4.4
1	B	143	LEU	4.3
1	A	403	THR	4.0
1	A	413	ARG	3.8
1	A	210	GLY	3.5
1	A	410	ARG	3.5
1	B	410	ARG	3.5
1	B	407	VAL	3.4
1	A	278	CYS	3.4
1	A	414	GLN	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	434	ASN	3.3
1	A	295	GLU	3.2
1	B	19	ARG	3.2
1	A	70	ARG	3.2
1	A	143	LEU	3.1
1	B	131	ARG	3.1
1	A	21	ALA	3.0
1	B	413	ARG	3.0
1	A	213	LYS	3.0
1	B	409	ASP	2.9
1	A	343	ASP	2.9
1	B	261	GLU	2.8
1	A	212	GLY	2.8
1	B	71	GLU	2.8
1	A	48	ASP	2.8
1	A	406	GLU	2.7
1	B	-1	HIS	2.7
1	B	428	LEU	2.7
1	B	33	ALA	2.7
1	A	404	ASP	2.7
1	A	167	THR	2.7
1	B	121	CYS	2.7
1	B	208	VAL	2.7
1	A	394	GLY	2.7
1	A	260	ALA	2.7
1	B	48	ASP	2.6
1	A	211	SER	2.6
1	A	294	PRO	2.6
1	A	289	LEU	2.6
1	B	164	LEU	2.6
1	A	344	LEU	2.5
1	B	308	VAL	2.5
1	A	456	ASN	2.5
1	B	72	ASP	2.5
1	B	239	ALA	2.5
1	A	11	ARG	2.5
1	A	261	GLU	2.4
1	A	262	LEU	2.4
1	B	408	HIS	2.4
1	B	55	LYS	2.4
1	A	398	GLU	2.3
1	A	273	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	295	GLU	2.3
1	B	343	ASP	2.3
1	B	170	THR	2.3
1	A	33	ALA	2.3
1	A	329	ALA	2.3
1	A	164	LEU	2.3
1	B	69	LEU	2.3
1	A	255	ASP	2.3
1	A	19	ARG	2.3
1	B	130	GLY	2.3
1	A	24	LEU	2.2
1	B	105	LEU	2.2
1	A	296	ASP	2.2
1	A	400	GLY	2.2
1	A	401	MET	2.2
1	A	131	ARG	2.2
1	B	173	ALA	2.2
1	B	465	VAL	2.1
1	B	394	GLY	2.1
1	A	417	LEU	2.1
1	A	236	ARG	2.1
1	A	67	ALA	2.1
1	A	207	ALA	2.1
1	B	166	GLU	2.1
1	B	12	VAL	2.1
1	B	278	CYS	2.1
1	B	58	ASP	2.0
1	A	284	GLU	2.0
1	B	75	GLU	2.0
1	B	456	ASN	2.0
1	A	372	ALA	2.0
1	B	260	ALA	2.0
1	B	264	ALA	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](#)

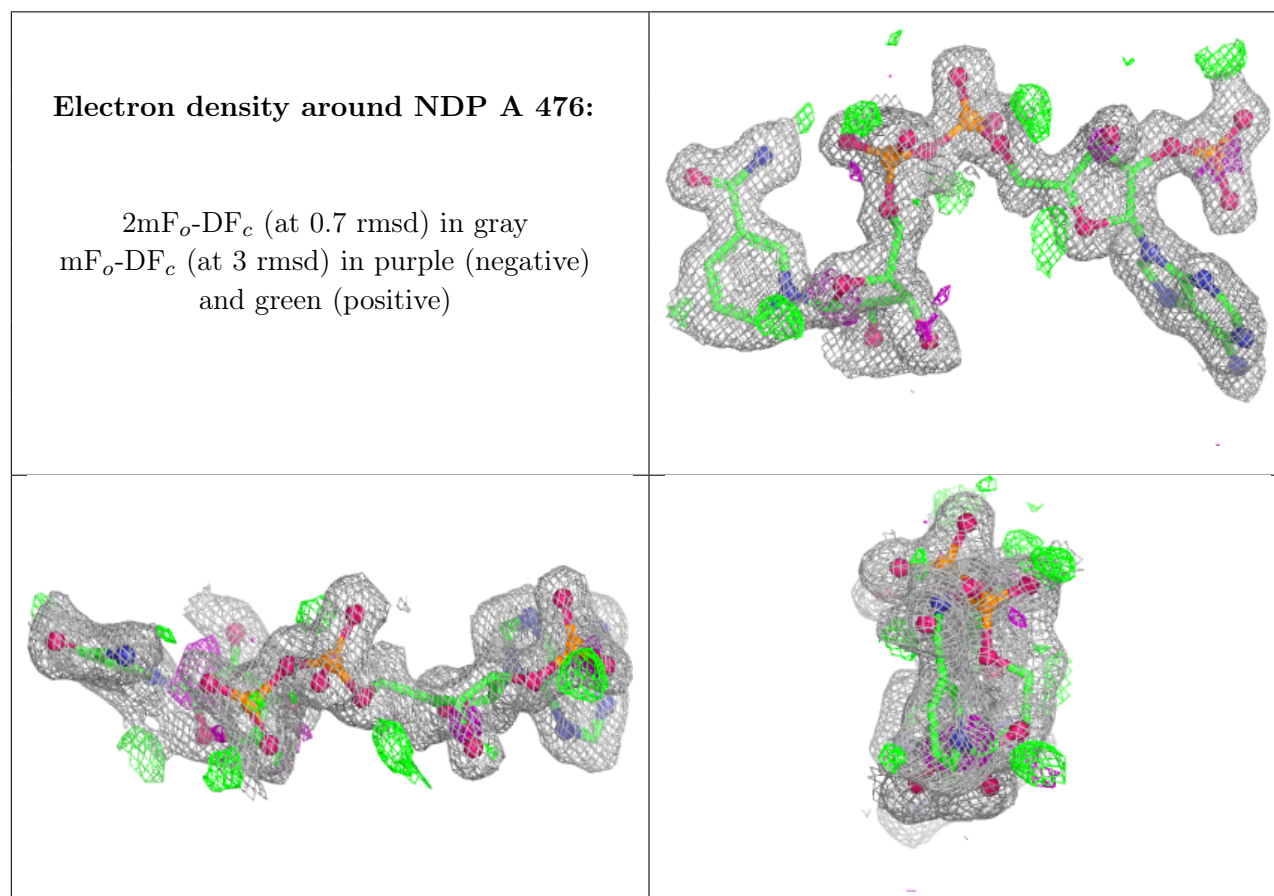
There are no oligosaccharides in this entry.

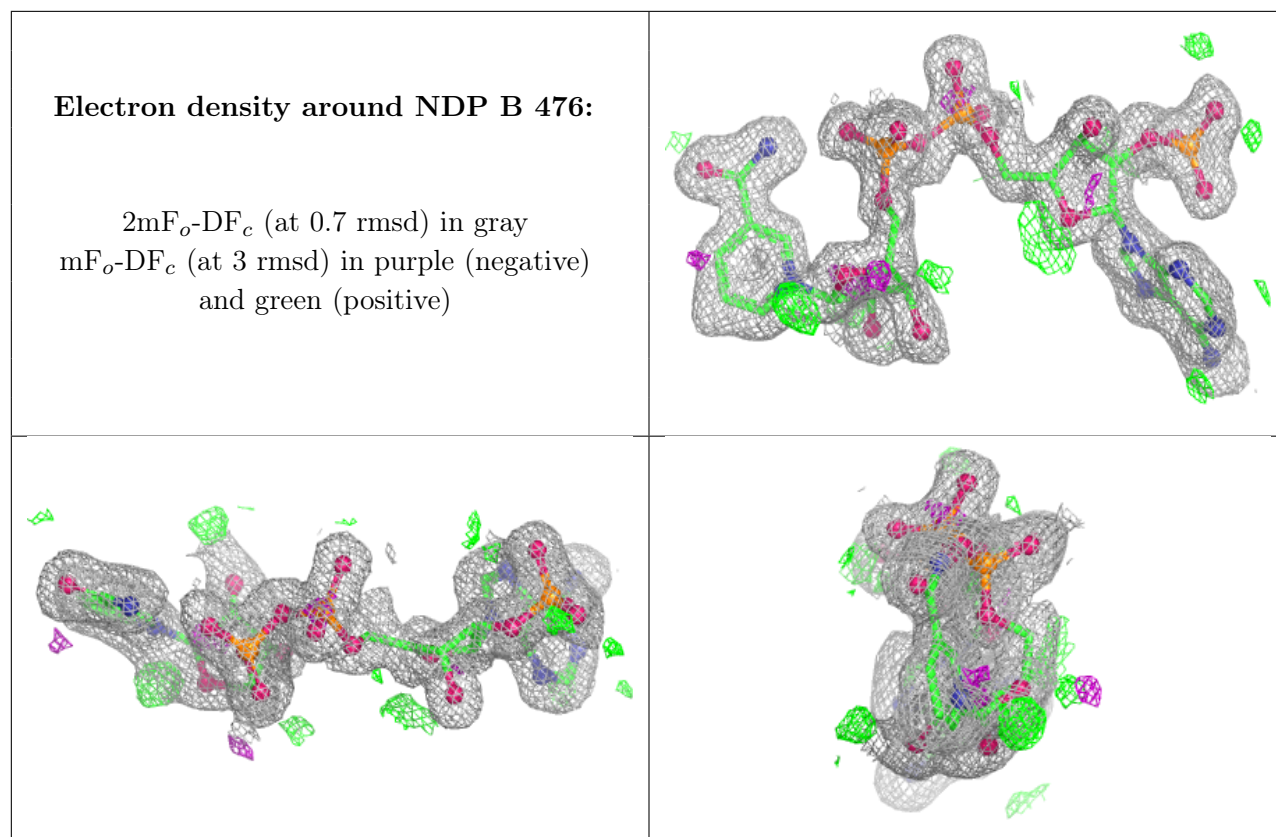
6.4 Ligands [i](#)

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
3	GOL	A	477	6/6	0.82	0.15	32,33,38,40	0
3	GOL	B	477	6/6	0.86	0.11	23,24,28,29	0
2	NDP	A	476	48/48	0.93	0.09	12,18,22,22	0
2	NDP	B	476	48/48	0.95	0.08	11,15,19,20	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.