



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:08 AM UTC

PDB ID : 8ICU / pdb_00008icu
Title : DNA POLYMERASE BETA (POL B) (E.C.2.7.7.7) COMPLEXED WITH SEVEN BASE PAIRS OF DNA; SOAKED IN THE PRESENCE OF DDATP (1 MILLIMOLAR) AND MNCL2 (5 MILLIMOLAR)
Authors : Pelletier, H.; Sawaya, M.R.
Deposited on : 1995-12-15
Resolution : 3.00 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
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

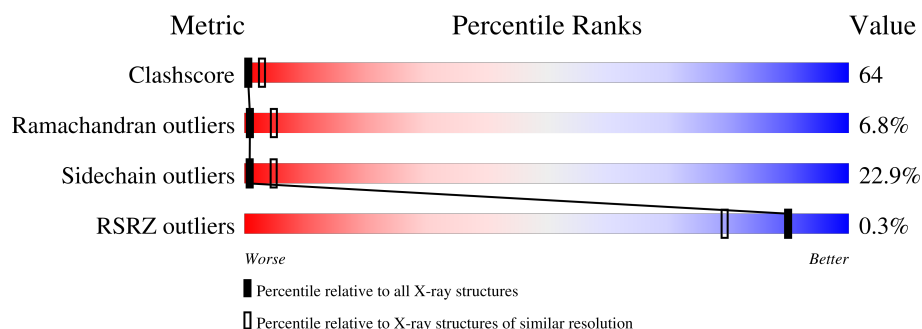
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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))
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	T	8	25% 25% 50%
2	P	8	12% 12% 88%
3	A	335	23% 41% 27% 8% .

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 3084 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 DNA chain called DNA (5'-D(*CP*AP*TP*TP*AP*GP*AP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	T	8	Total	C	N	O	P	0	0	0
			145	69	27	42	7			

- Molecule 2 is a DNA chain called DNA (5'-D(*TP*CP*TP*AP*AP*TP*GP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	8	Total	C	N	O	P	0	0	0
			164	79	29	48	8			

- Molecule 3 is a protein called PROTEIN (DNA POLYMERASE BETA (E.C.2.7.7.7)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	327	Total	C	N	O	S	18	0	0
			2623	1657	458	499	9			

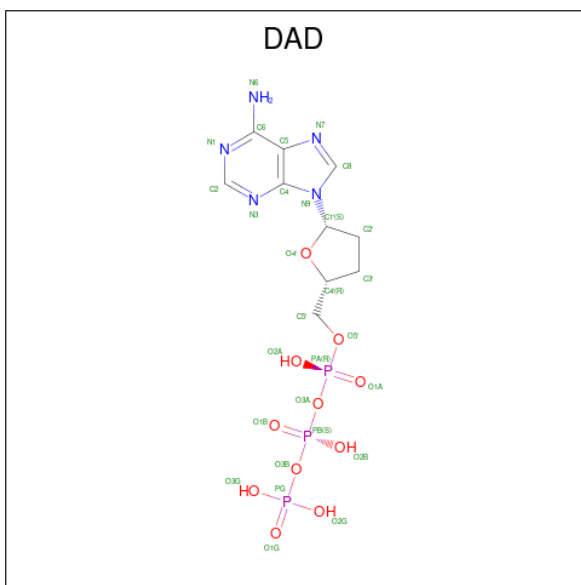
- Molecule 4 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mn	0	0
			2	2		

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Na	0	0
			2	2		

- Molecule 6 is 2',3'-DIDEOXYADENOSINE-5'-TRIPHOSPHATE (CCD ID: DAD) (formula: C₁₀H₁₆N₅O₁₁P₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	P	0	0
			13	10	3		

- Molecule 7 is water.

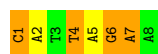
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	T	11	Total O 11 11	0	0
7	P	21	Total O 21 21	0	0
7	A	103	Total O 103 103	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

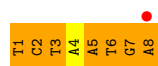
- Molecule 1: DNA (5'-D(*CP*AP*TP*TP*AP*GP*AP*A)-3')

Chain T: 

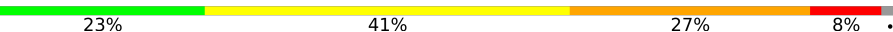


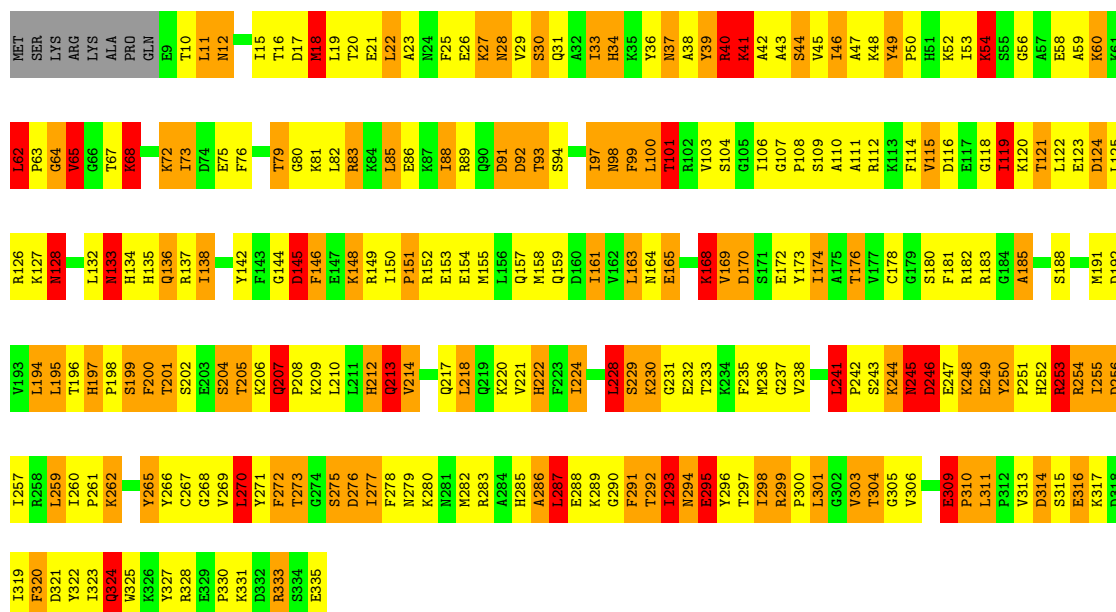
- Molecule 2: DNA (5'-D(*TP*CP*TP*AP*AP*TP*GP*A)-3')

Chain P: 



- Molecule 3: PROTEIN (DNA POLYMERASE BETA (E.C.2.7.7.7))

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	179.67Å 57.59Å 48.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.00 20.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	90.0 (20.00-3.00) 89.7 (20.00-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 2.70Å)	Xtriage
Refinement program	TNT 5-D	Depositor
R, R_{free}	0.175 , (Not available) (Not available) , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	36.8	Xtriage
Anisotropy	0.207	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.17 , 159.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3084	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.26% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NA, MN, DAD

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	T	0.99	0/162	4.08	14/249 (5.6%)
2	P	1.10	1/183 (0.5%)	4.23	19/278 (6.8%)
3	A	1.59	12/2672 (0.4%)	2.20	128/3590 (3.6%)
All	All	1.54	13/3017 (0.4%)	2.54	161/4117 (3.9%)

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
3	A	2	0

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	P	1	DT	OP3-P	6.78	1.62	1.48
3	A	138	ILE	CA-CB	-6.40	1.46	1.54
3	A	270	LEU	C-N	-6.33	1.26	1.33
3	A	165	GLU	N-CA	-6.06	1.39	1.46
3	A	213	GLN	CA-C	-5.74	1.45	1.52
3	A	97	ILE	CA-CB	-5.53	1.47	1.54
3	A	269	VAL	C-O	5.47	1.30	1.24
3	A	170	ASP	CA-C	-5.29	1.46	1.52
3	A	110	ALA	N-CA	-5.16	1.39	1.46
3	A	146	PHE	CA-C	-5.15	1.45	1.52
3	A	115	VAL	CA-C	-5.12	1.46	1.52
3	A	17	ASP	CG-OD1	5.02	1.34	1.25
3	A	63	PRO	C-O	5.00	1.29	1.23

All (161) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	T	7	DA	C4-N9-C1'	-33.68	76.53	127.05
1	T	7	DA	C8-N9-C1'	32.51	175.81	127.05
2	P	1	DT	C6-N1-C1'	-24.68	82.33	119.35
2	P	1	DT	C2-N1-C1'	24.43	155.99	119.35
2	P	6	DT	C2-N1-C1'	22.64	153.30	119.35
2	P	6	DT	C6-N1-C1'	-22.19	86.07	119.35
1	T	4	DT	C6-N1-C1'	-19.17	90.59	119.35
1	T	4	DT	C2-N1-C1'	18.36	146.88	119.35
2	P	3	DT	C2-N1-C1'	15.71	142.91	119.35
2	P	2	DC	C2-N1-C1'	14.95	142.12	119.70
2	P	3	DT	C6-N1-C1'	-14.85	97.08	119.35
2	P	7	DG	C4-N9-C1'	-14.52	105.22	127.00
1	T	6	DG	C4-N9-C1'	-14.47	105.29	127.00
2	P	8	DA	C8-N9-C1'	14.47	148.76	127.05
2	P	8	DA	C4-N9-C1'	-14.18	105.79	127.05
2	P	2	DC	C6-N1-C1'	-13.96	98.76	119.70
1	T	6	DG	C8-N9-C1'	13.56	147.34	127.00
2	P	7	DG	C8-N9-C1'	13.39	147.08	127.00
3	A	214	VAL	N-CA-CB	11.42	123.07	110.62
2	P	8	DA	P-O5'-C5'	-11.40	102.89	120.00
3	A	303	VAL	N-CA-C	11.04	123.20	110.62
2	P	5	DA	C8-N9-C1'	-9.80	112.34	127.05
2	P	5	DA	C4-N9-C1'	9.48	141.27	127.05
3	A	49	TYR	CA-C-N	9.13	131.25	119.84
3	A	49	TYR	C-N-CA	9.13	131.25	119.84
3	A	161	ILE	N-CA-C	9.06	119.05	110.53
3	A	49	TYR	O-C-N	8.77	128.57	121.38
3	A	168	LYS	N-CA-CB	8.65	123.04	110.06
1	T	1	DC	P-O3'-C3'	8.51	132.96	120.20
3	A	311	LEU	CA-C-N	8.47	128.80	119.90
3	A	311	LEU	C-N-CA	8.47	128.80	119.90
3	A	309	GLU	CA-C-N	8.43	130.38	119.84
3	A	309	GLU	C-N-CA	8.43	130.38	119.84
3	A	157	GLN	N-CA-CB	8.36	122.19	110.07
3	A	17	ASP	N-CA-C	8.36	120.01	111.07
3	A	266	TYR	CA-CB-CG	-8.27	99.02	113.90
2	P	3	DT	O4'-C1'-N1	7.99	120.38	108.40
3	A	146	PHE	N-CA-CB	7.85	121.96	110.26
3	A	286	ALA	CA-C-N	7.67	130.41	120.44
3	A	286	ALA	C-N-CA	7.67	130.41	120.44
2	P	3	DT	N1-C1'-C2'	7.64	124.95	113.50
3	A	41	LYS	N-CA-CB	7.56	120.97	110.01
3	A	83	ARG	N-CA-CB	7.42	120.74	109.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	151	PRO	N-CA-CB	7.39	109.84	103.19
3	A	40	ARG	N-CA-CB	7.23	120.97	110.20
3	A	272	PHE	CA-CB-CG	-7.16	106.64	113.80
3	A	235	PHE	N-CA-C	-7.04	99.51	110.42
3	A	73	ILE	N-CA-CB	6.97	117.94	110.62
3	A	157	GLN	CB-CA-C	6.96	121.90	110.90
3	A	64	GLY	CA-C-N	-6.92	113.86	122.93
3	A	64	GLY	C-N-CA	-6.92	113.86	122.93
3	A	291	PHE	N-CA-C	6.92	117.17	108.45
3	A	98	ASN	CA-CB-CG	-6.81	105.79	112.60
3	A	88	ILE	CB-CA-C	-6.80	103.31	111.81
3	A	116	ASP	N-CA-CB	6.80	119.83	109.91
3	A	214	VAL	N-CA-C	6.70	116.80	110.30
3	A	145	ASP	CA-CB-CG	6.65	119.25	112.60
3	A	320	PHE	CA-CB-CG	-6.58	107.22	113.80
1	T	6	DG	N9-C1'-C2'	-6.54	103.70	113.50
3	A	91	ASP	CA-CB-CG	-6.52	106.08	112.60
3	A	292	THR	CB-CA-C	6.49	120.84	109.80
3	A	17	ASP	N-CA-CB	6.48	119.41	110.01
3	A	134	HIS	CA-CB-CG	6.48	120.28	113.80
3	A	99	PHE	N-CA-C	6.42	117.97	110.97
1	T	4	DT	P-O3'-C3'	-6.39	110.61	120.20
3	A	256	ASP	CA-CB-CG	6.34	118.94	112.60
3	A	41	LYS	CB-CA-C	6.32	120.81	110.88
3	A	253	ARG	N-CA-C	6.32	119.80	109.24
3	A	116	ASP	CB-CA-C	6.28	120.62	110.96
3	A	200	PHE	N-CA-C	6.28	118.75	108.52
3	A	93	THR	N-CA-CB	6.25	119.04	109.98
3	A	195	LEU	N-CA-CB	-6.24	100.01	111.37
3	A	159	GLN	CB-CA-C	-6.23	100.08	110.68
3	A	241	LEU	CA-C-N	6.21	127.61	119.84
3	A	241	LEU	C-N-CA	6.21	127.61	119.84
3	A	207	GLN	CA-C-N	6.19	127.58	119.84
3	A	207	GLN	C-N-CA	6.19	127.58	119.84
3	A	314	ASP	N-CA-CB	6.18	119.61	110.53
3	A	207	GLN	C-N-CD	-6.14	99.84	125.00
3	A	30	SER	CA-C-N	-6.08	112.56	122.65
3	A	30	SER	C-N-CA	-6.08	112.56	122.65
3	A	54	LYS	N-CA-C	6.05	121.40	113.30
3	A	241	LEU	O-C-N	6.02	127.80	121.42
3	A	192	ASP	CB-CA-C	-6.00	102.93	111.76
3	A	228	LEU	N-CA-C	-5.95	104.17	112.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	119	ILE	CB-CA-C	-5.93	105.92	112.68
3	A	45	VAL	CA-C-O	-5.89	114.35	121.18
3	A	145	ASP	CB-CA-C	-5.89	98.69	110.42
3	A	250	TYR	N-CA-C	5.89	117.49	110.07
3	A	124	ASP	N-CA-C	-5.87	104.79	111.07
1	T	6	DG	O4'-C1'-N9	5.82	117.13	108.40
3	A	28	ASN	CA-CB-CG	-5.80	106.80	112.60
3	A	213	GLN	CB-CA-C	-5.79	101.55	110.81
3	A	164	ASN	CA-CB-CG	-5.75	106.85	112.60
1	T	2	DA	C8-N9-C1'	5.74	135.65	127.05
3	A	20	THR	CB-CA-C	5.73	121.68	110.67
3	A	68	LYS	N-CA-C	5.73	117.20	111.07
3	A	50	PRO	CB-CA-C	-5.72	102.12	111.56
3	A	249	GLU	N-CA-CB	5.72	119.73	110.41
3	A	101	THR	CB-CA-C	5.67	120.32	110.68
3	A	257	ILE	CB-CA-C	-5.65	102.08	110.33
3	A	18	MET	N-CA-C	-5.61	105.31	111.82
3	A	253	ARG	CA-C-O	5.60	127.10	120.66
3	A	148	LYS	N-CA-C	5.59	118.75	110.48
3	A	192	ASP	CA-C-N	-5.57	113.97	121.66
3	A	192	ASP	C-N-CA	-5.57	113.97	121.66
3	A	92	ASP	N-CA-CB	5.53	118.73	110.22
3	A	45	VAL	N-CA-C	5.52	115.61	110.42
1	T	2	DA	C4-N9-C1'	-5.52	118.77	127.05
3	A	333	ARG	N-CA-C	-5.52	106.13	112.92
3	A	293	ILE	CB-CA-C	-5.50	102.03	110.50
3	A	12	ASN	CB-CA-C	5.49	119.84	110.56
3	A	164	ASN	N-CA-C	5.44	117.02	111.14
3	A	197	HIS	CA-C-N	5.42	124.76	118.85
3	A	197	HIS	C-N-CA	5.42	124.76	118.85
3	A	212	HIS	N-CA-C	5.42	117.62	111.11
3	A	273	THR	N-CA-CB	5.40	118.65	110.28
3	A	254	ARG	CA-C-N	-5.40	114.93	122.71
3	A	254	ARG	C-N-CA	-5.40	114.93	122.71
3	A	195	LEU	N-CA-C	5.39	118.02	109.50
3	A	249	GLU	CB-CA-C	5.38	118.72	109.51
3	A	63	PRO	N-CA-C	-5.37	102.69	110.80
3	A	201	THR	N-CA-CB	-5.34	103.23	111.46
3	A	309	GLU	N-CA-C	5.34	121.62	109.81
3	A	34	HIS	CA-CB-CG	-5.34	108.46	113.80
3	A	65	VAL	CA-C-N	-5.32	117.70	122.43
3	A	65	VAL	C-N-CA	-5.32	117.70	122.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	45	VAL	N-CA-CB	5.31	116.65	110.65
3	A	205	THR	N-CA-C	5.31	122.10	110.80
3	A	58	GLU	O-C-N	5.28	127.58	122.09
3	A	133	ASN	OD1-CG-ND2	5.27	127.87	122.60
3	A	276	ASP	CA-CB-CG	5.26	117.86	112.60
3	A	49	TYR	CA-CB-CG	5.25	123.36	113.90
3	A	287	LEU	N-CA-C	5.24	116.68	111.07
3	A	163	LEU	O-C-N	5.24	127.68	122.12
3	A	115	VAL	CB-CA-C	-5.22	105.28	111.97
3	A	224	ILE	CB-CA-C	-5.21	104.66	110.96
3	A	11	LEU	CA-C-N	-5.18	114.36	122.65
3	A	11	LEU	C-N-CA	-5.18	114.36	122.65
3	A	62	LEU	N-CA-C	5.17	117.80	110.24
2	P	3	DT	O5'-C5'-C4'	-5.16	103.06	110.80
3	A	128	ASN	CA-CB-CG	-5.14	107.46	112.60
3	A	115	VAL	O-C-N	5.14	126.94	121.91
2	P	1	DT	O4'-C1'-N1	5.12	116.09	108.40
3	A	299	ARG	N-CA-C	5.12	117.59	109.40
3	A	168	LYS	CA-C-N	-5.08	114.52	120.72
3	A	168	LYS	C-N-CA	-5.08	114.52	120.72
3	A	246	ASP	N-CA-C	5.08	121.62	110.80
3	A	176	THR	CA-C-N	-5.07	116.04	122.43
3	A	176	THR	C-N-CA	-5.07	116.04	122.43
3	A	40	ARG	CA-C-N	-5.07	113.86	120.44
3	A	40	ARG	C-N-CA	-5.07	113.86	120.44
3	A	39	TYR	N-CA-CB	-5.06	101.93	110.49
3	A	313	VAL	N-CA-CB	-5.06	105.06	111.64
1	T	1	DC	C6-N1-C1'	-5.05	112.12	119.70
3	A	295	GLU	CB-CG-CD	5.04	121.17	112.60
3	A	97	ILE	CB-CA-C	-5.03	105.35	112.14
3	A	65	VAL	N-CA-CB	5.03	117.69	111.41
3	A	124	ASP	CA-CB-CG	5.02	117.62	112.60
1	T	4	DT	O4'-C1'-N1	5.02	115.93	108.40
3	A	245	ASN	CB-CA-C	5.01	119.66	112.05

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	157	GLN	CA
3	A	205	THR	CA

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	T	145	0	80	10	0
2	P	164	0	90	24	0
3	A	2623	0	2641	343	0
4	A	2	0	0	0	0
5	A	2	0	0	0	0
6	A	13	0	0	0	0
7	A	103	0	0	27	0
7	P	21	0	0	1	0
7	T	11	0	0	2	0
All	All	3084	0	2811	364	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 64.

All (364) 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)
2:P:6:DT:H2''	2:P:7:DG:H5''	1.30	1.11
3:A:18:MET:HE2	3:A:82:LEU:HD22	1.35	1.09
3:A:29:VAL:HG21	3:A:94:SER:HB2	1.32	1.08
2:P:5:DA:H2''	2:P:6:DT:H5''	1.39	1.04
2:P:1:DT:H2''	2:P:2:DC:H5'	1.39	1.04
3:A:245:ASN:N	3:A:245:ASN:HD22	1.54	1.00
3:A:150:ILE:HG21	3:A:158:MET:HE1	1.48	0.93
3:A:60:LYS:HA	3:A:65:VAL:HG23	1.52	0.91
3:A:165:GLU:HA	3:A:168:LYS:HG2	1.50	0.91
3:A:245:ASN:HD22	3:A:245:ASN:H	1.11	0.90
3:A:210:LEU:HB3	3:A:259:LEU:HD21	1.53	0.89
3:A:291:PHE:CD1	3:A:300:PRO:HA	2.09	0.88
3:A:31:GLN:NE2	3:A:112:ARG:HH12	1.72	0.88
3:A:114:PHE:HB3	3:A:119:ILE:HB	1.55	0.87
2:P:6:DT:C2'	2:P:7:DG:H5''	2.05	0.86
2:P:8:DA:H5''	3:A:272:PHE:HE2	1.38	0.86
3:A:18:MET:HG2	3:A:19:LEU:N	1.90	0.85
3:A:293:ILE:HD13	3:A:298:ILE:HG13	1.58	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:8:DA:H4'	3:A:272:PHE:CE2	2.14	0.83
3:A:31:GLN:HE21	3:A:112:ARG:HH12	1.26	0.83
3:A:103:VAL:HB	3:A:106:ILE:HD12	1.61	0.83
2:P:5:DA:H2''	2:P:6:DT:C5'	2.09	0.83
3:A:15:ILE:CG2	3:A:46:ILE:HD13	2.08	0.83
3:A:23:ALA:HB2	3:A:39:TYR:HB3	1.63	0.81
3:A:18:MET:HE2	3:A:82:LEU:CD2	2.11	0.81
3:A:23:ALA:HB2	3:A:39:TYR:CB	2.11	0.80
3:A:270:LEU:HD23	3:A:319:ILE:HD13	1.62	0.79
3:A:42:ALA:O	3:A:46:ILE:HG23	1.82	0.79
3:A:270:LEU:HD23	3:A:319:ILE:HG21	1.63	0.79
3:A:291:PHE:HD1	3:A:300:PRO:HA	1.48	0.79
3:A:128:ASN:N	3:A:128:ASN:ND2	2.30	0.78
3:A:180:SER:HB2	3:A:185:ALA:HB2	1.65	0.78
3:A:178:CYS:SG	3:A:194:LEU:HD22	2.23	0.78
3:A:27:LYS:HB3	3:A:36:TYR:CD1	2.19	0.78
3:A:128:ASN:N	3:A:128:ASN:HD22	1.81	0.78
3:A:29:VAL:CG2	3:A:94:SER:HB2	2.13	0.77
3:A:18:MET:HE1	3:A:76:PHE:HB2	1.66	0.77
3:A:19:LEU:HB3	3:A:43:ALA:HB2	1.63	0.77
1:T:4:DT:H5''	3:A:231:GLY:HA3	1.65	0.77
3:A:259:LEU:HD12	3:A:260:ILE:H	1.48	0.77
3:A:260:ILE:CG2	3:A:261:PRO:HD2	2.14	0.77
3:A:207:GLN:HB3	7:A:625:HOH:O	1.85	0.76
3:A:245:ASN:H	3:A:245:ASN:ND2	1.83	0.76
3:A:277:ILE:HD13	3:A:277:ILE:H	1.49	0.76
1:T:6:DG:H2''	1:T:7:DA:C8	2.20	0.76
3:A:123:GLU:O	3:A:127:LYS:HG2	1.86	0.75
3:A:181:PHE:HA	7:A:530:HOH:O	1.86	0.74
3:A:323:ILE:O	3:A:324:GLN:HG2	1.88	0.73
3:A:121:THR:HG23	3:A:124:ASP:OD1	1.88	0.73
3:A:243:SER:HB3	3:A:249:GLU:HA	1.69	0.73
3:A:114:PHE:CZ	3:A:132:LEU:HD23	2.24	0.72
2:P:1:DT:C2'	2:P:2:DC:H5'	2.17	0.71
3:A:197:HIS:CE1	3:A:199:SER:HB3	2.25	0.71
3:A:254:ARG:CZ	3:A:254:ARG:HB3	2.18	0.71
3:A:277:ILE:HG12	3:A:335:GLU:HA	1.71	0.70
3:A:18:MET:HE1	3:A:76:PHE:CB	2.21	0.70
3:A:294:ASN:HB2	3:A:295:GLU:OE1	1.92	0.70
3:A:18:MET:CE	3:A:82:LEU:HD22	2.20	0.70
2:P:8:DA:H5''	3:A:272:PHE:CE2	2.26	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:33:ILE:O	3:A:37:ASN:HB2	1.92	0.69
3:A:112:ARG:O	3:A:115:VAL:HB	1.92	0.69
3:A:60:LYS:HA	3:A:65:VAL:CG2	2.23	0.69
3:A:197:HIS:CD2	3:A:198:PRO:HD2	2.28	0.69
1:T:4:DT:H1'	7:T:527:HOH:O	1.93	0.69
3:A:11:LEU:HD23	3:A:11:LEU:H	1.56	0.69
3:A:270:LEU:CD2	3:A:319:ILE:HG21	2.22	0.69
3:A:142:TYR:CE2	3:A:238:VAL:HG11	2.28	0.69
3:A:180:SER:HB3	3:A:183:ARG:HH21	1.58	0.68
3:A:125:LEU:CD2	3:A:132:LEU:HD21	2.23	0.68
3:A:16:THR:HG21	3:A:47:ALA:HB2	1.76	0.68
3:A:122:LEU:HD21	3:A:126:ARG:CZ	2.25	0.67
2:P:5:DA:C2'	2:P:6:DT:H5''	2.20	0.67
3:A:46:ILE:HG13	3:A:47:ALA:N	2.05	0.67
2:P:8:DA:C5'	3:A:272:PHE:HE2	2.08	0.67
3:A:83:ARG:HA	3:A:86:GLU:HG2	1.75	0.67
3:A:182:ARG:NH2	3:A:316:GLU:OE1	2.28	0.66
3:A:68:LYS:HB2	3:A:68:LYS:NZ	2.11	0.65
3:A:11:LEU:H	3:A:11:LEU:CD2	2.10	0.65
3:A:271:TYR:HB2	7:A:592:HOH:O	1.94	0.65
3:A:233:THR:HB	7:A:538:HOH:O	1.95	0.65
2:P:6:DT:H2''	2:P:7:DG:C5'	2.19	0.65
3:A:41:LYS:HE2	3:A:64:GLY:HA3	1.79	0.65
3:A:182:ARG:NH1	3:A:273:THR:HG21	2.12	0.65
3:A:293:ILE:HD13	3:A:298:ILE:CG1	2.26	0.65
3:A:150:ILE:HG21	3:A:158:MET:CE	2.24	0.64
3:A:210:LEU:HB3	3:A:259:LEU:CD2	2.25	0.64
3:A:145:ASP:OD2	3:A:252:HIS:N	2.26	0.64
3:A:327:TYR:HD1	3:A:328:ARG:N	1.96	0.64
3:A:41:LYS:HE2	3:A:64:GLY:CA	2.28	0.64
3:A:180:SER:HA	3:A:183:ARG:HE	1.63	0.64
3:A:182:ARG:NH1	3:A:182:ARG:HG2	2.12	0.64
3:A:323:ILE:C	3:A:324:GLN:HG2	2.22	0.64
3:A:155:MET:HE2	3:A:188:SER:HB2	1.80	0.64
3:A:285:HIS:NE2	7:A:583:HOH:O	2.30	0.64
1:T:5:DA:H5'	7:T:527:HOH:O	1.98	0.64
3:A:218:LEU:HB2	3:A:224:ILE:HD12	1.78	0.64
3:A:15:ILE:HB	3:A:46:ILE:CD1	2.28	0.63
3:A:259:LEU:HD12	3:A:260:ILE:N	2.13	0.63
3:A:260:ILE:HG22	3:A:261:PRO:HD2	1.79	0.63
3:A:12:ASN:HB3	3:A:46:ILE:HD12	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:146:PHE:HZ	3:A:238:VAL:CG2	2.11	0.63
3:A:295:GLU:HA	7:A:592:HOH:O	1.97	0.63
3:A:204:SER:O	3:A:206:LYS:N	2.32	0.63
3:A:15:ILE:HB	3:A:46:ILE:HD13	1.80	0.62
3:A:15:ILE:CB	3:A:46:ILE:HD13	2.29	0.62
3:A:275:SER:OG	3:A:277:ILE:HD13	1.99	0.62
3:A:242:PRO:HG2	7:A:589:HOH:O	1.98	0.62
3:A:82:LEU:HB3	3:A:85:LEU:HB2	1.80	0.62
3:A:15:ILE:HD13	3:A:73:ILE:HD13	1.80	0.62
3:A:124:ASP:O	3:A:128:ASN:ND2	2.33	0.62
3:A:196:THR:OG1	3:A:197:HIS:N	2.32	0.62
3:A:289:LYS:HD3	3:A:289:LYS:N	2.14	0.62
3:A:23:ALA:HB2	3:A:39:TYR:HB2	1.82	0.61
3:A:11:LEU:HD23	3:A:11:LEU:N	2.16	0.61
3:A:182:ARG:HH12	3:A:273:THR:HG21	1.66	0.60
3:A:248:LYS:O	3:A:248:LYS:HG2	2.01	0.60
3:A:294:ASN:O	3:A:296:TYR:N	2.33	0.60
3:A:29:VAL:HA	3:A:97:ILE:HD12	1.82	0.60
3:A:260:ILE:HG23	3:A:261:PRO:HD2	1.83	0.60
3:A:209:LYS:HA	3:A:212:HIS:HB2	1.84	0.60
3:A:30:SER:HB3	7:A:561:HOH:O	2.02	0.60
3:A:155:MET:HE2	3:A:188:SER:CB	2.32	0.60
3:A:103:VAL:HA	7:A:600:HOH:O	2.02	0.60
3:A:18:MET:HG3	3:A:22:LEU:HD23	1.85	0.59
3:A:154:GLU:O	3:A:158:MET:HG3	2.01	0.59
3:A:125:LEU:HD22	3:A:132:LEU:HD21	1.84	0.59
3:A:286:ALA:O	3:A:291:PHE:HB2	2.02	0.59
3:A:210:LEU:CB	3:A:259:LEU:HD21	2.30	0.59
3:A:165:GLU:O	3:A:168:LYS:N	2.36	0.59
3:A:25:PHE:CG	3:A:88:ILE:HD13	2.37	0.59
3:A:122:LEU:HD21	3:A:126:ARG:NH2	2.17	0.58
3:A:148:LYS:HB2	7:A:620:HOH:O	2.03	0.58
3:A:79:THR:O	3:A:81:LYS:N	2.29	0.58
3:A:309:GLU:N	3:A:309:GLU:OE1	2.37	0.58
3:A:218:LEU:HB2	3:A:224:ILE:CD1	2.33	0.58
3:A:60:LYS:HG3	3:A:60:LYS:O	2.03	0.58
3:A:122:LEU:C	3:A:122:LEU:HD23	2.29	0.58
3:A:212:HIS:HB3	7:A:541:HOH:O	2.03	0.58
3:A:182:ARG:NH1	3:A:273:THR:OG1	2.37	0.58
3:A:111:ALA:O	3:A:115:VAL:HG23	2.04	0.57
3:A:31:GLN:HE21	3:A:112:ARG:NH1	2.01	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:165:GLU:HA	3:A:168:LYS:CG	2.30	0.57
3:A:169:VAL:HG22	3:A:170:ASP:N	2.20	0.56
3:A:213:GLN:HG3	7:A:644:HOH:O	2.04	0.56
3:A:15:ILE:HG22	3:A:46:ILE:HD13	1.87	0.56
3:A:68:LYS:O	3:A:72:LYS:HE3	2.06	0.56
3:A:328:ARG:O	3:A:333:ARG:NE	2.30	0.56
3:A:18:MET:O	3:A:21:GLU:HB2	2.05	0.56
3:A:241:LEU:O	3:A:250:TYR:HB2	2.06	0.56
3:A:128:ASN:HD22	3:A:128:ASN:H	1.55	0.55
3:A:183:ARG:NH1	3:A:275:SER:HA	2.20	0.55
3:A:182:ARG:HG2	3:A:182:ARG:HH11	1.69	0.55
3:A:127:LYS:HB2	3:A:128:ASN:ND2	2.22	0.55
3:A:245:ASN:N	3:A:245:ASN:ND2	2.30	0.55
3:A:292:THR:O	3:A:298:ILE:HA	2.07	0.55
3:A:188:SER:HB3	7:A:522:HOH:O	2.07	0.55
3:A:146:PHE:HZ	3:A:238:VAL:HG22	1.73	0.54
3:A:255:ILE:HG12	3:A:256:ASP:N	2.22	0.54
3:A:21:GLU:OE1	3:A:85:LEU:HG	2.08	0.54
3:A:200:PHE:CD2	3:A:261:PRO:HA	2.42	0.54
3:A:306:VAL:HG22	7:A:650:HOH:O	2.08	0.54
3:A:254:ARG:NH1	3:A:255:ILE:N	2.56	0.54
3:A:291:PHE:CE1	3:A:300:PRO:HA	2.43	0.54
3:A:18:MET:CE	3:A:82:LEU:HD13	2.38	0.54
3:A:317:LYS:HE3	3:A:321:ASP:OD1	2.07	0.54
3:A:270:LEU:C	3:A:270:LEU:HD12	2.32	0.53
3:A:292:THR:O	3:A:298:ILE:HG13	2.07	0.53
3:A:303:VAL:C	3:A:305:GLY:H	2.14	0.53
3:A:108:PRO:O	3:A:112:ARG:HG3	2.08	0.53
3:A:237:GLY:O	3:A:254:ARG:NH1	2.41	0.53
3:A:146:PHE:HB2	7:A:514:HOH:O	2.07	0.53
3:A:149:ARG:NH2	3:A:188:SER:HA	2.23	0.53
3:A:243:SER:O	3:A:244:LYS:O	2.27	0.53
3:A:291:PHE:O	3:A:301:LEU:HD22	2.09	0.53
3:A:169:VAL:HG22	3:A:173:TYR:HE2	1.74	0.53
3:A:103:VAL:CB	3:A:106:ILE:HD12	2.37	0.52
3:A:151:PRO:HG2	3:A:154:GLU:CD	2.35	0.52
3:A:173:TYR:OH	3:A:213:GLN:NE2	2.43	0.52
3:A:83:ARG:O	3:A:86:GLU:N	2.43	0.52
2:P:8:DA:C4'	3:A:272:PHE:CE2	2.91	0.51
3:A:244:LYS:HB3	3:A:245:ASN:HD22	1.75	0.51
3:A:100:LEU:HD21	3:A:119:ILE:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:145:ASP:OD2	3:A:252:HIS:HB2	2.10	0.51
3:A:146:PHE:CZ	3:A:238:VAL:HG22	2.46	0.51
3:A:268:GLY:O	3:A:271:TYR:HB3	2.09	0.51
3:A:330:PRO:HA	3:A:333:ARG:CG	2.40	0.51
2:P:5:DA:C2'	2:P:6:DT:H71	2.41	0.51
3:A:267:CYS:SG	3:A:297:THR:HA	2.49	0.51
3:A:322:TYR:C	3:A:324:GLN:H	2.18	0.51
3:A:48:LYS:HE3	7:A:607:HOH:O	2.10	0.51
3:A:319:ILE:O	3:A:322:TYR:HB2	2.11	0.51
3:A:259:LEU:O	3:A:260:ILE:HD13	2.11	0.51
3:A:295:GLU:OE1	3:A:295:GLU:N	2.44	0.50
1:T:4:DT:C5'	3:A:231:GLY:HA3	2.39	0.50
3:A:26:GLU:O	3:A:30:SER:O	2.29	0.50
3:A:38:ALA:O	3:A:41:LYS:HD3	2.11	0.50
3:A:183:ARG:HH11	3:A:275:SER:HA	1.76	0.50
3:A:254:ARG:HB3	3:A:254:ARG:NH1	2.26	0.50
3:A:41:LYS:O	3:A:44:SER:HB3	2.10	0.50
1:T:5:DA:H2''	1:T:6:DG:O5'	2.11	0.50
3:A:146:PHE:HZ	3:A:238:VAL:HG21	1.76	0.50
3:A:22:LEU:HD22	3:A:85:LEU:CD1	2.42	0.50
3:A:49:TYR:CD2	3:A:53:ILE:HG13	2.47	0.50
3:A:180:SER:HB2	3:A:185:ALA:CB	2.38	0.50
3:A:282:MET:HG2	7:A:555:HOH:O	2.11	0.50
3:A:228:LEU:O	3:A:229:SER:HB3	2.11	0.49
3:A:196:THR:HG22	3:A:265:TYR:CD1	2.48	0.49
3:A:197:HIS:ND1	3:A:199:SER:HB3	2.28	0.49
3:A:298:ILE:HG23	3:A:298:ILE:O	2.12	0.49
3:A:114:PHE:HZ	3:A:132:LEU:HD23	1.73	0.49
3:A:172:GLU:HB3	3:A:197:HIS:NE2	2.28	0.49
3:A:286:ALA:HA	3:A:323:ILE:CG2	2.43	0.49
2:P:5:DA:H2'	2:P:6:DT:H71	1.95	0.49
3:A:299:ARG:CD	3:A:310:PRO:HD3	2.42	0.49
3:A:218:LEU:CB	3:A:224:ILE:HD12	2.42	0.49
3:A:133:ASN:ND2	7:A:512:HOH:O	2.46	0.48
3:A:277:ILE:HG13	3:A:335:GLU:HB2	1.94	0.48
1:T:7:DA:H61	2:P:1:DT:H3	1.62	0.48
3:A:253:ARG:HH11	3:A:253:ARG:CG	2.26	0.48
3:A:26:GLU:HB2	3:A:36:TYR:HB2	1.96	0.48
3:A:123:GLU:HG3	7:A:623:HOH:O	2.14	0.48
3:A:16:THR:CG2	3:A:47:ALA:HB2	2.43	0.48
3:A:137:ARG:NH1	7:A:515:HOH:O	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:11:LEU:HD23	3:A:12:ASN:H	1.79	0.47
3:A:99:PHE:CD1	3:A:99:PHE:C	2.91	0.47
3:A:155:MET:HA	3:A:158:MET:HE3	1.96	0.47
3:A:327:TYR:CD1	3:A:328:ARG:N	2.79	0.47
3:A:12:ASN:HA	7:A:553:HOH:O	2.13	0.47
3:A:288:GLU:C	3:A:290:GLY:H	2.22	0.47
3:A:101:THR:HA	3:A:106:ILE:HG22	1.95	0.47
3:A:183:ARG:HH11	3:A:275:SER:CA	2.28	0.47
3:A:230:LYS:HE2	3:A:230:LYS:HB3	1.41	0.47
3:A:270:LEU:HD23	3:A:319:ILE:CD1	2.40	0.47
3:A:278:PHE:HB2	3:A:333:ARG:O	2.14	0.47
3:A:18:MET:CE	3:A:76:PHE:HB2	2.40	0.47
3:A:37:ASN:HB3	7:A:556:HOH:O	2.13	0.47
3:A:59:ALA:C	3:A:65:VAL:HG21	2.38	0.47
3:A:259:LEU:C	3:A:260:ILE:HG12	2.39	0.47
3:A:317:LYS:O	3:A:320:PHE:N	2.47	0.47
3:A:49:TYR:CE2	3:A:53:ILE:CG1	2.98	0.47
3:A:85:LEU:HD12	3:A:85:LEU:HA	1.68	0.47
3:A:330:PRO:HA	3:A:333:ARG:HG3	1.97	0.47
3:A:16:THR:HG23	3:A:46:ILE:HD11	1.96	0.47
3:A:121:THR:HG23	3:A:124:ASP:CG	2.39	0.47
3:A:250:TYR:HA	3:A:251:PRO:HD3	1.74	0.47
2:P:3:DT:H5"	7:P:615:HOH:O	2.15	0.47
3:A:241:LEU:HD12	3:A:250:TYR:CD1	2.50	0.47
3:A:294:ASN:N	3:A:294:ASN:HD22	2.13	0.47
3:A:133:ASN:H	3:A:136:GLN:HE21	1.61	0.46
1:T:4:DT:N3	2:P:4:DA:C2	2.78	0.46
3:A:97:ILE:HG23	3:A:111:ALA:HB1	1.98	0.46
3:A:75:GLU:O	3:A:79:THR:HG23	2.16	0.46
3:A:201:THR:HA	3:A:261:PRO:CB	2.46	0.46
3:A:22:LEU:HD22	3:A:85:LEU:HD11	1.96	0.46
3:A:133:ASN:ND2	3:A:135:HIS:H	2.13	0.46
3:A:127:LYS:HB3	3:A:127:LYS:HE3	1.58	0.46
3:A:176:THR:HG22	3:A:178:CYS:SG	2.56	0.46
3:A:327:TYR:CD1	3:A:327:TYR:C	2.94	0.46
3:A:22:LEU:HA	3:A:22:LEU:HD13	1.49	0.45
3:A:218:LEU:N	3:A:218:LEU:CD1	2.78	0.45
3:A:270:LEU:HD12	3:A:270:LEU:O	2.16	0.45
3:A:125:LEU:HD21	3:A:132:LEU:HD21	1.99	0.45
3:A:174:ILE:O	3:A:195:LEU:HD12	2.16	0.45
3:A:293:ILE:CD1	3:A:298:ILE:HD12	2.45	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:23:ALA:O	3:A:36:TYR:HD1	2.00	0.45
3:A:196:THR:CG2	3:A:265:TYR:CD1	2.99	0.45
3:A:276:ASP:O	3:A:280:LYS:HG3	2.16	0.45
3:A:316:GLU:O	3:A:320:PHE:HD2	1.99	0.45
3:A:56:GLY:O	3:A:59:ALA:HB3	2.17	0.45
3:A:182:ARG:HH11	3:A:182:ARG:CG	2.27	0.45
3:A:254:ARG:NH1	3:A:255:ILE:H	2.15	0.45
3:A:31:GLN:NE2	3:A:112:ARG:NH1	2.51	0.45
3:A:29:VAL:HG21	3:A:94:SER:CB	2.23	0.45
3:A:34:HIS:O	3:A:37:ASN:N	2.49	0.45
3:A:49:TYR:CE2	3:A:53:ILE:HG13	2.52	0.45
3:A:25:PHE:CG	3:A:88:ILE:CD1	3.00	0.45
3:A:125:LEU:HD23	3:A:125:LEU:HA	1.63	0.45
3:A:59:ALA:O	3:A:65:VAL:HG21	2.17	0.44
3:A:25:PHE:CD1	3:A:88:ILE:HD11	2.52	0.44
3:A:197:HIS:CG	3:A:198:PRO:HD2	2.52	0.44
3:A:25:PHE:CD2	3:A:88:ILE:HD13	2.52	0.44
3:A:197:HIS:O	3:A:262:LYS:HB2	2.17	0.44
3:A:146:PHE:CZ	3:A:238:VAL:CG2	2.98	0.44
3:A:182:ARG:NH1	3:A:273:THR:CG2	2.79	0.44
3:A:265:TYR:O	3:A:268:GLY:N	2.49	0.44
3:A:298:ILE:HA	7:A:593:HOH:O	2.17	0.44
3:A:62:LEU:HD12	3:A:62:LEU:HA	1.63	0.44
2:P:5:DA:P	3:A:109:SER:HB3	2.58	0.44
3:A:150:ILE:N	3:A:188:SER:O	2.44	0.44
3:A:228:LEU:HB2	3:A:236:MET:O	2.17	0.44
3:A:232:GLU:O	3:A:233:THR:HG23	2.18	0.44
3:A:16:THR:HG22	3:A:43:ALA:O	2.18	0.44
3:A:200:PHE:O	3:A:262:LYS:N	2.48	0.44
3:A:217:GLN:O	3:A:220:LYS:N	2.50	0.44
3:A:41:LYS:HD3	3:A:42:ALA:H	1.84	0.43
3:A:108:PRO:HA	3:A:111:ALA:HB3	1.99	0.43
3:A:144:GLY:O	3:A:146:PHE:N	2.51	0.43
3:A:49:TYR:CE2	3:A:53:ILE:HG12	2.53	0.43
2:P:2:DC:H6	2:P:2:DC:H2'	1.41	0.43
2:P:5:DA:O5'	3:A:107:GLY:HA3	2.18	0.43
3:A:12:ASN:HD21	3:A:53:ILE:H	1.66	0.43
3:A:163:LEU:HA	3:A:163:LEU:HD23	1.09	0.43
3:A:158:MET:HG2	3:A:241:LEU:HD21	2.01	0.43
3:A:287:LEU:HD13	3:A:287:LEU:HA	1.39	0.43
3:A:76:PHE:HB2	3:A:82:LEU:HD13	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:286:ALA:HA	3:A:323:ILE:HG22	2.00	0.43
2:P:8:DA:H4'	3:A:272:PHE:HE2	1.77	0.43
3:A:327:TYR:HD1	3:A:327:TYR:C	2.26	0.43
3:A:244:LYS:HB3	3:A:245:ASN:ND2	2.34	0.42
3:A:34:HIS:O	3:A:38:ALA:N	2.52	0.42
3:A:94:SER:O	3:A:98:ASN:OD1	2.37	0.42
3:A:28:ASN:HD22	3:A:28:ASN:HA	1.29	0.42
3:A:54:LYS:HD2	3:A:54:LYS:HA	1.42	0.42
3:A:151:PRO:HG2	3:A:154:GLU:HG3	2.01	0.42
3:A:260:ILE:HG23	3:A:260:ILE:HD12	1.53	0.42
3:A:277:ILE:CG1	3:A:335:GLU:HB2	2.49	0.42
3:A:280:LYS:O	3:A:283:ARG:HB2	2.20	0.42
3:A:22:LEU:CD2	3:A:85:LEU:HD13	2.50	0.42
3:A:127:LYS:C	3:A:128:ASN:ND2	2.78	0.42
3:A:149:ARG:NH2	3:A:188:SER:CA	2.82	0.42
3:A:127:LYS:HB2	3:A:128:ASN:HD21	1.84	0.42
3:A:170:ASP:HB3	3:A:173:TYR:CE2	2.55	0.42
3:A:289:LYS:HD2	3:A:289:LYS:HA	1.85	0.42
3:A:293:ILE:CD1	3:A:298:ILE:CD1	2.98	0.42
3:A:135:HIS:CD2	3:A:135:HIS:C	2.97	0.42
3:A:245:ASN:O	3:A:246:ASP:O	2.37	0.42
1:T:4:DT:O2	2:P:4:DA:H2	2.02	0.42
3:A:49:TYR:CD2	3:A:53:ILE:CG1	3.02	0.42
3:A:104:SER:N	7:A:600:HOH:O	2.27	0.42
3:A:118:GLY:HA2	3:A:120:LYS:HE3	2.02	0.42
3:A:200:PHE:CE2	3:A:261:PRO:N	2.88	0.42
3:A:253:ARG:NH2	7:A:620:HOH:O	2.28	0.42
3:A:315:SER:C	3:A:317:LYS:N	2.78	0.42
3:A:15:ILE:HG21	3:A:46:ILE:HD13	1.99	0.42
3:A:241:LEU:HA	3:A:242:PRO:HD2	1.69	0.42
3:A:97:ILE:HG23	3:A:111:ALA:CB	2.50	0.41
3:A:253:ARG:NH1	7:A:620:HOH:O	2.51	0.41
3:A:97:ILE:O	3:A:97:ILE:HG22	2.20	0.41
3:A:155:MET:HE3	3:A:181:PHE:HB2	2.01	0.41
3:A:221:VAL:O	3:A:222:HIS:HB2	2.20	0.41
2:P:4:DA:H2''	2:P:5:DA:H5'	2.03	0.41
3:A:270:LEU:HD21	3:A:282:MET:HE1	2.01	0.41
3:A:295:GLU:H	3:A:295:GLU:CD	2.29	0.41
3:A:40:ARG:HE	3:A:40:ARG:HB2	1.64	0.41
3:A:228:LEU:HD12	3:A:228:LEU:HA	1.64	0.41
3:A:79:THR:O	3:A:79:THR:OG1	2.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:204:SER:O	3:A:204:SER:OG	2.36	0.41
3:A:236:MET:HG3	3:A:256:ASP:OD1	2.21	0.41
3:A:277:ILE:CG1	3:A:335:GLU:CB	2.99	0.41
3:A:286:ALA:CB	3:A:293:ILE:CG1	2.99	0.41
1:T:1:DC:H6	1:T:1:DC:H2'	1.62	0.41
3:A:68:LYS:HB2	3:A:68:LYS:HZ1	1.82	0.41
3:A:183:ARG:HH11	3:A:275:SER:CB	2.33	0.41
3:A:114:PHE:CZ	3:A:132:LEU:CD2	2.99	0.40
3:A:197:HIS:CD2	3:A:198:PRO:CD	3.02	0.40
3:A:250:TYR:HB3	7:A:577:HOH:O	2.21	0.40
3:A:191:MET:HG2	3:A:255:ILE:HG13	2.03	0.40
3:A:210:LEU:HD23	3:A:210:LEU:HA	1.77	0.40
3:A:303:VAL:C	3:A:305:GLY:N	2.78	0.40
3:A:120:LYS:N	3:A:124:ASP:OD2	2.49	0.40
3:A:316:GLU:O	3:A:320:PHE:CD2	2.75	0.40
3:A:165:GLU:CA	3:A:168:LYS:HG2	2.37	0.40
3:A:299:ARG:HD2	3:A:310:PRO:HD3	2.02	0.40
3:A:172:GLU:HG3	3:A:198:PRO:HG2	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
3	A	325/335 (97%)	256 (79%)	47 (14%)	22 (7%)	1 5

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	185	ALA
3	A	205	THR
3	A	207	GLN

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Mol	Chain	Res	Type
3	A	244	LYS
3	A	246	ASP
3	A	295	GLU
3	A	202	SER
3	A	204	SER
3	A	247	GLU
3	A	265	TYR
3	A	316	GLU
3	A	91	ASP
3	A	145	ASP
3	A	304	THR
3	A	208	PRO
3	A	262	LYS
3	A	310	PRO
3	A	60	LYS
3	A	222	HIS
3	A	229	SER
3	A	324	GLN
3	A	80	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
3	A	288/295 (98%)	222 (77%)	66 (23%)	1 5

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	10	THR
3	A	18	MET
3	A	22	LEU
3	A	27	LYS
3	A	33	ILE
3	A	37	ASN
3	A	40	ARG

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Mol	Chain	Res	Type
3	A	41	LYS
3	A	44	SER
3	A	46	ILE
3	A	52	LYS
3	A	54	LYS
3	A	62	LEU
3	A	65	VAL
3	A	67	THR
3	A	68	LYS
3	A	72	LYS
3	A	79	THR
3	A	85	LEU
3	A	89	ARG
3	A	92	ASP
3	A	93	THR
3	A	100	LEU
3	A	101	THR
3	A	119	ILE
3	A	121	THR
3	A	128	ASN
3	A	133	ASN
3	A	136	GLN
3	A	138	ILE
3	A	152	ARG
3	A	153	GLU
3	A	161	ILE
3	A	168	LYS
3	A	169	VAL
3	A	174	ILE
3	A	194	LEU
3	A	199	SER
3	A	213	GLN
3	A	214	VAL
3	A	218	LEU
3	A	228	LEU
3	A	230	LYS
3	A	241	LEU
3	A	245	ASN
3	A	248	LYS
3	A	253	ARG
3	A	255	ILE
3	A	259	LEU

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Mol	Chain	Res	Type
3	A	270	LEU
3	A	275	SER
3	A	277	ILE
3	A	279	ASN
3	A	287	LEU
3	A	293	ILE
3	A	294	ASN
3	A	295	GLU
3	A	298	ILE
3	A	301	LEU
3	A	304	THR
3	A	309	GLU
3	A	311	LEU
3	A	314	ASP
3	A	324	GLN
3	A	325	TRP
3	A	331	LYS

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

Mol	Chain	Res	Type
3	A	12	ASN
3	A	28	ASN
3	A	31	GLN
3	A	37	ASN
3	A	98	ASN
3	A	128	ASN
3	A	133	ASN
3	A	134	HIS
3	A	136	GLN
3	A	157	GLN
3	A	212	HIS
3	A	213	GLN
3	A	217	GLN
3	A	245	ASN
3	A	264	GLN
3	A	281	ASN
3	A	294	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 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	DAD	A	338	4	10,12,31	1.63	2 (20%)	17,20,48	1.19	1 (5%)

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	DAD	A	338	4	-	2/12/12/31	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	338	DAD	PB-O3A	3.29	1.63	1.59
6	A	338	DAD	PB-O3B	-2.75	1.56	1.59

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	338	DAD	O2G-PG-O1G	2.12	119.09	110.83

There are no chirality outliers.

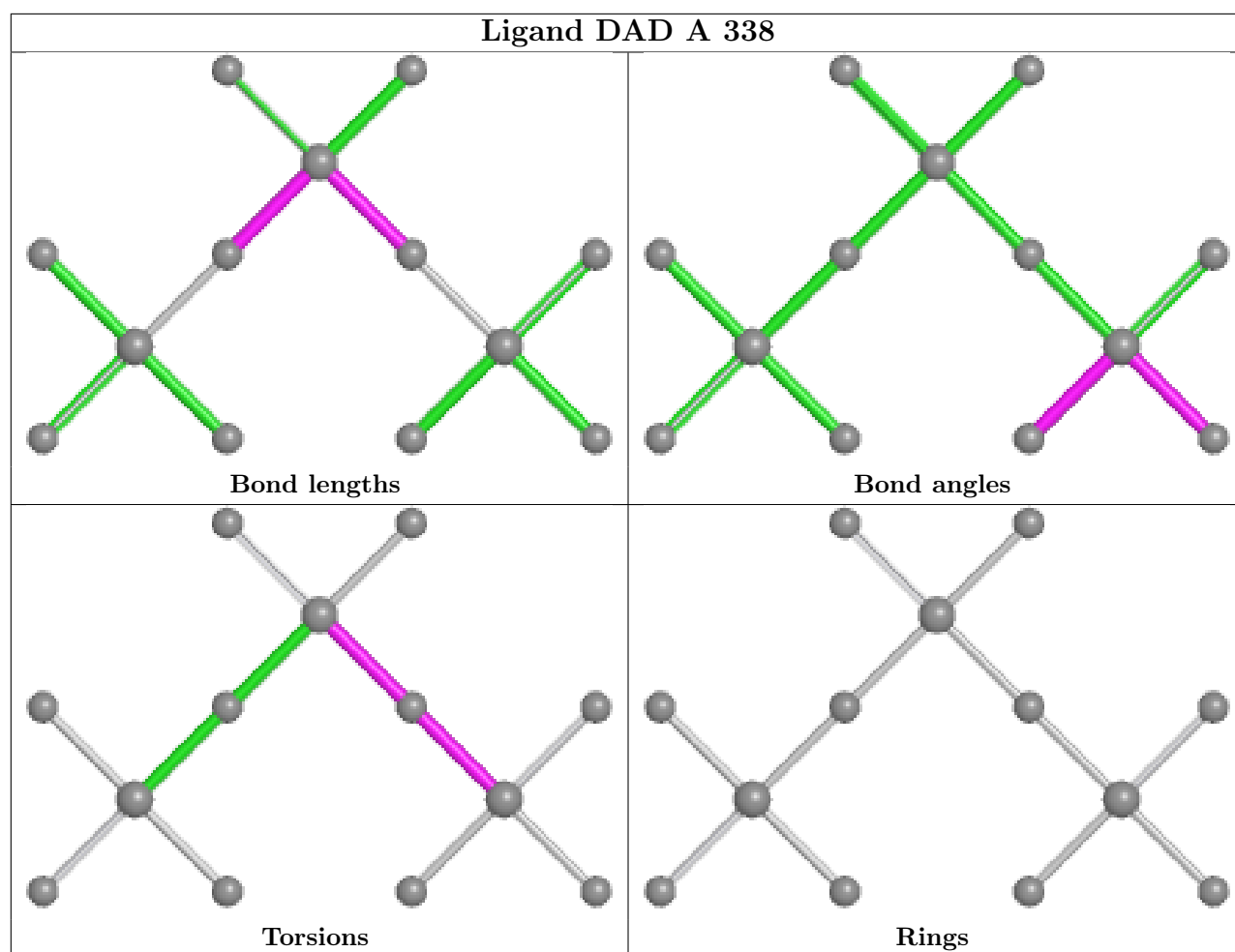
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	338	DAD	PB-O3B-PG-O3G
6	A	338	DAD	PG-O3B-PB-O2B

There are no ring outliers.

No monomer is involved in short contacts.

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	T	8/8 (100%)	-0.47	0 100 100	17, 42, 99, 99	0
2	P	8/8 (100%)	-0.42	1 (12%) 8 5	19, 36, 69, 100	0
3	A	325/335 (97%)	-0.93	0 100 100	4, 35, 85, 100	0
All	All	341/351 (97%)	-0.90	1 (0%) 90 79	4, 36, 88, 100	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	P	8	DA	3.0

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.

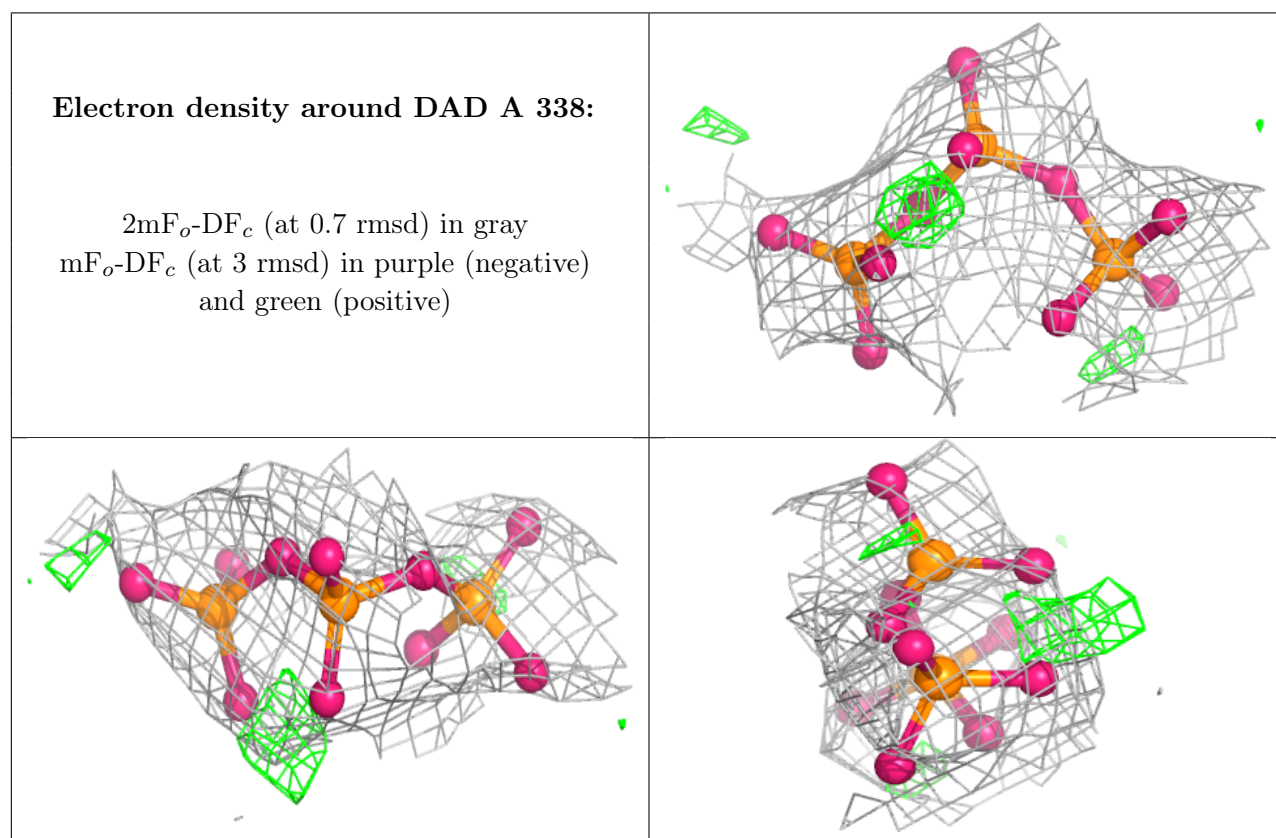
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	DAD	A	338	13/29	0.82	0.15	30,57,63,67	13
5	NA	A	342	1/1	0.83	0.10	57,57,57,57	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MN	A	340	1/1	0.88	0.08	30,30,30,30	1
4	MN	A	339	1/1	0.89	0.07	30,30,30,30	1
5	NA	A	341	1/1	0.99	0.03	17,17,17,17	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.