



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 10:43 AM UTC

PDB ID : 8VQ2 / pdb_00008vq2
Title : HSV1 polymerase ternary complex with dsDNA and compound 44
Authors : Hayes, R.P.; Heo, M.R.; Plotkin, M.
Deposited on : 2024-01-17
Resolution : 3.83 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:

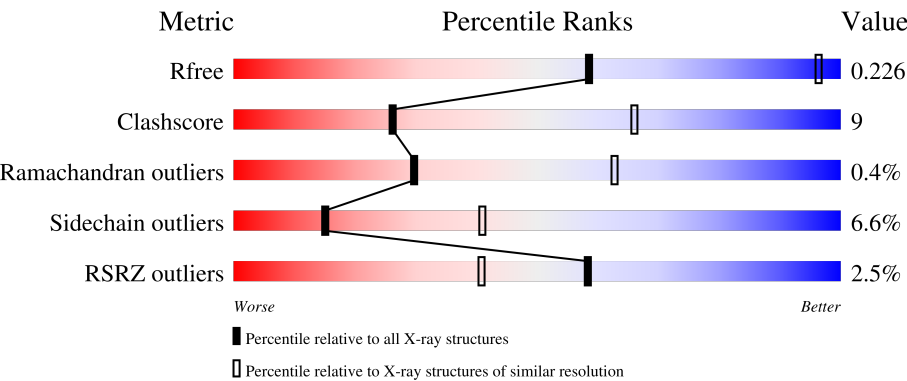
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal 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 i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.83 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.



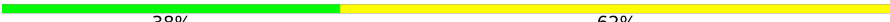
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	1043 (3.98-3.66)
Clashscore	190562	1075 (3.98-3.66)
Ramachandran outliers	187476	1029 (3.98-3.66)
Sidechain outliers	187428	1024 (3.98-3.66)
RSRZ outliers	180081	1042 (3.98-3.66)

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	1163	2% 69% 19% • 11%
1	B	1163	2% 68% 20% • 10%
2	C	15	7% 60% 40%
2	E	15	40% 53% 7%
3	D	13	8% 46% 54%

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Mol	Chain	Length	Quality of chain
3	F	13	 A horizontal bar chart showing the quality of chain F. The bar is divided into two segments: a green segment on the left representing 38% and a yellow segment on the right representing 62%. The percentages are labeled below the bar.

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 17099 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 DNA polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1037	Total	C	N	O	S	0	0	0
			7845	5029	1343	1434	39			
1	B	1046	Total	C	N	O	S	0	0	0
			8072	5188	1391	1453	40			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	35	LEU	-	expression tag	UNP I7GY94
A	36	GLU	-	expression tag	UNP I7GY94
A	37	VAL	-	expression tag	UNP I7GY94
A	38	LEU	-	expression tag	UNP I7GY94
A	39	PHE	-	expression tag	UNP I7GY94
A	40	GLN	-	expression tag	UNP I7GY94
A	41	GLY	-	expression tag	UNP I7GY94
A	42	PRO	-	expression tag	UNP I7GY94
A	370	ALA	GLU	engineered mutation	UNP I7GY94
B	35	LEU	-	expression tag	UNP I7GY94
B	36	GLU	-	expression tag	UNP I7GY94
B	37	VAL	-	expression tag	UNP I7GY94
B	38	LEU	-	expression tag	UNP I7GY94
B	39	PHE	-	expression tag	UNP I7GY94
B	40	GLN	-	expression tag	UNP I7GY94
B	41	GLY	-	expression tag	UNP I7GY94
B	42	PRO	-	expression tag	UNP I7GY94
B	370	ALA	GLU	engineered mutation	UNP I7GY94

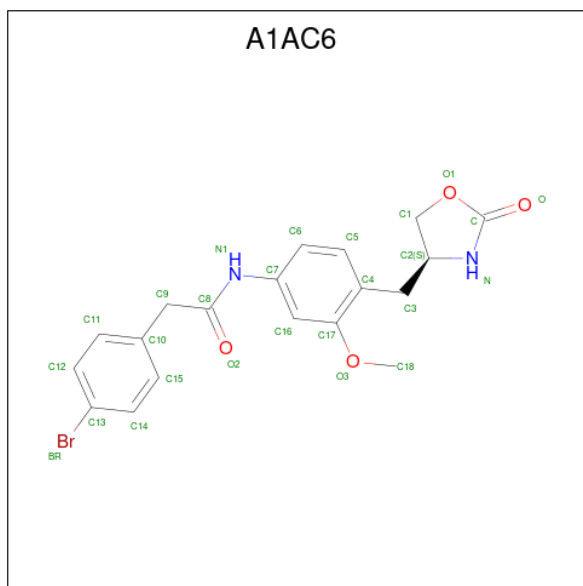
- Molecule 2 is a DNA chain called DNA (5'-D(P*TP*GP*GP*TP*AP*GP*GP*GP*GP*A P*AP*GP*GP*AP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	15	Total	C	N	O	P	0	0	0
			320	150	66	89	15			
2	E	14	Total	C	N	O	P	0	0	0
			300	140	64	82	14			

- Molecule 3 is a DNA chain called DNA (5'-D(P*AP*TP*CP*CP*TP*TP*CP*CP*CP*CP*TP*AP*C)-3').

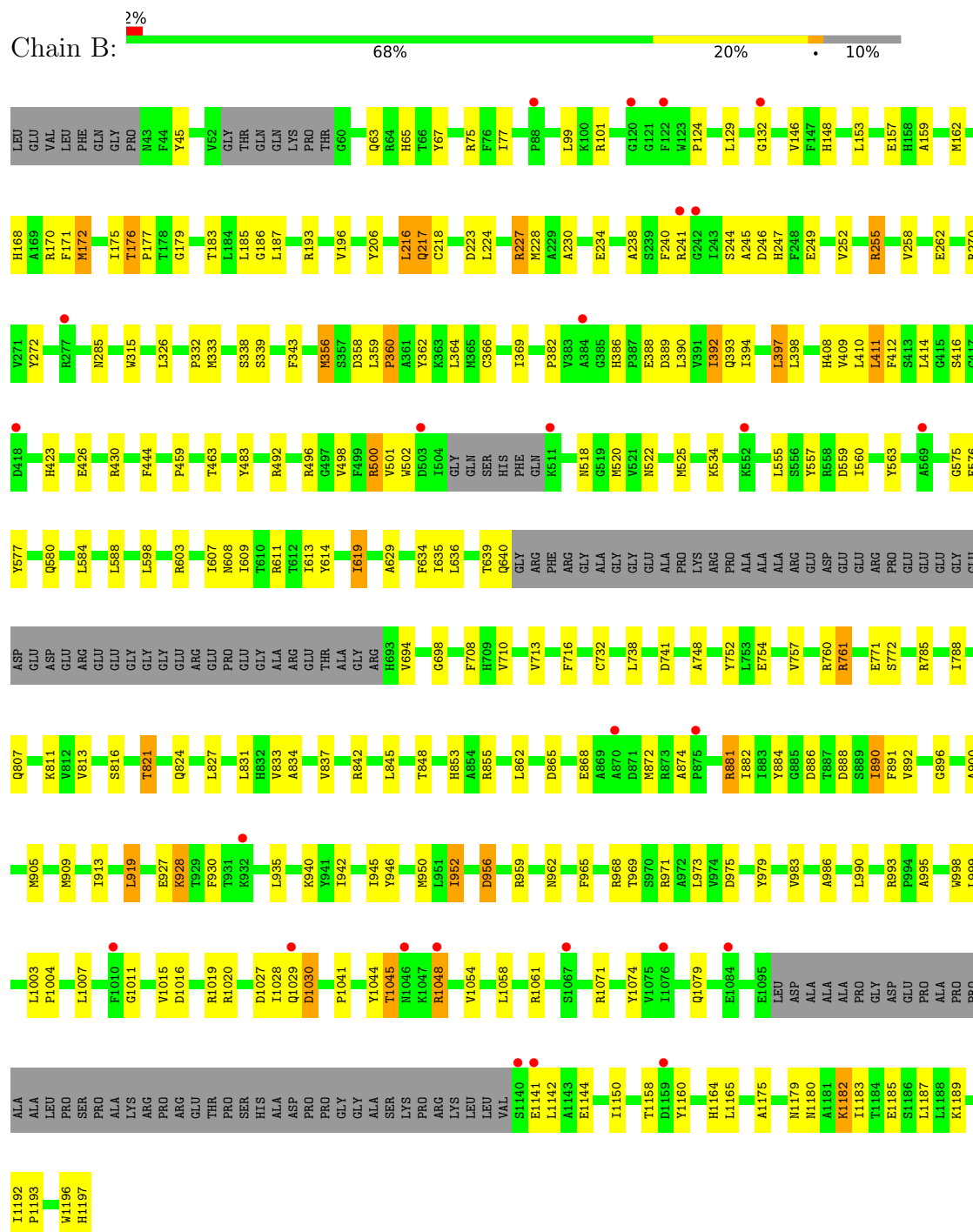
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	13	Total	C	N	O	P	0	0	0
			255	123	39	80	13			
3	F	13	Total	C	N	O	P	0	0	0
			255	123	39	80	13			

- Molecule 4 is 2-(4-bromophenyl)-N-(3-methoxy-4-{[(4S)-2-oxo-1,3-oxazolidin-4-yl]methyl}phenyl)acetamide (CCD ID: A1AC6) (formula: C₁₉H₁₉BrN₂O₄) (labeled as "Ligand of Interest" by depositor).

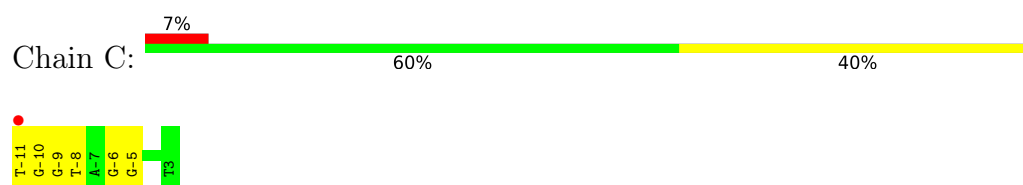


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	C	1	Total	Br	C	N	O	0	0
			26	1	19	2	4		
4	E	1	Total	Br	C	N	O	0	0
			26	1	19	2	4		

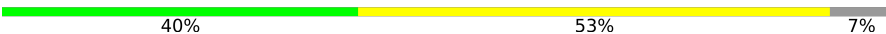
• Molecule 1: DNA polymerase



• Molecule 2: DNA (5'-D(P*TP*GP*GP*TP*AP*GP*GP*GP*GP*AP*AP*GP*GP*AP*T)-3')



- Molecule 2: DNA (5'-D(P*TP*GP*GP*TP*AP*GP*GP*GP*GP*AP*AP*GP*GP*AP*T)-3')

Chain E:  40% 53% 7%

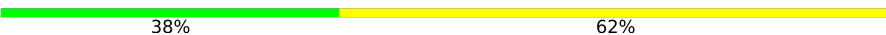


- Molecule 3: DNA (5'-D(P*AP*TP*CP*CP*TP*TP*CP*CP*CP*CP*TP*AP*C)-3')

Chain D:  8% 46% 54%



- Molecule 3: DNA (5'-D(P*AP*TP*CP*CP*TP*TP*CP*CP*CP*CP*TP*AP*C)-3')

Chain F:  38% 62%



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	180.63Å 180.63Å 231.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.81 – 3.83 43.81 – 3.83	Depositor EDS
% Data completeness (in resolution range)	98.6 (43.81-3.83) 98.6 (43.81-3.83)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.58 (at 3.88Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.209 , 0.227 0.211 , 0.226	Depositor DCC
R_{free} test set	3410 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å ²)	118.7	Xtriage
Anisotropy	0.177	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 80.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.064 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	17099	wwPDB-VP
Average B, all atoms (Å ²)	124.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.10% 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: A1AC6

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	0.16	0/8036	0.39	0/10957
1	B	0.16	0/8270	0.39	0/11249
2	C	0.25	0/361	0.50	0/558
2	E	0.25	0/339	0.48	0/524
3	D	0.38	0/282	0.60	0/430
3	F	0.36	0/282	0.53	0/430
All	All	0.18	0/17570	0.40	0/24148

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	7845	0	7490	133	0
1	B	8072	0	7870	156	0
2	C	320	0	169	10	0
2	E	300	0	157	10	0
3	D	255	0	148	8	0
3	F	255	0	148	10	0
4	C	26	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	E	26	0	0	0	0
All	All	17099	0	15982	311	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 9.

All (311) 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:258:VAL:HG21	1:A:636:LEU:HD23	1.54	0.90
1:B:788:ILE:HD11	1:B:807:GLN:HB3	1.52	0.89
1:B:258:VAL:HG21	1:B:636:LEU:HD23	1.57	0.86
1:A:1061:ARG:NH1	1:A:1144:GLU:OE2	2.10	0.84
1:B:392:ILE:HG22	1:B:393:GLN:HG3	1.64	0.78
1:B:754:GLU:OE2	1:B:761:ARG:NH2	2.15	0.78
1:A:934:LEU:HD22	1:A:936:ILE:HG12	1.64	0.78
2:E:-10:DG:H5''	2:E:-10:DG:H8	1.49	0.78
3:D:0:DA:H2''	3:D:1:DC:H5''	1.68	0.76
1:A:339:SER:OG	1:A:492:ARG:NH1	2.18	0.76
1:B:598:LEU:HD13	1:B:613:ILE:HG12	1.66	0.76
3:F:0:DA:H2''	3:F:1:DC:H5''	1.67	0.75
1:B:990:LEU:HD11	1:B:1007:LEU:HD11	1.69	0.74
1:B:639:THR:HG23	1:B:640:GLN:HG3	1.71	0.73
1:B:1003:LEU:HB3	1:B:1007:LEU:HD12	1.69	0.73
1:A:618:GLN:HB3	1:A:827:LEU:HD11	1.71	0.73
1:B:738:LEU:HD12	1:B:760:ARG:HH21	1.56	0.71
1:A:1067:SER:HB2	1:A:1070:ASP:OD2	1.90	0.70
1:A:478:LYS:HD3	1:A:482:ILE:HD12	1.74	0.69
1:A:598:LEU:HD13	1:A:613:ILE:HG12	1.75	0.69
1:B:716:PHE:HB2	1:B:890:ILE:HG22	1.75	0.68
1:A:998:TRP:HZ3	1:A:1004:PRO:HD3	1.58	0.68
1:B:339:SER:OG	1:B:492:ARG:NH1	2.27	0.68
2:E:-10:DG:H5''	2:E:-10:DG:C8	2.29	0.68
1:B:998:TRP:CZ3	1:B:1004:PRO:HD3	2.30	0.66
1:B:63:GLN:NE2	1:B:162:MET:SD	2.69	0.66
1:B:919:LEU:HD12	1:B:919:LEU:H	1.60	0.66
1:B:855:ARG:NH1	1:B:865:ASP:OD2	2.28	0.66
1:B:608:ASN:HD21	1:B:611:ARG:HG3	1.61	0.66
1:A:716:PHE:HB3	1:A:924:LEU:HD11	1.77	0.65
1:B:959:ARG:HH22	3:F:-3:DC:H1'	1.60	0.65
1:A:716:PHE:HB2	1:A:890:ILE:HG22	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:698:GLY:O	1:B:842:ARG:NH1	2.30	0.64
1:B:998:TRP:HZ3	1:B:1004:PRO:HD3	1.60	0.64
2:E:-6:DG:H2'	2:E:-5:DG:C8	2.33	0.64
1:B:959:ARG:NH1	3:F:-2:DC:O2	2.31	0.64
1:B:175:ILE:HG22	1:B:179:GLY:HA3	1.79	0.64
1:B:148:HIS:O	1:B:186:GLY:HA3	1.98	0.63
1:A:618:GLN:HE21	2:C:-11:DT:H2''	1.64	0.63
1:B:708:PHE:HE1	1:B:710:VAL:HG22	1.64	0.62
1:B:942:ILE:HG12	1:B:952:ILE:HG22	1.80	0.62
1:A:175:ILE:HG22	1:A:179:GLY:HA3	1.81	0.62
1:A:501:VAL:HG22	1:A:515:ILE:HD13	1.80	0.62
1:A:77:ILE:HG22	1:A:390:LEU:HD11	1.82	0.62
1:A:193:ARG:NH1	1:A:338:SER:O	2.33	0.62
1:B:940:LYS:NZ	2:E:-4:DG:H5'	2.15	0.61
1:A:739:ARG:NH2	1:A:741:ASP:OD2	2.34	0.61
1:B:611:ARG:HH21	1:B:619:ILE:HD11	1.63	0.61
1:B:397:LEU:HD23	1:B:409:VAL:HG13	1.83	0.61
1:A:557:TYR:O	1:A:560:ILE:HG22	2.00	0.61
1:A:337:THR:OG1	1:A:342:GLU:OE2	2.19	0.61
1:A:959:ARG:HH22	2:C:-5:DG:N2	1.99	0.60
1:B:358:ASP:OD2	1:B:358:ASP:N	2.33	0.60
1:B:1027:ASP:HB3	1:B:1030:ASP:HB2	1.84	0.60
2:E:-9:DG:H2''	2:E:-8:DT:H5'	1.82	0.59
1:B:223:ASP:O	1:B:227:ARG:HG2	2.02	0.59
1:B:1071:ARG:NH1	3:F:-2:DC:OP1	2.36	0.59
1:B:1041:PRO:HA	1:B:1044:TYR:HD1	1.67	0.59
1:A:618:GLN:NE2	2:C:-11:DT:H2''	2.17	0.59
1:A:364:LEU:HD11	1:A:463:THR:HG22	1.84	0.58
1:B:1061:ARG:NH1	1:B:1144:GLU:OE2	2.33	0.58
1:B:950:MET:SD	1:B:971:ARG:NH1	2.76	0.58
3:F:-2:DC:H2''	3:F:-1:DT:H6	1.68	0.58
1:B:821:THR:HG22	1:B:834:ALA:HB2	1.86	0.58
2:C:-9:DG:H2''	2:C:-8:DT:H5'	1.84	0.58
1:A:369:ILE:HG22	1:A:394:ILE:HG12	1.87	0.57
1:A:995:ALA:HA	1:A:998:TRP:CD1	2.39	0.57
1:B:159:ALA:HB2	1:B:176:THR:HA	1.86	0.57
1:A:1028:ILE:HD12	1:A:1150:ILE:HG12	1.86	0.57
1:B:1185:GLU:HG2	1:B:1189:LYS:HE3	1.86	0.57
1:B:397:LEU:C	1:B:398:LEU:HD23	2.29	0.57
1:A:998:TRP:CZ3	1:A:1004:PRO:HD3	2.37	0.57
1:A:176:THR:HB	1:A:177:PRO:HD3	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:-2:DC:H2''	3:D:-1:DT:H6	1.69	0.57
1:A:392:ILE:HG13	1:A:393:GLN:N	2.20	0.56
1:A:855:ARG:NH1	1:A:865:ASP:OD2	2.38	0.56
1:B:853:HIS:CD2	1:B:881:ARG:HD2	2.40	0.56
1:B:940:LYS:HZ3	2:E:-4:DG:H5'	1.70	0.56
1:A:1192:ILE:HG23	1:A:1196:TRP:HE3	1.71	0.56
1:B:216:LEU:HD13	1:B:227:ARG:HG3	1.88	0.56
1:B:965:PHE:HE1	1:B:1020:ARG:HH11	1.52	0.56
1:B:224:LEU:HG	1:B:228:MET:HE2	1.87	0.56
1:A:884:TYR:HD2	1:A:891:PHE:CD2	2.24	0.56
1:B:498:VAL:H	1:B:518:ASN:HD22	1.53	0.55
1:B:555:LEU:HB2	1:B:577:TYR:CD2	2.41	0.55
1:A:969:THR:HG21	1:A:1165:LEU:HD11	1.89	0.55
1:B:369:ILE:HG22	1:B:394:ILE:HG12	1.87	0.55
1:A:218:CYS:HB2	1:A:223:ASP:HB3	1.88	0.55
1:A:398:LEU:HD22	1:A:457:TYR:CD2	2.42	0.55
1:A:862:LEU:HD11	1:A:909:MET:HE2	1.88	0.55
1:A:924:LEU:HD12	1:A:925:GLU:N	2.22	0.55
1:B:741:ASP:OD1	1:B:741:ASP:N	2.40	0.55
1:B:1192:ILE:HG13	1:B:1197:HIS:HE1	1.72	0.55
1:A:366:CYS:HA	1:A:463:THR:O	2.07	0.54
1:A:1058:LEU:HD21	1:A:1145:ASP:OD2	2.06	0.54
1:B:218:CYS:HB2	1:B:223:ASP:HB3	1.88	0.54
1:B:1054:VAL:HG21	1:B:1074:TYR:HB3	1.89	0.54
2:C:-10:DG:O5'	2:C:-10:DG:H8	1.90	0.54
1:B:555:LEU:HB2	1:B:577:TYR:HD2	1.73	0.54
1:A:963:CYS:SG	1:A:965:PHE:HB3	2.48	0.54
1:A:1164:HIS:O	1:A:1167:GLY:N	2.39	0.54
1:A:1183:ILE:O	1:A:1187:LEU:HG	2.07	0.54
1:B:153:LEU:HB2	1:B:183:THR:HB	1.88	0.54
1:B:1016:ASP:O	1:B:1019:ARG:HB3	2.07	0.54
1:A:176:THR:CB	1:A:177:PRO:HD3	2.38	0.54
1:A:1014:LEU:O	1:A:1017:ALA:HB3	2.08	0.54
1:A:311:VAL:O	1:A:316:TYR:OH	2.21	0.53
1:A:880:MET:HE1	1:A:909:MET:HE1	1.88	0.53
2:C:-6:DG:H2'	2:C:-5:DG:C8	2.42	0.53
1:B:1003:LEU:HD21	1:B:1187:LEU:HD12	1.88	0.53
1:B:234:GLU:OE2	1:B:234:GLU:N	2.35	0.53
1:B:364:LEU:HD11	1:B:463:THR:HG22	1.91	0.53
3:F:-2:DC:H2''	3:F:-1:DT:C6	2.44	0.53
1:A:498:VAL:H	1:A:518:ASN:HD22	1.57	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:411:LEU:HD11	1:B:575:GLY:HA3	1.90	0.53
1:B:853:HIS:NE2	1:B:881:ARG:HD2	2.24	0.53
1:B:501:VAL:HG12	1:B:501:VAL:O	2.09	0.52
1:B:986:ALA:O	1:B:990:LEU:HD12	2.10	0.52
1:A:1071:ARG:NH2	3:D:-2:DC:OP1	2.41	0.52
1:B:132:GLY:HA2	1:B:333:MET:HE1	1.91	0.52
3:D:-2:DC:H2''	3:D:-1:DT:C6	2.44	0.52
1:B:1193:PRO:HD2	1:B:1196:TRP:CZ3	2.44	0.52
1:A:397:LEU:HD12	1:A:397:LEU:N	2.25	0.52
1:A:132:GLY:HA2	1:A:333:MET:HE1	1.92	0.52
1:B:886:ASP:O	1:B:888:ASP:N	2.42	0.51
1:A:824:GLN:HB3	1:A:831:LEU:HD11	1.93	0.51
1:B:153:LEU:HD11	1:B:185:LEU:HD11	1.93	0.51
1:B:629:ALA:HB1	1:B:634:PHE:HB2	1.91	0.51
1:A:696:TYR:HB2	1:A:835:ALA:HB2	1.93	0.51
1:A:223:ASP:O	1:A:227:ARG:HG2	2.10	0.51
1:A:618:GLN:HB3	1:A:827:LEU:CD1	2.39	0.50
1:B:1048:ARG:NE	2:E:0:DG:OP1	2.39	0.50
1:B:998:TRP:HZ3	1:B:1003:LEU:HA	1.76	0.50
1:B:1185:GLU:O	1:B:1189:LYS:HG2	2.11	0.50
1:A:286:PHE:CD2	1:A:287:CYS:HB2	2.46	0.50
3:F:-5:DC:H1'	3:F:-4:DC:H5'	1.92	0.50
1:B:500:ARG:HD2	1:B:502:TRP:CE3	2.46	0.50
1:B:168:HIS:HB2	1:B:171:PHE:CE1	2.47	0.50
1:B:708:PHE:CE1	1:B:710:VAL:HG22	2.46	0.50
1:A:365:MET:HB2	1:A:398:LEU:HD23	1.93	0.49
1:A:704:PRO:HB2	1:A:935:LEU:HD13	1.93	0.49
1:B:315:TRP:CD1	1:B:356:MET:HB3	2.46	0.49
1:A:1015:VAL:HG23	1:A:1196:TRP:CH2	2.46	0.49
1:A:1015:VAL:HG23	1:A:1196:TRP:HH2	1.77	0.49
1:B:1192:ILE:HG13	1:B:1197:HIS:CE1	2.48	0.49
1:B:77:ILE:HG22	1:B:390:LEU:HD11	1.93	0.49
1:B:1071:ARG:HH12	3:F:-2:DC:P	2.35	0.49
1:B:382:PRO:HB3	1:B:389:ASP:HB3	1.94	0.49
1:A:629:ALA:HB1	1:A:634:PHE:HB2	1.94	0.49
1:A:959:ARG:HH22	2:C:-5:DG:H21	1.59	0.49
1:B:99:LEU:HD21	1:B:101:ARG:HG2	1.95	0.49
1:B:874:ALA:HB3	1:B:896:GLY:C	2.38	0.48
1:A:124:PRO:HG3	1:A:408:HIS:CD2	2.49	0.48
1:A:184:LEU:HB2	1:A:196:VAL:HG13	1.96	0.48
1:A:444:PHE:HD1	1:A:483:TYR:CD2	2.31	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:597:GLU:HB2	1:A:805:LYS:HD2	1.95	0.48
1:A:830:CYS:SG	1:A:833:VAL:HG23	2.53	0.48
1:B:713:VAL:HG23	1:B:930:PHE:HB2	1.96	0.48
1:A:321:GLY:N	1:A:325:THR:O	2.45	0.48
1:A:325:THR:HG22	1:A:349:ASN:HD21	1.78	0.48
1:A:208:ASN:OD1	1:A:211:GLU:N	2.45	0.48
1:A:358:ASP:OD1	1:A:358:ASP:N	2.34	0.48
1:A:884:TYR:HB3	1:A:891:PHE:HB2	1.96	0.48
1:B:608:ASN:ND2	1:B:611:ARG:HG3	2.28	0.48
1:A:522:ASN:O	1:A:614:TYR:OH	2.32	0.48
1:B:423:HIS:HB2	1:B:576:GLU:OE1	2.14	0.48
1:A:153:LEU:HD11	1:A:185:LEU:HD11	1.96	0.48
1:B:187:LEU:HD23	1:B:193:ARG:HA	1.96	0.48
1:B:359:LEU:HD22	1:B:360:PRO:HD2	1.95	0.47
1:A:250:ALA:HA	1:A:270:ARG:O	2.14	0.47
1:A:1023:ASP:O	1:A:1026:ARG:HG3	2.14	0.47
1:B:1061:ARG:NH2	1:B:1144:GLU:OE2	2.44	0.47
1:A:525:MET:HE2	1:A:588:LEU:HB3	1.96	0.47
1:B:1175:ALA:HA	1:B:1179:ASN:OD1	2.15	0.47
1:B:500:ARG:HB3	1:B:502:TRP:CE3	2.49	0.47
1:A:1193:PRO:HD2	1:A:1196:TRP:CZ3	2.50	0.47
1:B:172:MET:HE3	1:B:175:ILE:HD11	1.97	0.47
1:B:998:TRP:CE3	1:B:1003:LEU:HD23	2.50	0.47
1:B:502:TRP:HE3	1:B:502:TRP:H	1.63	0.46
1:A:919:LEU:HD12	1:A:919:LEU:H	1.80	0.46
1:B:522:ASN:O	1:B:614:TYR:OH	2.26	0.46
1:B:1164:HIS:NE2	2:E:-2:DA:OP1	2.49	0.46
1:A:797:PRO:O	1:A:801:VAL:HG23	2.15	0.46
1:B:193:ARG:NH1	1:B:338:SER:O	2.48	0.46
1:B:884:TYR:HD2	1:B:891:PHE:CD2	2.34	0.46
2:C:-8:DT:H5'	2:C:-8:DT:C6	2.51	0.46
2:E:-8:DT:H5'	2:E:-8:DT:C6	2.50	0.46
1:A:129:LEU:HD11	1:A:332:PRO:O	2.16	0.46
1:A:148:HIS:O	1:A:186:GLY:HA3	2.16	0.46
1:B:813:VAL:HA	1:B:816:SER:HB3	1.97	0.46
1:B:995:ALA:HB1	1:B:1183:ILE:HG21	1.98	0.46
1:B:1160:TYR:CE1	1:B:1164:HIS:CE1	3.03	0.46
1:B:444:PHE:HB2	1:B:483:TYR:CE2	2.51	0.46
1:B:848:THR:HG23	1:B:913:ILE:HG21	1.97	0.46
1:A:65:HIS:HB3	1:A:67:TYR:CE2	2.51	0.45
1:A:1066:PRO:HG3	1:A:1072:ILE:HD11	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:824:GLN:HB3	1:B:831:LEU:HD11	1.97	0.45
1:B:244:SER:OG	1:B:245:ALA:N	2.50	0.45
1:B:969:THR:HG21	1:B:1165:LEU:HD11	1.99	0.45
1:A:77:ILE:HA	1:A:95:HIS:O	2.16	0.45
1:A:714:VAL:HG22	1:A:902:LEU:HD13	1.98	0.45
1:A:1035:ALA:HA	3:D:-2:DC:OP1	2.17	0.45
1:B:206:TYR:CE2	1:B:270:ARG:HG3	2.51	0.45
1:B:930:PHE:CE2	1:B:945:ILE:HG12	2.52	0.45
1:B:950:MET:HE3	1:B:979:TYR:HE2	1.82	0.45
1:B:1028:ILE:HG21	1:B:1150:ILE:HD12	1.98	0.45
1:A:383:VAL:HG12	1:A:385:GLY:H	1.82	0.45
3:F:-7:DT:H5'	3:F:-7:DT:H6	1.81	0.45
3:D:-3:DC:H5'	3:D:-3:DC:C6	2.51	0.45
1:A:804:ASP:O	1:A:807:GLN:HG3	2.16	0.45
1:A:444:PHE:HB2	1:A:483:TYR:CE2	2.51	0.44
1:A:454:VAL:HG13	1:A:520:MET:HE1	1.99	0.44
1:B:409:VAL:O	1:B:410:LEU:HD23	2.17	0.44
1:A:80:ARG:HG2	1:A:95:HIS:CD2	2.52	0.44
1:A:160:TYR:O	1:A:163:ARG:HG2	2.18	0.44
1:B:444:PHE:HD1	1:B:483:TYR:CD2	2.35	0.44
1:A:170:ARG:HH12	1:A:346:THR:HG22	1.82	0.44
1:A:395:SER:HB3	1:A:578:CYS:HB3	1.99	0.44
1:B:247:HIS:O	1:B:247:HIS:ND1	2.49	0.44
1:A:696:TYR:CZ	1:A:834:ALA:HB1	2.53	0.44
1:A:171:PHE:O	1:A:175:ILE:HG12	2.18	0.44
1:B:1079:GLN:HG2	1:B:1142:LEU:HD11	2.00	0.44
1:B:238:ALA:HB1	1:B:241:ARG:HD2	2.00	0.44
1:B:366:CYS:HA	1:B:463:THR:O	2.18	0.44
1:A:187:LEU:HD23	1:A:193:ARG:HA	2.00	0.44
1:B:75:ARG:HE	1:B:77:ILE:HD11	1.82	0.44
1:B:129:LEU:HD11	1:B:332:PRO:O	2.18	0.43
2:E:-9:DG:H2'	2:E:-8:DT:H71	2.00	0.43
1:A:397:LEU:HD13	1:A:582:SER:HB3	2.00	0.43
1:A:1054:VAL:HG21	1:A:1074:TYR:HB3	2.01	0.43
1:B:258:VAL:HG22	1:B:634:PHE:HB3	1.99	0.43
1:A:1012:ALA:O	1:A:1015:VAL:HG12	2.18	0.43
1:A:1164:HIS:O	1:A:1165:LEU:C	2.61	0.43
3:D:-5:DC:H1'	3:D:-4:DC:H5'	2.01	0.43
1:B:833:VAL:O	1:B:837:VAL:HG23	2.19	0.43
1:A:224:LEU:HD22	1:A:228:MET:HE2	2.01	0.43
1:A:322:ARG:HG2	1:B:771:GLU:OE2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:821:THR:HG22	1:A:834:ALA:HB2	2.00	0.43
1:B:459:PRO:HG2	1:B:520:MET:HE2	2.00	0.43
1:A:99:LEU:HD12	1:A:100:LYS:H	1.84	0.43
1:A:187:LEU:HG	1:A:362:TYR:CE2	2.53	0.43
1:A:555:LEU:HD13	1:A:556:SER:N	2.34	0.43
1:A:732:CYS:HB2	1:A:772:SER:OG	2.19	0.43
1:B:609:ILE:O	1:B:613:ILE:HG13	2.19	0.43
1:B:862:LEU:HD11	1:B:909:MET:HE2	2.00	0.43
1:A:234:GLU:H	1:A:234:GLU:HG2	1.65	0.43
1:A:898:THR:C	1:A:900:ALA:H	2.27	0.43
3:D:-3:DC:H5'	3:D:-3:DC:H6	1.84	0.43
1:B:157:GLU:O	1:B:175:ILE:HB	2.19	0.43
1:B:393:GLN:HA	1:B:412:PHE:O	2.19	0.43
1:B:218:CYS:HB2	1:B:223:ASP:CB	2.49	0.42
1:B:426:GLU:HG2	1:B:430:ARG:HD2	2.01	0.42
1:A:310:PHE:HA	1:A:352:ILE:HD11	2.01	0.42
1:A:199:TYR:OH	1:A:330:ARG:NH2	2.52	0.42
2:C:-11:DT:O2	2:C:-11:DT:H2'	2.18	0.42
1:A:398:LEU:HD22	1:A:457:TYR:CE2	2.55	0.42
1:A:788:ILE:HD12	1:A:806:GLN:HB3	2.00	0.42
1:B:230:ALA:O	1:B:234:GLU:OE2	2.37	0.42
1:A:80:ARG:HG2	1:A:95:HIS:NE2	2.35	0.42
1:B:249:GLU:HB3	1:B:272:TYR:HB2	2.01	0.42
1:A:611:ARG:HD2	1:A:615:ASP:OD1	2.20	0.42
1:A:732:CYS:SG	1:A:733:PHE:N	2.93	0.42
1:A:1189:LYS:HB3	1:A:1197:HIS:NE2	2.35	0.42
1:B:525:MET:HE2	1:B:588:LEU:HB3	2.01	0.42
1:B:65:HIS:HB3	1:B:67:TYR:CE2	2.55	0.42
1:B:905:MET:HG2	1:B:909:MET:HE3	2.02	0.42
1:B:965:PHE:HD1	1:B:968:ARG:HH21	1.68	0.42
1:B:1182:LYS:HG2	1:B:1183:ILE:HD12	2.02	0.42
1:A:1041:PRO:HA	1:A:1044:TYR:HD1	1.85	0.42
1:B:423:HIS:CG	1:B:576:GLU:HG3	2.54	0.42
1:B:927:GLU:C	1:B:928:LYS:HG3	2.44	0.42
1:B:956:ASP:OD2	1:B:962:ASN:ND2	2.41	0.42
1:A:153:LEU:HB2	1:A:183:THR:HB	2.02	0.42
1:A:1036:GLU:HA	1:A:1071:ARG:HA	2.01	0.42
1:B:315:TRP:CD2	1:B:356:MET:HG2	2.55	0.42
1:B:884:TYR:HB3	1:B:891:PHE:HB2	2.01	0.42
1:A:609:ILE:O	1:A:613:ILE:HG13	2.19	0.41
1:B:386:HIS:HB2	1:B:389:ASP:OD1	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1045:THR:HG23	3:F:-4:DC:P	2.60	0.41
1:A:103:PRO:HA	1:A:445:GLU:OE1	2.20	0.41
1:A:233:ARG:C	1:A:235:SER:H	2.29	0.41
2:C:-6:DG:H2''	2:C:-5:DG:H5'	2.03	0.41
1:B:559:ASP:O	1:B:563:TYR:HD1	2.04	0.41
1:B:868:GLU:O	1:B:872:MET:HE3	2.20	0.41
1:B:1011:GLY:O	1:B:1015:VAL:N	2.42	0.41
1:B:414:LEU:HD23	1:B:414:LEU:HA	1.90	0.41
1:A:714:VAL:HB	1:A:892:VAL:HG23	2.01	0.41
1:B:326:LEU:HD13	1:B:326:LEU:HA	1.90	0.41
1:A:931:THR:HG23	1:A:946:TYR:HA	2.02	0.41
1:B:993:ARG:CB	1:B:998:TRP:HE1	2.33	0.41
1:A:959:ARG:HE	1:A:959:ARG:HB3	1.67	0.41
1:B:187:LEU:HG	1:B:362:TYR:CE2	2.55	0.41
1:A:126:ARG:HH22	1:A:460:GLU:CD	2.28	0.41
1:A:253:VAL:HG21	1:A:255:ARG:HH21	1.86	0.41
1:A:987:ALA:HA	1:A:990:LEU:HD12	2.03	0.41
1:B:882:ILE:HG13	1:B:892:VAL:HG22	2.02	0.41
1:B:1180:ASN:ND2	1:B:1183:ILE:HD13	2.35	0.41
1:A:76:PHE:CE2	1:A:442:SER:HB3	2.56	0.41
1:A:319:LYS:HE3	1:A:348:ASP:O	2.21	0.41
1:A:966:ILE:H	1:A:966:ILE:HG13	1.56	0.41
1:B:63:GLN:OE1	1:B:496:ARG:NH2	2.53	0.41
1:B:557:TYR:HA	1:B:560:ILE:HG12	2.03	0.41
1:B:748:ALA:HA	1:B:752:TYR:CE1	2.56	0.41
1:A:184:LEU:HB2	1:A:196:VAL:CG1	2.51	0.40
1:B:255:ARG:HG2	1:B:635:ILE:HG23	2.03	0.40
1:B:124:PRO:HG3	1:B:408:HIS:CD2	2.55	0.40
1:B:580:GLN:O	1:B:584:LEU:HD23	2.22	0.40
1:B:196:VAL:HG22	1:B:343:PHE:HB2	2.03	0.40
1:B:785:ARG:NH2	1:B:811:LYS:HD2	2.36	0.40
1:B:881:ARG:HG3	1:B:881:ARG:HH11	1.86	0.40
1:B:732:CYS:HB2	1:B:772:SER:OG	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	1029/1163 (88%)	962 (94%)	66 (6%)	1 (0%)	48	79
1	B	1036/1163 (89%)	971 (94%)	58 (6%)	7 (1%)	18	52
All	All	2065/2326 (89%)	1933 (94%)	124 (6%)	8 (0%)	30	63

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	176	THR
1	B	900	ALA
1	B	946	TYR
1	B	217	GLN
1	B	177	PRO
1	B	416	SER
1	B	757	VAL
1	B	360	PRO

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	779/951 (82%)	722 (93%)	57 (7%)	13	38
1	B	819/951 (86%)	771 (94%)	48 (6%)	18	43
All	All	1598/1902 (84%)	1493 (93%)	105 (7%)	15	41

All (105) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	94	VAL
1	A	101	ARG
1	A	117	VAL
1	A	176	THR
1	A	178	THR
1	A	196	VAL
1	A	224	LEU
1	A	239	SER
1	A	240	PHE
1	A	258	VAL
1	A	285	ASN
1	A	287	CYS
1	A	322	ARG
1	A	358	ASP
1	A	397	LEU
1	A	398	LEU
1	A	424	LEU
1	A	426	GLU
1	A	515	ILE
1	A	555	LEU
1	A	607	ILE
1	A	609	ILE
1	A	694	VAL
1	A	746	LEU
1	A	750	LYS
1	A	765	VAL
1	A	821	THR
1	A	827	LEU
1	A	845	LEU
1	A	846	LEU
1	A	881	ARG
1	A	890	ILE
1	A	897	LEU
1	A	919	LEU
1	A	924	LEU
1	A	935	LEU
1	A	938	LYS
1	A	950	MET
1	A	951	LEU
1	A	956	ASP
1	A	963	CYS
1	A	966	ILE
1	A	975	ASP

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Mol	Chain	Res	Type
1	A	982	THR
1	A	999	LEU
1	A	1008	GLN
1	A	1039	ARG
1	A	1045	THR
1	A	1062	ARG
1	A	1064	GLN
1	A	1070	ASP
1	A	1075	VAL
1	A	1138	LEU
1	A	1139	VAL
1	A	1159	ASP
1	A	1171	VAL
1	A	1183	ILE
1	B	45	TYR
1	B	146	VAL
1	B	170	ARG
1	B	172	MET
1	B	176	THR
1	B	216	LEU
1	B	217	GLN
1	B	227	ARG
1	B	240	PHE
1	B	246	ASP
1	B	252	VAL
1	B	255	ARG
1	B	262	GLU
1	B	285	ASN
1	B	356	MET
1	B	388	GLU
1	B	392	ILE
1	B	397	LEU
1	B	411	LEU
1	B	500	ARG
1	B	534	LYS
1	B	603	ARG
1	B	607	ILE
1	B	619	ILE
1	B	694	VAL
1	B	761	ARG
1	B	821	THR
1	B	827	LEU

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Mol	Chain	Res	Type
1	B	845	LEU
1	B	881	ARG
1	B	890	ILE
1	B	919	LEU
1	B	928	LYS
1	B	935	LEU
1	B	952	ILE
1	B	956	ASP
1	B	973	LEU
1	B	975	ASP
1	B	983	VAL
1	B	999	LEU
1	B	1029	GLN
1	B	1030	ASP
1	B	1045	THR
1	B	1048	ARG
1	B	1058	LEU
1	B	1141	GLU
1	B	1158	THR
1	B	1182	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	168	HIS
1	A	192	HIS
1	A	408	HIS
1	A	587	GLN
1	A	745	HIS
1	A	832	HIS
1	B	587	GLN
1	B	693	HIS
1	B	745	HIS
1	B	787	GLN
1	B	832	HIS
1	B	1197	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

2 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	$\# Z > 2$
4	A1AC6	E	4000	-	28,28,28	0.14	0	38,38,38	0.33	0
4	A1AC6	C	4001	-	28,28,28	0.13	0	38,38,38	0.35	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
4	A1AC6	E	4000	-	-	2/14/23/23	0/3/3/3
4	A1AC6	C	4001	-	-	1/14/23/23	0/3/3/3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	E	4000	A1AC6	C1-C2-C3-C4
4	E	4000	A1AC6	N-C2-C3-C4

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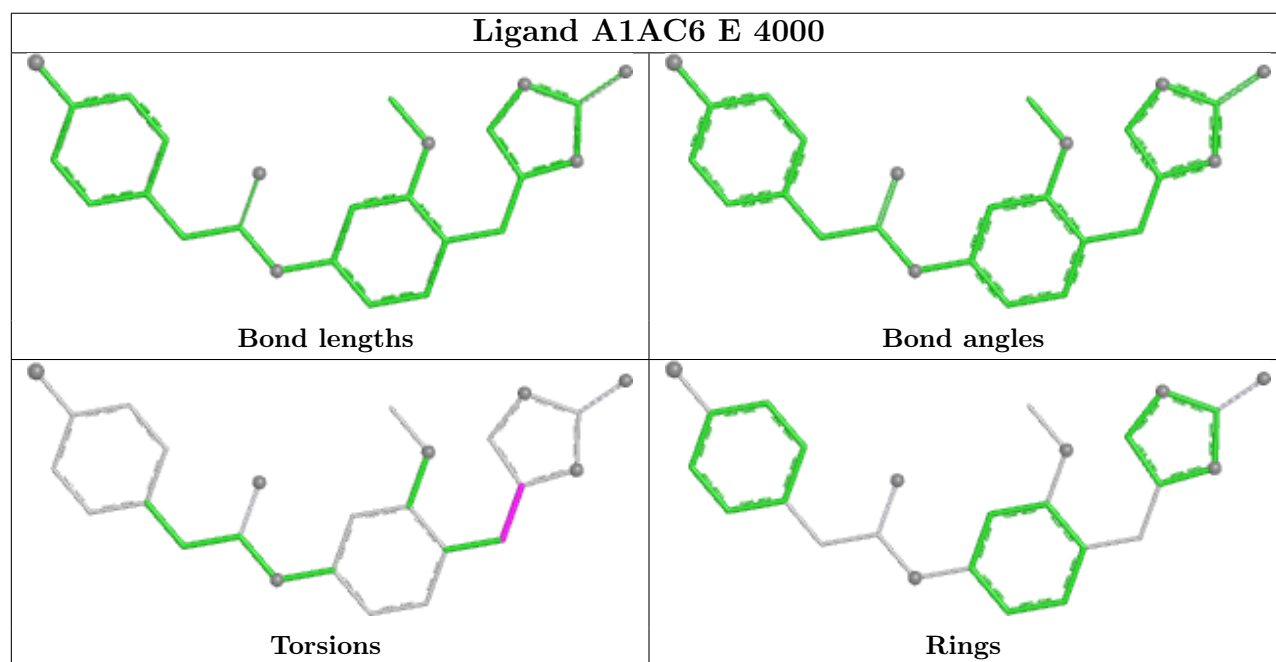
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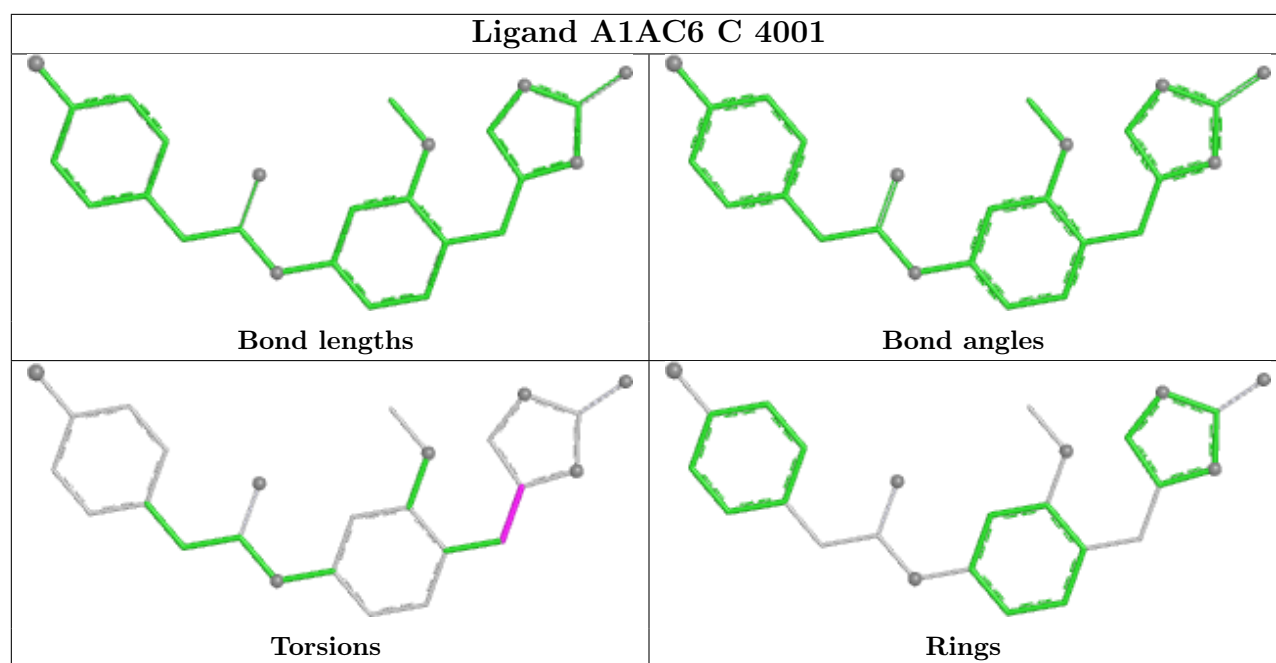
Mol	Chain	Res	Type	Atoms
4	C	4001	A1AC6	C1-C2-C3-C4

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

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	1037/1163 (89%)	0.12	25 (2%) 59 42	57, 123, 174, 223	0
1	B	1046/1163 (89%)	0.16	26 (2%) 58 41	65, 127, 170, 210	0
2	C	15/15 (100%)	0.57	1 (6%) 24 21	107, 149, 227, 241	0
2	E	14/15 (93%)	0.54	0 100 100	115, 146, 230, 239	0
3	D	13/13 (100%)	0.51	1 (7%) 19 18	115, 138, 236, 245	0
3	F	13/13 (100%)	0.35	0 100 100	118, 139, 228, 246	0
All	All	2138/2382 (89%)	0.15	53 (2%) 58 41	57, 126, 174, 246	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	503	ASP	6.4
1	B	1159	ASP	3.9
1	B	242	GLY	3.8
1	B	503	ASP	3.7
1	B	1140	SER	3.6
1	A	443	GLU	3.6
2	C	-11	DT	3.5
1	A	378	GLU	3.4
1	A	558	ARG	3.1
1	A	1095	GLU	2.8
1	B	418	ASP	2.8
1	A	1020	ARG	2.7
1	A	1048	ARG	2.7
1	B	1029	GLN	2.7
1	B	1046	ASN	2.6
1	B	384	ALA	2.5
1	A	954	GLY	2.5
1	A	1008	GLN	2.5
1	B	122	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	870	ALA	2.4
1	B	511	LYS	2.4
1	A	705	THR	2.4
1	B	1141	GLU	2.4
1	A	1139	VAL	2.3
1	B	241	ARG	2.3
1	B	120	GLY	2.3
1	B	569	ALA	2.3
1	B	1010	PHE	2.3
1	B	132	GLY	2.3
1	A	1047	LYS	2.3
1	A	1182	LYS	2.3
1	A	876	GLY	2.2
1	A	418	ASP	2.2
1	B	277	ARG	2.2
1	A	694	VAL	2.2
1	B	1048	ARG	2.2
1	B	552	LYS	2.2
1	B	1067	SER	2.2
1	B	88	PRO	2.2
1	A	999	LEU	2.2
3	D	-11	DA	2.1
1	A	939	LYS	2.1
1	A	491	GLY	2.1
1	A	552	LYS	2.1
1	B	875	PRO	2.1
1	A	875	PRO	2.0
1	A	1141	GLU	2.0
1	B	1076	ILE	2.0
1	B	1084	GLU	2.0
1	A	1009	ALA	2.0
1	B	932	LYS	2.0
1	A	1158	THR	2.0
1	A	1177	PHE	2.0

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

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](#)

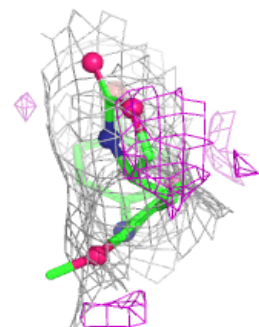
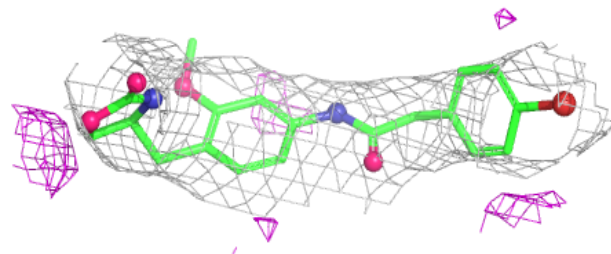
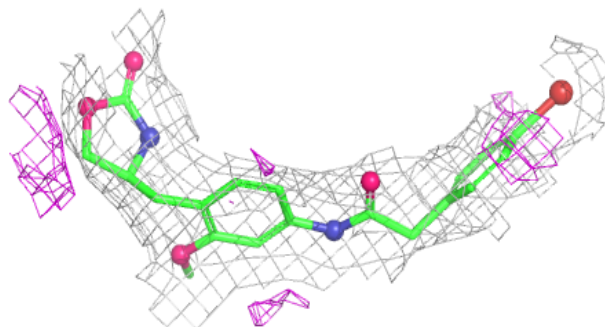
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists 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
4	A1AC6	C	4001	26/26	0.94	0.12	93,111,122,128	0
4	A1AC6	E	4000	26/26	0.96	0.11	97,113,120,133	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.

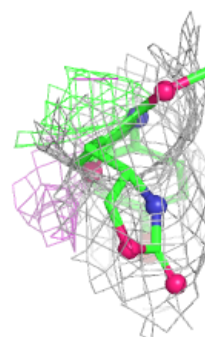
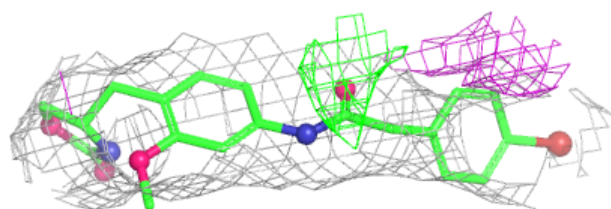
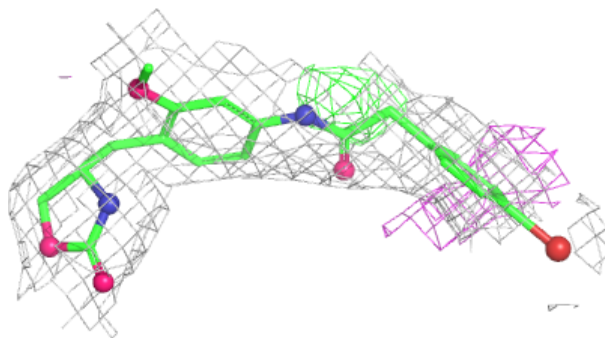
Electron density around A1AC6 C 4001:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around A1AC6 E 4000:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.