



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:52 PM UTC

PDB ID : 3QX3 / pdb_00003qx3
Title : Human topoisomerase IIbeta in complex with DNA and etoposide
Authors : Wu, C.C.; Li, T.K.; Farh, L.; Lin, L.Y.; Lin, T.S.; Yu, Y.J.; Yen, T.J.; Chiang, C.W.; Chan, N.L.
Deposited on : 2011-03-01
Resolution : 2.16 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

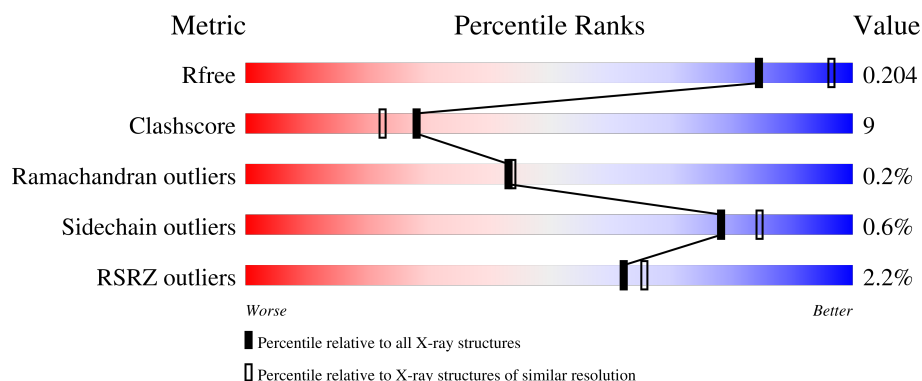
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.16 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	803	2% 69% 15% 16%
1	B	803	% 68% 15% 16%
2	C	8	88% 12%
2	E	8	100%
3	D	12	67% 33%

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Mol	Chain	Length	Quality of chain
3	F	12	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: green (58%), yellow (33%), and orange (8%). The percentages are labeled below the segments.

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 13012 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 topoisomerase 2-beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	671	Total	C	N	O	S	0	21	0
			5536	3535	956	1021	24			
1	B	671	Total	C	N	O	S	6	18	0
			5519	3526	956	1013	24			

There are 92 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	419	MET	-	expression tag	UNP Q02880
A	420	ALA	-	expression tag	UNP Q02880
A	421	SER	-	expression tag	UNP Q02880
A	422	TRP	-	expression tag	UNP Q02880
A	423	SER	-	expression tag	UNP Q02880
A	424	HIS	-	expression tag	UNP Q02880
A	425	PRO	-	expression tag	UNP Q02880
A	426	GLN	-	expression tag	UNP Q02880
A	427	PHE	-	expression tag	UNP Q02880
A	428	GLU	-	expression tag	UNP Q02880
A	429	LYS	-	expression tag	UNP Q02880
A	430	GLY	-	expression tag	UNP Q02880
A	431	ALA	-	expression tag	UNP Q02880
A	432	ASP	-	expression tag	UNP Q02880
A	433	ASP	-	expression tag	UNP Q02880
A	434	ASP	-	expression tag	UNP Q02880
A	435	ASP	-	expression tag	UNP Q02880
A	436	LYS	-	expression tag	UNP Q02880
A	437	VAL	-	expression tag	UNP Q02880
A	438	PRO	-	expression tag	UNP Q02880
A	439	ASP	-	expression tag	UNP Q02880
A	440	PRO	-	expression tag	UNP Q02880
A	441	THR	-	expression tag	UNP Q02880
A	442	SER	-	expression tag	UNP Q02880
A	443	VAL	-	expression tag	UNP Q02880

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Chain	Residue	Modelled	Actual	Comment	Reference
A	444	ASP	-	expression tag	UNP Q02880
A	1202	GLY	-	expression tag	UNP Q02880
A	1203	ALA	-	expression tag	UNP Q02880
A	1204	PRO	-	expression tag	UNP Q02880
A	1205	GLY	-	expression tag	UNP Q02880
A	1206	PHE	-	expression tag	UNP Q02880
A	1207	SER	-	expression tag	UNP Q02880
A	1208	SER	-	expression tag	UNP Q02880
A	1209	ILE	-	expression tag	UNP Q02880
A	1210	SER	-	expression tag	UNP Q02880
A	1211	ALA	-	expression tag	UNP Q02880
A	1212	HIS	-	expression tag	UNP Q02880
A	1213	HIS	-	expression tag	UNP Q02880
A	1214	HIS	-	expression tag	UNP Q02880
A	1215	HIS	-	expression tag	UNP Q02880
A	1216	HIS	-	expression tag	UNP Q02880
A	1217	HIS	-	expression tag	UNP Q02880
A	1218	HIS	-	expression tag	UNP Q02880
A	1219	HIS	-	expression tag	UNP Q02880
A	1220	HIS	-	expression tag	UNP Q02880
A	1221	HIS	-	expression tag	UNP Q02880
B	419	MET	-	expression tag	UNP Q02880
B	420	ALA	-	expression tag	UNP Q02880
B	421	SER	-	expression tag	UNP Q02880
B	422	TRP	-	expression tag	UNP Q02880
B	423	SER	-	expression tag	UNP Q02880
B	424	HIS	-	expression tag	UNP Q02880
B	425	PRO	-	expression tag	UNP Q02880
B	426	GLN	-	expression tag	UNP Q02880
B	427	PHE	-	expression tag	UNP Q02880
B	428	GLU	-	expression tag	UNP Q02880
B	429	LYS	-	expression tag	UNP Q02880
B	430	GLY	-	expression tag	UNP Q02880
B	431	ALA	-	expression tag	UNP Q02880
B	432	ASP	-	expression tag	UNP Q02880
B	433	ASP	-	expression tag	UNP Q02880
B	434	ASP	-	expression tag	UNP Q02880
B	435	ASP	-	expression tag	UNP Q02880
B	436	LYS	-	expression tag	UNP Q02880
B	437	VAL	-	expression tag	UNP Q02880
B	438	PRO	-	expression tag	UNP Q02880
B	439	ASP	-	expression tag	UNP Q02880

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Chain	Residue	Modelled	Actual	Comment	Reference
B	440	PRO	-	expression tag	UNP Q02880
B	441	THR	-	expression tag	UNP Q02880
B	442	SER	-	expression tag	UNP Q02880
B	443	VAL	-	expression tag	UNP Q02880
B	444	ASP	-	expression tag	UNP Q02880
B	1202	GLY	-	expression tag	UNP Q02880
B	1203	ALA	-	expression tag	UNP Q02880
B	1204	PRO	-	expression tag	UNP Q02880
B	1205	GLY	-	expression tag	UNP Q02880
B	1206	PHE	-	expression tag	UNP Q02880
B	1207	SER	-	expression tag	UNP Q02880
B	1208	SER	-	expression tag	UNP Q02880
B	1209	ILE	-	expression tag	UNP Q02880
B	1210	SER	-	expression tag	UNP Q02880
B	1211	ALA	-	expression tag	UNP Q02880
B	1212	HIS	-	expression tag	UNP Q02880
B	1213	HIS	-	expression tag	UNP Q02880
B	1214	HIS	-	expression tag	UNP Q02880
B	1215	HIS	-	expression tag	UNP Q02880
B	1216	HIS	-	expression tag	UNP Q02880
B	1217	HIS	-	expression tag	UNP Q02880
B	1218	HIS	-	expression tag	UNP Q02880
B	1219	HIS	-	expression tag	UNP Q02880
B	1220	HIS	-	expression tag	UNP Q02880
B	1221	HIS	-	expression tag	UNP Q02880

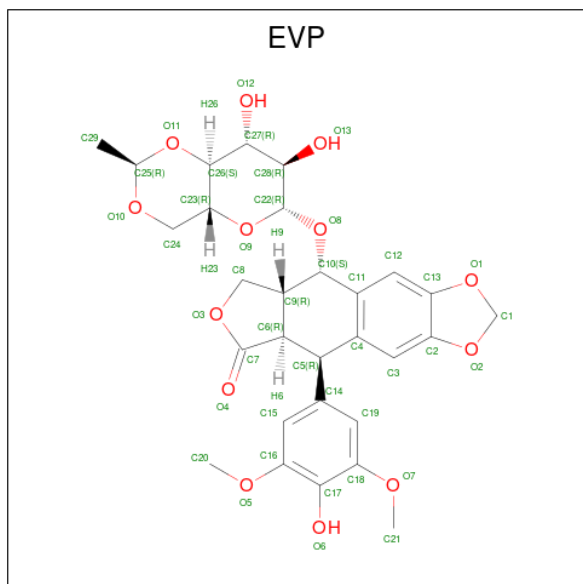
- Molecule 2 is a DNA chain called DNA (5'-D(P*AP*GP*CP*CP*GP*AP*GP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	8	Total 165	C 77	N 34	O 46	P 8	0	0	0
2	E	8	Total 165	C 77	N 34	O 46	P 8	0	0	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	12	Total 245	C 116	N 43	O 74	P 12	0	0	0
3	F	12	Total 245	C 116	N 43	O 74	P 12	0	0	0

- Molecule 4 is (5S,5aR,8aR,9R)-9-(4-hydroxy-3,5-dimethoxyphenyl)-8-oxo-5,5a,6,8,8a,9-hexahydrofuro[3',4':6,7]naphtho[2,3-d][1,3]dioxol -5-yl 4,6-O-[(1R)-ethylidene]-beta-D-glucopyranoside (CCD ID: EVP) (formula: C₂₉H₃₂O₁₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			42	29	13		
4	D	1	Total	C	O	0	0
			42	29	13		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Mg	0	0
			2	2		
5	B	2	Total	Mg	0	0
			2	2		
5	D	1	Total	Mg	0	0
			1	1		
5	F	1	Total	Mg	0	0
			1	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	422	Total	O	0	0
			422	422		
6	B	462	Total	O	0	0
			462	462		

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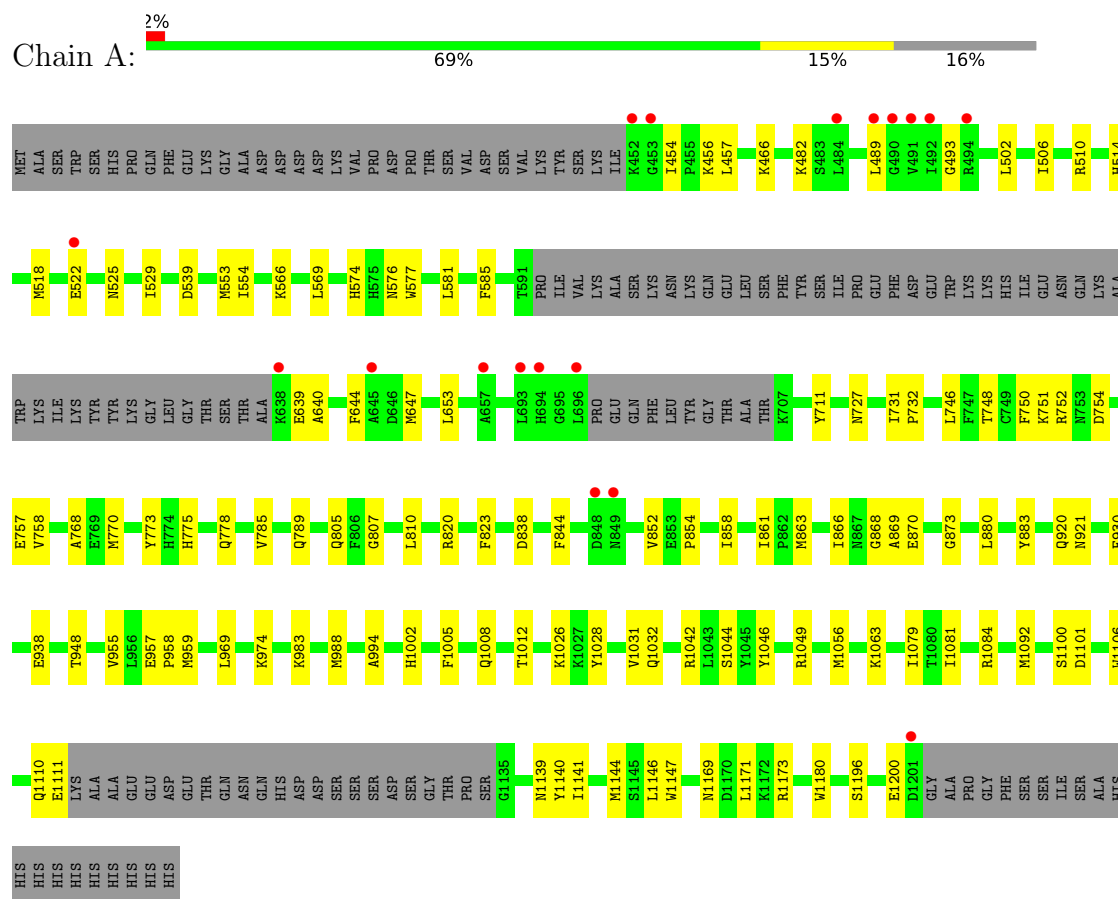
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	29	Total 29	O 29	0	0
6	D	48	Total 48	O 48	0	0
6	E	37	Total 37	O 37	0	0
6	F	49	Total 49	O 49	0	0

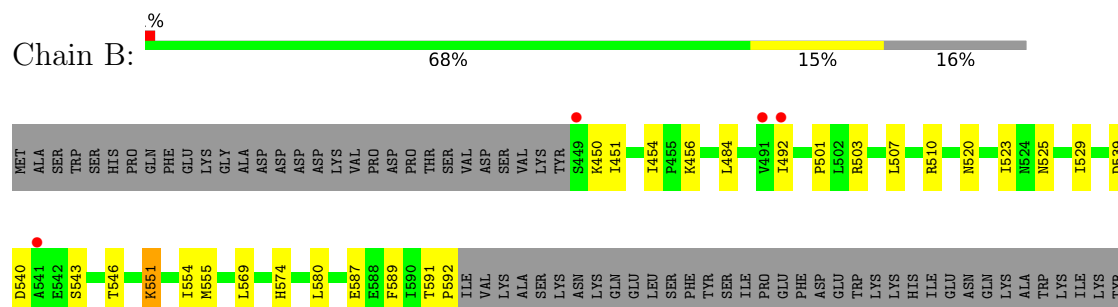
3 Residue-property plots [i](#)

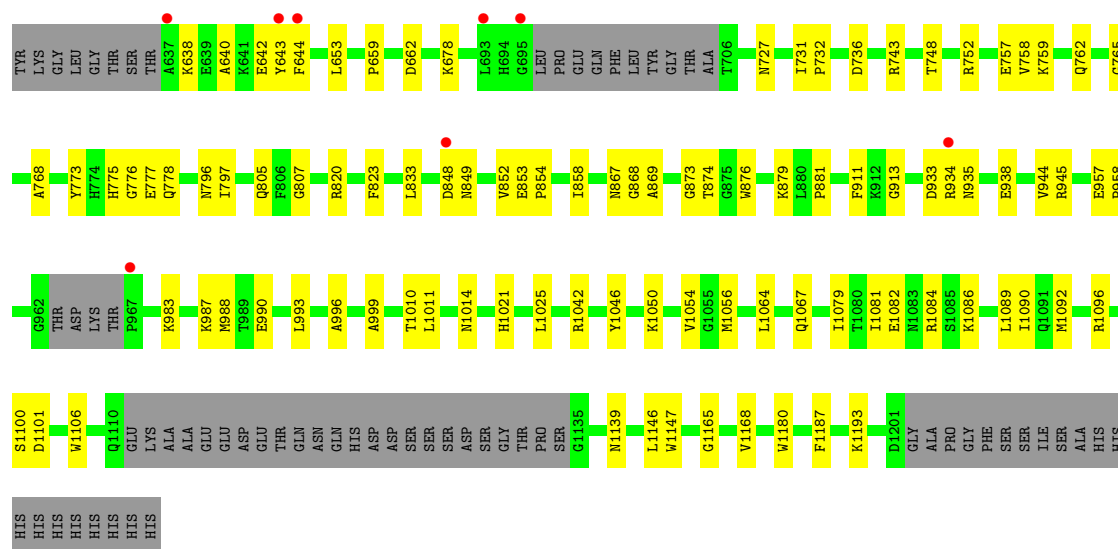
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

• Molecule 1: DNA topoisomerase 2-beta



• Molecule 1: DNA topoisomerase 2-beta





- Molecule 2: DNA (5'-D(P*AP*GP*CP*CP*GP*AP*GP*C)-3')

Chain C: 88% 12%



- Molecule 2: DNA (5'-D(P*AP*GP*CP*CP*GP*AP*GP*C)-3')

Chain E: 100%

There are no outlier residues recorded for this chain.

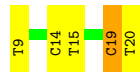
- Molecule 3: DNA (5'-D(P*TP*GP*CP*AP*GP*CP*TP*CP*GP*GP*CP*T)-3')

Chain D: 67% 33%



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

Chain F: 58% 33% 8%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	80.15Å 176.80Å 94.07Å 90.00° 111.58° 90.00°	Depositor
Resolution (Å)	29.54 – 2.16 29.54 – 2.16	Depositor EDS
% Data completeness (in resolution range)	90.8 (29.54-2.16) 90.7 (29.54-2.16)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.61 (at 2.16Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.168 , 0.207 0.167 , 0.204	Depositor DCC
R_{free} test set	6234 reflections (4.81%)	wwPDB-VP
Wilson B-factor (Å ²)	32.1	Xtriage
Anisotropy	0.081	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 43.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13012	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.00% 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: MG, EVP

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	0.47	0/5699	0.73	2/7672 (0.0%)
1	B	0.48	0/5676	0.73	2/7638 (0.0%)
2	C	0.30	0/185	0.91	0/283
2	E	0.35	0/185	0.90	0/283
3	D	0.35	0/273	0.99	1/419 (0.2%)
3	F	0.32	0/273	1.00	1/419 (0.2%)
All	All	0.47	0/12291	0.75	6/16714 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	19	DC	P-O5'-C5'	-5.97	111.05	120.00
3	D	18	DG	C4'-C3'-O3'	-5.29	102.06	110.00
1	B	454	ILE	CA-C-N	5.23	125.28	119.32
1	B	454	ILE	C-N-CA	5.23	125.28	119.32
1	A	454	ILE	CA-C-N	5.21	125.27	119.32
1	A	454	ILE	C-N-CA	5.21	125.27	119.32

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5536	0	5635	88	0
1	B	5519	0	5623	118	0
2	C	165	0	89	7	0
2	E	165	0	89	0	0
3	D	245	0	136	4	0
3	F	245	0	136	8	0
4	A	42	0	30	0	0
4	D	42	0	30	1	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
5	D	1	0	0	0	0
5	F	1	0	0	0	0
6	A	422	0	0	12	0
6	B	462	0	0	14	0
6	C	29	0	0	0	0
6	D	48	0	0	1	0
6	E	37	0	0	0	0
6	F	49	0	0	0	0
All	All	13012	0	11768	211	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 (211) 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:501:PRO:HG2	1:B:503[B]:ARG:HH22	1.27	0.98
1:A:805:GLN:HE21	1:A:807:GLY:H	1.09	0.96
1:B:456:LYS:HD3	1:B:503[B]:ARG:HH21	1.31	0.95
1:B:456:LYS:HD3	1:B:503[B]:ARG:NH2	1.82	0.95
1:B:510:ARG:HH22	1:B:574:HIS:HD2	1.13	0.89
2:C:1:DA:H2	3:F:20:DT:H3	1.22	0.87
1:A:510:ARG:HH22	1:A:574:HIS:HD2	1.24	0.86
1:B:988:MET:HE3	1:B:993:LEU:HD13	1.58	0.86
1:B:944:VAL:HG12	1:B:945:ARG:HG2	1.61	0.83
1:B:805:GLN:HE21	1:B:807:GLY:H	1.27	0.82
1:A:938:GLU:OE2	1:A:983:LYS:HE3	1.80	0.82
1:B:797:ILE:H	1:B:867:ASN:HD21	1.26	0.81
1:A:805:GLN:NE2	1:A:807:GLY:H	1.77	0.81
1:B:868:GLY:HA2	6:B:1439:HOH:O	1.83	0.78
1:B:1096:ARG:HD3	6:B:1464:HOH:O	1.84	0.77
1:B:501:PRO:HG2	1:B:503[B]:ARG:NH2	2.01	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:1:DA:C8	2:C:1:DA:H5'	2.21	0.75
1:B:501:PRO:HB2	1:B:503[B]:ARG:HH12	1.53	0.74
1:B:765:GLY:HA2	1:B:777:GLU:OE2	1.90	0.72
1:A:868:GLY:HA2	6:A:1432:HOH:O	1.90	0.72
1:B:776:GLY:HA3	1:B:778[B]:GLN:NE2	2.05	0.72
1:A:778[B]:GLN:HE22	3:D:9:DT:P	2.13	0.72
1:B:456:LYS:CD	1:B:503[B]:ARG:HH21	2.03	0.71
1:A:1079:ILE:HD11	1:A:1092:MET:HE3	1.72	0.71
1:B:1056:MET:HE2	1:B:1106:TRP:CZ3	2.26	0.70
2:C:1:DA:H8	2:C:1:DA:P	2.14	0.70
1:B:456:LYS:HE2	6:B:1471:HOH:O	1.92	0.69
1:B:1086:LYS:O	1:B:1090:ILE:HG12	1.92	0.69
1:A:1026[A]:LYS:HE2	6:A:1455:HOH:O	1.93	0.69
1:B:510:ARG:HH22	1:B:574:HIS:CD2	2.05	0.68
1:A:868:GLY:CA	6:A:1432:HOH:O	2.41	0.68
1:B:752:ARG:HH12	1:B:758:VAL:HG21	1.58	0.68
1:A:553:MET:HE1	1:A:647:MET:HE3	1.76	0.66
1:B:938:GLU:OE2	1:B:983:LYS:HE3	1.96	0.66
1:B:554[B]:ILE:CD1	1:B:569:LEU:HD23	2.27	0.65
1:B:554[B]:ILE:HG12	1:B:569:LEU:HD23	1.78	0.65
1:B:1079:ILE:HD11	1:B:1092:MET:HE3	1.79	0.65
1:B:778[B]:GLN:H	1:B:778[B]:GLN:CD	2.05	0.65
1:B:554[B]:ILE:HG12	1:B:569:LEU:CD2	2.26	0.64
1:B:501:PRO:CG	1:B:503[B]:ARG:HH22	2.05	0.64
1:A:868:GLY:N	6:A:1432:HOH:O	2.31	0.64
1:B:868:GLY:CA	6:B:1439:HOH:O	2.43	0.63
1:A:861:ILE:HD12	1:A:866[B]:ILE:HD11	1.81	0.62
1:B:935:ASN:HA	1:B:988:MET:O	1.99	0.62
2:C:1:DA:H5'	2:C:1:DA:H8	1.66	0.61
1:B:879:LYS:HE2	6:D:118:HOH:O	2.00	0.61
1:A:732:PRO:HG2	1:A:869:ALA:HB1	1.82	0.61
1:A:554[A]:ILE:HD13	1:A:566:LYS:HA	1.82	0.60
1:A:554[B]:ILE:HD11	1:A:569:LEU:HD12	1.82	0.60
1:B:993:LEU:HD23	6:B:1251:HOH:O	2.00	0.60
1:A:820[A]:ARG:NH2	3:F:9:DT:OP1	2.35	0.60
1:B:520:ASN:HB3	1:B:523[B]:ILE:HG12	1.84	0.60
1:A:639:GLU:H	1:A:639:GLU:CD	2.10	0.59
1:A:930[B]:PHE:CZ	1:A:938:GLU:OE1	2.55	0.59
1:B:757:GLU:H	1:B:757:GLU:CD	2.09	0.59
1:B:805:GLN:NE2	1:B:807:GLY:H	1.98	0.59
1:B:555:MET:HE1	1:B:644:PHE:HZ	1.67	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:555:MET:HE1	1:B:644:PHE:CZ	2.38	0.58
1:A:1056:MET:HE2	1:A:1106:TRP:CZ3	2.38	0.58
1:A:778[A]:GLN:HG3	3:D:9:DT:H71	1.85	0.58
1:B:934:ARG:HB3	1:B:990:GLU:HG2	1.85	0.58
1:B:1081:ILE:HA	1:B:1084:ARG:HD3	1.85	0.58
1:A:510:ARG:HH22	1:A:574:HIS:CD2	2.13	0.57
1:A:1139:ASN:HD22	1:B:1147:TRP:HE1	1.51	0.57
1:B:1089:LEU:HD12	1:B:1092:MET:HE2	1.87	0.56
2:C:1:DA:H2	3:F:20:DT:N3	1.99	0.56
1:A:525:ASN:O	1:A:529:ILE:HG13	2.05	0.56
1:A:457:LEU:HD22	1:A:529:ILE:HG12	1.88	0.56
1:B:934:ARG:O	1:B:935:ASN:HB2	2.05	0.56
1:B:732:PRO:HG2	1:B:869:ALA:HB1	1.88	0.55
1:A:1169:ASN:HB3	1:A:1173[A]:ARG:HH12	1.70	0.55
1:A:1169:ASN:O	1:A:1173[A]:ARG:HG2	2.06	0.55
1:A:522:GLU:CD	6:A:1236:HOH:O	2.49	0.54
1:B:591:THR:HB	1:B:592:PRO:CD	2.37	0.54
1:A:1081:ILE:HA	1:A:1084:ARG:HD3	1.90	0.54
1:B:540:ASP:HB2	1:B:543:SER:H	1.71	0.54
1:A:883:TYR:CZ	1:A:1031[A]:VAL:HG21	2.42	0.54
1:A:1141:ILE:HA	1:A:1144:MET:HE2	1.90	0.54
1:B:748:THR:O	1:B:752:ARG:HG3	2.07	0.54
1:A:858:ILE:HG13	1:A:1042:ARG:HD2	1.90	0.53
1:A:514:HIS:O	1:A:518:MET:HG2	2.09	0.53
1:A:820[B]:ARG:NH2	1:B:777:GLU:OE1	2.42	0.53
1:B:554[B]:ILE:CG1	1:B:569:LEU:HD23	2.38	0.53
1:B:778[A]:GLN:HG3	3:F:9:DT:H71	1.90	0.53
1:B:510:ARG:NH2	1:B:574:HIS:HD2	1.94	0.52
2:C:1:DA:C2	3:F:20:DT:N3	2.64	0.52
1:A:1056:MET:HE2	1:A:1106:TRP:HZ3	1.75	0.52
1:A:994:ALA:HB2	6:A:1441:HOH:O	2.10	0.52
1:B:1010:THR:HB	6:B:1234:HOH:O	2.09	0.51
1:A:585:PHE:N	6:A:1422:HOH:O	2.44	0.51
1:A:748:THR:O	1:A:752:ARG:HG3	2.11	0.51
1:A:581:LEU:HD22	1:A:711:TYR:OH	2.11	0.51
1:B:587:GLU:HG2	1:B:653[B]:LEU:HD22	1.92	0.51
1:B:957:GLU:HB2	1:B:958:PRO:HD3	1.91	0.51
1:A:1026[A]:LYS:HD3	1:A:1028:TYR:CZ	2.46	0.50
1:B:1100:SER:O	1:B:1101:ASP:C	2.54	0.50
1:A:1147:TRP:HE1	1:B:1139:ASN:HD22	1.60	0.50
1:A:778[A]:GLN:HG3	3:D:9:DT:C7	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1196:SER:O	1:A:1200:GLU:HG2	2.12	0.50
1:B:1165:GLY:O	1:B:1168:VAL:HG12	2.11	0.50
1:B:520:ASN:HB3	1:B:523[B]:ILE:CG1	2.41	0.50
1:A:752:ARG:NH1	1:A:754:ASP:OD2	2.45	0.50
1:A:820[A]:ARG:NH2	3:F:9:DT:H71	2.27	0.49
1:B:911:PHE:CZ	1:B:913:GLY:HA3	2.48	0.49
1:A:1100:SER:O	1:A:1101:ASP:C	2.54	0.49
1:B:1193[A]:LYS:HE2	6:B:1256:HOH:O	2.11	0.49
1:B:820[B]:ARG:NH2	6:B:1371:HOH:O	2.45	0.49
1:B:934:ARG:O	1:B:935:ASN:CB	2.60	0.49
1:A:863:MET:SD	1:A:866[B]:ILE:HD12	2.52	0.49
1:B:881:PRO:HB3	1:B:1011:LEU:HD21	1.94	0.49
1:A:861:ILE:CD1	1:A:866[B]:ILE:HD11	2.43	0.48
1:B:868:GLY:N	6:B:1439:HOH:O	2.45	0.48
1:B:848:ASP:O	1:B:849:ASN:HB2	2.12	0.48
1:B:1021:HIS:HD2	6:B:1337:HOH:O	1.94	0.48
1:B:833:LEU:HD21	1:B:1187:PHE:CE1	2.48	0.48
1:B:752:ARG:NH1	1:B:758:VAL:HG21	2.26	0.48
1:B:796:ASN:H	1:B:867:ASN:HD22	1.61	0.48
1:B:727:ASN:ND2	1:B:873:GLY:HA3	2.29	0.47
1:A:1146:LEU:HD23	1:B:1146:LEU:HD23	1.96	0.47
1:B:736:ASP:OD2	1:B:743:ARG:HD3	2.13	0.47
1:A:785:VAL:O	1:A:789:GLN:HG3	2.15	0.47
1:B:450:LYS:O	1:B:451:ILE:HD13	2.14	0.47
1:B:659:PRO:HA	1:B:662:ASP:OD1	2.15	0.47
1:B:759:LYS:HE2	1:B:823:PHE:CZ	2.49	0.47
1:A:751:LYS:HD3	1:A:770:MET:HE3	1.98	0.46
3:F:14:DC:H2''	3:F:15:DT:H5'	1.97	0.46
1:B:678:LYS:HE3	1:B:876:TRP:CD1	2.50	0.46
1:A:746[B]:LEU:HG	1:A:750:PHE:CE2	2.50	0.46
1:B:484[B]:LEU:CD1	1:B:640:ALA:HB2	2.46	0.46
1:B:555:MET:HE2	1:B:589:PHE:CD1	2.50	0.46
1:B:574:HIS:HE1	1:B:662:ASP:OD2	1.97	0.46
1:A:858:ILE:O	1:A:858:ILE:HG23	2.15	0.46
1:A:1002:HIS:HB3	1:A:1008:GLN:HG3	1.96	0.46
1:B:503[A]:ARG:HD2	6:B:1467:HOH:O	2.14	0.46
1:A:1012:THR:HB	3:F:19:DC:H5''	1.98	0.46
1:B:638:LYS:O	1:B:642:GLU:HG3	2.14	0.46
1:B:752:ARG:HH12	1:B:758:VAL:HG11	1.81	0.46
1:A:456:LYS:O	1:A:482:LYS:NZ	2.45	0.46
1:A:844:PHE:HA	1:A:854:PRO:HA	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:757:GLU:HB2	1:B:823:PHE:HB3	1.97	0.46
1:A:1169:ASN:HB3	1:A:1173[A]:ARG:NH1	2.32	0.45
1:B:935:ASN:O	1:B:987:LYS:HA	2.16	0.45
1:A:757:GLU:H	1:A:757:GLU:CD	2.24	0.45
1:B:580:LEU:HD23	1:B:580:LEU:HA	1.87	0.45
6:A:1246:HOH:O	1:B:762[A]:GLN:HG2	2.16	0.45
2:C:1:DA:C8	2:C:1:DA:P	3.02	0.45
3:D:14:DC:H2''	3:D:15:DT:H5'	1.98	0.45
1:B:507:LEU:HB3	1:B:523[A]:ILE:HD11	1.98	0.44
1:A:810:LEU:HD12	1:A:948:THR:HB	1.99	0.44
1:A:838:ASP:OD1	1:A:1049:ARG:HD2	2.17	0.44
1:A:757:GLU:HB2	1:A:823:PHE:HB3	2.00	0.44
1:B:768:ALA:HA	1:B:773:TYR:HB3	2.00	0.44
1:B:858:ILE:O	1:B:858:ILE:HG23	2.18	0.43
1:A:955:VAL:O	1:A:959:MET:HG3	2.18	0.43
1:B:933:ASP:OD2	1:B:934:ARG:O	2.36	0.43
1:B:993:LEU:HD12	1:B:993:LEU:O	2.18	0.43
1:A:1046:TYR:CE2	1:A:1180:TRP:CG	3.07	0.43
1:B:507:LEU:HB3	1:B:523[A]:ILE:CD1	2.47	0.43
1:B:757:GLU:CD	1:B:757:GLU:N	2.77	0.43
1:A:731:ILE:HA	1:A:732:PRO:HD3	1.85	0.43
1:B:727:ASN:HD22	1:B:873:GLY:HA3	1.83	0.43
1:A:773:TYR:CZ	1:A:775:HIS:HB2	2.54	0.43
1:A:957:GLU:HB2	1:A:958:PRO:HD3	2.01	0.43
1:B:551:LYS:N	1:B:551:LYS:HD3	2.33	0.43
1:B:1046:TYR:CZ	1:B:1180:TRP:HA	2.53	0.43
1:B:759:LYS:HE3	1:B:759:LYS:HB2	1.86	0.43
1:B:1064:LEU:HD23	1:B:1067:GLN:NE2	2.34	0.43
1:A:805:GLN:NE2	1:A:807:GLY:N	2.58	0.43
1:A:1110:GLN:O	1:A:1111:GLU:HB3	2.18	0.43
1:B:1014:ASN:ND2	6:B:1445:HOH:O	2.52	0.42
1:B:1081:ILE:HA	1:B:1084:ARG:CD	2.49	0.42
1:A:489:LEU:HB3	1:A:493:GLY:HA3	2.01	0.42
1:B:944:VAL:O	1:B:945:ARG:HB2	2.19	0.42
1:A:1063:LYS:HG3	1:A:1140:TYR:CD1	2.54	0.42
1:A:653[B]:LEU:HG	6:A:1424:HOH:O	2.18	0.42
1:B:587:GLU:HG2	1:B:653[B]:LEU:CD2	2.49	0.42
1:A:502:LEU:HD13	1:A:506:ILE:HD11	2.00	0.42
1:A:640:ALA:O	1:A:644:PHE:HD2	2.02	0.42
1:B:996:ALA:O	1:B:999:ALA:O	2.37	0.42
1:A:969:LEU:HG	1:A:988:MET:HE3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1032[A]:GLN:HG2	6:A:1438:HOH:O	2.18	0.42
1:A:752:ARG:HH12	1:A:758[B]:VAL:HG21	1.85	0.42
1:A:820[B]:ARG:NH1	6:B:1371:HOH:O	2.52	0.42
1:B:773:TYR:CZ	1:B:775:HIS:HB2	2.54	0.42
1:A:727:ASN:ND2	1:A:873:GLY:HA3	2.35	0.42
1:A:959:MET:HE1	1:A:1005:PHE:HA	2.02	0.42
1:B:520:ASN:OD1	1:B:523[B]:ILE:HG12	2.19	0.42
1:B:731:ILE:HG21	1:B:1025:LEU:HD11	2.02	0.42
1:B:858:ILE:HG13	1:B:1042:ARG:HD2	2.01	0.42
1:B:727:ASN:ND2	1:B:874:THR:H	2.18	0.42
1:A:1106:TRP:O	1:A:1110:GLN:HG2	2.20	0.41
1:B:503[B]:ARG:HB3	4:D:1:EVP:H3	2.03	0.41
1:B:539:ASP:OD2	1:B:540:ASP:N	2.49	0.41
1:A:974:LYS:HE3	6:A:1447:HOH:O	2.19	0.41
1:A:1173[A]:ARG:NH2	6:A:1458:HOH:O	2.51	0.41
1:A:576:ASN:HB2	1:A:577:TRP:CE3	2.56	0.41
1:A:1171:LEU:C	1:A:1171:LEU:HD13	2.45	0.41
1:B:546:THR:HG23	6:B:1428:HOH:O	2.20	0.41
1:B:525:ASN:O	1:B:529:ILE:HG13	2.20	0.41
1:B:1050:LYS:O	1:B:1054:VAL:HG23	2.21	0.41
1:B:731:ILE:HA	1:B:732:PRO:HD3	1.84	0.41
1:B:592:PRO:HG2	1:B:643:TYR:OH	2.21	0.41
1:B:1082:GLU:OE1	1:B:1082:GLU:N	2.44	0.41
1:B:1056:MET:HE2	1:B:1106:TRP:CE3	2.56	0.41
1:A:768:ALA:HA	1:A:773:TYR:HB3	2.03	0.40
1:B:554[B]:ILE:HD11	1:B:569:LEU:HD23	2.02	0.40
1:B:853:GLU:HB2	1:B:854:PRO:HD2	2.03	0.40
1:A:727:ASN:HD22	1:A:873:GLY:HA3	1.86	0.40
1:A:920:GLN:O	1:A:921:ASN:HB2	2.21	0.40
1:A:751:LYS:HD3	1:A:770:MET:CE	2.51	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	684/803 (85%)	664 (97%)	19 (3%)	1 (0%)	48	50
1	B	679/803 (85%)	661 (97%)	17 (2%)	1 (0%)	48	50
All	All	1363/1606 (85%)	1325 (97%)	36 (3%)	2 (0%)	43	50

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	852	VAL
1	B	852	VAL

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	610/704 (87%)	605 (99%)	5 (1%)	73	79
1	B	606/704 (86%)	604 (100%)	2 (0%)	86	91
All	All	1216/1408 (86%)	1209 (99%)	7 (1%)	78	84

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	466	LYS
1	A	539	ASP
1	A	870	GLU
1	A	880	LEU
1	A	1044	SER
1	B	492	ILE
1	B	551	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	574	HIS
1	A	576	ASN
1	A	712	ASN
1	A	727	ASN
1	A	805	GLN
1	A	882	ASN
1	A	922	GLN
1	A	1002	HIS
1	A	1021	HIS
1	A	1067	GLN
1	A	1083	ASN
1	A	1139	ASN
1	A	1160	GLN
1	B	533	GLN
1	B	574	HIS
1	B	576	ASN
1	B	583	HIS
1	B	712	ASN
1	B	727	ASN
1	B	786	ASN
1	B	805	GLN
1	B	867	ASN
1	B	882	ASN
1	B	920	GLN
1	B	935	ASN
1	B	949	GLN
1	B	1014	ASN
1	B	1021	HIS
1	B	1067	GLN
1	B	1076	GLN
1	B	1095	GLN
1	B	1139	ASN
1	B	1169	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 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	EVP	D	1	-	48,48,48	4.31	24 (50%)	71,73,73	2.34	25 (35%)
4	EVP	A	1	-	48,48,48	4.48	25 (52%)	71,73,73	2.31	24 (33%)

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	EVP	D	1	-	-	0/12/76/76	0/7/7/7
4	EVP	A	1	-	-	0/12/76/76	0/7/7/7

All (49) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1	EVP	C22-C28	-10.06	1.23	1.52
4	D	1	EVP	C22-C28	-9.73	1.24	1.52
4	A	1	EVP	C19-C14	8.69	1.52	1.39
4	A	1	EVP	C12-C13	8.61	1.53	1.38
4	D	1	EVP	C12-C13	8.38	1.53	1.38
4	D	1	EVP	C15-C14	8.27	1.51	1.39
4	A	1	EVP	C15-C14	8.16	1.51	1.39
4	A	1	EVP	C3-C2	8.02	1.52	1.38
4	D	1	EVP	C16-C17	7.95	1.51	1.40
4	A	1	EVP	C15-C16	7.93	1.52	1.38
4	D	1	EVP	C15-C16	7.92	1.52	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1	EVP	C24-C23	-7.81	1.38	1.51
4	D	1	EVP	C19-C14	7.77	1.50	1.39
4	A	1	EVP	C18-C17	7.74	1.51	1.40
4	A	1	EVP	C16-C17	7.72	1.51	1.40
4	D	1	EVP	C3-C2	7.63	1.51	1.38
4	A	1	EVP	C24-C23	-7.58	1.39	1.51
4	A	1	EVP	C19-C18	7.56	1.51	1.38
4	D	1	EVP	C19-C18	7.15	1.51	1.38
4	D	1	EVP	C18-C17	7.11	1.50	1.40
4	A	1	EVP	C4-C5	6.62	1.61	1.51
4	A	1	EVP	C3-C4	5.83	1.49	1.39
4	D	1	EVP	C4-C5	5.59	1.59	1.51
4	D	1	EVP	C3-C4	5.55	1.48	1.39
4	A	1	EVP	C12-C11	5.26	1.48	1.39
4	D	1	EVP	C12-C11	5.17	1.47	1.39
4	A	1	EVP	O3-C7	5.10	1.46	1.35
4	D	1	EVP	O3-C7	4.91	1.45	1.35
4	A	1	EVP	O11-C26	-4.83	1.36	1.44
4	D	1	EVP	O8-C22	4.82	1.55	1.41
4	A	1	EVP	O8-C22	4.78	1.55	1.41
4	D	1	EVP	O11-C26	-4.71	1.36	1.44
4	D	1	EVP	C6-C9	-4.54	1.45	1.54
4	A	1	EVP	C6-C9	-4.22	1.46	1.54
4	A	1	EVP	C6-C5	-4.10	1.48	1.56
4	D	1	EVP	C6-C5	-3.68	1.49	1.56
4	A	1	EVP	C14-C5	3.30	1.56	1.52
4	A	1	EVP	O9-C22	-3.03	1.34	1.41
4	D	1	EVP	O9-C22	-2.81	1.34	1.41
4	A	1	EVP	O7-C18	2.59	1.41	1.37
4	A	1	EVP	C11-C10	2.51	1.55	1.50
4	D	1	EVP	O7-C18	2.35	1.41	1.37
4	A	1	EVP	C11-C4	2.32	1.44	1.40
4	D	1	EVP	C11-C10	2.18	1.54	1.50
4	A	1	EVP	O10-C25	2.07	1.46	1.41
4	A	1	EVP	C28-C27	-2.07	1.47	1.52
4	D	1	EVP	C14-C5	2.05	1.55	1.52
4	D	1	EVP	O12-C27	-2.04	1.37	1.43
4	D	1	EVP	O2-C1	-2.02	1.39	1.43

All (49) bond angle outliers are listed below:

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1	EVP	O11-C26-C27	6.59	121.26	109.72
4	D	1	EVP	O11-C26-C27	6.57	121.23	109.72
4	A	1	EVP	O3-C7-O4	5.95	127.75	121.43
4	D	1	EVP	O8-C22-O9	-5.75	95.57	110.69
4	A	1	EVP	O8-C22-O9	-5.61	95.94	110.69
4	D	1	EVP	O5-C16-C17	5.23	120.00	114.53
4	A	1	EVP	O5-C16-C17	4.96	119.72	114.53
4	D	1	EVP	O3-C7-O4	4.92	126.66	121.43
4	D	1	EVP	O7-C18-C17	4.91	119.67	114.53
4	A	1	EVP	C27-C26-C23	-4.90	100.07	110.93
4	D	1	EVP	C4-C5-C6	4.70	114.31	106.57
4	D	1	EVP	C27-C26-C23	-4.42	101.13	110.93
4	A	1	EVP	C4-C5-C6	4.29	113.63	106.57
4	A	1	EVP	O13-C28-C27	-4.22	100.43	110.38
4	A	1	EVP	O9-C22-C28	4.09	118.77	110.37
4	A	1	EVP	O7-C18-C17	3.92	118.63	114.53
4	D	1	EVP	O9-C23-C26	3.82	117.61	109.72
4	D	1	EVP	O13-C28-C27	-3.57	101.97	110.38
4	A	1	EVP	C22-O9-C23	-3.43	107.01	113.72
4	D	1	EVP	O9-C22-C28	3.40	117.35	110.37
4	D	1	EVP	C22-O9-C23	-3.35	107.19	113.72
4	A	1	EVP	C24-O10-C25	3.31	114.91	111.67
4	D	1	EVP	O8-C22-C28	-3.30	99.96	108.09
4	A	1	EVP	O8-C22-C28	-3.20	100.22	108.09
4	A	1	EVP	O9-C23-C26	3.09	116.10	109.72
4	A	1	EVP	O11-C25-O10	2.97	113.85	110.86
4	D	1	EVP	C8-C9-C6	2.95	106.31	101.73
4	A	1	EVP	O4-C7-C6	-2.91	125.68	129.38
4	D	1	EVP	O1-C13-C12	2.87	131.68	127.86
4	A	1	EVP	C21-O7-C18	2.86	121.71	117.51
4	D	1	EVP	C24-O10-C25	2.69	114.30	111.67
4	D	1	EVP	O4-C7-C6	-2.68	125.97	129.38
4	D	1	EVP	O2-C2-C3	2.66	131.40	127.86
4	A	1	EVP	C9-C6-C7	2.63	107.03	103.14
4	A	1	EVP	O1-C13-C12	2.59	131.30	127.86
4	D	1	EVP	O7-C18-C19	-2.56	119.67	124.08
4	A	1	EVP	C8-C9-C6	2.55	105.68	101.73
4	D	1	EVP	C14-C5-C4	-2.52	109.06	112.95
4	D	1	EVP	O2-C1-O1	-2.38	104.36	108.09
4	D	1	EVP	C9-C6-C7	2.22	106.42	103.14
4	D	1	EVP	O5-C16-C15	-2.21	120.27	124.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1	EVP	O2-C2-C3	2.21	130.80	127.86
4	A	1	EVP	O5-C16-C15	-2.21	120.28	124.08
4	D	1	EVP	C8-O3-C7	-2.17	108.16	110.31
4	D	1	EVP	O3-C8-C9	2.14	107.95	104.70
4	D	1	EVP	O11-C25-C29	2.13	111.39	108.50
4	A	1	EVP	O3-C8-C9	2.06	107.83	104.70
4	A	1	EVP	C28-C27-C26	2.01	114.23	109.68
4	A	1	EVP	O3-C7-C6	-2.00	106.51	109.50

There are no chirality outliers.

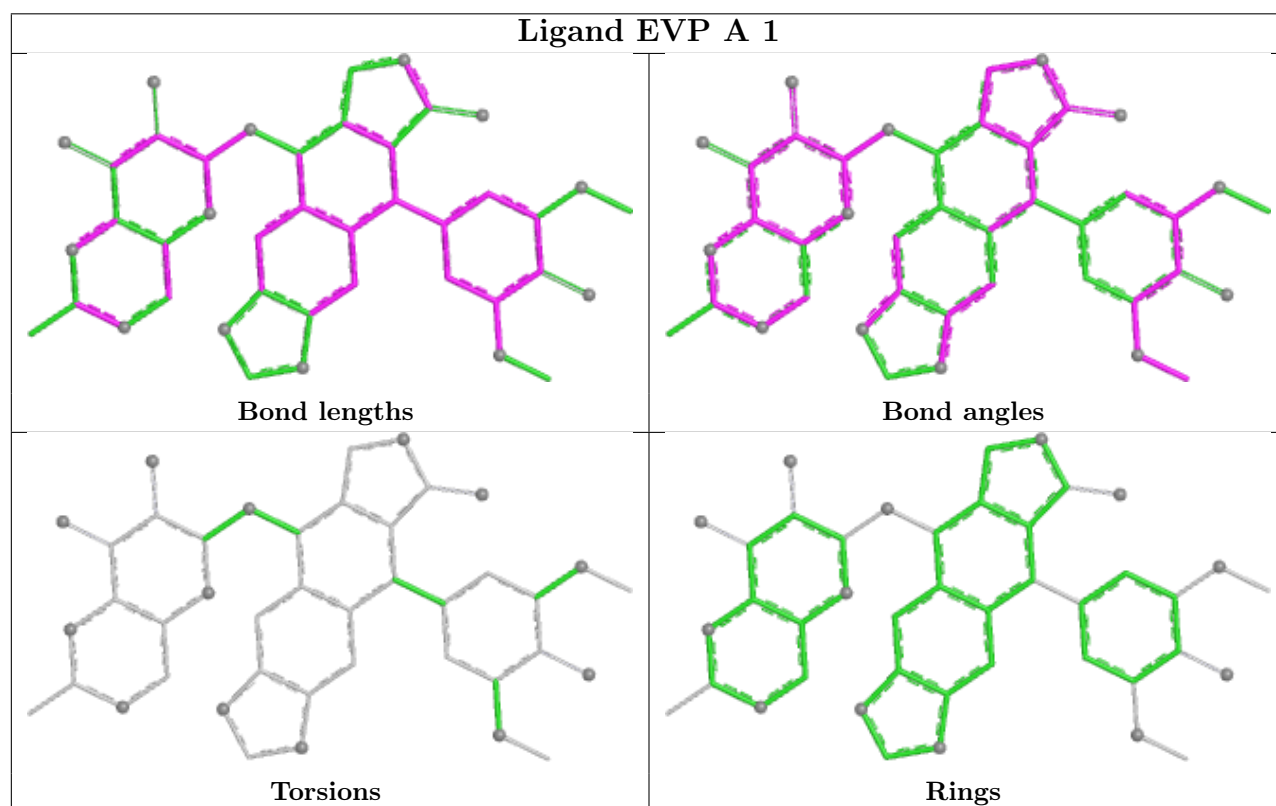
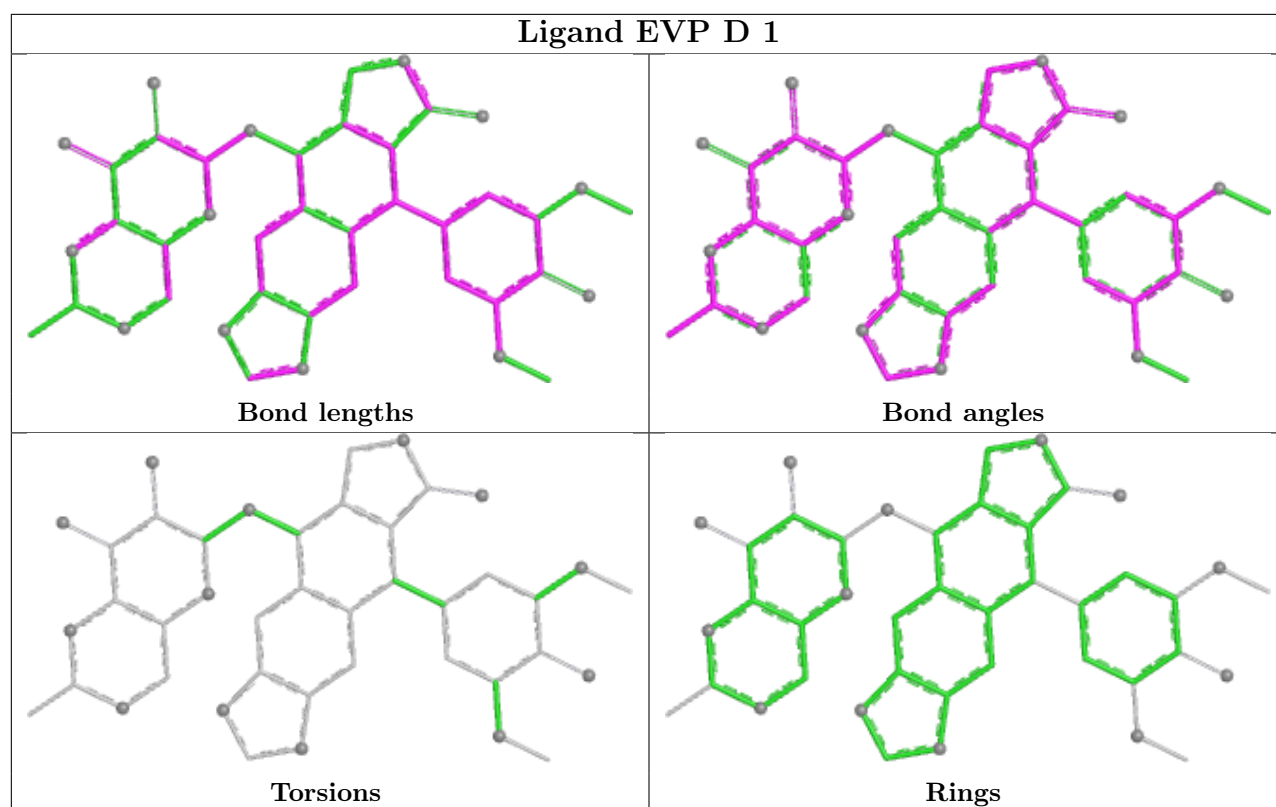
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	1	EVP	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 ⓘ

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	671/803 (83%)	-0.17	18 (2%) 56 60	11, 34, 71, 102	21 (3%)
1	B	671/803 (83%)	-0.22	12 (1%) 67 71	17, 31, 65, 91	18 (2%)
2	C	8/8 (100%)	-0.74	0 100 100	21, 24, 46, 67	0
2	E	8/8 (100%)	-0.77	0 100 100	19, 22, 43, 68	0
3	D	12/12 (100%)	-1.03	0 100 100	19, 28, 34, 34	0
3	F	12/12 (100%)	-0.96	0 100 100	23, 30, 39, 56	0
All	All	1382/1646 (83%)	-0.22	30 (2%) 62 66	11, 32, 67, 102	39 (2%)

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	452	LYS	4.4
1	B	695	GLY	4.3
1	B	491	VAL	4.3
1	B	637	ALA	4.1
1	A	490	GLY	4.0
1	A	491	VAL	3.9
1	A	696	LEU	3.4
1	B	848	ASP	3.4
1	A	453	GLY	3.1
1	A	657	ALA	2.9
1	B	967	PRO	2.9
1	A	522	GLU	2.8
1	B	541	ALA	2.7
1	A	693	LEU	2.6
1	A	638	LYS	2.6
1	A	848	ASP	2.6
1	B	449	SER	2.6
1	B	934	ARG	2.5
1	B	693	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	492	ILE	2.5
1	A	849	ASN	2.4
1	A	484	LEU	2.3
1	A	494	ARG	2.3
1	B	644	PHE	2.3
1	B	492	ILE	2.3
1	B	643	TYR	2.2
1	A	489	LEU	2.2
1	A	1201	ASP	2.1
1	A	694	HIS	2.0
1	A	645	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](#)

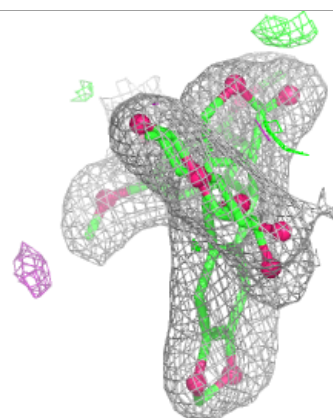
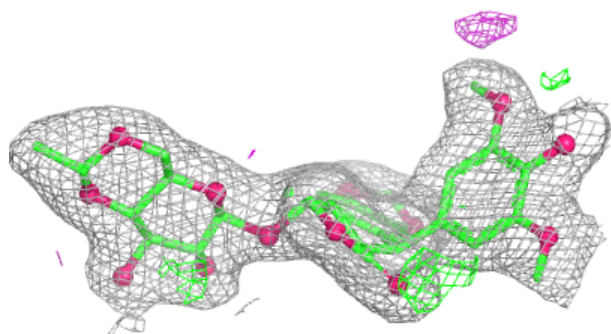
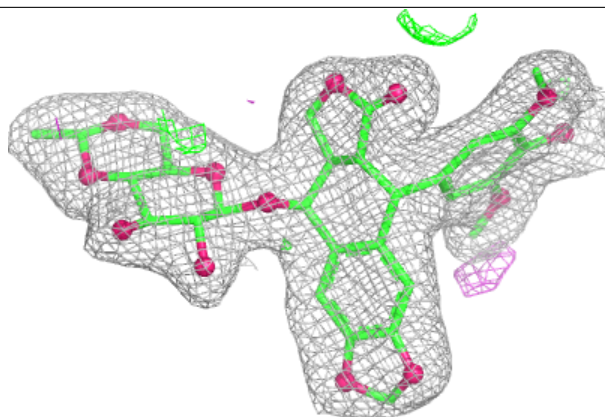
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists 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
5	MG	B	1222	1/1	0.87	0.12	48,48,48,48	0
5	MG	A	1223	1/1	0.93	0.07	45,45,45,45	0
5	MG	D	21	1/1	0.94	0.07	44,44,44,44	0
4	EVP	A	1	42/42	0.96	0.07	21,29,33,36	0
4	EVP	D	1	42/42	0.97	0.06	19,25,34,38	0
5	MG	F	1	1/1	0.97	0.05	50,50,50,50	0
5	MG	B	1	1/1	0.99	0.06	30,30,30,30	0
5	MG	A	1222	1/1	0.99	0.09	31,31,31,31	0

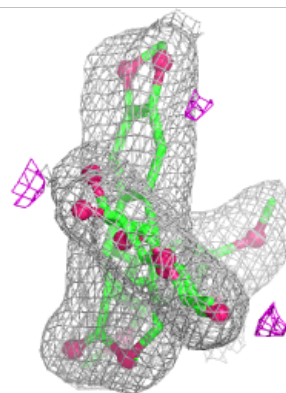
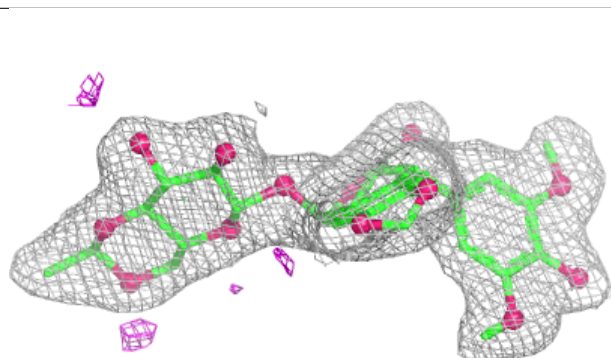
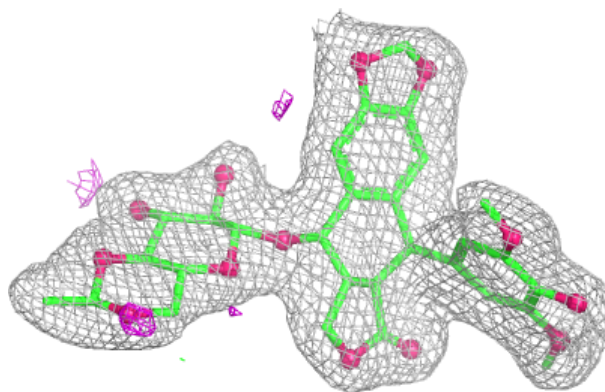
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around EVP A 1:

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

**Electron density around EVP D 1:**

$2mF_o-DF_c$ (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.