



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 29, 2021 – 12:07 PM EDT

PDB ID : 7LUF
Title : HSV1 polymerase ternary complex with dsDNA and PNU-183792
Authors : Hayes, R.P.; Klein, D.
Deposited on : 2021-02-22
Resolution : 3.50 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.18
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.18

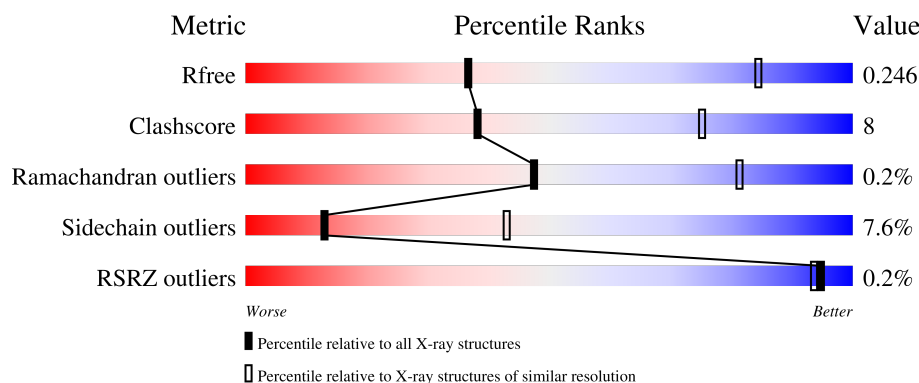
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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}	130704	1659 (3.60-3.40)
Clashscore	141614	1036 (3.58-3.42)
Ramachandran outliers	138981	1005 (3.58-3.42)
Sidechain outliers	138945	1006 (3.58-3.42)
RSRZ outliers	127900	1559 (3.60-3.40)

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	<div> <div>71%</div> <div>17%</div> <div>• 10%</div> </div>
1	B	1163	<div> <div>70%</div> <div>19%</div> <div>• 9%</div> </div>
2	C	27	<div> <div>7%</div> <div>56%</div> <div>11%</div> <div>33%</div> </div>
2	E	27	<div> <div>48%</div> <div>19%</div> <div>33%</div> </div>
3	D	23	<div> <div>35%</div> <div>30%</div> <div>35%</div> </div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	F	23	<div><div><div></div><div></div><div></div></div><div><div>4%</div><div>13%</div><div>52%</div><div>35%</div></div></div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 17531 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	1044	Total	C	N	O	S	0	0	0
			7946	5093	1365	1449	39			
1	B	1054	Total	C	N	O	S	0	0	0
			8175	5249	1409	1477	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	conflict	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	conflict	UNP I7GY94

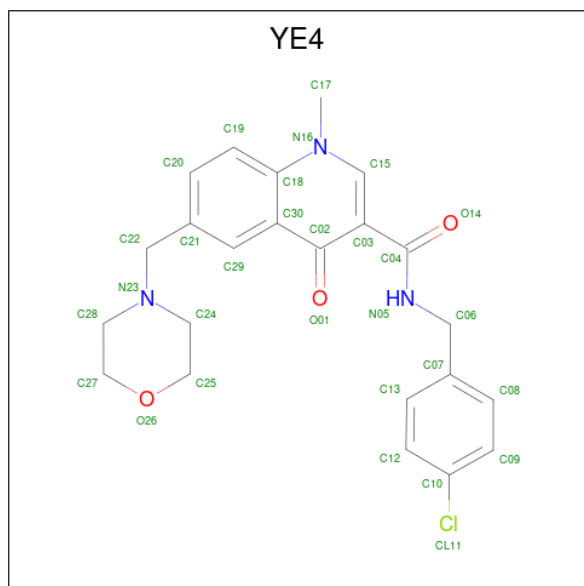
- Molecule 2 is a DNA chain called DNA (5'-D(P*AP*TP*GP*GP*TP*AP*GP*GP*GP*GP*AP*AP*GP*GP*AP*TP*CP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	18	Total 382	C 179	N 79	O 106	P 18	0	0	0
2	E	18	Total 382	C 179	N 79	O 106	P 18	0	0	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	15	Total 293	C 142	N 47	O 90	P 14	0	0	0
3	F	15	Total 293	C 142	N 47	O 90	P 14	0	0	0

- Molecule 4 is N-(4-chlorobenzyl)-1-methyl-6-(morpholinomethyl)-4-oxo-1,4-dihydroquinolin e-3-carboxamide (three-letter code: YE4) (formula: C₂₃H₂₄ClN₃O₃) (labeled as "Ligand of Interest" by depositor).

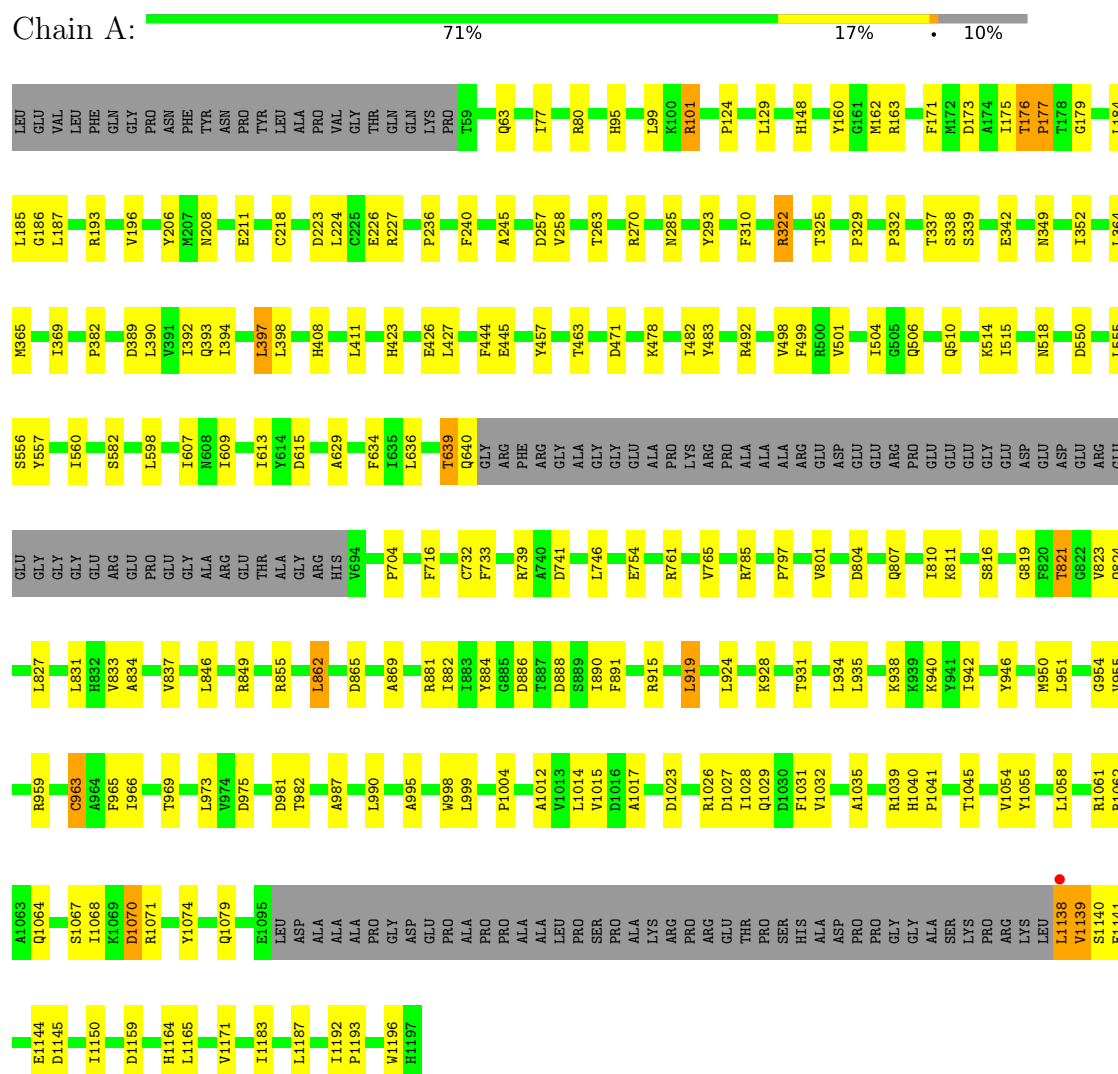


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	C	1	Total 30	C 23	Cl 1	N 3	O 3	0	0
4	E	1	Total 30	C 23	Cl 1	N 3	O 3	0	0

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 polymerase

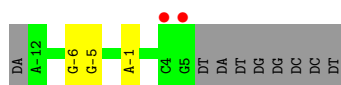


• Molecule 1: DNA polymerase





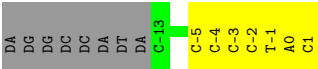
Chain C:  7% 56% 11% 26%



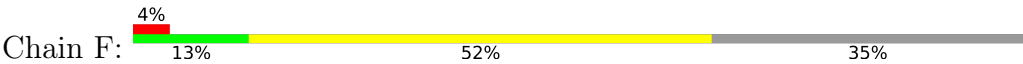
Chain E:  48% 19% 33%



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	181.30Å 181.30Å 233.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	57.33 – 3.50 57.33 – 3.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (57.33-3.50) 99.9 (57.33-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 3.49Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.226 , 0.245 0.226 , 0.246	Depositor DCC
R_{free} test set	4710 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	104.9	Xtriage
Anisotropy	0.198	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 65.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.034 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	17531	wwPDB-VP
Average B, all atoms (Å ²)	109.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.29% 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: YE4

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	0.25	0/8140	0.44	0/11087
1	B	0.25	0/8376	0.45	0/11387
2	C	0.48	0/431	0.84	0/666
2	E	0.46	0/431	0.84	0/666
3	D	0.60	0/325	1.02	0/497
3	F	0.59	0/325	1.01	0/497
All	All	0.29	0/18028	0.51	0/24800

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	7946	0	7640	114	0
1	B	8175	0	7991	123	0
2	C	382	0	202	5	0
2	E	382	0	202	5	0
3	D	293	0	171	10	0
3	F	293	0	171	14	0
4	C	30	0	0	1	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	E	30	0	0	0	0
All	All	17531	0	16377	257	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 8.

All (257) 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:788:ILE:HD11	1:B:807:GLN:HB3	1.50	0.90
3:D:0:DA:H2"	3:D:1:DC:H5"	1.54	0.90
3:F:0:DA:H2"	3:F:1:DC:H5"	1.57	0.85
1:A:258:VAL:HG21	1:A:636:LEU:HD23	1.65	0.78
1:B:990:LEU:HD11	1:B:1007:LEU:HD11	1.67	0.77
1:B:392:ILE:HG22	1:B:393:GLN:HG3	1.67	0.76
1:B:234:GLU:OE2	1:B:234:GLU:N	2.16	0.75
1:A:176:THR:HG23	1:A:177:PRO:HD3	1.69	0.74
1:B:598:LEU:HD13	1:B:613:ILE:HG12	1.71	0.72
1:B:1003:LEU:HB3	1:B:1007:LEU:HD12	1.70	0.72
1:B:608:ASN:HD21	1:B:611:ARG:HG3	1.56	0.70
1:A:1067:SER:HB2	1:A:1070:ASP:OD2	1.92	0.70
1:B:950:MET:SD	1:B:971:ARG:NH1	2.66	0.69
1:A:598:LEU:HD13	1:A:613:ILE:HG12	1.74	0.69
1:B:698:GLY:O	1:B:842:ARG:NH1	2.26	0.69
1:B:998:TRP:HZ3	1:B:1004:PRO:HD3	1.59	0.68
1:A:1061:ARG:NH1	1:A:1144:GLU:OE2	2.21	0.67
1:B:1041:PRO:HA	1:B:1044:TYR:HD1	1.57	0.67
1:B:998:TRP:CZ3	1:B:1004:PRO:HD3	2.29	0.67
1:B:1027:ASP:HB3	1:B:1030:ASP:HB2	1.76	0.67
1:A:397:LEU:HD11	1:A:582:SER:HB3	1.76	0.67
1:B:223:ASP:O	1:B:227:ARG:HG2	1.95	0.67
1:A:855:ARG:NH1	1:A:865:ASP:OD2	2.29	0.66
1:B:153:LEU:HB2	1:B:183:THR:HB	1.77	0.65
1:B:63:GLN:NE2	1:B:162:MET:SD	2.69	0.65
1:A:339:SER:OG	1:A:492:ARG:NH1	2.29	0.65
1:B:611:ARG:HH21	1:B:619:ILE:HD11	1.63	0.64
1:A:998:TRP:HZ3	1:A:1004:PRO:HD3	1.63	0.63
1:A:175:ILE:HG22	1:A:179:GLY:HA3	1.80	0.63
1:B:716:PHE:HB2	1:B:890:ILE:HG22	1.82	0.62
1:A:959:ARG:HH22	2:C:-5:DG:H21	1.47	0.61
1:B:339:SER:OG	1:B:492:ARG:NH1	2.33	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:995:ALA:HA	1:A:998:TRP:CD1	2.36	0.61
1:A:1058:LEU:HD21	1:A:1145:ASP:OD2	2.01	0.60
1:A:557:TYR:O	1:A:560:ILE:HG22	2.01	0.60
1:B:738:LEU:HD12	1:B:760:ARG:HH21	1.66	0.60
1:A:824:GLN:HB3	1:A:831:LEU:HD11	1.84	0.59
1:A:101:ARG:NH1	1:A:445:GLU:OE1	2.34	0.59
1:B:1061:ARG:NH1	1:B:1144:GLU:OE2	2.31	0.59
1:A:1183:ILE:O	1:A:1187:LEU:HG	2.02	0.59
1:B:986:ALA:O	1:B:990:LEU:HD12	2.03	0.59
1:A:739:ARG:NH2	1:A:741:ASP:OD2	2.36	0.59
1:B:855:ARG:NH1	1:B:865:ASP:OD2	2.36	0.59
3:D:0:DA:C2'	3:D:1:DC:H5''	2.31	0.58
1:B:824:GLN:HB3	1:B:831:LEU:HD11	1.85	0.58
1:B:1048:ARG:NE	2:E:0:DG:OP1	2.31	0.58
1:A:193:ARG:NH1	1:A:338:SER:O	2.37	0.57
1:A:963:CYS:SG	1:A:965:PHE:HB3	2.44	0.57
1:B:175:ILE:HG22	1:B:179:GLY:HA3	1.86	0.57
1:B:728:ALA:HB1	1:B:919:LEU:HD13	1.87	0.57
1:A:716:PHE:HB2	1:A:890:ILE:HG22	1.86	0.57
1:B:206:TYR:HB2	1:B:293:TYR:HB2	1.87	0.57
1:A:392:ILE:HG13	1:A:393:GLN:N	2.20	0.57
1:A:1192:ILE:HG23	1:A:1196:TRP:HE3	1.69	0.57
1:A:325:THR:HG22	1:A:349:ASN:HD21	1.70	0.56
1:B:713:VAL:HG23	1:B:930:PHE:HB2	1.88	0.56
1:B:821:THR:HG22	1:B:834:ALA:HB2	1.88	0.56
1:A:969:THR:HG21	1:A:1165:LEU:HD11	1.88	0.55
1:B:159:ALA:HB2	1:B:176:THR:HA	1.89	0.55
1:A:1028:ILE:HD12	1:A:1150:ILE:HG12	1.89	0.55
1:A:77:ILE:HG22	1:A:390:LEU:HD11	1.89	0.55
1:A:337:THR:OG1	1:A:342:GLU:OE2	2.22	0.55
1:A:998:TRP:CZ3	1:A:1004:PRO:HD3	2.42	0.55
1:B:193:ARG:NH1	1:B:338:SER:O	2.41	0.54
1:B:148:HIS:O	1:B:186:GLY:HA3	2.08	0.54
1:B:258:VAL:HG13	1:B:634:PHE:HB3	1.90	0.53
1:B:629:ALA:HB1	1:B:634:PHE:HB2	1.91	0.53
1:B:785:ARG:HB2	1:B:810:ILE:HG21	1.91	0.53
1:B:1054:VAL:HG21	1:B:1074:TYR:HB3	1.91	0.53
1:B:184:LEU:HB2	1:B:196:VAL:HG13	1.90	0.53
1:B:187:LEU:HD23	1:B:193:ARG:HA	1.91	0.53
1:A:223:ASP:O	1:A:227:ARG:HG2	2.08	0.53
1:A:173:ASP:O	1:A:176:THR:HG22	2.08	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:VAL:HG22	1:A:634:PHE:HB3	1.91	0.52
3:F:0:DA:C2'	3:F:1:DC:H5''	2.34	0.52
3:F:-5:DC:H1'	3:F:-4:DC:H5'	1.91	0.52
1:B:942:ILE:HD13	1:B:952:ILE:HG23	1.92	0.52
1:A:124:PRO:HG3	1:A:408:HIS:CD2	2.45	0.52
1:A:206:TYR:HB2	1:A:293:TYR:HB2	1.90	0.52
1:A:959:ARG:HH22	2:C:-5:DG:N2	2.07	0.52
1:B:886:ASP:O	1:B:888:ASP:N	2.42	0.52
1:A:218:CYS:HB2	1:A:223:ASP:HB3	1.92	0.51
1:B:959:ARG:HH22	3:F:-3:DC:H1'	1.75	0.51
1:B:249:GLU:HB3	1:B:272:TYR:HB2	1.91	0.51
1:B:940:LYS:NZ	2:E:-4:DG:H5'	2.25	0.51
1:B:965:PHE:HE1	1:B:1020:ARG:HH11	1.58	0.51
1:B:247:HIS:O	1:B:247:HIS:ND1	2.44	0.51
1:B:230:ALA:O	1:B:234:GLU:OE2	2.28	0.51
1:A:732:CYS:SG	1:A:733:PHE:N	2.84	0.51
1:B:398:LEU:HD22	1:B:408:HIS:HB2	1.93	0.51
1:B:732:CYS:SG	1:B:733:PHE:N	2.84	0.50
1:A:187:LEU:HD23	1:A:193:ARG:HA	1.93	0.50
1:A:504:ILE:HD13	1:A:1068:ILE:HB	1.92	0.50
1:B:500:ARG:HD2	1:B:502:TRP:CE3	2.46	0.50
1:B:741:ASP:OD1	1:B:741:ASP:N	2.42	0.50
1:B:1192:ILE:HG13	1:B:1197:HIS:HE1	1.76	0.50
1:B:1045:THR:HG23	3:F:-4:DC:P	2.51	0.50
1:A:208:ASN:OD1	1:A:211:GLU:N	2.43	0.50
1:A:80:ARG:HG2	1:A:95:HIS:CD2	2.46	0.50
1:B:848:THR:HG23	1:B:913:ILE:HG21	1.93	0.50
1:B:747:GLU:HB3	1:B:750:LYS:HG2	1.93	0.49
1:B:1185:GLU:HG2	1:B:1189:LYS:HE3	1.95	0.49
1:B:426:GLU:HG2	1:B:430:ARG:HD2	1.94	0.49
3:D:-2:DC:H2''	3:D:-1:DT:C6	2.48	0.49
1:A:501:VAL:HG22	1:A:515:ILE:HD13	1.95	0.49
1:A:506:GLN:HG3	1:A:1040:HIS:CD2	2.48	0.49
2:E:-6:DG:H2'	2:E:-5:DG:C8	2.48	0.49
1:A:1014:LEU:O	1:A:1017:ALA:HB3	2.13	0.49
3:D:-3:DC:H5'	3:D:-3:DC:C6	2.48	0.48
1:B:555:LEU:HB2	1:B:577:TYR:CD2	2.48	0.48
1:B:382:PRO:HB3	1:B:389:ASP:HB3	1.95	0.48
1:B:411:LEU:HD11	1:B:575:GLY:HA3	1.94	0.48
1:A:398:LEU:HD22	1:A:457:TYR:CD2	2.47	0.48
1:B:945:ILE:HD11	1:B:951:LEU:HD13	1.94	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1185:GLU:O	1:B:1189:LYS:HG2	2.13	0.48
1:B:423:HIS:O	1:B:427:LEU:HG	2.14	0.48
1:A:148:HIS:O	1:A:186:GLY:HA3	2.14	0.48
1:A:629:ALA:HB1	1:A:634:PHE:HB2	1.96	0.48
1:B:310:PHE:HA	1:B:352:ILE:HD11	1.96	0.48
1:B:611:ARG:O	1:B:615:ASP:HB2	2.14	0.48
1:B:1193:PRO:HD2	1:B:1196:TRP:CZ3	2.49	0.47
1:A:365:MET:HB2	1:A:398:LEU:HD23	1.96	0.47
1:B:216:LEU:HD13	1:B:227:ARG:HG3	1.96	0.47
1:A:184:LEU:HB2	1:A:196:VAL:HG13	1.97	0.47
1:B:99:LEU:HD21	1:B:101:ARG:HG2	1.96	0.47
1:B:911:SER:O	1:B:915:ARG:HG3	2.15	0.47
1:B:77:ILE:HG22	1:B:390:LEU:HD11	1.95	0.47
1:A:969:THR:O	1:A:973:LEU:HG	2.15	0.47
3:F:-12:DG:H2"	3:F:-11:DA:C8	2.50	0.47
1:B:358:ASP:OD2	1:B:358:ASP:N	2.36	0.47
1:B:882:ILE:HG13	1:B:892:VAL:HG22	1.96	0.47
1:A:754:GLU:OE1	1:A:761:ARG:NH2	2.43	0.47
1:B:206:TYR:CE2	1:B:270:ARG:HG3	2.49	0.47
1:B:959:ARG:NH1	3:F:-2:DC:O2	2.48	0.47
1:A:423:HIS:O	1:A:427:LEU:HG	2.15	0.46
1:B:500:ARG:HB3	1:B:502:TRP:CE3	2.51	0.46
1:B:1192:ILE:HG13	1:B:1197:HIS:CE1	2.51	0.46
1:A:99:LEU:HD21	1:A:101:ARG:HD2	1.97	0.46
1:A:478:LYS:HD3	1:A:482:ILE:HD12	1.98	0.46
1:B:1082:GLU:O	1:B:1086:THR:OG1	2.32	0.46
1:A:704:PRO:HB2	1:A:935:LEU:HD13	1.96	0.46
1:A:310:PHE:HA	1:A:352:ILE:HD11	1.98	0.46
1:A:804:ASP:O	1:A:807:GLN:HG3	2.15	0.46
1:A:257:ASP:HB3	1:A:263:THR:HG21	1.98	0.46
1:B:423:HIS:HB2	1:B:576:GLU:OE1	2.16	0.46
1:A:940:LYS:HG2	1:A:954:GLY:HA3	1.98	0.46
1:B:998:TRP:HZ3	1:B:1003:LEU:HA	1.81	0.45
1:A:987:ALA:HA	1:A:990:LEU:HD12	1.98	0.45
1:A:1023:ASP:O	1:A:1026:ARG:HG3	2.16	0.45
3:D:-2:DC:H2"	3:D:-1:DT:H6	1.80	0.45
1:B:319:LYS:HE3	1:B:348:ASP:O	2.17	0.45
1:A:206:TYR:CE2	1:A:270:ARG:HG3	2.52	0.45
1:A:1193:PRO:HD2	1:A:1196:TRP:CZ3	2.51	0.45
1:A:816:SER:OG	4:C:1201:YE4:CL11	2.72	0.45
1:A:886:ASP:O	1:A:888:ASP:N	2.49	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:-6:DG:H2''	2:C:-5:DG:H5'	1.98	0.45
1:B:580:GLN:O	1:B:584:LEU:HD23	2.16	0.44
1:B:969:THR:HG21	1:B:1165:LEU:HD11	1.99	0.44
1:B:956:ASP:OD2	1:B:962:ASN:ND2	2.41	0.44
3:F:-10:DT:H2''	3:F:-9:DC:C6	2.53	0.44
1:B:301:THR:O	1:B:305:LEU:HD13	2.17	0.44
1:B:995:ALA:HB1	1:B:1183:ILE:HG21	1.99	0.44
2:E:-6:DG:N2	3:F:-1:DT:O2	2.50	0.44
1:A:80:ARG:HG2	1:A:95:HIS:NE2	2.33	0.44
1:B:998:TRP:CE3	1:B:1003:LEU:HD23	2.53	0.44
1:A:171:PHE:O	1:A:175:ILE:HG12	2.18	0.44
1:B:498:VAL:H	1:B:518:ASN:HD22	1.64	0.44
1:A:884:TYR:HD2	1:A:891:PHE:CD2	2.36	0.44
1:B:785:ARG:NH2	1:B:811:LYS:HD2	2.32	0.44
1:A:741:ASP:N	1:A:741:ASP:OD1	2.51	0.44
1:A:849:ARG:HD3	1:A:882:ILE:HG21	2.00	0.44
1:B:555:LEU:HB2	1:B:577:TYR:HD2	1.82	0.44
1:B:609:ILE:O	1:B:613:ILE:HG13	2.17	0.44
1:A:819:GLY:O	1:A:823:VAL:HG23	2.18	0.43
1:B:608:ASN:ND2	1:B:611:ARG:HG3	2.29	0.43
1:A:1192:ILE:HG23	1:A:1196:TRP:CE3	2.51	0.43
3:F:-7:DT:H5'	3:F:-7:DT:H6	1.83	0.43
1:A:919:LEU:HD12	1:A:919:LEU:H	1.82	0.43
3:F:-2:DC:H2''	3:F:-1:DT:C6	2.53	0.43
1:A:1012:ALA:O	1:A:1015:VAL:HG12	2.19	0.43
3:D:-5:DC:H1'	3:D:-4:DC:H5'	2.00	0.43
1:B:1028:ILE:HG21	1:B:1150:ILE:HD12	2.00	0.43
1:B:1061:ARG:NH2	1:B:1144:GLU:OE2	2.49	0.43
1:A:382:PRO:HB3	1:A:389:ASP:HB3	2.01	0.43
1:A:1068:ILE:H	1:A:1068:ILE:HG13	1.64	0.43
1:A:364:LEU:HD11	1:A:463:THR:HG22	2.00	0.43
1:A:639:THR:HG22	1:A:640:GLN:H	1.83	0.43
1:A:931:THR:HG23	1:A:946:TYR:HA	2.01	0.43
1:B:511:LYS:HD3	1:B:511:LYS:HA	1.54	0.43
1:B:639:THR:HG22	1:B:640:GLN:H	1.83	0.43
1:A:369:ILE:HG22	1:A:394:ILE:HG12	2.01	0.42
1:A:1139:VAL:O	1:A:1141:GLU:N	2.45	0.42
1:A:1139:VAL:HG22	2:C:-1:DA:P	2.59	0.42
1:B:315:TRP:CD1	1:B:356:MET:HB3	2.53	0.42
1:A:1015:VAL:HG23	1:A:1196:TRP:HH2	1.85	0.42
1:B:956:ASP:O	1:B:962:ASN:ND2	2.42	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:-3:DC:H5'	3:F:-3:DC:C6	2.53	0.42
1:A:862:LEU:HD12	1:A:869:ALA:HB2	2.00	0.42
1:A:1071:ARG:NH2	3:D:-2:DC:OP1	2.48	0.42
1:A:322:ARG:HG2	1:B:771:GLU:OE2	2.20	0.42
1:B:373:ALA:HA	1:B:389:ASP:OD2	2.19	0.42
1:B:994:PRO:O	1:B:998:TRP:HD1	2.01	0.42
3:F:-2:DC:H2''	3:F:-1:DT:H6	1.85	0.42
1:B:129:LEU:HD11	1:B:332:PRO:O	2.19	0.42
1:B:369:ILE:HG22	1:B:394:ILE:HG12	2.02	0.42
2:E:-8:DT:H5'	2:E:-8:DT:C6	2.54	0.42
1:A:444:PHE:HD1	1:A:483:TYR:CD2	2.38	0.42
1:A:942:ILE:HD11	1:A:955:VAL:HG21	2.02	0.42
1:B:238:ALA:HB1	1:B:241:ARG:HD2	2.02	0.42
1:B:739:ARG:HB2	1:B:741:ASP:OD1	2.20	0.42
1:B:1180:ASN:ND2	1:B:1183:ILE:HD13	2.34	0.42
1:A:63:GLN:NE2	1:A:162:MET:SD	2.93	0.42
1:A:797:PRO:O	1:A:801:VAL:HG23	2.20	0.42
1:B:526:TYR:O	1:B:530:THR:HG23	2.20	0.42
1:B:536:SER:OG	1:B:960:LYS:O	2.31	0.42
1:B:804:ASP:O	1:B:807:GLN:HG3	2.20	0.42
1:A:884:TYR:HB3	1:A:891:PHE:HB2	2.02	0.41
1:A:1041:PRO:HG3	1:A:1055:TYR:CE2	2.55	0.41
1:B:204:TYR:HA	1:B:271:VAL:O	2.18	0.41
1:A:821:THR:HG22	1:A:834:ALA:HB2	2.01	0.41
1:A:1015:VAL:HG23	1:A:1196:TRP:CH2	2.55	0.41
1:A:1027:ASP:OD2	1:A:1029:GLN:HB2	2.20	0.41
1:B:198:VAL:HA	1:B:345:CYS:O	2.21	0.41
1:B:423:HIS:CG	1:B:576:GLU:HG3	2.55	0.41
1:B:890:ILE:HD12	1:B:890:ILE:HA	1.89	0.41
1:B:1035:ALA:HA	3:F:-2:DC:OP1	2.20	0.41
1:A:77:ILE:HA	1:A:95:HIS:O	2.20	0.41
1:A:498:VAL:H	1:A:518:ASN:HD22	1.67	0.41
1:A:510:GLN:HE22	1:A:514:LYS:HE2	1.86	0.41
3:D:-3:DC:H5'	3:D:-3:DC:H6	1.84	0.41
1:A:785:ARG:HB2	1:A:810:ILE:HG21	2.02	0.41
1:A:185:LEU:HD13	1:A:492:ARG:HG2	2.02	0.41
1:A:555:LEU:HD13	1:A:556:SER:N	2.36	0.41
1:A:963:CYS:SG	1:A:1031:PHE:HA	2.61	0.41
1:A:1029:GLN:O	1:A:1032:VAL:HG22	2.21	0.41
1:B:63:GLN:OE1	1:B:496:ARG:NH2	2.54	0.41
1:B:337:THR:OG1	1:B:342:GLU:OE2	2.32	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:930:PHE:CE2	1:B:945:ILE:HG12	2.56	0.41
1:A:160:TYR:O	1:A:163:ARG:HG2	2.21	0.41
1:B:1175:ALA:HA	1:B:1179:ASN:OD1	2.21	0.41
1:A:1035:ALA:HA	3:D:-2:DC:OP1	2.21	0.41
2:C:-6:DG:N2	3:D:-1:DT:O2	2.54	0.40
1:B:399:TYR:CE2	1:B:406:LEU:HD13	2.56	0.40
1:B:713:VAL:CG2	1:B:930:PHE:HB2	2.50	0.40
1:A:226:GLU:HA	1:A:245:ALA:HB2	2.03	0.40
1:B:1003:LEU:HD21	1:B:1187:LEU:HD12	2.03	0.40
1:A:176:THR:CG2	1:A:177:PRO:HD3	2.46	0.40
1:A:478:LYS:O	1:A:483:TYR:HD1	2.03	0.40
1:A:609:ILE:O	1:A:613:ILE:HG13	2.21	0.40
1:A:129:LEU:HD11	1:A:332:PRO:O	2.21	0.40
1:A:785:ARG:NH2	1:A:811:LYS:HD2	2.36	0.40
1:A:1079:GLN:NE2	1:A:1138:LEU:HB2	2.37	0.40
1:B:879:SER:OG	1:B:895:ARG:HB3	2.22	0.40
1:A:833:VAL:O	1:A:837:VAL:HG23	2.21	0.40
1:A:1054:VAL:HG21	1:A:1074:TYR:HB3	2.04	0.40
1:B:168:HIS:HB2	1:B:171:PHE:CE1	2.56	0.40
1:B:1188:LEU:HD13	1:B:1188:LEU:HA	1.88	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	1038/1163 (89%)	974 (94%)	61 (6%)	3 (0%)	41	75
1	B	1046/1163 (90%)	975 (93%)	69 (7%)	2 (0%)	47	81
All	All	2084/2326 (90%)	1949 (94%)	130 (6%)	5 (0%)	47	81

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	217	GLN
1	A	177	PRO
1	A	1140	SER
1	B	177	PRO
1	A	236	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	796/951 (84%)	749 (94%)	47 (6%)	19	53
1	B	836/951 (88%)	759 (91%)	77 (9%)	9	36
All	All	1632/1902 (86%)	1508 (92%)	124 (8%)	13	43

All (124) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	101	ARG
1	A	176	THR
1	A	224	LEU
1	A	240	PHE
1	A	285	ASN
1	A	322	ARG
1	A	329	PRO
1	A	397	LEU
1	A	411	LEU
1	A	426	GLU
1	A	471	ASP
1	A	499	PHE
1	A	550	ASP
1	A	607	ILE
1	A	615	ASP
1	A	639	THR
1	A	746	LEU
1	A	765	VAL
1	A	821	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	827	LEU
1	A	846	LEU
1	A	862	LEU
1	A	881	ARG
1	A	915	ARG
1	A	919	LEU
1	A	924	LEU
1	A	928	LYS
1	A	934	LEU
1	A	938	LYS
1	A	950	MET
1	A	951	LEU
1	A	963	CYS
1	A	966	ILE
1	A	975	ASP
1	A	981	ASP
1	A	982	THR
1	A	999	LEU
1	A	1039	ARG
1	A	1045	THR
1	A	1062	ARG
1	A	1064	GLN
1	A	1070	ASP
1	A	1138	LEU
1	A	1139	VAL
1	A	1159	ASP
1	A	1164	HIS
1	A	1171	VAL
1	B	45	TYR
1	B	170	ARG
1	B	172	MET
1	B	176	THR
1	B	178	THR
1	B	196	VAL
1	B	214	ARG
1	B	217	GLN
1	B	224	LEU
1	B	240	PHE
1	B	246	ASP
1	B	252	VAL
1	B	255	ARG
1	B	258	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	262	GLU
1	B	285	ASN
1	B	292	LYS
1	B	305	LEU
1	B	322	ARG
1	B	325	THR
1	B	329	PRO
1	B	356	MET
1	B	388	GLU
1	B	397	LEU
1	B	398	LEU
1	B	424	LEU
1	B	434	THR
1	B	471	ASP
1	B	499	PHE
1	B	500	ARG
1	B	501	VAL
1	B	510	GLN
1	B	511	LYS
1	B	512	ARG
1	B	514	LYS
1	B	552	LYS
1	B	603	ARG
1	B	615	ASP
1	B	619	ILE
1	B	639	THR
1	B	693	HIS
1	B	705	THR
1	B	710	VAL
1	B	714	VAL
1	B	756	GLU
1	B	761	ARG
1	B	765	VAL
1	B	791	ARG
1	B	821	THR
1	B	827	LEU
1	B	845	LEU
1	B	862	LEU
1	B	871	ASP
1	B	881	ARG
1	B	919	LEU
1	B	924	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	942	ILE
1	B	952	ILE
1	B	956	ASP
1	B	958	VAL
1	B	959	ARG
1	B	963	CYS
1	B	973	LEU
1	B	975	ASP
1	B	981	ASP
1	B	983	VAL
1	B	996	GLU
1	B	999	LEU
1	B	1029	GLN
1	B	1030	ASP
1	B	1039	ARG
1	B	1045	THR
1	B	1058	LEU
1	B	1086	THR
1	B	1159	ASP
1	B	1164	HIS
1	B	1188	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	YE4	E	1201	-	31,33,33	4.14	17 (54%)	41,46,46	1.77	6 (14%)
4	YE4	C	1201	-	31,33,33	4.15	17 (54%)	41,46,46	1.83	7 (17%)

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	YE4	E	1201	-	-	0/13/21/21	0/4/4/4
4	YE4	C	1201	-	-	0/13/21/21	0/4/4/4

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	1201	YE4	C15-N16	12.88	1.49	1.33
4	E	1201	YE4	C15-N16	12.83	1.49	1.33
4	C	1201	YE4	C22-N23	-9.76	1.28	1.47
4	E	1201	YE4	C22-N23	-9.71	1.28	1.47
4	C	1201	YE4	C15-C03	7.24	1.51	1.39
4	E	1201	YE4	C15-C03	7.19	1.51	1.39
4	E	1201	YE4	C30-C18	6.53	1.46	1.41
4	C	1201	YE4	C30-C18	6.42	1.46	1.41
4	E	1201	YE4	C18-N16	5.67	1.52	1.39
4	C	1201	YE4	C18-N16	5.62	1.52	1.39
4	E	1201	YE4	C04-N05	5.35	1.45	1.33
4	C	1201	YE4	C04-N05	5.33	1.45	1.33
4	E	1201	YE4	C22-C21	4.70	1.59	1.51
4	C	1201	YE4	C22-C21	4.66	1.59	1.51
4	C	1201	YE4	O01-C02	-3.75	1.16	1.23
4	E	1201	YE4	O01-C02	-3.73	1.17	1.23
4	C	1201	YE4	C20-C21	3.58	1.46	1.38
4	E	1201	YE4	C20-C21	3.57	1.46	1.38
4	E	1201	YE4	C24-N23	-3.40	1.37	1.46

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	1201	YE4	C24-N23	-3.39	1.37	1.46
4	E	1201	YE4	C19-C20	3.02	1.43	1.36
4	C	1201	YE4	C10-CL11	3.00	1.81	1.74
4	C	1201	YE4	C03-C04	2.99	1.55	1.50
4	C	1201	YE4	C19-C20	2.98	1.42	1.36
4	E	1201	YE4	C10-CL11	2.95	1.80	1.74
4	E	1201	YE4	C03-C04	2.94	1.55	1.50
4	E	1201	YE4	C09-C10	2.60	1.43	1.38
4	C	1201	YE4	C09-C10	2.57	1.42	1.38
4	C	1201	YE4	C12-C13	2.47	1.43	1.38
4	E	1201	YE4	C12-C13	2.34	1.43	1.38
4	E	1201	YE4	C28-N23	-2.34	1.40	1.46
4	C	1201	YE4	C28-N23	-2.33	1.40	1.46
4	C	1201	YE4	C06-C07	2.23	1.56	1.51
4	E	1201	YE4	C06-C07	2.15	1.56	1.51

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	1201	YE4	C25-C24-N23	6.18	119.47	110.10
4	E	1201	YE4	C25-C24-N23	5.88	119.02	110.10
4	E	1201	YE4	C28-N23-C24	4.73	119.48	108.83
4	C	1201	YE4	C28-N23-C24	4.56	119.09	108.83
4	E	1201	YE4	C27-C28-N23	4.08	116.29	110.10
4	C	1201	YE4	C27-C28-N23	4.05	116.24	110.10
4	C	1201	YE4	C03-C04-N05	3.51	121.05	116.30
4	E	1201	YE4	O26-C25-C24	3.13	118.69	111.80
4	C	1201	YE4	O26-C25-C24	3.07	118.56	111.80
4	E	1201	YE4	C03-C04-N05	2.95	120.28	116.30
4	E	1201	YE4	C27-O26-C25	2.12	116.96	109.89
4	C	1201	YE4	C27-O26-C25	2.09	116.87	109.89
4	C	1201	YE4	C15-N16-C18	-2.04	119.80	121.59

There are no chirality outliers.

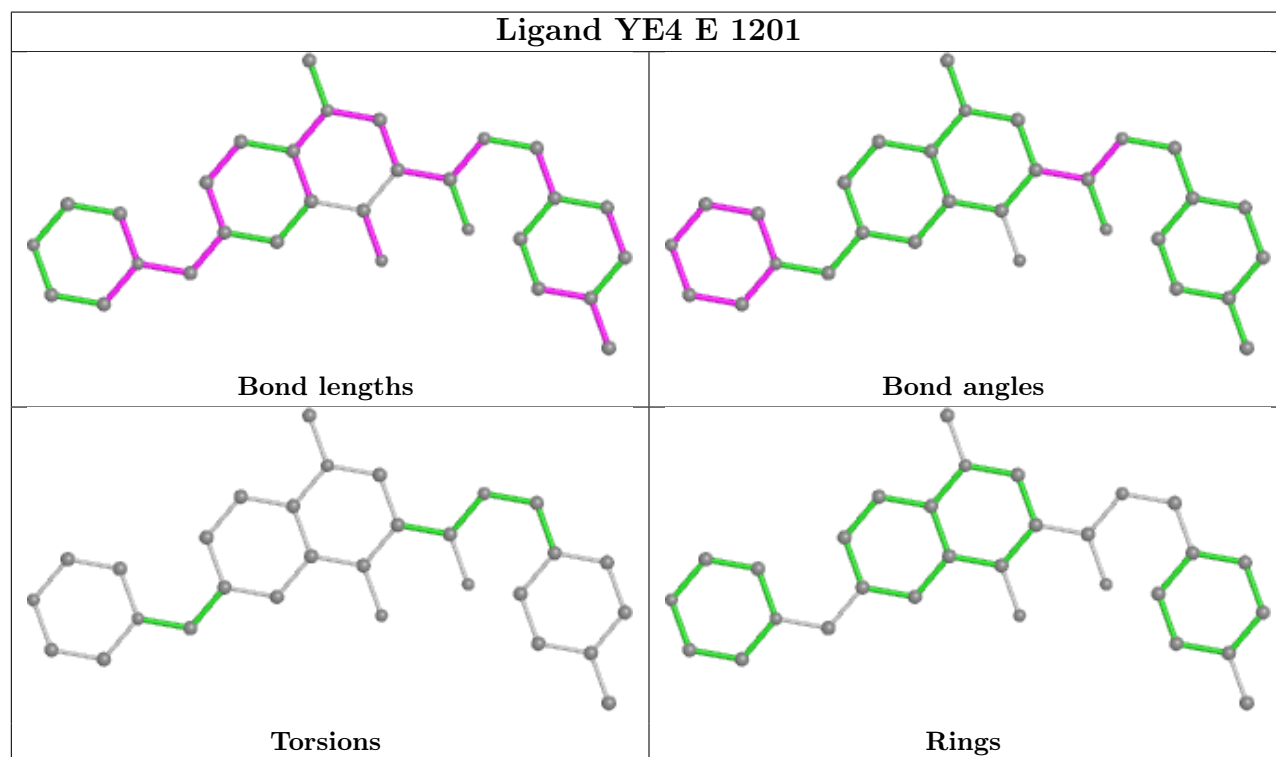
There are no torsion outliers.

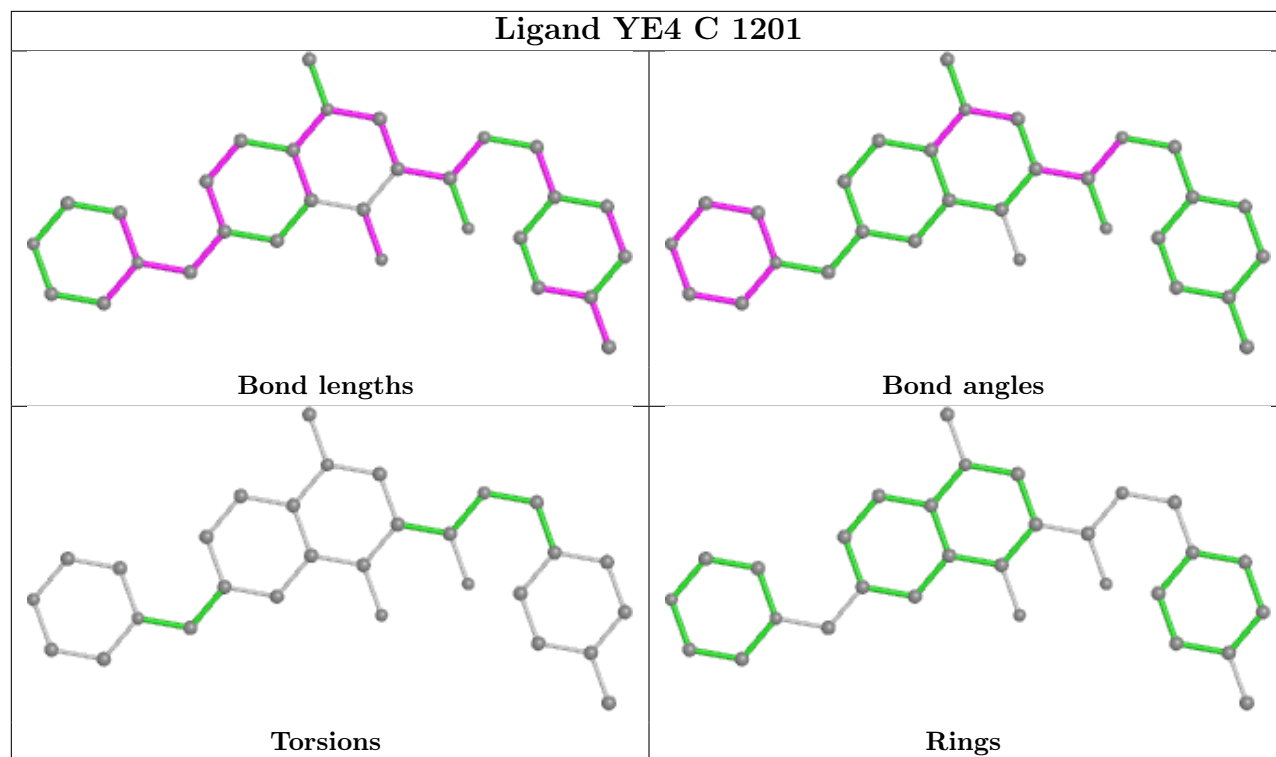
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	1201	YE4	1	0

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	1044/1163 (89%)	-0.26	1 (0%) 95 95	46, 107, 156, 210	0
1	B	1054/1163 (90%)	-0.24	1 (0%) 95 95	52, 109, 154, 196	0
2	C	18/27 (66%)	-0.19	2 (11%) 5 6	88, 132, 243, 251	0
2	E	18/27 (66%)	-0.29	0 100 100	96, 126, 236, 242	0
3	D	15/23 (65%)	-0.12	0 100 100	96, 136, 238, 240	0
3	F	15/23 (65%)	-0.17	1 (6%) 17 16	100, 128, 236, 238	0
All	All	2164/2426 (89%)	-0.25	5 (0%) 95 93	46, 109, 158, 251	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	5	DG	3.1
3	F	-13	DC	2.8
2	C	4	DC	2.2
1	A	1138	LEU	2.1
1	B	122	PHE	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](#)

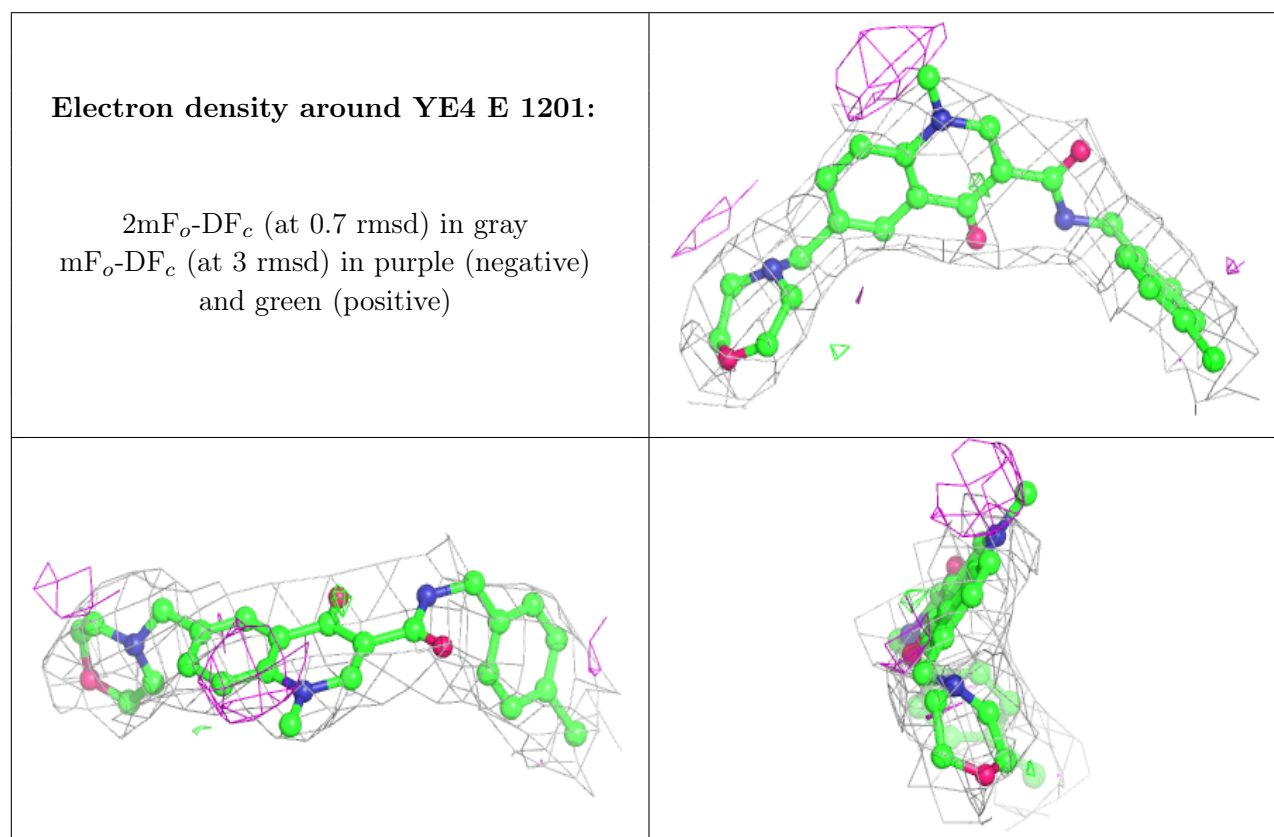
There are no monosaccharides in this entry.

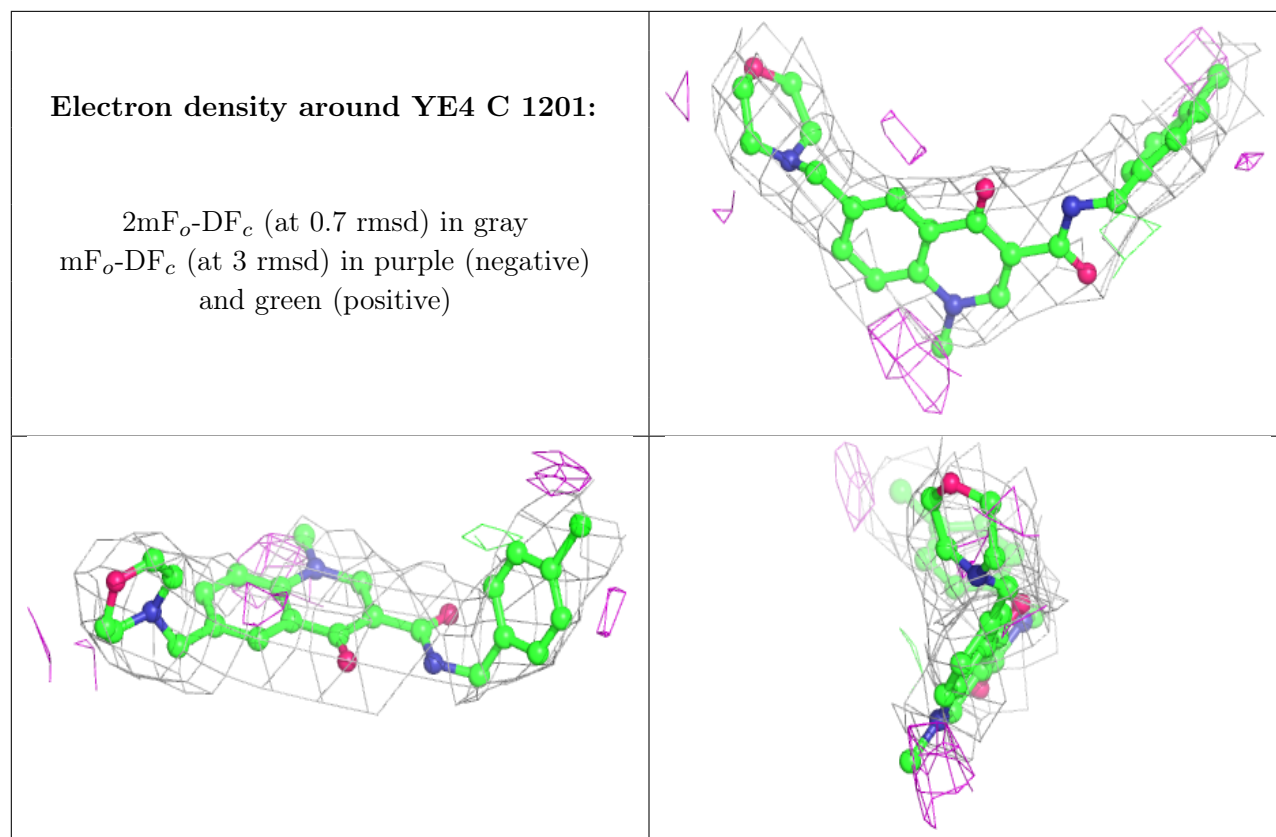
6.4 Ligands [i](#)

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(\AA^2)	Q<0.9
4	YE4	E	1201	30/30	0.93	0.25	70,87,95,99	0
4	YE4	C	1201	30/30	0.95	0.26	63,79,88,95	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.