Full wwPDB X-ray Structure Validation Report (i)

Sep 10, 2023 – 03:44 AM EDT

PDB ID	:	4H4O
Title	:	Crystal Structure of HIV-1 Reverse Transcriptase (RT) in Complex with (E)-3-(3-(2-(2-(2,4-dioxo-3,4-dihydropyrimidin-1(2H)-yl)ethoxy)- 4-fluorophenoxy)-5-fluorophenyl)acrylonitrile (JLJ506), A Non-nucleoside inhibitor
Authors	:	Frey, K.M.; Anderson, K.S.
Deposited on	:	2012-09-17
Resolution	:	2.90 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 (i) 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 (i)) were used in the production of this report:

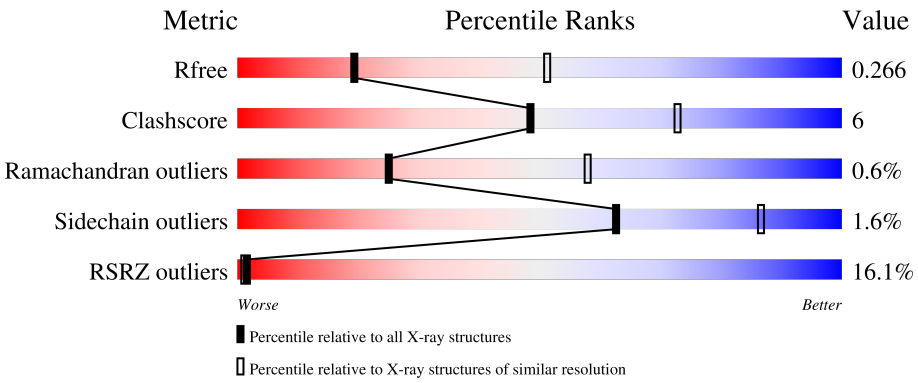
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 <=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	557	16% 85% 13% •
2	B	428	16% 82% 17% •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 8082 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 Reverse transcriptase/ribonuclease H, Exoribonuclease H, p66 RT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	549	4465	2891	742	825	7	0	0	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	initiating methionine	UNP P03366
A	0	VAL	-	expression tag	UNP P03366
A	172	ALA	LYS	engineered mutation	UNP P03366
A	173	ALA	LYS	engineered mutation	UNP P03366
A	280	SER	CYS	engineered mutation	UNP P03366

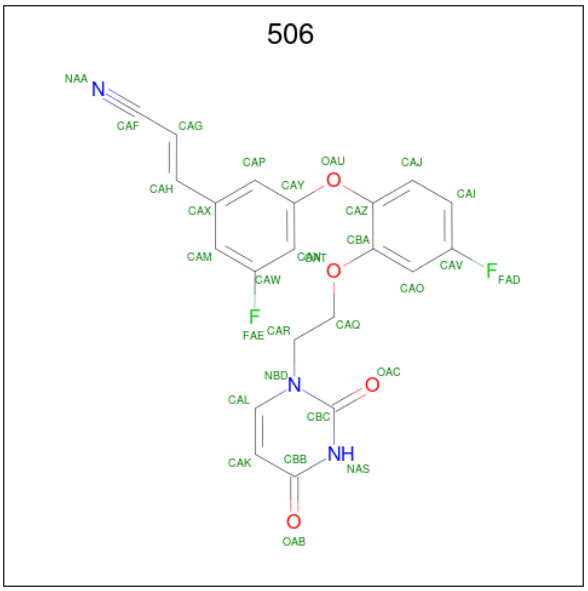
- Molecule 2 is a protein called Reverse transcriptase/ribonuclease H, Exoribonuclease H, p51 RT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	428	3538	2305	586	640	7	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	280	SER	CYS	engineered mutation	UNP P03366

- Molecule 3 is (2E)-3-(3-{2-[2-(2,4-dioxo-3,4-dihydropyrimidin-1(2H)-yl)ethoxy]-4-fluorophenoxy}-5-fluorophenyl)prop-2-enenitrile (three-letter code: 506) (formula: $C_{21}H_{15}F_2N_3O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	F	N	O		
3	A	1	30	21	2	3	4	0	0

- Molecule 4 is water.

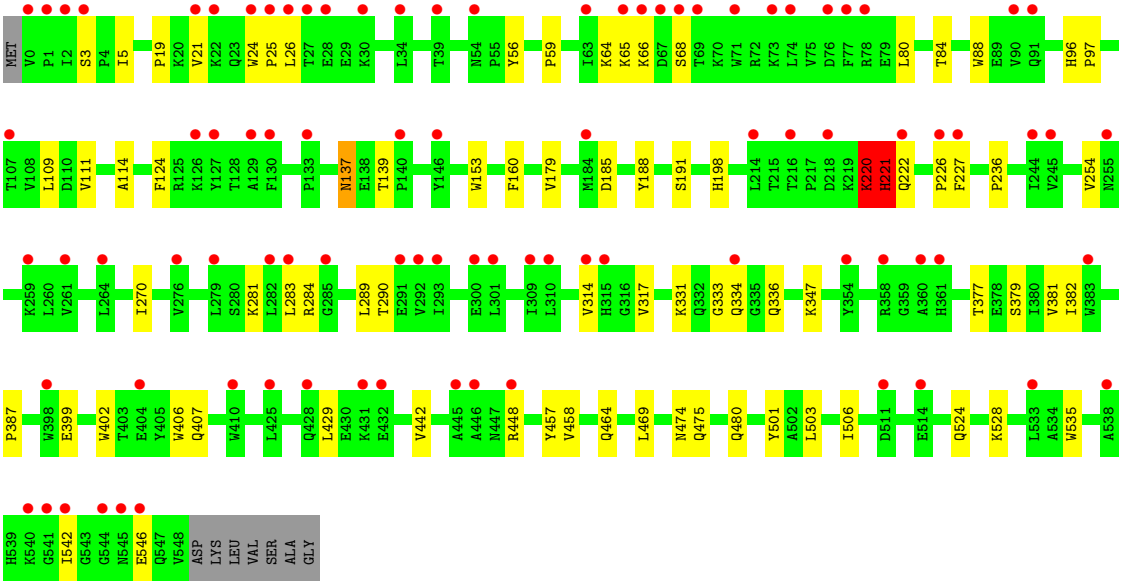
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	22	22	22	0	0
			Total	O		
4	B	27	27	27	0	0

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.

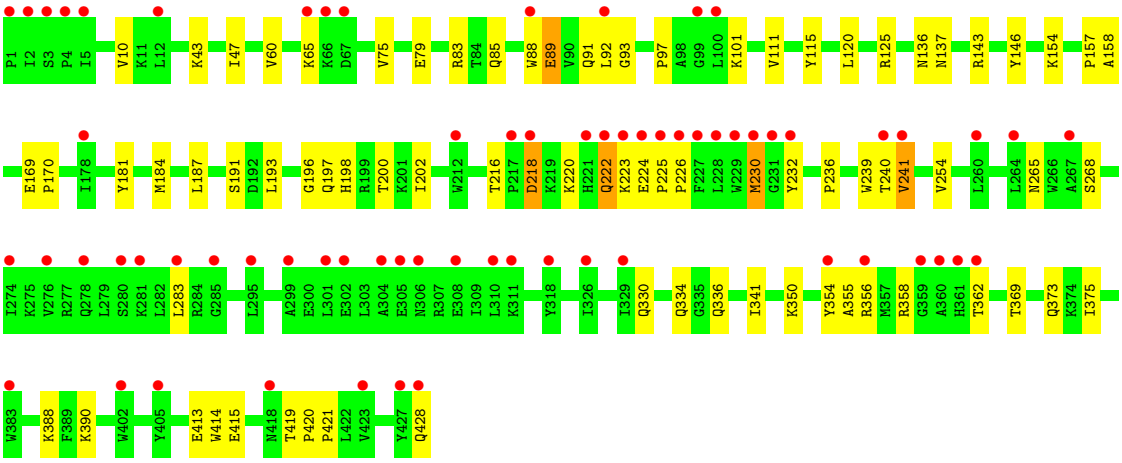
• Molecule 1: Reverse transcriptase/ribonuclease H, Exoribonuclease H, p66 RT

Chain A:



• Molecule 2: Reverse transcriptase/ribonuclease H, Exoribonuclease H, p51 RT

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	225.45Å 69.58Å 104.40Å 90.00° 106.21° 90.00°	Depositor
Resolution (Å)	41.72 – 2.90 41.72 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.8 (41.72-2.90) 99.8 (41.72-2.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	2.27 (at 2.90Å)	Xtriage
Refinement program	PHENIX 1.8_1069	Depositor
R, R_{free}	0.232 , 0.265 0.234 , 0.266	Depositor DCC
R_{free} test set	2000 reflections (5.75%)	wwPDB-VP
Wilson B-factor (Å^2)	79.0	Xtriage
Anisotropy	0.256	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.27 , 113.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	8082	wwPDB-VP
Average B, all atoms (Å^2)	164.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: 506

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	0.23	0/4583	0.45	0/6231
2	B	0.24	0/3643	0.48	1/4952 (0.0%)
All	All	0.24	0/8226	0.46	1/11183 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
2	B	0	1
All	All	0	2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	89	GLU	N-CA-C	-5.05	97.36	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	220	LYS	Peptide
2	B	419	THR	Peptide

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	4465	0	4516	49	0
2	B	3538	0	3576	52	0
3	A	30	0	15	2	0
4	A	22	0	0	0	1
4	B	27	0	0	1	0
All	All	8082	0	8107	96	1

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 6.

All (96) 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:220:LYS:HB3	1:A:222:GLN:H	1.51	0.76
1:A:331:LYS:HE2	1:A:333:GLY:HA2	1.68	0.76
1:A:191:SER:OG	1:A:198:HIS:ND1	2.23	0.68
1:A:64:LYS:HD2	1:A:68:SER:HA	1.75	0.67
1:A:226:PRO:HA	1:A:227:PHE:HB2	1.77	0.66
1:A:317:VAL:HG11	1:A:347:LYS:HD3	1.79	0.65
2:B:137:ASN:O	4:B:505:HOH:O	2.15	0.64
2:B:336:GLN:HG2	2:B:355:ALA:HB2	1.80	0.63
1:A:84:THR:HG21	1:A:153:TRP:HE1	1.61	0.63
2:B:390:LYS:NZ	2:B:415:GLU:OE2	2.34	0.61
1:A:88:TRP:CE2	2:B:143:ARG:HD2	2.37	0.59
1:A:334:GLN:HB2	1:A:336:GLN:HE21	1.68	0.58
2:B:157:PRO:HG3	2:B:184:MET:HA	1.86	0.58
2:B:354:TYR:HB3	2:B:356:ARG:HH12	1.71	0.56
1:A:220:LYS:O	1:A:221:HIS:HB2	2.04	0.55
2:B:191:SER:OG	2:B:198:HIS:ND1	2.30	0.54
1:A:220:LYS:CB	1:A:222:GLN:H	2.20	0.54
1:A:270:ILE:HG13	1:A:314:VAL:HG13	1.90	0.54
2:B:91:GLN:HG2	2:B:92:LEU:HD12	1.88	0.53
2:B:111:VAL:HG11	2:B:187:LEU:HD22	1.90	0.52
1:A:399:GLU:HA	1:A:402:TRP:CD1	2.45	0.52
1:A:84:THR:HG22	1:A:124:PHE:HZ	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:236:PRO:HA	2:B:239:TRP:CG	2.46	0.51
1:A:88:TRP:CD2	2:B:143:ARG:HD2	2.46	0.51
2:B:79:GLU:OE2	2:B:83:ARG:NH1	2.44	0.50
1:A:281:LYS:HG3	1:A:284:ARG:HE	1.76	0.50
2:B:89:GLU:O	2:B:93:GLY:HA3	2.12	0.50
2:B:47:ILE:HG22	2:B:146:TYR:HA	1.93	0.50
2:B:97:PRO:HG3	2:B:181:TYR:HB2	1.94	0.49
2:B:239:TRP:O	2:B:350:LYS:HE2	2.13	0.49
2:B:92:LEU:HB3	2:B:158:ALA:HB1	1.95	0.48
1:A:114:ALA:HB1	1:A:160:PHE:CE1	2.49	0.48
1:A:220:LYS:HA	1:A:220:LYS:HD2	1.55	0.48
2:B:336:GLN:OE1	2:B:428:GLN:NE2	2.46	0.48
2:B:240:THR:OG1	2:B:241:VAL:N	2.46	0.48
2:B:254:VAL:HG13	2:B:283:LEU:HD22	1.96	0.48
2:B:334:GLN:O	2:B:358:ARG:NH1	2.47	0.47
1:A:19:PRO:O	1:A:56:TYR:HB3	2.14	0.47
1:A:254:VAL:HG13	1:A:283:LEU:HD11	1.96	0.47
1:A:475:GLN:HB3	1:A:501:TYR:CE2	2.49	0.47
1:A:236:PRO:HA	3:A:601:506:CBB	2.44	0.47
1:A:399:GLU:HA	1:A:402:TRP:NE1	2.29	0.47
2:B:88:TRP:CD1	2:B:88:TRP:N	2.84	0.46
2:B:120:LEU:HD23	2:B:125:ARG:HG2	1.98	0.46
2:B:354:TYR:HB3	2:B:356:ARG:NH1	2.29	0.46
1:A:111:VAL:HG23	1:A:114:ALA:HB2	1.98	0.46
1:A:65:LYS:HG2	1:A:66:LYS:H	1.80	0.46
2:B:193:LEU:HB3	2:B:197:GLN:HG3	1.98	0.45
2:B:222:GLN:HB3	2:B:223:LYS:H	1.59	0.45
1:A:21:VAL:HB	1:A:59:PRO:HD3	1.98	0.45
2:B:420:PRO:HA	2:B:421:PRO:HD3	1.65	0.45
1:A:448:ARG:HE	1:A:474:ASN:ND2	2.15	0.45
1:A:25:PRO:HA	1:A:26:LEU:HA	1.72	0.45
1:A:442:VAL:HG12	1:A:457:TYR:HB3	1.99	0.45
1:A:406:TRP:CH2	2:B:420:PRO:HD2	2.52	0.45
1:A:406:TRP:CZ3	1:A:407:GLN:HG3	2.52	0.45
2:B:115:TYR:OH	2:B:184:MET:O	2.33	0.45
1:A:379:SER:CB	1:A:387:PRO:HD3	2.47	0.44
1:A:377:THR:O	1:A:381:VAL:HG23	2.17	0.44
1:A:111:VAL:HG22	1:A:185:ASP:O	2.17	0.44
1:A:254:VAL:HB	1:A:289:LEU:O	2.17	0.44
2:B:220:LYS:HA	2:B:220:LYS:HD3	1.68	0.44
2:B:265:ASN:O	2:B:268:SER:OG	2.27	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:PRO:HG3	1:A:80:LEU:HB2	1.98	0.44
2:B:10:VAL:HG22	2:B:88:TRP:CH2	2.52	0.44
2:B:236:PRO:HA	2:B:239:TRP:CD2	2.53	0.44
2:B:220:LYS:C	2:B:222:GLN:H	2.21	0.44
2:B:43:LYS:HE3	2:B:43:LYS:HB2	1.86	0.44
2:B:101:LYS:O	2:B:236:PRO:HB2	2.17	0.44
1:A:3:SER:HB3	1:A:5:ILE:HG13	2.00	0.43
1:A:188:TYR:CZ	3:A:601:506:H14	2.53	0.43
1:A:96:HIS:HA	1:A:97:PRO:HD3	1.92	0.43
2:B:341:ILE:HD11	2:B:375:ILE:HG23	2.00	0.43
1:A:469:LEU:HD11	1:A:480:GLN:HG2	2.00	0.43
2:B:60:VAL:HG12	2:B:75:VAL:HG22	1.99	0.43
2:B:369:THR:HG22	2:B:373:GLN:NE2	2.34	0.43
1:A:524:GLN:O	1:A:528:LYS:HG2	2.19	0.42
1:A:458:VAL:HA	1:A:464:GLN:HB3	2.01	0.42
2:B:198:HIS:O	2:B:202:ILE:HG12	2.20	0.42
1:A:429:LEU:HD11	1:A:506:ILE:HG22	2.02	0.42
1:A:24:TRP:N	1:A:25:PRO:HD2	2.35	0.42
1:A:503:LEU:HD22	1:A:535:TRP:HB2	2.01	0.42
2:B:224:GLU:HA	2:B:225:PRO:HD2	1.96	0.42
1:A:382:ILE:O	2:B:136:ASN:HB2	2.20	0.41
2:B:187:LEU:HD12	2:B:187:LEU:HA	1.90	0.41
1:A:458:VAL:HG22	1:A:464:GLN:HB3	2.01	0.41
1:A:406:TRP:CZ3	2:B:420:PRO:HD2	2.56	0.41
2:B:85:GLN:HG3	2:B:154:LYS:HB3	2.02	0.41
2:B:388:LYS:HD2	2:B:413:GLU:HB2	2.02	0.41
2:B:216:THR:O	2:B:218:ASP:N	2.49	0.41
2:B:169:GLU:HB3	2:B:170:PRO:HD3	2.03	0.41
2:B:196:GLY:O	2:B:200:THR:HG23	2.21	0.40
2:B:236:PRO:O	2:B:239:TRP:HB2	2.21	0.40
1:A:109:LEU:HD21	1:A:221:HIS:HE1	1.87	0.40
2:B:369:THR:HG22	2:B:373:GLN:HE21	1.85	0.40
2:B:220:LYS:HG3	2:B:230:MET:HG3	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:718:HOH:O	4:A:718:HOH:O[2_756]	2.16	0.04

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	547/557 (98%)	519 (95%)	24 (4%)	4 (1%)	22	54
2	B	426/428 (100%)	397 (93%)	27 (6%)	2 (0%)	29	61
All	All	973/985 (99%)	916 (94%)	51 (5%)	6 (1%)	25	58

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	221	HIS
1	A	220	LYS
1	A	137	ASN
2	B	226	PRO
1	A	542	ILE
2	B	65	LYS

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	489/495 (99%)	483 (99%)	6 (1%)	71	91
2	B	390/390 (100%)	382 (98%)	8 (2%)	53	81
All	All	879/885 (99%)	865 (98%)	14 (2%)	62	86

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	137	ASN
1	A	139	THR
1	A	179	VAL
1	A	221	HIS
1	A	290	THR
1	A	546	GLU
2	B	218	ASP
2	B	222	GLN
2	B	230	MET
2	B	232	TYR
2	B	241	VAL
2	B	330	GLN
2	B	362	THR
2	B	414	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	361	HIS
1	A	474	ASN
2	B	197	GLN
2	B	278	GLN

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)

1 ligand is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	506	A	601	-	32,32,32	1.58	5 (15%)	42,43,43	1.79	9 (21%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	506	A	601	-	-	5/13/14/14	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	506	CBC-NBD	4.48	1.44	1.37
3	A	601	506	CAL-NBD	2.48	1.40	1.37
3	A	601	506	CAZ-CBA	2.42	1.45	1.40
3	A	601	506	CAM-CAW	2.19	1.41	1.37
3	A	601	506	CBB-NAS	2.00	1.42	1.38

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	601	506	OAT-CBA-CAZ	5.81	127.74	115.73
3	A	601	506	OAT-CBA-CAO	-3.97	114.45	123.58
3	A	601	506	NAS-CBC-NBD	3.73	118.27	114.86
3	A	601	506	CAX-CAP-CAY	3.06	122.62	119.91
3	A	601	506	CBA-CAO-CAV	2.88	122.31	116.45
3	A	601	506	CAL-NBD-CBC	-2.73	118.62	121.04
3	A	601	506	OAU-CAZ-CBA	2.30	124.08	116.87
3	A	601	506	CAI-CAV-CAO	-2.08	120.59	123.29
3	A	601	506	CAI-CAJ-CAZ	2.07	124.00	120.06

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	601	506	CAZ-CBA-OAT-CAQ
3	A	601	506	CAO-CBA-OAT-CAQ
3	A	601	506	CAG-CAH-CAX-CAM
3	A	601	506	CAG-CAH-CAX-CAP
3	A	601	506	OAT-CAQ-CAR-NBD

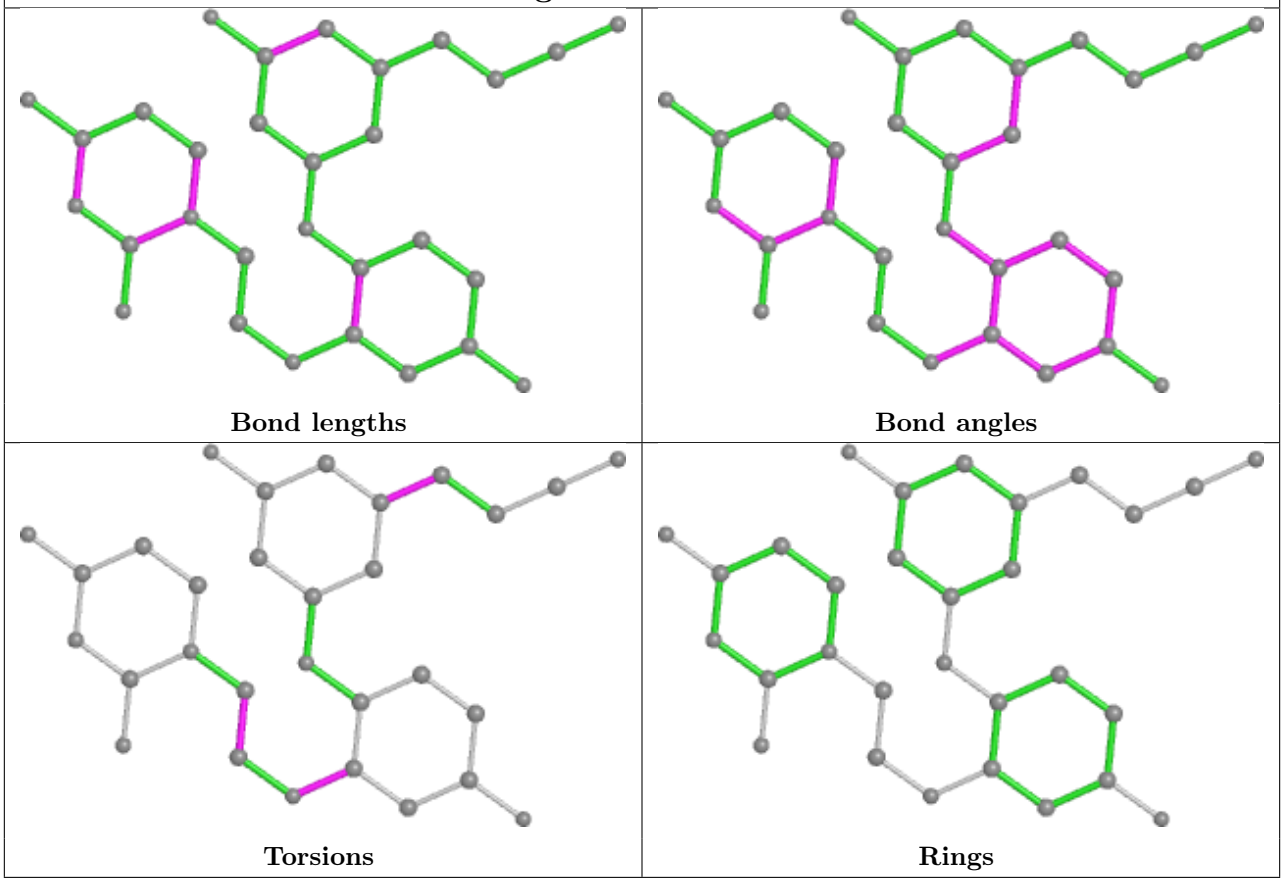
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	601	506	2	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 then 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.

Ligand 506 A 601



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	549/557 (98%)	1.05	90 (16%)	1	1	112, 157, 242, 303	1 (0%)
2	B	428/428 (100%)	1.22	67 (15%)	2	1	114, 148, 248, 369	1 (0%)
All	All	977/985 (99%)	1.12	157 (16%)	1	1	112, 154, 246, 369	2 (0%)

All (157) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	1	PRO	37.0
2	B	2	ILE	18.6
1	A	541	GLY	17.2
2	B	224	GLU	16.2
2	B	231	GLY	15.9
1	A	25	PRO	15.0
1	A	545	ASN	12.6
1	A	24	TRP	12.4
2	B	4	PRO	12.1
2	B	67	ASP	11.0
1	A	67	ASP	10.2
1	A	542	ILE	9.1
1	A	2	ILE	9.1
2	B	226	PRO	8.7
2	B	225	PRO	7.9
2	B	217	PRO	7.6
2	B	221	HIS	7.5
2	B	66	LYS	7.4
2	B	3	SER	7.4
2	B	229	TRP	7.3
2	B	227	PHE	6.6
2	B	228	LEU	6.6
2	B	428	GLN	6.4
2	B	361	HIS	6.1

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Mol	Chain	Res	Type	RSRZ
2	B	360	ALA	5.9
2	B	418	ASN	5.9
1	A	26	LEU	5.8
1	A	546	GLU	5.8
1	A	448	ARG	5.7
2	B	301	LEU	5.7
2	B	310	LEU	5.5
2	B	354	TYR	5.5
2	B	230	MET	5.4
1	A	360	ALA	5.3
1	A	264	LEU	5.1
1	A	68	SER	4.9
1	A	107	THR	4.8
1	A	22	LYS	4.8
1	A	354	TYR	4.8
1	A	28	GLU	4.7
1	A	21	VAL	4.7
2	B	88	TRP	4.7
1	A	133	PRO	4.6
1	A	227	PHE	4.6
2	B	5	ILE	4.6
1	A	226	PRO	4.5
1	A	63	ILE	4.4
2	B	311	LYS	4.4
1	A	130	PHE	4.4
2	B	218	ASP	4.3
1	A	255	ASN	4.2
1	A	0	VAL	4.2
2	B	283	LEU	4.2
2	B	299	ALA	4.2
2	B	295	LEU	4.1
1	A	309	ILE	4.1
1	A	514	GLU	4.1
1	A	77	PHE	4.1
1	A	283	LEU	4.0
2	B	304	ALA	4.0
1	A	358	ARG	4.0
1	A	54	ASN	4.0
1	A	261	VAL	3.9
1	A	129	ALA	3.9
2	B	326	ILE	3.8
1	A	540	LYS	3.7

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Mol	Chain	Res	Type	RSRZ
2	B	285	GLY	3.6
1	A	291	GLU	3.5
2	B	329	ILE	3.5
2	B	306	ASN	3.5
1	A	74	LEU	3.5
1	A	222	GLN	3.5
1	A	446	ALA	3.5
1	A	216	THR	3.5
1	A	71	TRP	3.4
2	B	318	TYR	3.4
1	A	361	HIS	3.4
2	B	65	LYS	3.4
2	B	302	GLU	3.3
1	A	245	VAL	3.3
1	A	78	ARG	3.3
2	B	305	GLU	3.3
1	A	293	ILE	3.2
2	B	232	TYR	3.2
1	A	30	LYS	3.2
1	A	282	LEU	3.2
1	A	445	ALA	3.1
1	A	91	GLN	3.1
1	A	431	LYS	3.1
1	A	544	GLY	3.0
2	B	359	GLY	3.0
1	A	140	PRO	2.9
1	A	538	ALA	2.9
1	A	218	ASP	2.9
1	A	410	TRP	2.9
1	A	127	TYR	2.9
2	B	427	TYR	2.8
1	A	3	SER	2.8
2	B	362	THR	2.8
1	A	76	ASP	2.8
2	B	264	LEU	2.7
2	B	356	ARG	2.7
2	B	423	VAL	2.7
1	A	300	GLU	2.7
1	A	34	LEU	2.7
1	A	292	VAL	2.7
1	A	279	LEU	2.7
2	B	405	TYR	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	334	GLN	2.6
1	A	404	GLU	2.6
1	A	90	VAL	2.6
2	B	100	LEU	2.6
1	A	398	TRP	2.6
1	A	66	LYS	2.5
2	B	241	VAL	2.5
1	A	276	VAL	2.5
2	B	280	SER	2.5
1	A	126	LYS	2.5
1	A	315	HIS	2.5
1	A	285	GLY	2.5
1	A	244	ILE	2.5
1	A	259	LYS	2.5
1	A	27	THR	2.5
2	B	276	VAL	2.4
1	A	428	GLN	2.4
1	A	511	ASP	2.4
1	A	73	LYS	2.4
1	A	533	LEU	2.4
1	A	310	LEU	2.4
2	B	281	LYS	2.4
2	B	274	ILE	2.3
2	B	278	GLN	2.3
2	B	308	GLU	2.3
1	A	69	THR	2.3
2	B	383	TRP	2.3
2	B	267	ALA	2.3
1	A	432	GLU	2.2
1	A	1	PRO	2.2
1	A	214	LEU	2.2
1	A	39	THR	2.2
1	A	425	LEU	2.2
2	B	402	TRP	2.2
2	B	178	ILE	2.2
2	B	12	LEU	2.2
2	B	260	LEU	2.1
1	A	65	LYS	2.1
1	A	383	TRP	2.1
1	A	184	MET	2.1
2	B	212	TRP	2.1
2	B	92	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
2	B	222	GLN	2.1
2	B	99	GLY	2.1
2	B	240	THR	2.1
1	A	146	TYR	2.0
2	B	223	LYS	2.0
1	A	314	VAL	2.0
1	A	301	LEU	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)

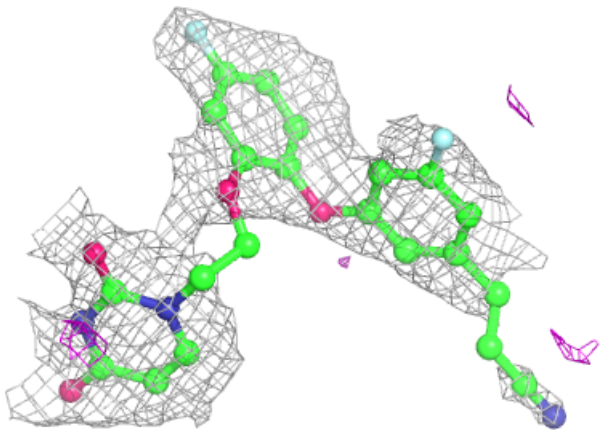
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

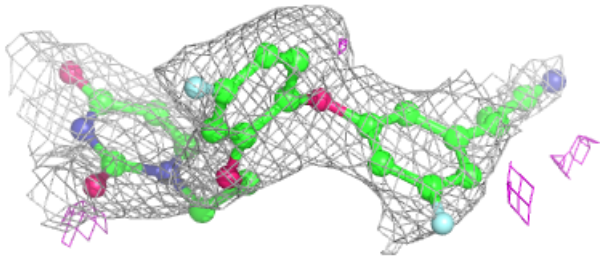
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	506	A	601	30/30	0.80	0.39	125,145,153,160	0

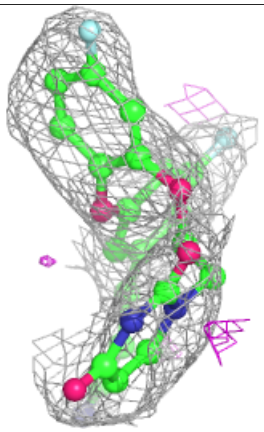
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 506 A 601:

$2mF_o$-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.