



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 03:45 AM UTC

PDB ID : 7OTZ / pdb_00007otz
Title : HIV-1 REVERSE TRANSCRIPTASE COMPLEX WITH DNA AND INHIBITOR RMC-259
Authors : Martinez, S.E.; Singh, A.K.; Gu, W.; Das, K.
Deposited on : 2021-06-10
Resolution : 3.10 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

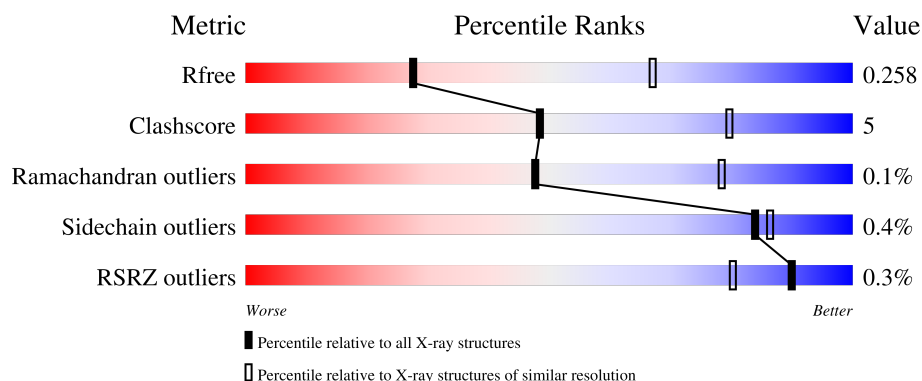
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	556	82% 18%
1	C	556	86% 14%
2	B	428	83% 12%
2	D	428	80% 15%
3	E	27	52% 33%

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Mol	Chain	Length	Quality of chain
3	T	27	 63% 22% 11%
4	F	21	 62% 33% 5%
4	P	21	 81% 14% 5%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 17619 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	555	Total	C	N	O	S	0	0	0
			4509	2919	750	832	8			
1	A	556	Total	C	N	O	S	0	0	0
			4521	2927	752	833	9			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	MET	-	initiating methionine	UNP P03366
C	0	VAL	-	expression tag	UNP P03366
C	258	CYS	GLN	conflict	UNP P03366
C	280	SER	CYS	conflict	UNP P03366
C	498	ASN	ASP	conflict	UNP P03366
A	-1	MET	-	initiating methionine	UNP P03366
A	0	VAL	-	expression tag	UNP P03366
A	258	CYS	GLN	conflict	UNP P03366
A	280	SER	CYS	conflict	UNP P03366
A	498	ASN	ASP	conflict	UNP P03366

- Molecule 2 is a protein called Reverse transcriptase/ribonuclease H.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	411	Total	C	N	O	S	0	0	0
			3398	2213	563	616	6			
2	B	412	Total	C	N	O	S	0	0	0
			3391	2206	560	619	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	280	SER	CYS	conflict	UNP P03366
B	280	SER	CYS	conflict	UNP P03366

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	23	Total	C	N	O	P	0	0	0
			477	223	95	136	23			
3	T	24	Total	C	N	O	P	0	0	0
			494	233	97	141	23			

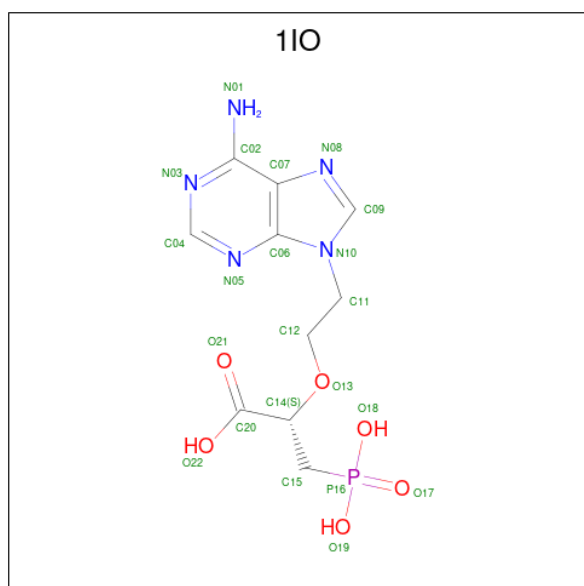
- Molecule 4 is a DNA chain called DNA (5'-D(*CP*AP*GP*TP*CP*CP*CP*TP*GP*TP*TP*CP*GP*GP*(MRG)P*CP*GP*CP*CP*(DDG))-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	20	Total	C	N	O	P	0	0	0
			403	192	72	120	19			
4	P	20	Total	C	N	O	P	0	0	0
			403	192	72	120	19			

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mg	0	0
			1	1		

- Molecule 6 is (S)-2-(2-(6-amino-9H-purin-9-yl)ethoxy)-3-phosphonopropanoic acid (CCD ID: 1IO) (formula: C₁₀H₁₄N₅O₆P) (labeled as "Ligand of Interest" by depositor).

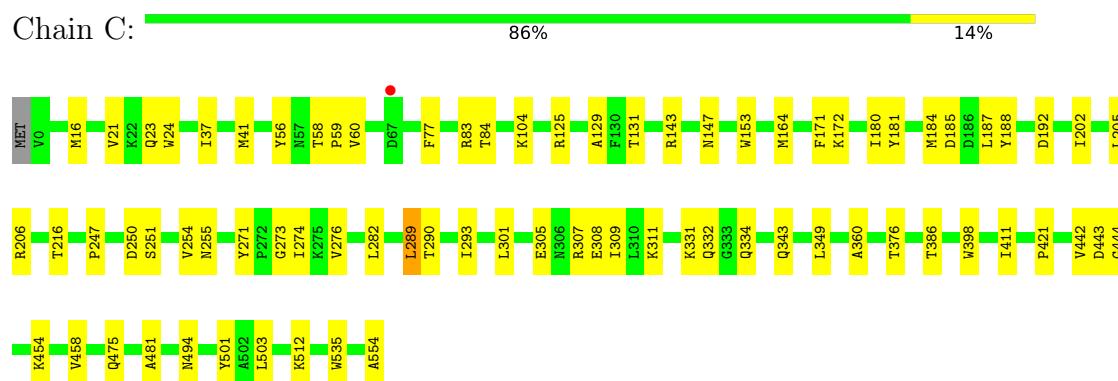


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	P	0	0
			22	10	5	6	1		

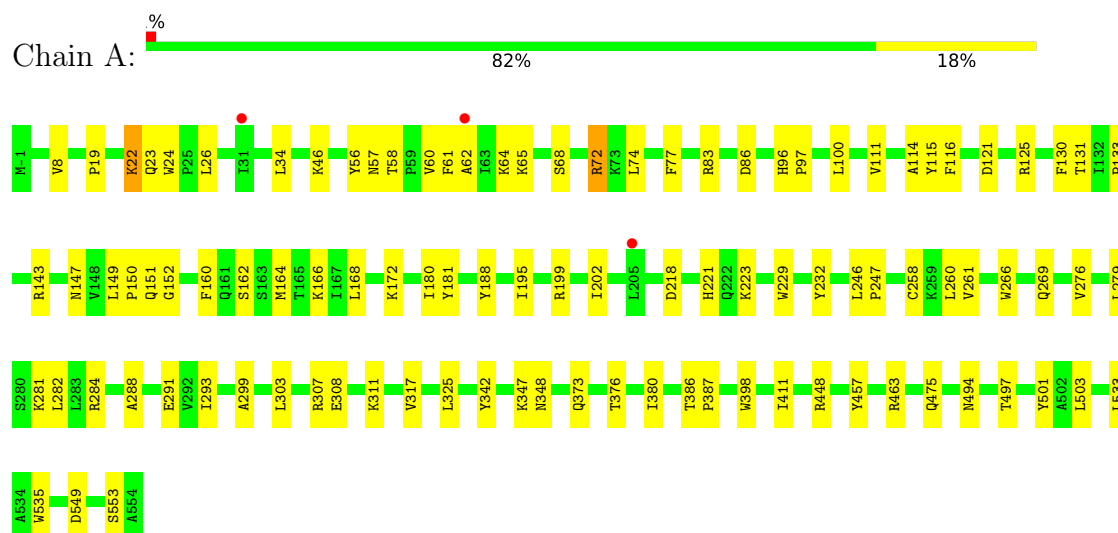
3 Residue-property plots

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.

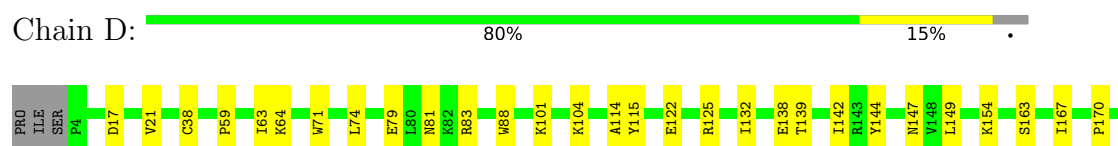
• Molecule 1: Reverse transcriptase/ribonuclease H

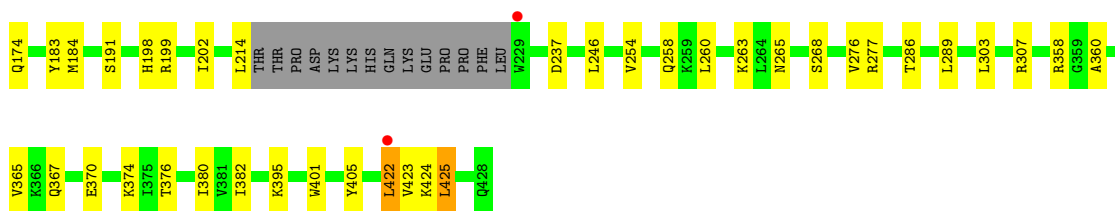


• Molecule 1: Reverse transcriptase/ribonuclease H



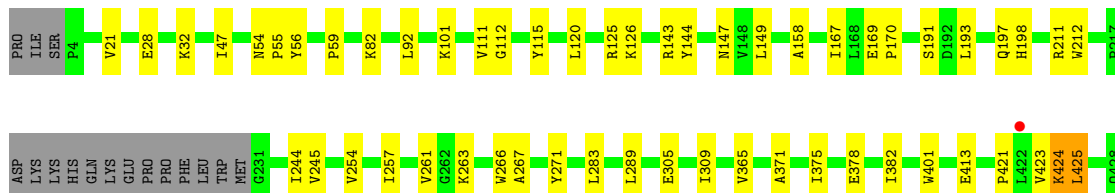
• Molecule 2: Reverse transcriptase/ribonuclease H





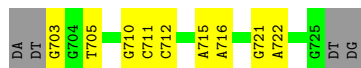
- Molecule 2: Reverse transcriptase/ribonuclease H

Chain B: 83% 12% .



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

Chain E: 52% 33% 15%



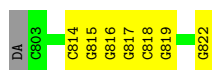
- Molecule 3: DNA (5'-D(P*GP*GP*TP*CP*GP*GP*CP*GP*CP*CP*CP*GP*AP*AP*CP*A P*GP*GP*GP*AP*CP*TP*G)-3')

Chain T: 63% 22% . 11%



- Molecule 4: DNA (5'-D(*CP*AP*GP*TP*CP*CP*CP*TP*GP*TP*TP*CP*GP*GP*(MRG) P*CP*GP*CP*CP*(DDG))-3')

Chain F: 62% 33% 5%



- Molecule 4: DNA (5'-D(*CP*AP*GP*TP*CP*CP*CP*TP*GP*TP*TP*CP*GP*GP*(MRG) P*CP*GP*CP*CP*(DDG))-3')

Chain P: 81% 14% 5%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	89.75Å 133.13Å 140.37Å 90.00° 97.59° 90.00°	Depositor
Resolution (Å)	96.19 – 3.10 96.19 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (96.19-3.10) 100.0 (96.19-3.10)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.60 (at 3.13Å)	Xtriage
Refinement program	PHENIX 1.19.1_4122	Depositor
R, R_{free}	0.210 , 0.257 0.216 , 0.258	Depositor DCC
R_{free} test set	1780 reflections (3.00%)	wwPDB-VP
Wilson B-factor (Å ²)	94.6	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 66.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	17619	wwPDB-VP
Average B, all atoms (Å ²)	115.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.89% 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: MRG, DDG, MG, IIO

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	0.31	0/4639	0.48	0/6303
1	C	0.30	0/4627	0.48	3/6289 (0.0%)
2	B	0.27	0/3488	0.49	1/4740 (0.0%)
2	D	0.26	1/3496 (0.0%)	0.39	1/4749 (0.0%)
3	E	0.37	0/536	0.59	0/826
3	T	0.35	0/555	0.65	1/856 (0.1%)
4	F	0.39	0/400	0.69	0/612
4	P	0.41	0/400	0.73	0/612
All	All	0.30	1/18141 (0.0%)	0.49	6/24987 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	395	LYS	CA-C	-6.65	1.44	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	424	LYS	N-CA-C	-13.48	92.95	111.56
1	C	289	LEU	N-CA-C	-6.05	105.75	113.01
2	D	425	LEU	N-CA-C	5.88	117.77	111.36
1	C	360	ALA	N-CA-C	-5.68	105.45	112.90
1	C	290	THR	N-CA-C	-5.53	104.97	112.26
3	T	711	DC	C2'-C3'-O3'	5.04	119.05	111.50

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	4521	0	4587	58	0
1	C	4509	0	4567	47	0
2	B	3391	0	3413	32	0
2	D	3398	0	3424	41	0
3	E	477	0	256	5	0
3	T	494	0	269	7	0
4	F	403	0	224	4	0
4	P	403	0	224	3	0
5	A	1	0	0	0	0
6	A	22	0	0	2	0
All	All	17619	0	16964	189	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 5.

All (189) 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:72:ARG:HD3	1:A:74:LEU:HD21	1.63	0.81
1:A:64:LYS:NZ	1:A:68:SER:O	2.14	0.79
1:C:84:THR:HG21	1:C:153:TRP:HE1	1.55	0.71
1:C:206:ARG:NH2	1:C:216:THR:O	2.23	0.71
3:T:704:DG:OP2	3:T:704:DG:N2	2.19	0.69
1:C:334:GLN:HB3	1:C:512:LYS:HE2	1.75	0.68
1:A:125:ARG:HE	1:A:147:ASN:HA	1.55	0.68
1:A:34:LEU:HD21	1:A:62:ALA:HB2	1.77	0.66
2:D:125:ARG:HE	2:D:147:ASN:HA	1.59	0.66
2:B:56:TYR:HE2	2:B:126:LYS:HE2	1.61	0.66
1:C:21:VAL:HG23	1:C:58:THR:HA	1.78	0.66
2:D:122:GLU:OE1	2:D:125:ARG:NH1	2.30	0.65
1:A:181:TYR:HB2	1:A:188:TYR:HB3	1.78	0.65
1:A:23:GLN:HG2	1:A:133:PRO:HD3	1.78	0.64
2:B:112:GLY:HA2	2:B:115:TYR:CD2	2.33	0.64
1:C:172:LYS:HE2	1:C:180:ILE:HB	1.79	0.64
1:C:308:GLU:HA	1:C:311:LYS:HE2	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:164:MET:HE2	1:C:187:LEU:HD11	1.80	0.62
1:C:181:TYR:HB2	1:C:188:TYR:HB3	1.82	0.61
2:D:199:ARG:HA	2:D:202:ILE:HD12	1.81	0.61
1:A:115:TYR:HD2	1:A:151:GLN:HG2	1.65	0.61
1:C:125:ARG:HD3	1:C:147:ASN:HA	1.82	0.61
1:C:56:TYR:O	1:C:143:ARG:NH2	2.34	0.61
1:A:503:LEU:HD12	1:A:533:LEU:HD13	1.83	0.60
6:A:602:1IO:C04	3:T:705:DT:H3	2.14	0.60
1:A:317:VAL:HG11	1:A:347:LYS:HB3	1.84	0.60
1:C:185:ASP:OD1	4:F:822:DDG:H3'1	2.02	0.59
2:B:112:GLY:HA2	2:B:115:TYR:HD2	1.67	0.59
2:B:254:VAL:HG13	2:B:283:LEU:HD22	1.85	0.59
1:C:454:LYS:NZ	1:C:554:ALA:HB3	2.18	0.59
2:D:81:ASN:HB3	2:D:154:LYS:HD2	1.83	0.58
1:A:56:TYR:O	1:A:143:ARG:NH2	2.36	0.58
2:D:114:ALA:HB2	2:D:214:LEU:HD13	1.86	0.58
2:B:101:LYS:HD3	2:B:382:ILE:HG23	1.86	0.57
1:C:254:VAL:HB	1:C:289:LEU:HA	1.88	0.56
1:A:281:LYS:HG2	1:A:284:ARG:NH2	2.20	0.56
1:C:131:THR:HG22	1:C:143:ARG:HG2	1.88	0.56
2:D:88:TRP:CZ2	2:D:154:LYS:HD3	2.41	0.55
1:C:58:THR:HG21	1:C:77:PHE:CE1	2.41	0.55
2:D:260:LEU:HD21	2:D:303:LEU:HD13	1.88	0.55
1:C:21:VAL:HG23	1:C:59:PRO:HD3	1.88	0.55
2:D:365:VAL:HG11	2:D:401:TRP:HB2	1.87	0.55
1:A:503:LEU:HD22	1:A:535:TRP:HB2	1.88	0.54
2:B:245:VAL:O	2:B:263:LYS:NZ	2.39	0.54
2:B:191:SER:OG	2:B:198:HIS:ND1	2.31	0.54
1:C:503:LEU:HD22	1:C:535:TRP:HB2	1.89	0.54
2:B:305:GLU:O	2:B:309:ILE:HG13	2.08	0.53
4:P:818:DC:H2'	4:P:819:DG:C8	2.42	0.53
1:A:164:MET:HE3	1:A:168:LEU:HD11	1.90	0.53
2:B:378:GLU:O	2:B:382:ILE:HG13	2.09	0.53
1:A:261:VAL:HG13	1:A:276:VAL:HG11	1.90	0.53
1:C:24:TRP:HD1	1:C:59:PRO:HB3	1.74	0.53
2:D:246:LEU:HD12	2:D:307:ARG:HG2	1.91	0.53
1:C:301:LEU:O	1:C:305:GLU:HG3	2.09	0.52
2:D:21:VAL:HB	2:D:59:PRO:HD3	1.91	0.52
2:D:358:ARG:NH2	2:D:405:TYR:O	2.31	0.52
1:C:21:VAL:CG2	1:C:59:PRO:HD3	2.40	0.52
2:B:421:PRO:HB2	2:B:423:VAL:HG22	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:ALA:HB1	1:A:160:PHE:CZ	2.44	0.52
2:D:276:VAL:HG22	2:D:277:ARG:HH21	1.75	0.51
1:A:162:SER:O	1:A:166:LYS:HG3	2.10	0.51
2:B:125:ARG:HD3	2:B:147:ASN:HA	1.92	0.51
2:D:163:SER:O	2:D:167:ILE:HG13	2.10	0.50
1:A:172:LYS:HE2	1:A:180:ILE:HB	1.93	0.50
4:P:818:DC:H2'	4:P:819:DG:H8	1.77	0.50
2:B:115:TYR:O	2:B:149:LEU:HB2	2.12	0.50
2:D:191:SER:OG	2:D:198:HIS:ND1	2.30	0.50
1:A:199:ARG:HA	1:A:202:ILE:HD12	1.93	0.50
3:E:703:DG:H21	3:E:705:DT:H72	1.77	0.50
1:A:58:THR:HG21	1:A:77:PHE:CD1	2.47	0.49
1:A:199:ARG:NH1	1:A:223:LYS:HB3	2.28	0.49
2:D:254:VAL:O	2:D:258:GLN:HG3	2.13	0.49
2:B:371:ALA:O	2:B:375:ILE:HG13	2.13	0.49
4:F:814:DC:H2''	4:F:815:DG:C8	2.48	0.49
4:F:816:DG:H2'	4:F:817:MRG:H8	1.95	0.49
1:C:343:GLN:HG3	1:C:349:LEU:HD11	1.95	0.48
1:C:458:VAL:HG12	2:D:286:THR:HG21	1.93	0.48
1:A:195:ILE:HD11	1:A:199:ARG:NH2	2.29	0.48
1:C:58:THR:HG22	1:C:129:ALA:O	2.13	0.48
2:B:266:TRP:CZ3	2:B:425:LEU:HD22	2.48	0.48
6:A:602:1IO:N03	3:T:705:DT:N3	2.61	0.48
2:B:193:LEU:HB3	2:B:197:GLN:HG3	1.94	0.48
2:B:21:VAL:HB	2:B:59:PRO:HD3	1.96	0.48
1:C:171:PHE:CD1	1:C:205:LEU:HD13	2.49	0.48
1:C:276:VAL:O	1:C:276:VAL:HG12	2.14	0.48
1:A:288:ALA:HB3	1:A:291:GLU:HB2	1.96	0.47
1:A:448:ARG:NH2	3:T:723:DC:H1'	2.29	0.47
1:A:100:LEU:HD11	1:A:229:TRP:CZ3	2.49	0.47
2:B:244:ILE:HD12	2:B:271:TYR:HE1	1.79	0.47
1:C:494:ASN:HB3	2:D:289:LEU:HD12	1.96	0.47
2:B:28:GLU:HG2	2:B:32:LYS:HE2	1.95	0.47
2:B:167:ILE:HG23	2:B:212:TRP:CD1	2.49	0.47
1:A:549:ASP:O	1:A:553:SER:OG	2.20	0.46
1:C:331:LYS:HB3	1:C:421:PRO:HG2	1.96	0.46
1:A:46:LYS:HD3	1:A:116:PHE:HB3	1.98	0.46
1:A:65:LYS:HE2	1:A:65:LYS:HA	1.96	0.46
2:B:257:ILE:O	2:B:261:VAL:HG23	2.16	0.46
2:D:170:PRO:O	2:D:174:GLN:HG3	2.15	0.46
1:A:266:TRP:O	1:A:269:GLN:HG2	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:342:TYR:HB3	1:A:348:ASN:HA	1.97	0.46
1:C:255:ASN:HB2	1:C:289:LEU:HB3	1.97	0.46
1:C:202:ILE:O	1:C:206:ARG:HG3	2.16	0.46
2:B:47:ILE:HD12	2:B:144:TYR:CD2	2.50	0.46
1:C:305:GLU:O	1:C:309:ILE:HG13	2.17	0.45
1:A:26:LEU:HD21	1:A:60:VAL:O	2.17	0.45
1:A:86:ASP:O	2:B:55:PRO:HB3	2.16	0.45
1:C:443:ASP:OD1	1:C:444:GLY:N	2.49	0.45
1:A:457:TYR:HE1	1:A:463:ARG:HG2	1.82	0.45
1:A:22:LYS:HA	1:A:22:LYS:HD2	1.48	0.45
1:A:325:LEU:HB3	1:A:387:PRO:HB3	1.98	0.45
1:A:19:PRO:HA	1:A:83:ARG:NH1	2.32	0.45
1:A:398:TRP:CZ2	1:A:411:ILE:HG13	2.52	0.45
1:C:398:TRP:CH2	1:C:411:ILE:HG13	2.51	0.44
2:B:169:GLU:HB3	2:B:170:PRO:HD3	1.99	0.44
1:C:247:PRO:O	1:C:307:ARG:NH2	2.51	0.44
1:C:475:GLN:HB3	1:C:501:TYR:CE2	2.53	0.44
1:A:303:LEU:HD23	1:A:307:ARG:HG3	1.99	0.44
1:C:37:ILE:HG22	1:C:41:MET:HE2	1.98	0.44
1:C:454:LYS:HZ1	1:C:554:ALA:HB3	1.83	0.44
2:B:92:LEU:HB2	2:B:158:ALA:HB1	1.99	0.44
1:A:376:THR:HG23	1:A:386:THR:HG22	2.00	0.44
1:A:97:PRO:HD3	1:A:232:TYR:CE2	2.52	0.44
2:B:125:ARG:NH1	2:B:147:ASN:O	2.51	0.44
1:C:376:THR:HG23	1:C:386:THR:HG23	2.00	0.43
1:C:16:MET:SD	1:C:83:ARG:HG2	2.59	0.43
1:A:475:GLN:HB3	1:A:501:TYR:CE2	2.53	0.43
3:E:715:DA:H2'	3:E:716:DA:C8	2.54	0.43
3:T:706:DC:H42	4:P:822:DDG:H1	1.65	0.43
2:D:17:ASP:O	2:D:83:ARG:HD3	2.19	0.43
2:D:101:LYS:HD3	2:D:382:ILE:HG23	2.00	0.43
2:D:423:VAL:HG22	2:D:423:VAL:O	2.19	0.43
1:A:246:LEU:HD22	1:A:260:LEU:HD12	2.01	0.43
2:D:265:ASN:O	2:D:268:SER:OG	2.27	0.43
3:E:721:DG:H2'	3:E:722:DA:C8	2.53	0.43
1:A:497:THR:O	1:A:535:TRP:HA	2.18	0.43
1:C:104:LYS:HB2	1:C:192:ASP:HA	2.01	0.43
2:D:183:TYR:CE2	2:D:184:MET:HG2	2.54	0.43
2:D:376:THR:O	2:D:380:ILE:HG13	2.19	0.43
2:B:54:ASN:HB3	2:B:143:ARG:HH21	1.84	0.43
4:F:818:DC:H2'	4:F:819:DG:H8	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:T:711:DC:H2'	3:T:712:DC:C6	2.54	0.43
2:B:365:VAL:HG11	2:B:401:TRP:HB2	2.01	0.43
2:B:120:LEU:HD23	2:B:125:ARG:HG2	2.00	0.43
1:C:250:ASP:OD1	1:C:251:SER:N	2.51	0.42
1:C:273:GLY:HA2	1:C:332:GLN:NE2	2.34	0.42
2:D:64:LYS:HE3	2:D:71:TRP:CE2	2.54	0.42
2:D:138:GLU:HG2	2:D:139:THR:HG23	2.01	0.42
2:B:82:LYS:HD3	2:B:413:GLU:OE2	2.19	0.42
3:E:710:DG:H2'	3:E:711:DC:C6	2.54	0.42
2:D:263:LYS:HE3	2:D:425:LEU:HA	2.00	0.42
1:A:279:LEU:HD23	1:A:299:ALA:HB1	2.00	0.42
1:A:494:ASN:HB3	2:B:289:LEU:HD12	2.01	0.42
3:E:711:DC:H2'	3:E:712:DC:C6	2.54	0.42
1:C:23:GLN:HE22	1:C:60:VAL:H	1.66	0.42
2:D:142:ILE:HG22	2:D:144:TYR:CE1	2.54	0.42
2:D:423:VAL:O	2:D:424:LYS:C	2.63	0.42
1:C:376:THR:HG21	2:D:401:TRP:CH2	2.55	0.42
2:D:38:CYS:SG	2:D:132:ILE:HD11	2.59	0.42
1:A:96:HIS:NE2	1:A:269:GLN:OE1	2.49	0.42
1:A:24:TRP:O	1:A:26:LEU:HD12	2.20	0.42
1:C:184:MET:HE2	1:C:184:MET:HB2	1.92	0.42
2:D:104:LYS:HD3	2:D:237:ASP:OD2	2.20	0.42
2:D:360:ALA:HA	2:D:367:GLN:CD	2.45	0.42
3:T:706:DC:H2'	3:T:707:DG:C8	2.55	0.42
2:D:115:TYR:O	2:D:149:LEU:HB2	2.20	0.41
1:A:247:PRO:O	1:A:307:ARG:NH2	2.53	0.41
1:A:111:VAL:HG12	1:A:114:ALA:HB2	2.02	0.41
1:A:308:GLU:HA	1:A:311:LYS:HE2	2.02	0.41
2:D:125:ARG:NE	2:D:147:ASN:HA	2.32	0.41
1:A:57:ASN:HD21	1:A:131:THR:HG22	1.84	0.41
2:B:267:ALA:HB2	2:B:425:LEU:HD11	2.01	0.41
1:C:282:LEU:HB3	1:C:293:ILE:HD13	2.03	0.41
1:A:60:VAL:HG21	1:A:130:PHE:CD2	2.55	0.41
1:A:218:ASP:HA	1:A:221:HIS:HB2	2.03	0.41
1:A:373:GLN:HE21	1:A:373:GLN:HB2	1.65	0.41
2:D:422:LEU:H	2:D:422:LEU:HG	1.55	0.41
2:D:63:ILE:HD13	2:D:74:LEU:HD22	2.02	0.41
2:D:198:HIS:CD2	2:D:202:ILE:HD11	2.56	0.41
1:C:271:TYR:O	1:C:274:ILE:HD13	2.22	0.40
2:D:183:TYR:CD2	2:D:184:MET:HG2	2.55	0.40
1:C:442:VAL:HB	1:C:481:ALA:HB1	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:370:GLU:O	2:D:374:LYS:HG3	2.21	0.40
1:A:8:VAL:O	1:A:121:ASP:HB2	2.21	0.40
1:A:282:LEU:HB3	1:A:293:ILE:HD13	2.02	0.40
1:A:376:THR:O	1:A:380:ILE:HG13	2.21	0.40
2:D:79:GLU:HG3	2:D:83:ARG:HE	1.86	0.40
1:A:61:PHE:O	1:A:74:LEU:HB2	2.21	0.40
2:B:211:ARG:HG3	2:B:212:TRP:CD1	2.56	0.40
1:A:149:LEU:HA	1:A:150:PRO:HD3	1.90	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	554/556 (100%)	535 (97%)	18 (3%)	1 (0%)	43	73
1	C	553/556 (100%)	537 (97%)	16 (3%)	0	100	100
2	B	408/428 (95%)	391 (96%)	17 (4%)	0	100	100
2	D	407/428 (95%)	390 (96%)	17 (4%)	0	100	100
All	All	1922/1968 (98%)	1853 (96%)	68 (4%)	1 (0%)	48	78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	152	GLY

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	497/497 (100%)	494 (99%)	3 (1%)	78	83
1	C	495/497 (100%)	495 (100%)	0	100	100
2	B	372/390 (95%)	369 (99%)	3 (1%)	73	81
2	D	372/390 (95%)	371 (100%)	1 (0%)	86	87
All	All	1736/1774 (98%)	1729 (100%)	7 (0%)	84	86

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

Mol	Chain	Res	Type
2	D	422	LEU
1	A	22	LYS
1	A	72	ARG
1	A	258	CYS
2	B	111	VAL
2	B	424	LYS
2	B	425	LEU

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

Mol	Chain	Res	Type
1	C	222	GLN
1	C	255	ASN
2	D	269	GLN
2	D	361	HIS
2	D	428	GLN
1	A	147	ASN
1	A	255	ASN
1	A	373	GLN
1	A	407	GLN
1	A	524	GLN
1	A	545	ASN
2	B	175	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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	MRG	F	817	3,4	21,24,29	0.32	0	30,35,42	0.26	0
4	DDG	F	822	3,4	20,23,24	0.37	0	27,33,36	0.44	0
4	MRG	P	817	3,4	21,24,29	0.32	0	30,35,42	0.29	0
4	DDG	P	822	3,4	20,23,24	0.47	0	27,33,36	0.41	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	MRG	F	817	3,4	-	0/7/21/27	0/3/3/3
4	DDG	F	822	3,4	-	0/7/18/19	0/3/3/3
4	MRG	P	817	3,4	-	3/7/21/27	0/3/3/3
4	DDG	P	822	3,4	-	0/7/18/19	0/3/3/3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	P	817	MRG	O4'-C4'-C5'-O5'
4	P	817	MRG	C3'-C4'-C5'-O5'
4	P	817	MRG	C4'-C5'-O5'-P

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	F	817	MRG	1	0
4	F	822	DDG	1	0
4	P	822	DDG	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
6	1IO	A	602	5	22,23,23	2.06	8 (36%)	29,33,33	1.72	8 (27%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	1IO	A	602	5	-	2/14/15/15	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	602	1IO	O13-C14	4.58	1.50	1.43
6	A	602	1IO	C14-C20	3.70	1.56	1.52
6	A	602	1IO	C09-N10	3.29	1.42	1.36
6	A	602	1IO	C07-C06	3.07	1.44	1.39
6	A	602	1IO	C04-N03	2.42	1.38	1.33
6	A	602	1IO	C11-C12	2.31	1.57	1.50
6	A	602	1IO	C09-N08	2.16	1.35	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	602	1IO	C06-N05	2.07	1.38	1.34

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	602	1IO	O17-P16-C15	-4.92	102.35	111.56
6	A	602	1IO	O13-C14-C20	-3.60	106.66	112.17
6	A	602	1IO	O19-P16-O18	2.81	115.96	107.96
6	A	602	1IO	O22-C20-C14	2.50	119.22	112.71
6	A	602	1IO	O21-C20-C14	-2.50	117.44	122.85
6	A	602	1IO	C15-C14-C20	2.40	115.99	110.43
6	A	602	1IO	N05-C06-N10	2.30	129.72	126.87
6	A	602	1IO	C07-C06-N05	-2.24	123.63	126.72

There are no chirality outliers.

All (2) torsion outliers are listed below:

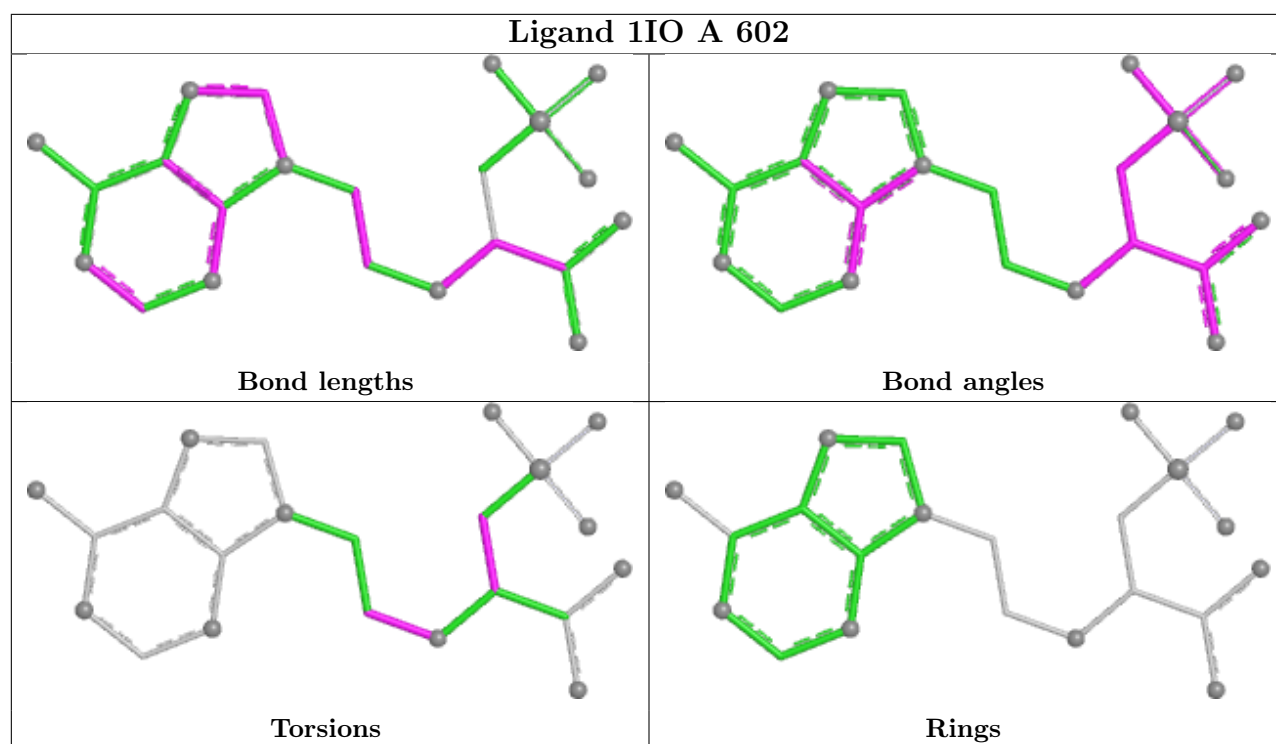
Mol	Chain	Res	Type	Atoms
6	A	602	1IO	C20-C14-C15-P16
6	A	602	1IO	C11-C12-O13-C14

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	602	1IO	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 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	556/556 (100%)	-0.24	3 (0%) 87 73	45, 116, 202, 235	0
1	C	555/556 (99%)	-0.23	1 (0%) 91 83	44, 118, 211, 260	0
2	B	412/428 (96%)	-0.31	1 (0%) 91 83	54, 102, 157, 187	0
2	D	411/428 (96%)	-0.41	2 (0%) 87 73	41, 88, 142, 171	0
3	E	23/27 (85%)	-0.22	0 100 100	109, 156, 212, 275	0
3	T	24/27 (88%)	-0.36	0 100 100	99, 141, 243, 266	0
4	F	18/21 (85%)	-0.42	0 100 100	105, 142, 194, 196	0
4	P	18/21 (85%)	-0.70	0 100 100	86, 129, 196, 197	0
All	All	2017/2064 (97%)	-0.29	7 (0%) 90 80	41, 107, 190, 275	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	67	ASP	3.3
2	B	422	LEU	2.9
1	A	205	LEU	2.7
1	A	31	ILE	2.3
2	D	229	TRP	2.3
1	A	62	ALA	2.2
2	D	422	LEU	2.0

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MRG	P	817	22/27	0.70	0.09	122,148,160,183	0
4	MRG	F	817	22/27	0.80	0.07	145,160,177,190	0
4	DDG	F	822	21/22	0.84	0.09	106,115,138,147	0
4	DDG	P	822	21/22	0.90	0.09	101,128,138,143	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

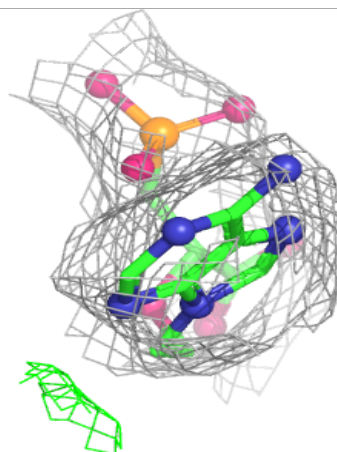
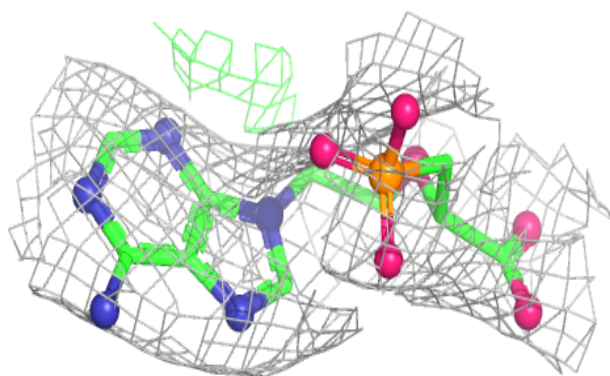
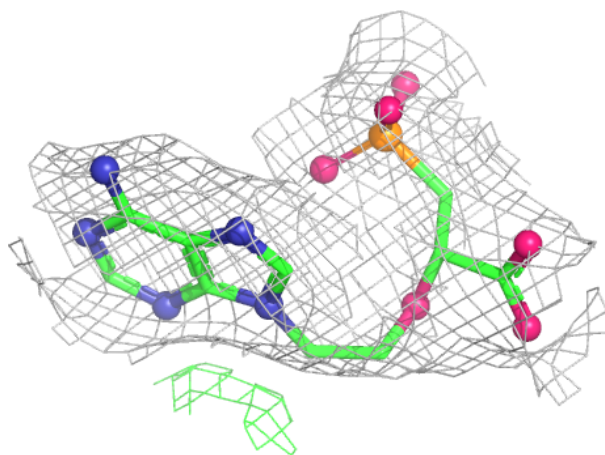
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
6	1IO	A	602	22/22	0.53	0.10	163,170,199,209	0
5	MG	A	601	1/1	0.91	0.09	156,156,156,156	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 1IO A 602:

$2mF_o-DF_c$ (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.