



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 05:44 PM UTC

PDB ID : 7AIF / pdb_00007aif
Title : HIV-1 REVERSE TRANSCRIPTASE COMPLEX WITH DNA AND L-GLUTAMATE TENOFOVIR WITH BOUND MANGANESE
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Deposited on : 2020-09-27
Resolution : 2.75 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

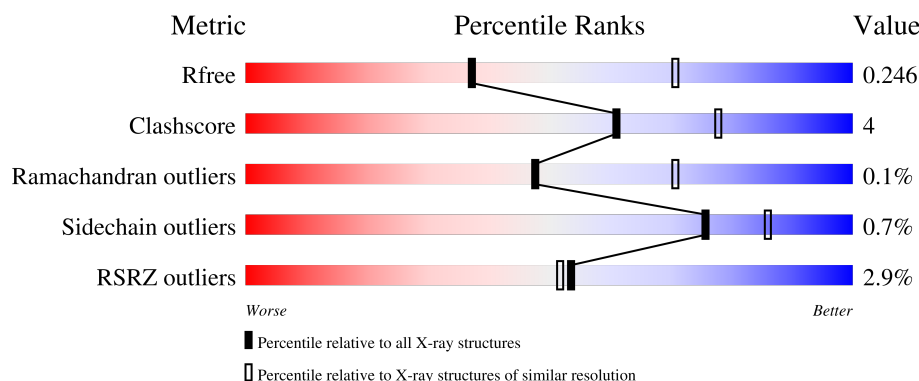
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	 3% 88% 12%
1	C	556	 3% 88% 12%
2	B	428	 2% 87% 9% .
2	D	428	 4% 82% 14% .

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Mol	Chain	Length	Quality of chain
3	E	27	 52% 37% 11%
3	T	27	 63% 22% 15%
4	F	21	 48% 48% 5%
4	P	21	 76% 19% 5%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 17898 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 Gag-Pol polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	555	Total	C	N	O	S	0	0	0
			4513	2922	751	832	8			
1	C	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
A	-1	MET	-	initiating methionine	UNP P03366
A	0	VAL	-	expression tag	UNP P03366
A	258	CYS	GLN	engineered mutation	UNP P03366
A	280	SER	CYS	engineered mutation	UNP P03366
A	498	ASN	ASP	engineered mutation	UNP P03366
C	-1	MET	-	initiating methionine	UNP P03366
C	0	VAL	-	expression tag	UNP P03366
C	258	CYS	GLN	engineered mutation	UNP P03366
C	280	SER	CYS	engineered mutation	UNP P03366
C	498	ASN	ASP	engineered mutation	UNP P03366

- Molecule 2 is a protein called Gag-Pol polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	411	Total	C	N	O	S	0	0	0
			3401	2215	563	616	7			
2	D	412	Total	C	N	O	S	0	0	0
			3400	2212	563	619	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	280	SER	CYS	engineered mutation	UNP P03366
D	280	SER	CYS	engineered mutation	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	T	23	Total	C	N	O	P	0	0	0
			477	223	95	136	23			
3	E	24	Total	C	N	O	P	0	0	0
			494	233	97	141	23			

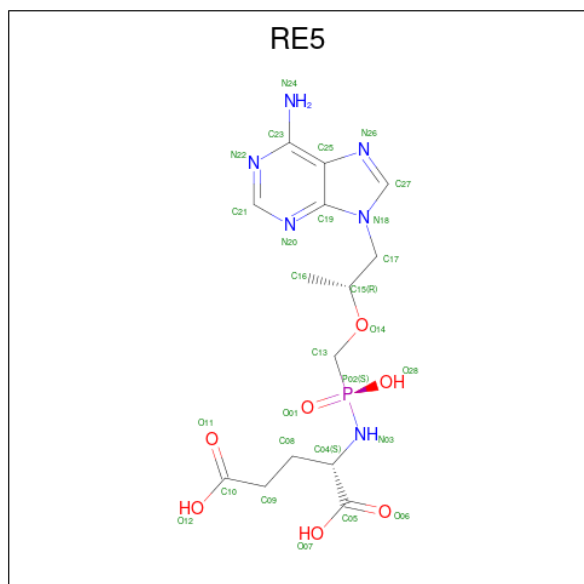
- Molecule 4 is a DNA chain called DNA (5'-D(*CP*AP*GP*TP*CP*CP*CP*TP*GP*TP*TP*CP*GP*GP*(MRG)P*CP*GP*CP*CP*(DDG))-3').

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

- Molecule 5 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Mn	0	0
			2	2		
5	C	2	Total	Mn	0	0
			2	2		

- Molecule 6 is L-Glutamate Tenofovir (CCD ID: RE5) (formula: C₁₄H₂₁N₆O₇P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	P	0	0
			28	14	6	7	1		
6	C	1	Total	C	N	O	P	0	0
			28	14	6	7	1		

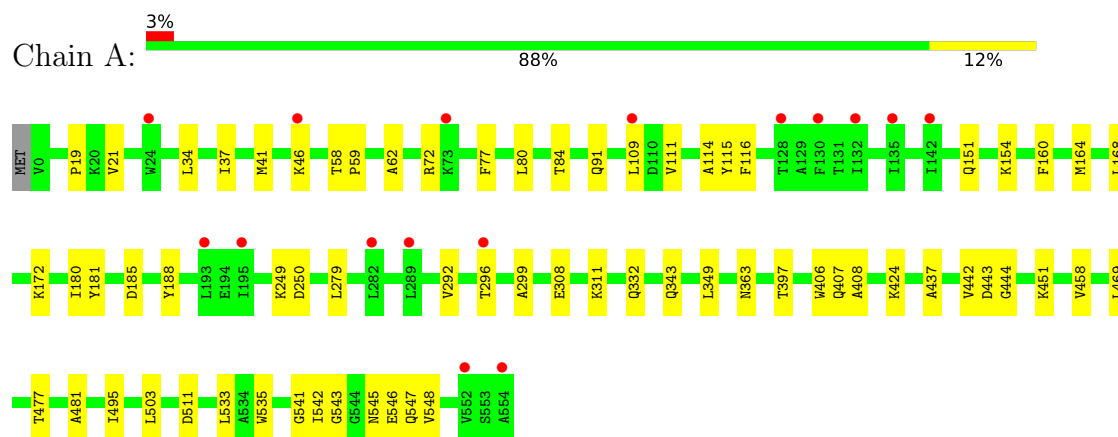
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	70	Total	O	0	0
			70	70		
7	B	64	Total	O	0	0
			64	64		
7	C	44	Total	O	0	0
			44	44		
7	D	30	Total	O	0	0
			30	30		
7	T	5	Total	O	0	0
			5	5		
7	P	6	Total	O	0	0
			6	6		
7	E	3	Total	O	0	0
			3	3		
7	F	4	Total	O	0	0
			4	4		

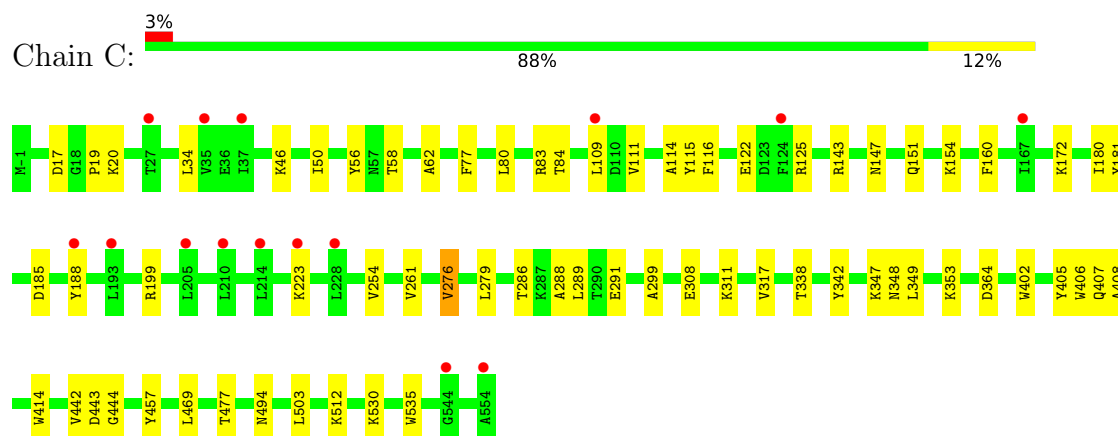
3 Residue-property plots [i](#)

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.

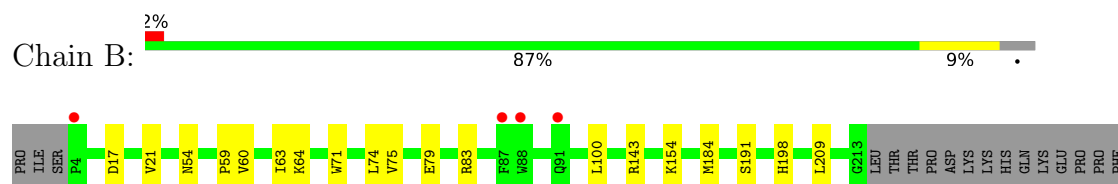
• Molecule 1: Gag-Pol polyprotein



• Molecule 1: Gag-Pol polyprotein

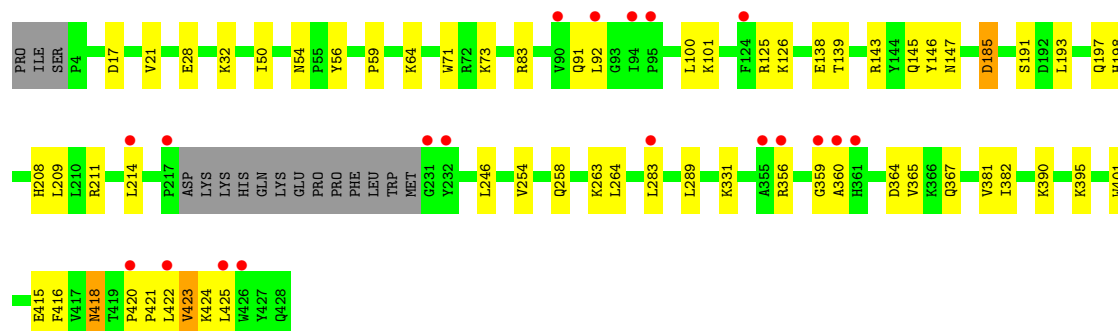
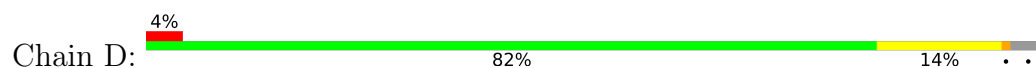


• Molecule 2: Gag-Pol polyprotein





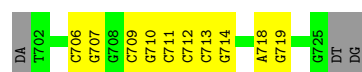
• Molecule 2: Gag-Pol polyprotein



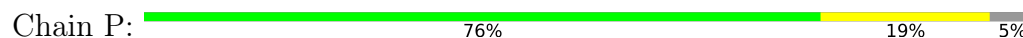
• Molecule 3: DNA (5'-D(P*GP*GP*TP*CP*GP*GP*CP*GP*CP*CP*CP*GP*AP*AP*CP*A P*GP*GP*GP*AP*CP*TP*G)-3')



• 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')



• 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')



• 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')



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	89.65Å 132.70Å 138.77Å 90.00° 98.65° 90.00°	Depositor
Resolution (Å)	80.14 – 2.75 80.14 – 2.75	Depositor EDS
% Data completeness (in resolution range)	98.3 (80.14-2.75) 98.3 (80.14-2.75)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.75 (at 2.73Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.218 , 0.249 0.218 , 0.246	Depositor DCC
R_{free} test set	2570 reflections (3.08%)	wwPDB-VP
Wilson B-factor (Å ²)	63.4	Xtriage
Anisotropy	0.163	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 49.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	17898	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.07% 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: MN, RE5, DDG, MRG

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	0.07	0/4631	0.22	0/6293
1	C	0.07	0/4639	0.22	0/6303
2	B	0.07	0/3499	0.23	0/4752
2	D	0.07	0/3497	0.22	0/4751
3	E	0.16	0/555	0.30	0/856
3	T	0.15	0/536	0.27	0/826
4	F	0.22	0/400	0.40	0/612
4	P	0.20	0/400	0.37	0/612
All	All	0.09	0/18157	0.24	0/25005

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	4513	0	4578	38	0
1	C	4521	0	4587	40	0
2	B	3401	0	3431	25	0
2	D	3400	0	3433	33	0
3	E	494	0	269	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	T	477	0	256	4	0
4	F	403	0	224	8	0
4	P	403	0	224	3	0
5	A	2	0	0	0	0
5	C	2	0	0	0	0
6	A	28	0	0	1	0
6	C	28	0	0	0	0
7	A	70	0	0	2	0
7	B	64	0	0	0	0
7	C	44	0	0	2	0
7	D	30	0	0	1	0
7	E	3	0	0	0	0
7	F	4	0	0	1	0
7	P	6	0	0	0	0
7	T	5	0	0	0	0
All	All	17898	0	17002	145	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 4.

All (145) 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:308:GLU:HA	1:A:311:LYS:HZ3	1.57	0.67
1:C:172:LYS:HE2	1:C:180:ILE:HB	1.80	0.64
2:B:390:LYS:NZ	2:B:415:GLU:OE2	2.29	0.63
1:C:125:ARG:HE	1:C:147:ASN:HA	1.64	0.63
1:A:181:TYR:HB2	1:A:188:TYR:HB3	1.79	0.62
2:D:209:LEU:HB3	2:D:214:LEU:HB2	1.82	0.62
1:A:343:GLN:HG3	1:A:349:LEU:HD11	1.82	0.61
1:C:181:TYR:HB2	1:C:188:TYR:HB3	1.83	0.61
2:D:390:LYS:NZ	2:D:415:GLU:OE2	2.32	0.59
1:A:503:LEU:HD22	1:A:535:TRP:HB2	1.86	0.58
1:C:503:LEU:HD22	1:C:535:TRP:HB2	1.86	0.57
2:B:209:LEU:HB3	2:B:228:LEU:HD22	1.86	0.57
1:A:172:LYS:HE2	1:A:180:ILE:HB	1.86	0.56
1:C:122:GLU:OE1	1:C:125:ARG:NH1	2.39	0.56
1:C:50:ILE:HD11	1:C:143:ARG:HB2	1.87	0.55
2:D:56:TYR:HE2	2:D:126:LYS:HE2	1.72	0.55
1:C:342:TYR:HB3	1:C:348:ASN:HA	1.89	0.55
2:D:246:LEU:HD11	2:D:264:LEU:HD21	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:21:VAL:HB	2:B:59:PRO:HD3	1.89	0.54
1:A:443:ASP:OD1	1:A:444:GLY:N	2.39	0.54
1:C:84:THR:HB	1:C:154:LYS:HD3	1.89	0.54
2:D:73:LYS:NZ	2:D:146:TYR:OH	2.37	0.54
2:D:28:GLU:HG2	2:D:32:LYS:HE2	1.89	0.53
2:D:359:GLY:O	2:D:367:GLN:NE2	2.41	0.53
3:E:718:DA:H1'	3:E:719:DG:C8	2.44	0.53
2:D:263:LYS:HG3	2:D:425:LEU:HG	1.91	0.52
2:D:125:ARG:NH1	2:D:147:ASN:O	2.43	0.52
1:A:21:VAL:HB	1:A:59:PRO:HD3	1.90	0.52
1:A:46:LYS:HD3	1:A:116:PHE:HB3	1.91	0.52
2:D:356:ARG:HD2	2:D:360:ALA:HB1	1.92	0.52
2:D:421:PRO:HB2	2:D:423:VAL:HG23	1.92	0.51
2:B:79:GLU:HG3	2:B:83:ARG:HE	1.74	0.51
1:C:80:LEU:O	1:C:84:THR:OG1	2.28	0.51
1:C:402:TRP:O	2:D:331:LYS:NZ	2.35	0.51
2:D:254:VAL:HG13	2:D:283:LEU:HD22	1.93	0.51
1:A:437:ALA:O	7:A:701:HOH:O	2.19	0.50
2:B:422:LEU:HD12	2:B:423:VAL:H	1.77	0.50
1:C:469:LEU:HD12	1:C:477:THR:HG22	1.94	0.50
2:B:63:ILE:HD13	2:B:74:LEU:HD22	1.94	0.50
2:B:229:TRP:CG	2:B:230:MET:H	2.30	0.50
1:C:414:TRP:NE1	7:C:701:HOH:O	2.39	0.50
1:A:84:THR:HB	1:A:154:LYS:HE2	1.94	0.49
1:A:469:LEU:HD12	1:A:477:THR:HG22	1.94	0.49
1:A:72:ARG:NH2	6:A:603:RE5:O11	2.46	0.49
1:A:397:THR:HG21	1:A:424:LYS:HA	1.94	0.49
2:D:258:GLN:NE2	7:D:504:HOH:O	2.43	0.49
2:D:21:VAL:HB	2:D:59:PRO:HD3	1.95	0.48
2:D:100:LEU:HG	2:D:381:VAL:HG13	1.96	0.48
1:C:114:ALA:HB1	1:C:160:PHE:CZ	2.49	0.48
1:A:111:VAL:HB	1:A:185:ASP:HB2	1.96	0.48
2:B:282:LEU:HD21	2:B:295:LEU:HA	1.96	0.48
2:D:54:ASN:HB3	2:D:143:ARG:HH21	1.79	0.48
1:C:442:VAL:HG12	1:C:457:TYR:HB3	1.96	0.47
2:B:54:ASN:HB3	2:B:143:ARG:HH21	1.80	0.47
1:A:19:PRO:HD3	1:A:80:LEU:HD13	1.95	0.47
1:A:363:ASN:HA	1:A:511:ASP:OD1	2.14	0.47
1:C:46:LYS:HD3	1:C:116:PHE:HB3	1.96	0.47
4:F:818:DC:H2'	4:F:819:DG:C8	2.50	0.47
1:A:543:GLY:N	2:B:283:LEU:O	2.41	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:T:713:DC:H2'	3:T:714:DG:C8	2.50	0.47
2:D:208:HIS:CD2	2:D:211:ARG:HH22	2.33	0.47
1:A:58:THR:HG21	1:A:77:PHE:CD1	2.50	0.47
1:C:19:PRO:HG3	1:C:80:LEU:HB2	1.98	0.46
2:D:101:LYS:HD3	2:D:382:ILE:HG23	1.97	0.46
4:F:820:DC:N4	7:F:902:HOH:O	2.48	0.46
2:B:60:VAL:HG22	2:B:75:VAL:HG22	1.97	0.46
2:B:365:VAL:HG11	2:B:401:TRP:HB2	1.98	0.46
1:C:364:ASP:OD1	1:C:512:LYS:NZ	2.43	0.46
3:T:711:DC:H2'	3:T:712:DC:C6	2.51	0.46
2:D:91:GLN:HG2	2:D:92:LEU:HG	1.98	0.45
2:D:191:SER:OG	2:D:198:HIS:ND1	2.39	0.45
1:C:19:PRO:HD3	1:C:80:LEU:HD13	1.97	0.45
1:C:405:TYR:CE2	1:C:407:GLN:HB2	2.51	0.45
3:E:711:DC:H2'	3:E:712:DC:C6	2.52	0.45
1:C:261:VAL:HG13	1:C:276:VAL:HG11	1.99	0.45
1:A:442:VAL:HB	1:A:481:ALA:HB1	1.98	0.44
2:B:64:LYS:HE3	2:B:71:TRP:CE2	2.52	0.44
2:B:191:SER:OG	2:B:198:HIS:ND1	2.44	0.44
2:B:154:LYS:HG2	2:B:184:MET:SD	2.56	0.44
1:C:199:ARG:HH12	1:C:223:LYS:HB3	1.82	0.44
1:A:80:LEU:O	1:A:84:THR:OG1	2.24	0.44
1:C:288:ALA:HB3	1:C:291:GLU:HB2	2.00	0.44
3:E:713:DC:H2'	3:E:714:DG:C8	2.53	0.44
2:D:365:VAL:HG11	2:D:401:TRP:HB2	2.00	0.44
4:P:818:DC:H2'	4:P:819:DG:C8	2.52	0.44
1:A:406:TRP:CD1	1:A:407:GLN:HG2	2.53	0.44
1:C:115:TYR:HD2	1:C:151:GLN:HG2	1.82	0.44
4:F:814:DC:H2''	4:F:815:DG:H8	1.83	0.43
4:F:804:DA:H4'	4:F:805:DG:OP1	2.18	0.43
1:A:279:LEU:HD23	1:A:299:ALA:HB1	2.00	0.43
4:F:805:DG:H2''	4:F:806:DT:O5'	2.18	0.43
1:A:451:LYS:NZ	4:P:808:DC:OP1	2.51	0.43
1:C:276:VAL:HG23	1:C:353:LYS:HE2	2.01	0.43
1:A:408:ALA:O	2:B:393:ILE:HG13	2.18	0.43
2:D:138:GLU:HG2	2:D:139:THR:HG23	2.01	0.43
1:A:495:ILE:HB	1:A:533:LEU:HD23	2.00	0.43
1:C:17:ASP:O	1:C:83:ARG:HD3	2.19	0.43
1:A:91:GLN:O	3:T:708:DG:H4'	2.19	0.43
1:A:114:ALA:HB1	1:A:160:PHE:CZ	2.54	0.43
1:A:541:GLY:HA2	1:A:546:GLU:HB2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:58:THR:HG21	1:C:77:PHE:CD1	2.54	0.43
1:C:289:LEU:HD11	4:F:817:MRG:H4'	2.01	0.43
1:C:34:LEU:HD21	1:C:62:ALA:HB2	2.00	0.42
2:B:358:ARG:NH2	2:B:405:TYR:O	2.37	0.42
2:B:260:LEU:HD21	2:B:303:LEU:HD13	2.02	0.42
2:D:17:ASP:O	2:D:83:ARG:HD3	2.19	0.42
1:A:164:MET:HE3	1:A:168:LEU:HD11	2.01	0.42
2:B:100:LEU:HG	2:B:381:VAL:HG13	2.02	0.42
2:B:229:TRP:CG	2:B:230:MET:N	2.88	0.42
4:F:817:MRG:H2'	4:F:818:DC:C6	2.54	0.42
1:A:542:ILE:HG23	2:B:283:LEU:HD13	2.02	0.42
3:E:709:DC:H2'	3:E:710:DG:C8	2.54	0.42
4:F:818:DC:H2'	4:F:819:DG:H8	1.84	0.42
2:B:17:ASP:O	2:B:83:ARG:HD3	2.20	0.42
1:C:254:VAL:HG21	1:C:286:THR:HG21	2.01	0.42
1:C:317:VAL:HG23	1:C:349:LEU:HD23	2.02	0.42
2:B:422:LEU:H	2:B:422:LEU:HG	1.57	0.41
1:C:443:ASP:OD1	1:C:444:GLY:N	2.52	0.41
1:A:250:ASP:N	1:A:250:ASP:OD1	2.54	0.41
2:B:332:GLN:HB2	2:B:336:GLN:HB2	2.02	0.41
1:C:494:ASN:HB3	2:D:289:LEU:HD12	2.01	0.41
1:C:408:ALA:HB1	2:D:364:ASP:HB3	2.02	0.41
2:D:185:ASP:N	2:D:185:ASP:OD1	2.53	0.41
1:A:115:TYR:HD2	1:A:151:GLN:HG2	1.85	0.41
3:E:706:DC:H2'	3:E:707:DG:C8	2.55	0.41
1:C:308:GLU:HA	1:C:311:LYS:HE2	2.03	0.41
2:D:193:LEU:HB3	2:D:197:GLN:HG3	2.01	0.41
1:A:545:ASN:HA	1:A:548:VAL:HG12	2.02	0.41
2:D:64:LYS:HE3	2:D:71:TRP:CE2	2.55	0.41
1:A:34:LEU:HD21	1:A:62:ALA:HB2	2.02	0.41
1:A:37:ILE:HG22	1:A:41:MET:HE2	2.03	0.41
1:C:111:VAL:HB	1:C:185:ASP:HB2	2.02	0.41
1:C:406:TRP:CZ2	2:D:420:PRO:HG3	2.56	0.41
1:C:279:LEU:HD23	1:C:299:ALA:HB1	2.03	0.41
1:C:317:VAL:HG11	1:C:347:LYS:HB3	2.03	0.40
1:A:547:GLN:OE1	1:A:547:GLN:N	2.54	0.40
2:D:50:ILE:HD13	2:D:145:GLN:HB3	2.02	0.40
2:D:395:LYS:HG3	2:D:416:PHE:CE2	2.56	0.40
2:D:418:ASN:OD1	2:D:418:ASN:N	2.55	0.40
1:A:249:LYS:NZ	7:A:714:HOH:O	2.49	0.40
1:C:530:LYS:NZ	7:C:703:HOH:O	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:T:707:DG:H2'	3:T:708:DG:C8	2.57	0.40
1:A:458:VAL:HG12	2:B:286:THR:HG21	2.03	0.40
1:C:20:LYS:HG2	1:C:56:TYR:HA	2.02	0.40
4:P:818:DC:H2'	4:P:819:DG:H8	1.86	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	553/556 (100%)	539 (98%)	13 (2%)	1 (0%)	43	64
1	C	554/556 (100%)	537 (97%)	17 (3%)	0	100	100
2	B	407/428 (95%)	388 (95%)	19 (5%)	0	100	100
2	D	408/428 (95%)	388 (95%)	20 (5%)	0	100	100
All	All	1922/1968 (98%)	1852 (96%)	69 (4%)	1 (0%)	48	71

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	332	GLN

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	496/497 (100%)	493 (99%)	3 (1%)	78	88
1	C	497/497 (100%)	494 (99%)	3 (1%)	78	88
2	B	373/390 (96%)	372 (100%)	1 (0%)	86	93
2	D	374/390 (96%)	369 (99%)	5 (1%)	61	76
All	All	1740/1774 (98%)	1728 (99%)	12 (1%)	76	86

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	109	LEU
1	A	292	VAL
1	A	296	THR
2	B	422	LEU
1	C	109	LEU
1	C	276	VAL
1	C	338	THR
2	D	185	ASP
2	D	418	ASN
2	D	422	LEU
2	D	423	VAL
2	D	424	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	332	GLN
1	A	348	ASN
1	A	373	GLN
1	A	407	GLN
1	A	498	ASN
1	A	545	ASN
2	B	361	HIS
2	B	428	GLN
1	C	407	GLN
1	C	487	GLN
1	C	520	GLN
1	C	545	ASN
2	D	182	GLN
2	D	278	GLN
2	D	394	GLN

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	DDG	F	822	4,3	20,23,24	0.81	0	27,33,36	2.10	9 (33%)
4	DDG	P	822	4,3	20,23,24	0.82	0	27,33,36	2.10	9 (33%)
4	MRG	P	817	4,3	21,24,29	0.40	0	30,35,42	0.35	0
4	MRG	F	817	4,3	21,24,29	0.39	0	30,35,42	0.36	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	DDG	F	822	4,3	-	0/7/18/19	0/3/3/3
4	DDG	P	822	4,3	-	1/7/18/19	0/3/3/3
4	MRG	P	817	4,3	-	0/7/21/27	0/3/3/3
4	MRG	F	817	4,3	-	0/7/21/27	0/3/3/3

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	P	822	DDG	C5-C4-N3	-5.42	119.77	128.39
4	F	822	DDG	C5-C4-N3	-5.39	119.82	128.39
4	P	822	DDG	C2-N3-C4	5.10	121.08	112.30
4	F	822	DDG	C2-N3-C4	5.10	121.08	112.30
4	P	822	DDG	N9-C4-N3	4.08	134.11	125.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	822	DDG	N9-C4-N3	4.05	134.04	125.95
4	P	822	DDG	C8-N7-C5	2.84	109.32	104.26
4	F	822	DDG	C8-N7-C5	2.81	109.26	104.26
4	F	822	DDG	C2-N1-C6	-2.64	120.32	125.11
4	P	822	DDG	C2-N1-C6	-2.63	120.34	125.11
4	F	822	DDG	O6-C6-C5	-2.21	120.71	126.53
4	F	822	DDG	C6-C5-N7	2.19	134.28	130.29
4	P	822	DDG	O6-C6-C5	-2.18	120.78	126.53
4	P	822	DDG	C6-C5-N7	2.18	134.25	130.29
4	P	822	DDG	C4-C5-N7	-2.17	107.23	110.67
4	F	822	DDG	C4-C5-N7	-2.16	107.25	110.67
4	F	822	DDG	C5-C6-N1	2.02	118.39	113.25
4	P	822	DDG	C5-C6-N1	2.01	118.37	113.25

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	P	822	DDG	O4'-C4'-C5'-O5'

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	F	817	MRG	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	RE5	A	603	5	26,29,29	1.71	6 (23%)	32,41,41	1.06	1 (3%)
6	RE5	C	603	5	26,29,29	1.67	6 (23%)	32,41,41	1.06	1 (3%)

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	RE5	A	603	5	-	10/21/24/24	0/2/2/2
6	RE5	C	603	5	-	10/21/24/24	0/2/2/2

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	603	RE5	C27-N18	3.26	1.42	1.36
6	C	603	RE5	C27-N18	3.24	1.42	1.36
6	C	603	RE5	C21-N22	2.81	1.38	1.33
6	A	603	RE5	C21-N22	2.79	1.38	1.33
6	A	603	RE5	P02-O28	-2.54	1.50	1.56
6	C	603	RE5	P02-O28	-2.53	1.50	1.56
6	C	603	RE5	C27-N26	2.44	1.36	1.31
6	A	603	RE5	C27-N26	2.41	1.36	1.31
6	C	603	RE5	C25-C19	2.35	1.43	1.39
6	A	603	RE5	C25-C19	2.29	1.43	1.39
6	A	603	RE5	O14-C15	-2.18	1.40	1.44
6	C	603	RE5	O14-C15	-2.10	1.40	1.44

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	603	RE5	O28-P02-O01	2.85	119.77	112.36
6	A	603	RE5	O28-P02-O01	2.74	119.48	112.36

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	603	RE5	N03-C04-C05-O06
6	A	603	RE5	N03-C04-C05-O07

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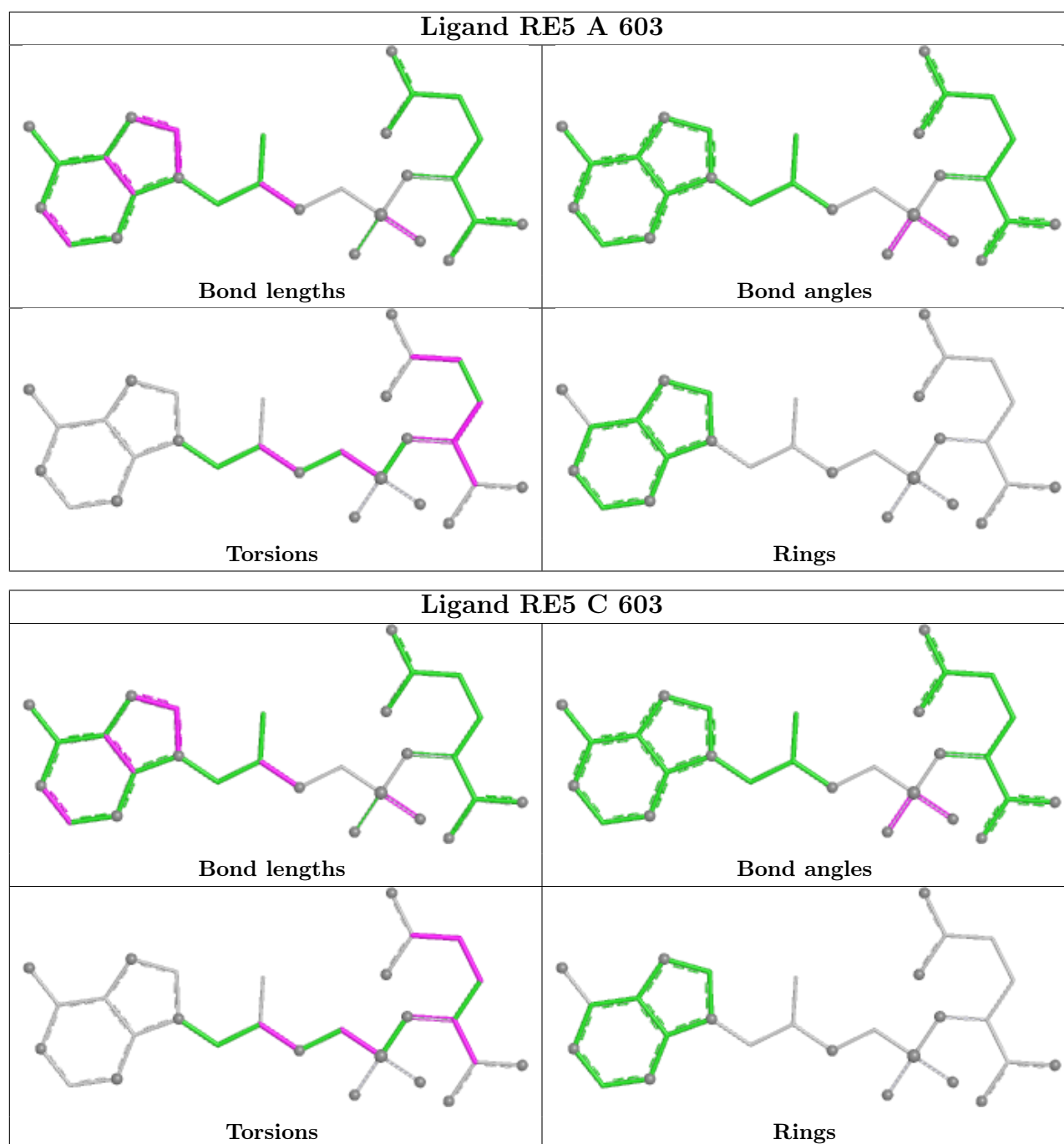
Mol	Chain	Res	Type	Atoms
6	A	603	RE5	O14-C13-P02-O28
6	A	603	RE5	C05-C04-C08-C09
6	A	603	RE5	N03-C04-C08-C09
6	C	603	RE5	N03-C04-C05-O06
6	C	603	RE5	N03-C04-C05-O07
6	A	603	RE5	O14-C13-P02-O01
6	C	603	RE5	O14-C13-P02-O01
6	C	603	RE5	C08-C04-N03-P02
6	C	603	RE5	C08-C04-C05-O06
6	C	603	RE5	C08-C04-C05-O07
6	A	603	RE5	C05-C04-N03-P02
6	A	603	RE5	C08-C09-C10-O12
6	C	603	RE5	C08-C09-C10-O11
6	A	603	RE5	C08-C09-C10-O11
6	C	603	RE5	C04-C08-C09-C10
6	C	603	RE5	C08-C09-C10-O12
6	A	603	RE5	C16-C15-O14-C13
6	C	603	RE5	C16-C15-O14-C13

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	603	RE5	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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	555/556 (99%)	0.10	16 (2%) 53 52	31, 72, 138, 187	0
1	C	556/556 (100%)	0.26	15 (2%) 56 54	33, 97, 163, 198	0
2	B	411/428 (96%)	-0.05	8 (1%) 66 64	27, 63, 119, 154	0
2	D	412/428 (96%)	0.27	19 (4%) 37 36	35, 80, 133, 167	0
3	E	24/27 (88%)	-0.25	0 100 100	75, 126, 194, 203	0
3	T	23/27 (85%)	-0.19	0 100 100	62, 107, 172, 191	0
4	F	18/21 (85%)	-0.44	0 100 100	69, 108, 172, 189	0
4	P	18/21 (85%)	-0.35	0 100 100	62, 92, 162, 164	0
All	All	2017/2064 (97%)	0.13	58 (2%) 53 52	27, 76, 149, 203	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	205	LEU	4.6
2	D	425	LEU	4.6
1	A	554	ALA	4.4
2	D	422	LEU	4.2
2	D	359	GLY	3.9
1	C	167	ILE	3.8
1	A	73	LYS	3.8
1	A	296	THR	3.5
1	C	210	LEU	3.5
2	D	426	TRP	3.4
1	A	128	THR	3.3
2	D	420	PRO	3.3
1	A	552	VAL	3.2
2	B	87	PHE	3.0
1	A	109	LEU	3.0
2	B	88	TRP	2.9

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Mol	Chain	Res	Type	RSRZ
2	D	90	VAL	2.9
2	D	355	ALA	2.9
2	D	361	HIS	2.9
1	A	142	ILE	2.9
2	D	232	TYR	2.8
1	C	124	PHE	2.8
2	D	95	PRO	2.8
2	D	360	ALA	2.8
1	C	214	LEU	2.6
2	B	422	LEU	2.6
2	B	4	PRO	2.6
2	B	229	TRP	2.6
2	D	92	LEU	2.5
2	D	217	PRO	2.5
1	A	132	ILE	2.5
2	B	266	TRP	2.5
1	C	188	TYR	2.4
1	C	37	ILE	2.4
1	C	554	ALA	2.4
1	C	228	LEU	2.4
2	D	214	LEU	2.3
1	C	27	THR	2.3
1	A	193	LEU	2.3
1	C	109	LEU	2.3
2	D	94	ILE	2.2
1	C	223	LYS	2.2
1	A	24	TRP	2.2
2	B	285	GLY	2.2
1	A	46	LYS	2.1
1	C	193	LEU	2.1
1	A	135	ILE	2.1
2	D	124	PHE	2.1
1	A	195	ILE	2.1
1	A	282	LEU	2.0
1	A	130	PHE	2.0
1	A	289	LEU	2.0
2	B	91	GLN	2.0
2	D	283	LEU	2.0
1	C	544	GLY	2.0
2	D	231	GLY	2.0
1	C	35	VAL	2.0
2	D	356	ARG	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(Å ²)	Q<0.9
4	DDG	F	822	21/22	0.78	0.10	79,97,112,118	0
4	MRG	F	817	22/27	0.84	0.07	97,124,160,168	0
4	MRG	P	817	22/27	0.88	0.08	70,91,120,138	0
4	DDG	P	822	21/22	0.94	0.07	48,64,69,73	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

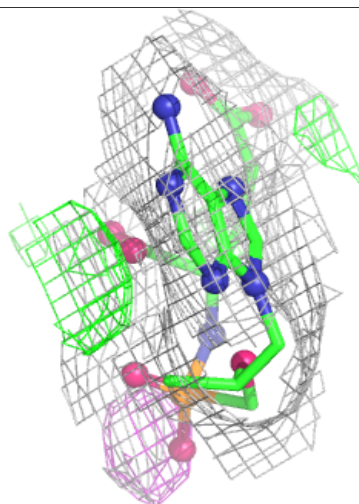
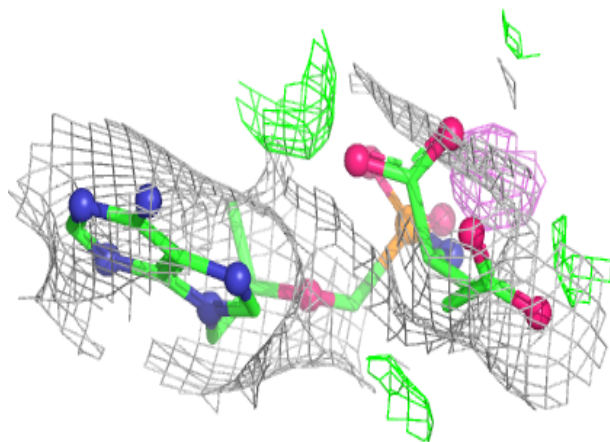
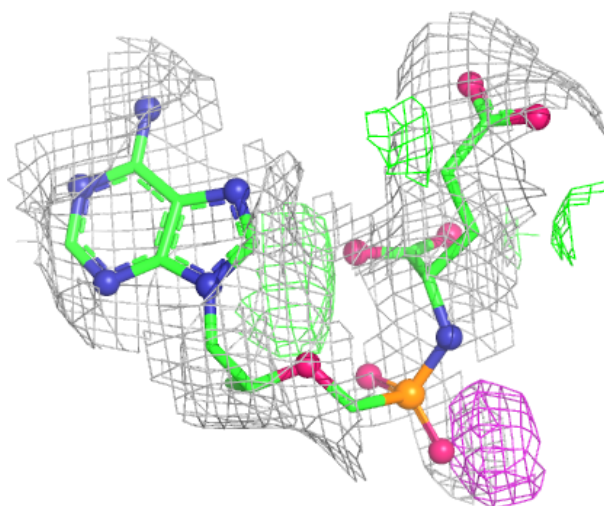
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	MN	A	602	1/1	0.74	0.23	155,155,155,155	0
6	RE5	C	603	28/28	0.75	0.11	91,111,133,138	0
6	RE5	A	603	28/28	0.89	0.10	57,81,101,117	0
5	MN	C	602	1/1	0.90	0.11	85,85,85,85	0
5	MN	C	601	1/1	0.94	0.05	133,133,133,133	0
5	MN	A	601	1/1	0.99	0.03	83,83,83,83	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.

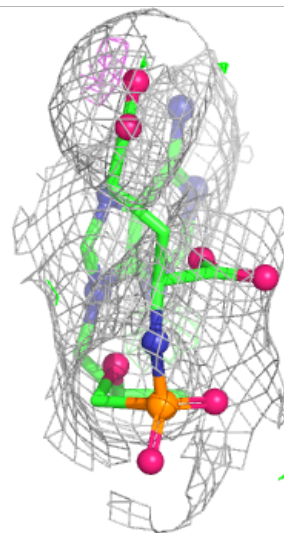
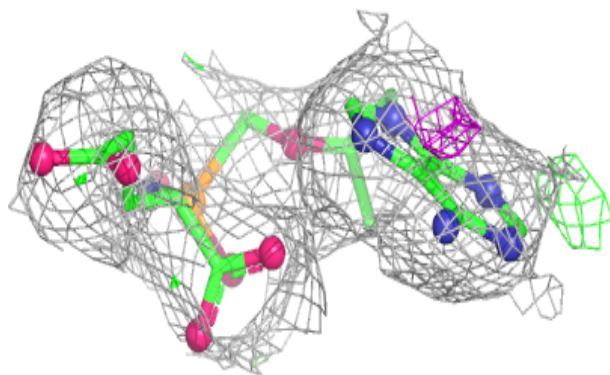
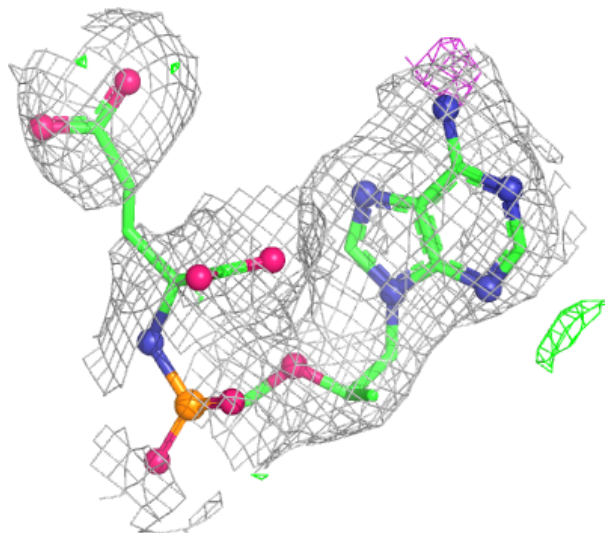
Electron density around RE5 C 603:

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



Electron density around RE5 A 603:

$2mF_o-DF_c$ (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.