



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 04:49 PM EDT

PDB ID : 5HV1 / pdb_00005hv1
Title : Rifampin phosphotransferase in complex with AMPPNP and rifampin from *Listeria monocytogenes*
Authors : Zhang, P.; Qi, X.
Deposited on : 2016-01-28
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

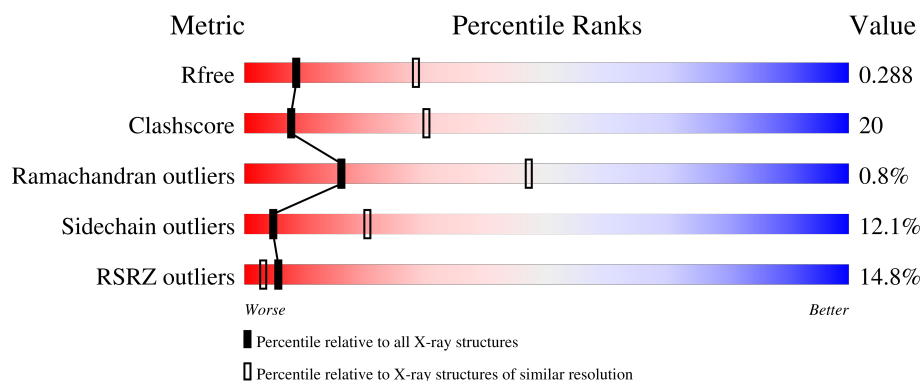
1 Overall quality at a glance

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	883	14% 57% 30% 7% ••

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6757 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.

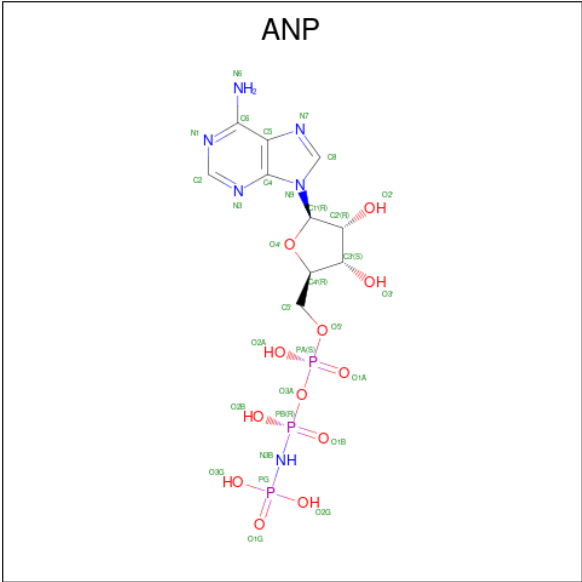
- Molecule 1 is a protein called Phosphoenolpyruvate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	847	6663	4229	1119	1285	30	0	0	0

There are 16 discrepancies between the modelled and reference sequences:

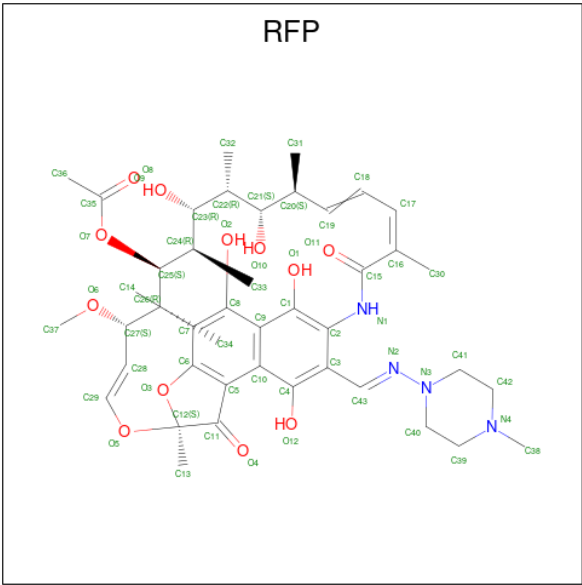
Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	MET	-	expression tag	UNP A0A0S2YLC8
A	-14	ARG	-	expression tag	UNP A0A0S2YLC8
A	-13	GLY	-	expression tag	UNP A0A0S2YLC8
A	-12	SER	-	expression tag	UNP A0A0S2YLC8
A	-11	HIS	-	expression tag	UNP A0A0S2YLC8
A	-10	HIS	-	expression tag	UNP A0A0S2YLC8
A	-9	HIS	-	expression tag	UNP A0A0S2YLC8
A	-8	HIS	-	expression tag	UNP A0A0S2YLC8
A	-7	HIS	-	expression tag	UNP A0A0S2YLC8
A	-6	HIS	-	expression tag	UNP A0A0S2YLC8
A	-5	GLY	-	expression tag	UNP A0A0S2YLC8
A	-4	SER	-	expression tag	UNP A0A0S2YLC8
A	-3	ALA	-	expression tag	UNP A0A0S2YLC8
A	-2	CYS	-	expression tag	UNP A0A0S2YLC8
A	-1	GLU	-	expression tag	UNP A0A0S2YLC8
A	0	LEU	-	expression tag	UNP A0A0S2YLC8

- Molecule 2 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	31	10	6	12	3	0	0

- Molecule 3 is RIFAMPICIN (CCD ID: RFP) (formula: C₄₃H₅₈N₄O₁₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O			
3	A	1	59	43	4	12		0	0

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	3	Total	O	0	0
			3	3		



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	151.09Å 151.09Å 191.65Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.02 – 3.10 30.02 – 3.10	Depositor EDS
% Data completeness (in resolution range)	95.0 (30.02-3.10) 88.5 (30.02-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 3.11Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.250 , 0.286 0.252 , 0.288	Depositor DCC
R_{free} test set	2000 reflections (8.32%)	wwPDB-VP
Wilson B-factor (Å ²)	56.1	Xtriage
Anisotropy	0.044	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 31.7	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	0.87	EDS
Total number of atoms	6757	wwPDB-VP
Average B, all atoms (Å ²)	56.0	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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ANP, RFP, MG

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	0.99	9/6782 (0.1%)	1.27	53/9174 (0.6%)

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	2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	314	THR	CA-C	7.85	1.62	1.52
1	A	315	THR	CA-C	6.01	1.60	1.52
1	A	274	GLN	CD-OE1	-5.92	1.12	1.23
1	A	233	GLU	CA-C	-5.84	1.48	1.52
1	A	108	GLY	C-N	-5.48	1.25	1.33
1	A	586	GLU	C-O	-5.36	1.17	1.24
1	A	586	GLU	CA-C	-5.22	1.45	1.52
1	A	585	GLY	C-O	-5.15	1.17	1.23
1	A	468	VAL	CA-CB	-5.03	1.47	1.54

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	842	GLU	CA-C-N	28.80	176.55	121.54
1	A	842	GLU	C-N-CA	28.80	176.55	121.54
1	A	842	GLU	N-CA-C	-17.39	85.02	110.23
1	A	232	ARG	N-CA-C	14.24	126.52	111.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	842	GLU	CB-CA-C	-13.46	87.37	109.72
1	A	382	MET	N-CA-C	-13.15	82.79	110.80
1	A	108	GLY	N-CA-C	-12.80	92.36	112.58
1	A	382	MET	CB-CA-C	9.93	130.18	110.42
1	A	109	GLY	N-CA-C	-9.40	86.04	113.30
1	A	752	ASN	N-CA-C	-9.39	95.02	109.95
1	A	770	VAL	N-CA-C	-8.91	90.81	109.34
1	A	735	PHE	N-CA-C	-7.74	97.86	109.86
1	A	221	SER	N-CA-C	6.95	122.73	112.94
1	A	108	GLY	CA-C-N	6.74	133.82	121.70
1	A	108	GLY	C-N-CA	6.74	133.82	121.70
1	A	770	VAL	CB-CA-C	6.43	121.83	111.29
1	A	664	GLY	N-CA-C	-6.40	105.07	112.50
1	A	380	ARG	CA-C-N	6.17	132.81	121.70
1	A	380	ARG	C-N-CA	6.17	132.81	121.70
1	A	231	VAL	N-CA-CB	-6.15	104.71	112.15
1	A	324	GLU	C-N-CD	-6.15	99.78	125.00
1	A	391	ASP	CA-C-N	6.13	127.50	119.84
1	A	391	ASP	C-N-CA	6.13	127.50	119.84
1	A	792	ASP	N-CA-C	6.05	118.38	111.11
1	A	788	ALA	CA-C-N	-5.98	114.77	122.79
1	A	788	ALA	C-N-CA	-5.98	114.77	122.79
1	A	524	SER	N-CA-C	-5.96	104.72	112.23
1	A	866	LEU	CB-CG-CD2	-5.95	92.85	110.70
1	A	750	MET	CA-C-N	-5.77	115.14	123.11
1	A	750	MET	C-N-CA	-5.77	115.14	123.11
1	A	546	TYR	N-CA-C	-5.75	104.92	111.14
1	A	223	LEU	CA-C-N	5.74	132.03	121.70
1	A	223	LEU	C-N-CA	5.74	132.03	121.70
1	A	841	VAL	N-CA-C	5.72	116.65	107.73
1	A	607	SER	CA-C-N	-5.67	113.57	122.65
1	A	607	SER	C-N-CA	-5.67	113.57	122.65
1	A	804	SER	CB-CA-C	-5.58	99.31	110.42
1	A	860	GLU	N-CA-C	-5.47	102.18	110.28
1	A	119	SER	CB-CA-C	5.38	118.27	110.79
1	A	287	GLU	N-CA-C	-5.30	105.41	111.14
1	A	508	LYS	N-CA-C	5.26	118.00	108.69
1	A	577	GLN	N-CA-C	-5.24	106.39	112.89
1	A	756	LYS	N-CA-C	5.16	118.06	107.69
1	A	108	GLY	O-C-N	-5.15	116.63	122.91
1	A	72	ASP	N-CA-C	-5.14	99.62	108.56
1	A	741	PRO	CA-C-O	-5.12	115.83	122.13

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	495	ALA	N-CA-C	-5.12	105.61	111.14
1	A	518	ALA	CA-C-N	-5.08	115.12	120.66
1	A	518	ALA	C-N-CA	-5.08	115.12	120.66
1	A	602	ILE	CA-C-N	-5.05	114.41	119.56
1	A	602	ILE	C-N-CA	-5.05	114.41	119.56
1	A	107	VAL	N-CA-C	5.04	119.83	109.34
1	A	119	SER	N-CA-C	-5.04	98.84	107.61

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	107	VAL	Peptide
1	A	770	VAL	Peptide

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	6663	0	6704	264	2
2	A	31	0	13	5	0
3	A	59	0	55	13	0
4	A	1	0	0	0	0
5	A	3	0	0	3	0
All	All	6757	0	6772	271	2

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 20.

All (271) 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)
3:A:902:RFP:H311	3:A:902:RFP:H323	1.32	1.09
1:A:583:CYS:O	1:A:586:GLU:HG3	1.58	1.02
1:A:581:MET:HE3	1:A:671:TYR:CE1	2.01	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:THR:HG22	1:A:46:GLU:H	1.29	0.95
1:A:328:ARG:NH2	1:A:371:THR:OG1	2.02	0.92
1:A:582:ARG:HD3	1:A:586:GLU:CD	1.93	0.92
1:A:393:LEU:HD11	1:A:755:TYR:CZ	2.06	0.91
1:A:381:ASP:H	1:A:384:VAL:HG22	1.35	0.90
1:A:578:LYS:O	1:A:581:MET:HE1	1.70	0.89
1:A:322:VAL:HG12	1:A:324:GLU:H	1.36	0.87
1:A:642:LEU:HG	1:A:643:PRO:HD2	1.60	0.84
1:A:745:THR:HG22	1:A:747:ASP:H	1.43	0.83
3:A:902:RFP:O11	3:A:902:RFP:O1	1.97	0.83
1:A:353:MET:HE3	1:A:471:LEU:HD11	1.61	0.82
1:A:696:ILE:CD1	1:A:720:VAL:HG22	2.11	0.81
1:A:581:MET:CE	1:A:671:TYR:CE1	2.63	0.81
1:A:380:ARG:HH22	1:A:408:LEU:HD23	1.44	0.80
1:A:232:ARG:HG2	1:A:232:ARG:HH11	1.46	0.80
1:A:233:GLU:C	1:A:233:GLU:OE1	2.27	0.78
1:A:287:GLU:O	1:A:291:GLY:N	2.14	0.77
1:A:581:MET:O	1:A:593:ARG:HG3	1.85	0.77
1:A:762:LYS:NZ	1:A:763:ASP:OD2	2.19	0.76
1:A:488:MET:HE1	3:A:902:RFP:H362	1.67	0.75
1:A:550:THR:HG22	1:A:552:ASN:H	1.53	0.74
1:A:823:MET:H	1:A:831:ARG:HH22	1.35	0.73
1:A:328:ARG:H	1:A:746:SER:HG	1.36	0.73
3:A:902:RFP:O4	3:A:902:RFP:O12	2.06	0.73
1:A:841:VAL:HG12	1:A:842:GLU:O	1.89	0.73
1:A:328:ARG:N	1:A:746:SER:OG	2.17	0.72
1:A:855:ARG:HB2	1:A:866:LEU:CD1	2.20	0.71
1:A:20:GLY:HA2	1:A:121:THR:HG23	1.70	0.71
1:A:137:TYR:HB2	1:A:140:ILE:HD11	1.71	0.71
1:A:3:PRO:HG2	1:A:6:LEU:HD11	1.72	0.70
1:A:64:ARG:HB2	1:A:79:ILE:HD11	1.72	0.70
1:A:753:GLY:N	5:A:1001:HOH:O	2.08	0.70
1:A:60:GLN:O	1:A:64:ARG:HG2	1.92	0.69
1:A:696:ILE:HD11	1:A:720:VAL:HG22	1.73	0.69
1:A:286:ILE:O	1:A:290:PHE:HB2	1.91	0.69
1:A:65:LEU:HD12	1:A:68:LEU:HD11	1.75	0.68
1:A:221:SER:N	1:A:222:GLY:HA2	2.08	0.68
1:A:345:PRO:HG2	1:A:705:PHE:HA	1.76	0.68
1:A:855:ARG:HB2	1:A:866:LEU:HD12	1.74	0.68
1:A:329:VAL:H	1:A:371:THR:HG22	1.59	0.67
1:A:696:ILE:HD12	1:A:720:VAL:HG22	1.76	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:902:RFP:H311	3:A:902:RFP:C32	2.17	0.66
1:A:530:LEU:HA	1:A:533:VAL:HG12	1.76	0.65
1:A:581:MET:CE	1:A:671:TYR:HE1	2.08	0.65
1:A:136:THR:OG1	2:A:901:ANP:O2A	2.06	0.65
1:A:504:TRP:HB3	1:A:642:LEU:HD21	1.78	0.64
1:A:582:ARG:HD3	1:A:586:GLU:OE1	1.96	0.64
1:A:702:ASP:OD1	1:A:729:LYS:NZ	2.31	0.64
1:A:385:ASN:OD1	1:A:386:SER:N	2.31	0.64
1:A:233:GLU:OE1	1:A:233:GLU:O	2.17	0.63
1:A:487:ILE:HD11	1:A:669:PRO:HG2	1.81	0.62
1:A:218:ALA:HB1	1:A:224:VAL:H	1.64	0.62
1:A:131:ALA:HA	2:A:901:ANP:O3G	2.00	0.62
1:A:841:VAL:C	1:A:842:GLU:O	2.34	0.62
1:A:80:SER:OG	1:A:84:ARG:NH2	2.32	0.62
1:A:377:LYS:O	1:A:378:VAL:HG12	2.00	0.62
1:A:393:LEU:HD11	1:A:755:TYR:CE1	2.35	0.61
1:A:465:ASP:OD2	1:A:712:ARG:NH2	2.33	0.61
1:A:514:LEU:HD13	1:A:656:ILE:HD13	1.81	0.61
1:A:769:PRO:O	1:A:770:VAL:C	2.43	0.61
1:A:22:LYS:NZ	2:A:901:ANP:O2G	2.31	0.61
1:A:757:ARG:O	1:A:757:ARG:HG3	2.00	0.61
1:A:340:THR:O	1:A:365:ARG:NH1	2.34	0.61
1:A:380:ARG:O	1:A:384:VAL:HG13	2.02	0.60
1:A:752:ASN:CA	5:A:1001:HOH:O	2.48	0.60
1:A:772:SER:HA	1:A:858:GLY:O	2.01	0.60
1:A:659:LEU:O	1:A:663:ILE:HG12	2.01	0.59
1:A:359:MET:HE3	1:A:366:LEU:HB3	1.83	0.59
1:A:451:LEU:O	1:A:455:ILE:HD12	2.03	0.59
1:A:841:VAL:O	1:A:842:GLU:C	2.43	0.59
1:A:404:LYS:NZ	1:A:404:LYS:HA	2.18	0.59
1:A:286:ILE:HG22	1:A:290:PHE:CD1	2.38	0.59
1:A:593:ARG:HG2	1:A:678:PHE:CD1	2.38	0.58
1:A:752:ASN:HA	5:A:1001:HOH:O	2.02	0.58
1:A:741:PRO:HB3	1:A:750:MET:HB3	1.84	0.58
1:A:368:VAL:HG11	3:A:902:RFP:H133	1.86	0.57
1:A:455:ILE:C	1:A:457:THR:H	2.13	0.57
1:A:44:THR:HG22	1:A:46:GLU:N	2.11	0.57
1:A:581:MET:HE3	1:A:671:TYR:CZ	2.39	0.57
1:A:548:GLU:HA	1:A:606:LEU:HD21	1.87	0.56
1:A:584:ALA:HA	1:A:804:SER:CB	2.35	0.56
1:A:286:ILE:HG22	1:A:290:PHE:HD1	1.69	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:806:THR:O	1:A:808:ALA:N	2.38	0.56
1:A:368:VAL:HG12	1:A:370:ILE:HG23	1.86	0.56
1:A:782:ILE:HG21	1:A:788:ALA:HB2	1.87	0.56
1:A:648:LYS:O	1:A:652:THR:HG23	2.05	0.56
1:A:72:ASP:O	1:A:76:ILE:HG13	2.06	0.55
1:A:101:ASP:OD2	1:A:143:LYS:NZ	2.36	0.55
1:A:369:ASP:OD1	1:A:371:THR:HG23	2.06	0.55
1:A:50:ARG:HG2	1:A:96:ILE:HD11	1.88	0.55
1:A:537:ILE:HB	1:A:609:ILE:HD11	1.88	0.55
3:A:902:RFP:O9	3:A:902:RFP:O10	2.12	0.55
1:A:329:VAL:H	1:A:371:THR:CG2	2.18	0.55
3:A:902:RFP:H323	3:A:902:RFP:C31	2.14	0.55
1:A:352:LEU:HD23	1:A:359:MET:HE2	1.89	0.54
1:A:534:ALA:HB1	1:A:609:ILE:HG13	1.89	0.54
1:A:589:LEU:O	1:A:681:LYS:HD2	2.07	0.54
1:A:696:ILE:HD12	1:A:720:VAL:CG2	2.36	0.54
1:A:721:ASP:O	1:A:724:LEU:HB3	2.08	0.54
1:A:246:LEU:HD13	1:A:258:THR:HG21	1.90	0.54
1:A:324:GLU:HB2	1:A:328:ARG:HE	1.71	0.54
1:A:802:ASP:HB2	1:A:803:PRO:HD2	1.88	0.54
1:A:324:GLU:CB	1:A:328:ARG:HH11	2.21	0.54
1:A:322:VAL:HG21	1:A:330:TYR:OH	2.07	0.54
1:A:801:THR:HG21	1:A:809:PHE:HZ	1.73	0.54
1:A:227:ASP:OD2	1:A:241:ILE:HA	2.08	0.54
1:A:457:THR:HG22	1:A:458:LYS:HD3	1.89	0.54
1:A:471:LEU:O	1:A:475:LYS:HG3	2.08	0.53
1:A:735:PHE:O	1:A:736:GLU:HB2	2.08	0.53
1:A:356:PRO:HG2	3:A:902:RFP:H402	1.91	0.53
1:A:569:LYS:O	1:A:573:GLU:HG3	2.08	0.53
1:A:335:HIS:CD2	1:A:819:VAL:HG22	2.44	0.53
1:A:516:GLU:HA	1:A:832:GLU:HA	1.90	0.52
1:A:351:TYR:CD2	1:A:359:MET:HE1	2.43	0.52
1:A:284:ARG:HG3	1:A:284:ARG:HH11	1.74	0.52
1:A:547:LEU:HG	1:A:602:ILE:HD12	1.91	0.52
1:A:232:ARG:NH1	1:A:232:ARG:CG	2.73	0.52
1:A:350:PHE:O	1:A:354:THR:HG22	2.08	0.52
1:A:119:SER:O	1:A:179:ALA:N	2.41	0.52
1:A:343:MET:HB2	1:A:348:LEU:HD13	1.91	0.52
1:A:351:TYR:HD2	1:A:359:MET:HE1	1.76	0.51
1:A:685:LEU:HD22	1:A:700:HIS:CE1	2.46	0.51
1:A:232:ARG:HH11	1:A:232:ARG:CG	2.17	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:398:LEU:O	1:A:401:VAL:HG12	2.11	0.51
1:A:539:PRO:O	1:A:541:PRO:HD3	2.11	0.51
1:A:521:ASN:O	1:A:523:THR:N	2.44	0.51
1:A:715:VAL:HG23	1:A:716:ARG:N	2.25	0.51
1:A:354:THR:HG21	1:A:677:TYR:OH	2.11	0.51
1:A:494:SER:HB3	1:A:659:LEU:CD1	2.41	0.50
1:A:204:ARG:HG3	1:A:204:ARG:HH11	1.76	0.50
1:A:110:TYR:HE2	1:A:146:LEU:HD12	1.75	0.50
1:A:530:LEU:HA	1:A:533:VAL:CG1	2.41	0.50
1:A:232:ARG:HG2	1:A:232:ARG:NH1	2.19	0.50
1:A:330:TYR:HB3	1:A:367:PHE:HB3	1.94	0.50
1:A:247:ALA:HB2	1:A:261:LEU:HD11	1.92	0.50
1:A:322:VAL:HG12	1:A:324:GLU:N	2.17	0.49
1:A:380:ARG:HB3	1:A:381:ASP:HB2	1.93	0.49
1:A:508:LYS:HB3	1:A:510:VAL:HG13	1.94	0.49
1:A:521:ASN:OD1	1:A:523:THR:HG22	2.12	0.49
1:A:546:TYR:O	1:A:550:THR:OG1	2.28	0.49
1:A:584:ALA:HA	1:A:804:SER:HB2	1.94	0.49
1:A:287:GLU:HG3	1:A:288:ALA:N	2.28	0.49
1:A:494:SER:HB3	1:A:659:LEU:HD13	1.94	0.49
1:A:589:LEU:HD12	1:A:677:TYR:HB3	1.95	0.49
1:A:104:LEU:O	1:A:108:GLY:HA2	2.13	0.49
1:A:530:LEU:HD12	1:A:533:VAL:HG11	1.95	0.48
1:A:286:ILE:HD13	1:A:296:ILE:HD13	1.96	0.48
1:A:398:LEU:O	1:A:402:ILE:HG23	2.13	0.48
1:A:523:THR:CG2	1:A:667:GLU:HG2	2.43	0.48
1:A:404:LYS:HB2	1:A:407:PHE:HB2	1.95	0.48
1:A:441:VAL:HG13	1:A:668:TYR:CZ	2.48	0.48
1:A:802:ASP:OD1	1:A:804:SER:HB3	2.14	0.48
1:A:392:PRO:HG2	1:A:768:LEU:HD21	1.94	0.48
1:A:393:LEU:HD12	1:A:393:LEU:H	1.79	0.48
1:A:757:ARG:HH11	1:A:757:ARG:HG2	1.77	0.48
1:A:66:SER:HB2	1:A:174:ARG:HH21	1.78	0.48
1:A:685:LEU:HD22	1:A:700:HIS:HE1	1.79	0.47
1:A:203:ASN:OD1	1:A:205:LYS:HB2	2.15	0.47
1:A:507:GLU:C	1:A:508:LYS:HG2	2.39	0.47
1:A:287:GLU:O	1:A:290:PHE:N	2.47	0.47
1:A:433:ASP:OD1	1:A:434:SER:N	2.48	0.47
1:A:227:ASP:HB3	1:A:229:TYR:CZ	2.49	0.47
1:A:770:VAL:CG1	1:A:827:ALA:HB1	2.45	0.47
1:A:203:ASN:C	1:A:205:LYS:H	2.21	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:359:MET:HG3	1:A:368:VAL:HG22	1.97	0.47
1:A:478:LEU:HD21	3:A:902:RFP:H17C	1.97	0.47
1:A:62:LEU:HD11	1:A:158:PHE:CE1	2.50	0.47
1:A:511:ALA:O	1:A:515:SER:OG	2.25	0.47
1:A:223:LEU:N	1:A:223:LEU:HD22	2.30	0.47
1:A:253:GLU:N	1:A:253:GLU:OE1	2.48	0.47
1:A:269:GLN:HG2	1:A:271:LEU:O	2.14	0.47
1:A:505:LEU:HA	1:A:505:LEU:HD13	1.42	0.47
1:A:202:SER:OG	1:A:316:LEU:HD23	2.15	0.46
1:A:405:LYS:HB2	1:A:405:LYS:NZ	2.30	0.46
1:A:556:MET:O	1:A:569:LYS:HD2	2.15	0.46
1:A:583:CYS:O	1:A:586:GLU:CG	2.47	0.46
1:A:363:GLY:HA2	1:A:735:PHE:CE1	2.51	0.46
1:A:224:VAL:HG22	1:A:225:SER:N	2.30	0.46
1:A:502:TYR:O	1:A:506:GLY:HA2	2.14	0.46
1:A:559:LEU:O	1:A:565:GLY:HA3	2.16	0.46
1:A:117:ARG:NH2	2:A:901:ANP:O1A	2.48	0.45
1:A:491:MET:HE1	3:A:902:RFP:H361	1.97	0.45
1:A:584:ALA:HB2	1:A:804:SER:OG	2.16	0.45
1:A:328:ARG:HA	1:A:371:THR:HG21	1.96	0.45
1:A:824:THR:O	1:A:825:HIS:C	2.59	0.45
1:A:16:GLU:O	1:A:121:THR:HG21	2.16	0.45
1:A:198:ASP:O	1:A:202:SER:HA	2.17	0.45
1:A:348:LEU:HD12	1:A:366:LEU:HD21	1.97	0.45
1:A:506:GLY:O	1:A:508:LYS:HG2	2.17	0.45
1:A:283:GLY:HA2	1:A:286:ILE:CD1	2.46	0.45
1:A:583:CYS:H	1:A:586:GLU:HG2	1.81	0.44
1:A:339:MET:CE	1:A:341:ASP:HB2	2.46	0.44
1:A:455:ILE:HG13	1:A:466:PHE:CE1	2.52	0.44
3:A:902:RFP:H18C	3:A:902:RFP:H211	1.85	0.44
1:A:201:THR:C	1:A:203:ASN:H	2.24	0.44
1:A:438:PHE:N	1:A:438:PHE:CD2	2.85	0.44
1:A:245:LYS:H	1:A:245:LYS:HD2	1.81	0.44
1:A:374:LEU:HD12	1:A:374:LEU:HA	1.70	0.44
1:A:439:GLU:HG2	1:A:442:ARG:NH2	2.32	0.44
1:A:239:LYS:HB3	1:A:239:LYS:HE2	1.74	0.44
1:A:337:GLN:HG2	1:A:587:ILE:HD13	1.99	0.44
1:A:406:GLY:O	1:A:407:PHE:C	2.60	0.44
1:A:491:MET:HE3	1:A:491:MET:HB3	1.87	0.44
1:A:646:GLU:O	1:A:650:MET:HG3	2.17	0.44
1:A:583:CYS:HB3	1:A:593:ARG:HA	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:ARG:CB	1:A:64:ARG:HH11	2.31	0.44
1:A:324:GLU:CG	1:A:328:ARG:HH11	2.31	0.44
1:A:283:GLY:HA2	1:A:286:ILE:HD12	2.00	0.43
1:A:311:ARG:HA	1:A:312:PRO:HD2	1.90	0.43
1:A:584:ALA:CA	1:A:804:SER:HB2	2.48	0.43
1:A:313:ILE:HG22	1:A:314:THR:N	2.32	0.43
1:A:670:LYS:O	1:A:674:ILE:HG12	2.18	0.43
1:A:536:VAL:O	1:A:539:PRO:HD2	2.19	0.43
1:A:582:ARG:CD	1:A:586:GLU:CD	2.80	0.43
1:A:765:ILE:HB	1:A:863:ILE:HB	2.00	0.43
1:A:769:PRO:HB3	1:A:859:THR:O	2.19	0.43
1:A:194:LEU:HD21	1:A:207:LEU:HD21	2.01	0.43
1:A:829:ILE:HD13	1:A:829:ILE:HA	1.78	0.43
1:A:22:LYS:HB2	1:A:22:LYS:HE3	1.85	0.43
1:A:404:LYS:O	1:A:406:GLY:N	2.51	0.43
1:A:324:GLU:HB3	1:A:328:ARG:HH11	1.83	0.42
1:A:405:LYS:HE3	1:A:405:LYS:O	2.19	0.42
1:A:324:GLU:HA	1:A:325:PRO:HD3	1.53	0.42
1:A:356:PRO:HD2	3:A:902:RFP:HC43	1.99	0.42
1:A:581:MET:HG2	1:A:582:ARG:N	2.33	0.42
1:A:94:SER:O	1:A:98:SER:OG	2.37	0.42
1:A:81:GLU:O	1:A:85:THR:HG23	2.19	0.42
1:A:768:LEU:HA	1:A:769:PRO:HD2	1.86	0.42
1:A:551:LYS:O	1:A:553:PRO:HD3	2.19	0.42
1:A:642:LEU:CG	1:A:643:PRO:HD2	2.41	0.42
1:A:237:THR:C	1:A:238:ASN:HD22	2.27	0.42
1:A:359:MET:HE3	1:A:366:LEU:HD22	2.01	0.42
1:A:365:ARG:HD2	1:A:740:PRO:HB3	2.02	0.42
1:A:440:LEU:HA	1:A:440:LEU:HD23	1.84	0.42
1:A:735:PHE:O	1:A:736:GLU:CB	2.67	0.42
1:A:370:ILE:HD12	1:A:370:ILE:O	2.20	0.42
1:A:62:LEU:HD23	1:A:174:ARG:HA	2.01	0.41
1:A:339:MET:HE3	1:A:341:ASP:HB2	2.02	0.41
1:A:451:LEU:CD1	1:A:455:ILE:HD11	2.49	0.41
1:A:7:LYS:NZ	1:A:106:ASP:OD2	2.49	0.41
1:A:392:PRO:HD2	1:A:768:LEU:HD21	2.03	0.41
1:A:768:LEU:HA	1:A:768:LEU:HD23	1.81	0.41
1:A:798:THR:OG1	1:A:799:ALA:N	2.53	0.41
1:A:406:GLY:C	1:A:408:LEU:N	2.75	0.41
1:A:762:LYS:O	1:A:763:ASP:HB2	2.19	0.41
1:A:398:LEU:HA	1:A:401:VAL:HG12	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:768:LEU:H	1:A:840:GLY:H	1.69	0.41
1:A:5:VAL:C	1:A:6:LEU:HD12	2.46	0.41
1:A:345:PRO:HG2	1:A:705:PHE:CA	2.49	0.41
1:A:674:ILE:O	1:A:677:TYR:N	2.53	0.41
1:A:210:ASP:HA	1:A:227:ASP:O	2.20	0.41
1:A:380:ARG:NH2	1:A:408:LEU:HD23	2.23	0.41
1:A:404:LYS:HA	1:A:404:LYS:CE	2.51	0.41
1:A:652:THR:O	1:A:653:LYS:C	2.62	0.41
1:A:246:LEU:HD13	1:A:258:THR:CG2	2.51	0.41
1:A:854:ILE:HG21	1:A:854:ILE:HD13	1.67	0.41
1:A:101:ASP:O	1:A:105:LEU:HD13	2.22	0.40
1:A:339:MET:O	1:A:800:TYR:CD2	2.74	0.40
1:A:397:ALA:O	1:A:400:THR:HG22	2.21	0.40
1:A:474:LEU:HD23	1:A:474:LEU:C	2.46	0.40
1:A:783:LEU:HD23	1:A:783:LEU:HA	1.82	0.40
1:A:117:ARG:NE	2:A:901:ANP:O1A	2.48	0.40
1:A:741:PRO:HG3	1:A:750:MET:SD	2.62	0.40
1:A:64:ARG:HH11	1:A:64:ARG:HB3	1.87	0.40
1:A:18:LEU:HG	1:A:44:THR:HG23	2.04	0.40
1:A:218:ALA:HB2	1:A:248:ILE:HG13	2.04	0.40
1:A:696:ILE:CD1	1:A:720:VAL:CG2	2.92	0.40
1:A:781:VAL:C	1:A:782:ILE:HG13	2.46	0.40

All (2) 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)
1:A:225:SER:OG	1:A:225:SER:OG[9_555]	2.08	0.12
1:A:50:ARG:NH2	1:A:613:ASP:OD2[8_565]	2.17	0.03

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	843/883 (96%)	778 (92%)	58 (7%)	7 (1%)	16	47

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	378	VAL
1	A	804	SER
1	A	770	VAL
1	A	325	PRO
1	A	807	PRO
1	A	643	PRO
1	A	2	LYS

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	728/759 (96%)	640 (88%)	88 (12%)	5	20

All (88) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	26	LEU
1	A	44	THR
1	A	62	LEU
1	A	64	ARG
1	A	68	LEU
1	A	88	GLN
1	A	98	SER
1	A	200	ILE
1	A	202	SER
1	A	204	ARG
1	A	205	LYS
1	A	206	SER
1	A	207	LEU

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Mol	Chain	Res	Type
1	A	208	SER
1	A	209	ILE
1	A	220	VAL
1	A	223	LEU
1	A	225	SER
1	A	228	SER
1	A	230	THR
1	A	233	GLU
1	A	238	ASN
1	A	239	LYS
1	A	256	THR
1	A	292	LYS
1	A	297	GLU
1	A	308	VAL
1	A	314	THR
1	A	323	ASN
1	A	328	ARG
1	A	331	ILE
1	A	348	LEU
1	A	359	MET
1	A	370	ILE
1	A	372	GLN
1	A	374	LEU
1	A	375	SER
1	A	378	VAL
1	A	383	MET
1	A	386	SER
1	A	394	ILE
1	A	400	THR
1	A	404	LYS
1	A	405	LYS
1	A	408	LEU
1	A	430	SER
1	A	442	ARG
1	A	449	LYS
1	A	457	THR
1	A	491	MET
1	A	503	GLN
1	A	505	LEU
1	A	507	GLU
1	A	508	LYS
1	A	510	VAL

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Mol	Chain	Res	Type
1	A	513	LYS
1	A	547	LEU
1	A	556	MET
1	A	578	LYS
1	A	581	MET
1	A	586	GLU
1	A	591	LYS
1	A	606	LEU
1	A	609	ILE
1	A	625	GLU
1	A	642	LEU
1	A	656	ILE
1	A	696	ILE
1	A	700	HIS
1	A	709	GLU
1	A	716	ARG
1	A	720	VAL
1	A	723	GLU
1	A	737	LYS
1	A	745	THR
1	A	746	SER
1	A	757	ARG
1	A	766	LEU
1	A	772	SER
1	A	784	GLU
1	A	804	SER
1	A	815	LEU
1	A	819	VAL
1	A	822	LEU
1	A	835	LEU
1	A	849	LYS
1	A	866	LEU

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	89	HIS
1	A	148	GLN
1	A	238	ASN
1	A	500	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is 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	$\# Z > 2$
3	RFP	A	902	-	63,63,63	1.78	13 (20%)	94,94,94	2.15	28 (29%)
2	ANP	A	901	4	33,33,33	1.01	2 (6%)	45,52,52	0.94	3 (6%)

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	RFP	A	902	-	-	11/60/85/85	0/5/5/5
2	ANP	A	901	4	-	8/18/38/38	0/3/3/3

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	902	RFP	C2-N1	-6.02	1.32	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	902	RFP	C6-C7	3.65	1.46	1.39
3	A	902	RFP	O7-C35	3.37	1.42	1.35
3	A	902	RFP	C15-N1	-3.37	1.28	1.35
3	A	902	RFP	C2-C1	3.25	1.46	1.38
3	A	902	RFP	O7-C25	-3.17	1.40	1.44
3	A	902	RFP	N3-N2	-3.11	1.30	1.38
2	A	901	ANP	PG-N3B	2.86	1.70	1.63
3	A	902	RFP	O5-C29	2.77	1.46	1.39
2	A	901	ANP	PG-O1G	2.58	1.50	1.46
3	A	902	RFP	C33-C24	-2.45	1.48	1.53
3	A	902	RFP	O3-C12	-2.45	1.37	1.45
3	A	902	RFP	C12-C11	-2.37	1.45	1.54
3	A	902	RFP	C30-C16	-2.07	1.45	1.50
3	A	902	RFP	C41-N3	-2.05	1.42	1.46

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	902	RFP	C3-C43-N2	-6.53	110.91	121.58
3	A	902	RFP	O3-C6-C7	5.83	131.04	121.16
3	A	902	RFP	C42-C41-N3	-5.24	102.01	110.51
3	A	902	RFP	C5-C10-C4	-5.04	119.18	124.08
3	A	902	RFP	C31-C20-C19	-4.85	98.93	109.91
3	A	902	RFP	O7-C35-C36	4.20	118.58	111.09
3	A	902	RFP	C43-N2-N3	4.16	126.50	120.55
3	A	902	RFP	O4-C11-C12	4.02	128.65	120.56
3	A	902	RFP	O4-C11-C5	-3.85	124.53	131.63
3	A	902	RFP	C41-C42-N4	-3.65	104.98	110.86
3	A	902	RFP	C38-N4-C42	-3.56	103.85	110.63
3	A	902	RFP	O10-C21-C22	-3.36	102.11	109.56
3	A	902	RFP	C3-C2-N1	-3.34	113.60	119.33
3	A	902	RFP	O9-C23-C22	-3.26	102.32	109.56
3	A	902	RFP	C17-C18-C19	-3.13	117.16	124.43
3	A	902	RFP	C3-C4-C10	-3.03	118.73	121.19
3	A	902	RFP	O11-C15-C16	-3.01	115.95	121.49
3	A	902	RFP	O5-C12-C13	2.99	114.74	106.98
3	A	902	RFP	C13-C12-C11	-2.76	107.11	113.90
3	A	902	RFP	C12-O5-C29	2.75	124.59	117.84
3	A	902	RFP	C42-N4-C39	2.71	113.85	109.54
3	A	902	RFP	C32-C22-C23	-2.59	106.34	111.43
2	A	901	ANP	O2A-PA-O3A	-2.58	100.29	107.27
3	A	902	RFP	C5-C6-C7	-2.24	119.89	125.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	902	RFP	O6-C27-C26	-2.18	103.47	107.97
2	A	901	ANP	O3G-PG-O1G	-2.17	108.00	113.45
2	A	901	ANP	O2B-PB-O3A	2.14	111.78	104.64
3	A	902	RFP	C12-O3-C6	2.13	111.24	107.69
3	A	902	RFP	C2-C3-C43	-2.10	121.76	123.99
3	A	902	RFP	C38-N4-C39	-2.05	106.72	110.63
3	A	902	RFP	C1-C2-N1	2.00	125.78	119.67

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	901	ANP	PB-N3B-PG-O1G
2	A	901	ANP	PG-N3B-PB-O1B
2	A	901	ANP	PA-O3A-PB-O2B
2	A	901	ANP	PB-O3A-PA-O5'
2	A	901	ANP	C5'-O5'-PA-O3A
3	A	902	RFP	C19-C20-C21-C22
3	A	902	RFP	C19-C20-C21-O10
3	A	902	RFP	C31-C20-C21-C22
3	A	902	RFP	C31-C20-C21-O10
3	A	902	RFP	C43-N2-N3-C41
3	A	902	RFP	C3-C2-N1-C15
3	A	902	RFP	C4-C3-C43-N2
2	A	901	ANP	O4'-C4'-C5'-O5'
3	A	902	RFP	C43-N2-N3-C40
2	A	901	ANP	C3'-C4'-C5'-O5'
3	A	902	RFP	C1-C2-N1-C15
2	A	901	ANP	C5'-O5'-PA-O1A
3	A	902	RFP	C2-C3-C43-N2
3	A	902	RFP	O7-C25-C26-C34

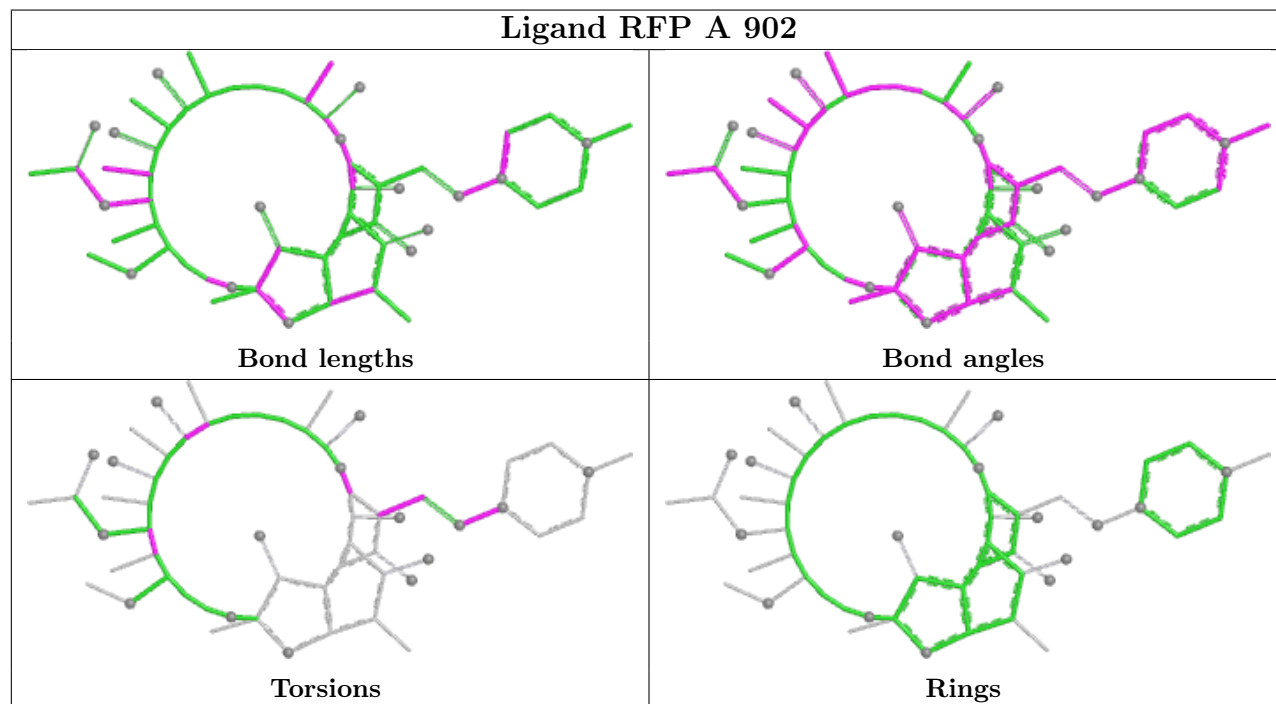
There are no ring outliers.

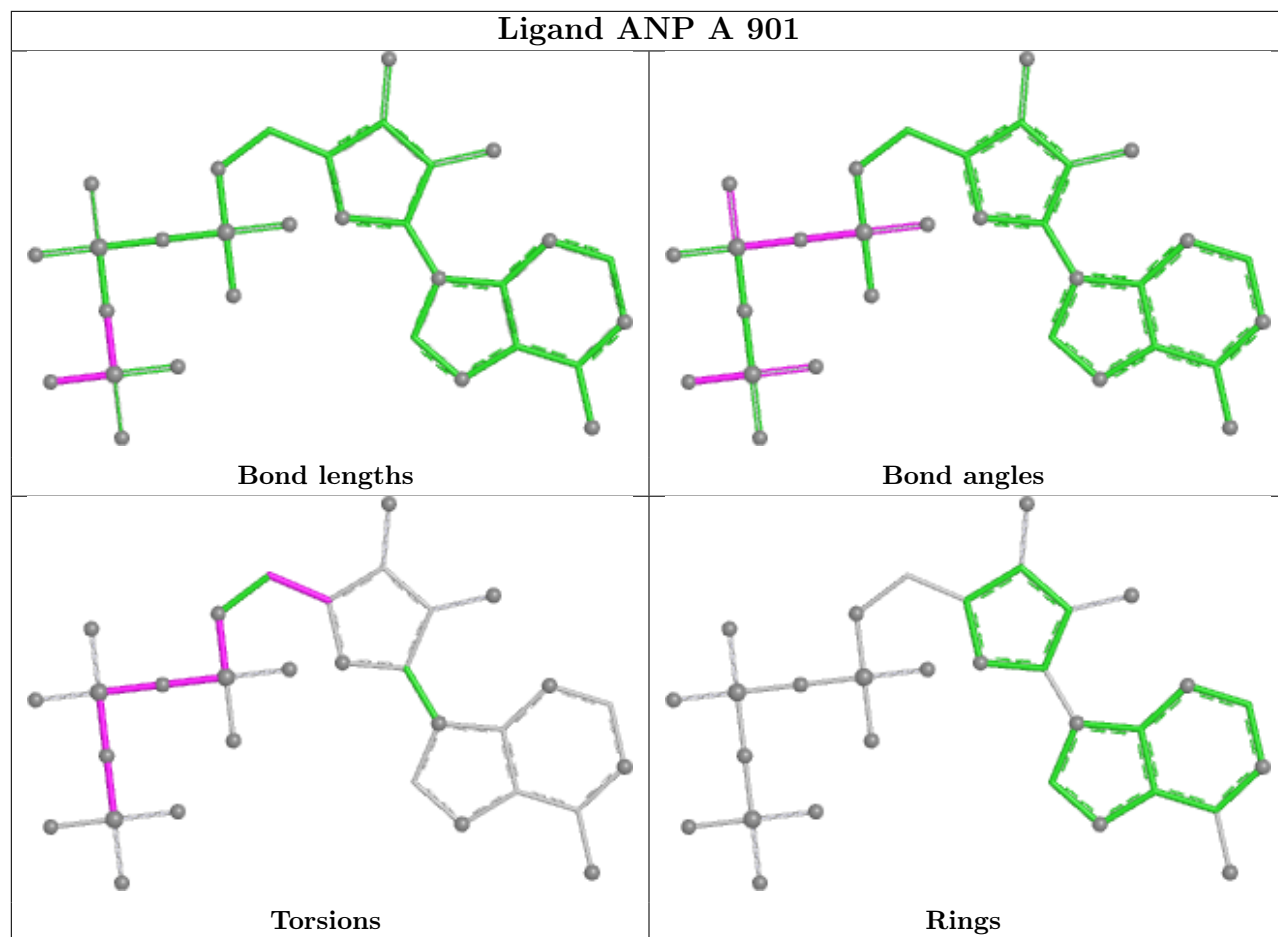
2 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	902	RFP	13	0
2	A	901	ANP	5	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	847/883 (95%)	0.49	125 (14%) 5 3	2, 27, 155, 168	0

All (125) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	48	TYR	6.4
1	A	176	VAL	6.2
1	A	52	LEU	5.5
1	A	507	GLU	5.4
1	A	53	ALA	5.0
1	A	42	CYS	4.9
1	A	62	LEU	4.9
1	A	96	ILE	4.8
1	A	214	GLY	4.8
1	A	68	LEU	4.7
1	A	219	LEU	4.6
1	A	58	PHE	4.6
1	A	129	SER	4.6
1	A	100	MET	4.3
1	A	75	ALA	4.2
1	A	16	GLU	4.2
1	A	76	ILE	4.2
1	A	65	LEU	4.1
1	A	303	GLY	4.1
1	A	279	LEU	4.0
1	A	134	HIS	4.0
1	A	223	LEU	3.9
1	A	97	ALA	3.8
1	A	308	VAL	3.7
1	A	178	LEU	3.7
1	A	220	VAL	3.6
1	A	6	LEU	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	180	VAL	3.5
1	A	26	LEU	3.4
1	A	130	PHE	3.4
1	A	19	VAL	3.4
1	A	246	LEU	3.4
1	A	67	SER	3.4
1	A	57	GLU	3.4
1	A	61	LEU	3.4
1	A	307	ILE	3.2
1	A	218	ALA	3.2
1	A	59	THR	3.2
1	A	34	GLY	3.1
1	A	126	PRO	3.1
1	A	186	ILE	3.1
1	A	222	GLY	3.1
1	A	173	HIS	3.1
1	A	18	LEU	3.0
1	A	171	PHE	3.0
1	A	215	LEU	3.0
1	A	306	TYR	3.0
1	A	147	LEU	3.0
1	A	80	SER	3.0
1	A	127	HIS	2.9
1	A	163	ILE	2.9
1	A	3	PRO	2.9
1	A	113	PRO	2.9
1	A	118	SER	2.9
1	A	177	GLN	2.9
1	A	247	ALA	2.9
1	A	128	ALA	2.8
1	A	71	SER	2.8
1	A	181	VAL	2.8
1	A	51	THR	2.8
1	A	32	ILE	2.8
1	A	60	GLN	2.7
1	A	409	PRO	2.7
1	A	842	GLU	2.7
1	A	167	ILE	2.7
1	A	25	ASN	2.7
1	A	83	ILE	2.7
1	A	117	ARG	2.7
1	A	94	SER	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	248	ILE	2.6
1	A	43	LEU	2.6
1	A	138	LEU	2.6
1	A	90	THR	2.6
1	A	224	VAL	2.6
1	A	125	LEU	2.6
1	A	154	TRP	2.6
1	A	13	PRO	2.6
1	A	288	ALA	2.6
1	A	164	ILE	2.5
1	A	185	MET	2.5
1	A	175	LYS	2.5
1	A	116	VAL	2.5
1	A	11	ILE	2.5
1	A	277	ILE	2.5
1	A	256	THR	2.4
1	A	225	SER	2.4
1	A	17	ALA	2.4
1	A	299	CYS	2.4
1	A	107	VAL	2.4
1	A	122	ALA	2.4
1	A	271	LEU	2.4
1	A	49	LYS	2.3
1	A	211	ALA	2.3
1	A	23	GLY	2.3
1	A	55	ASN	2.3
1	A	157	LEU	2.3
1	A	92	ILE	2.3
1	A	4	TYR	2.3
1	A	29	CYS	2.3
1	A	86	LEU	2.3
1	A	221	SER	2.2
1	A	153	CYS	2.2
1	A	2	LYS	2.2
1	A	187	SER	2.2
1	A	140	ILE	2.2
1	A	150	ILE	2.2
1	A	56	ASN	2.2
1	A	79	ILE	2.2
1	A	9	GLN	2.1
1	A	1	MET	2.1
1	A	8	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	172	ASP	2.1
1	A	179	ALA	2.1
1	A	216	GLY	2.1
1	A	276	ILE	2.1
1	A	174	ARG	2.1
1	A	406	GLY	2.1
1	A	259	ARG	2.1
1	A	110	TYR	2.1
1	A	141	ILE	2.0
1	A	112	MET	2.0
1	A	234	ASN	2.0
1	A	272	THR	2.0
1	A	280	ALA	2.0
1	A	22	LYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

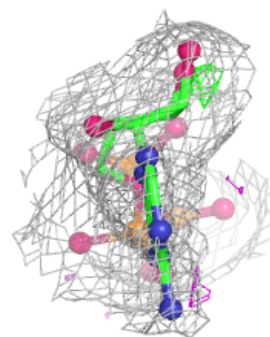
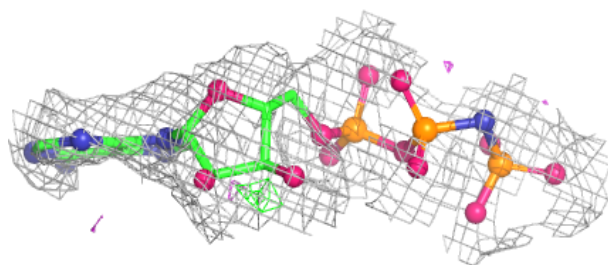
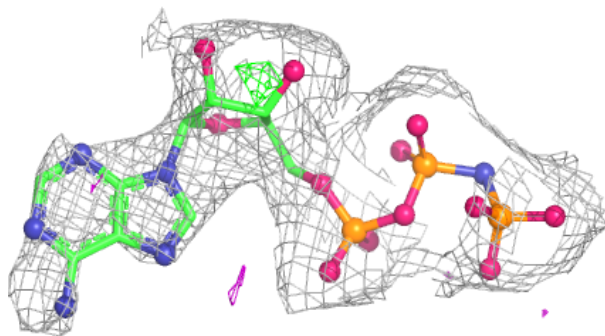
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MG	A	903	1/1	0.76	0.11	103,103,103,103	0
2	ANP	A	901	31/31	0.79	0.14	99,106,117,118	0
3	RFP	A	902	59/59	0.93	0.10	0,7,36,37	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.

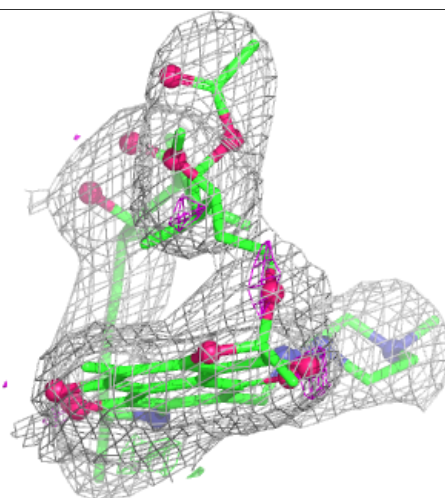
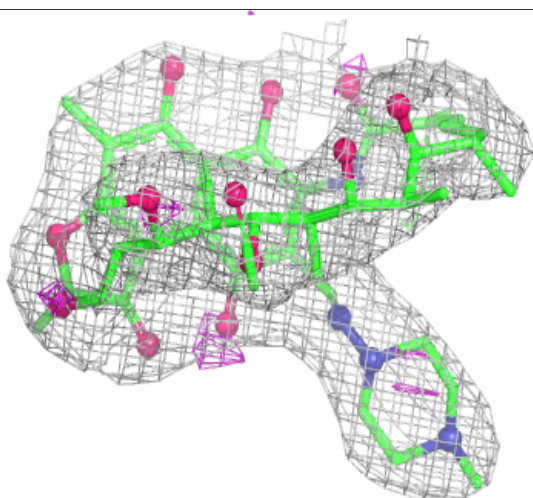
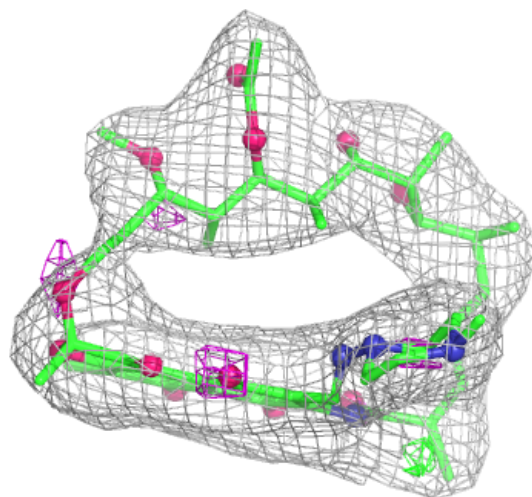
Electron density around ANP A 901:

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



Electron density around RFP A 902:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.