



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:50 PM UTC

PDB ID : 1HOO / pdb_00001hoo
Title : STRUCTURE OF GUANINE NUCLEOTIDE (GPPCP) COMPLEX OF ADENYLOSUCCINATE SYNTHETASE FROM E. COLI AT PH 6.5 AND 25 DEGREES CELSIUS
Authors : Poland, B.W.; Hou, Z.; Bruns, C.; Fromm, H.J.; Honzatko, R.B.
Deposited on : 1996-04-26
Resolution : 2.30 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

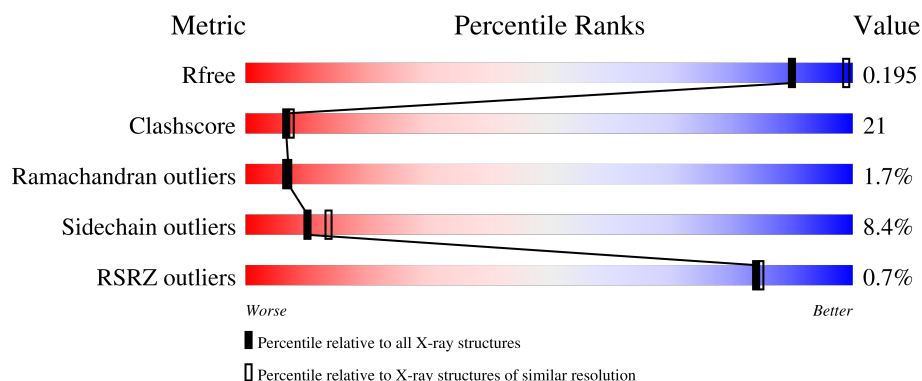
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	431	
1	B	431	

2 Entry composition [i](#)

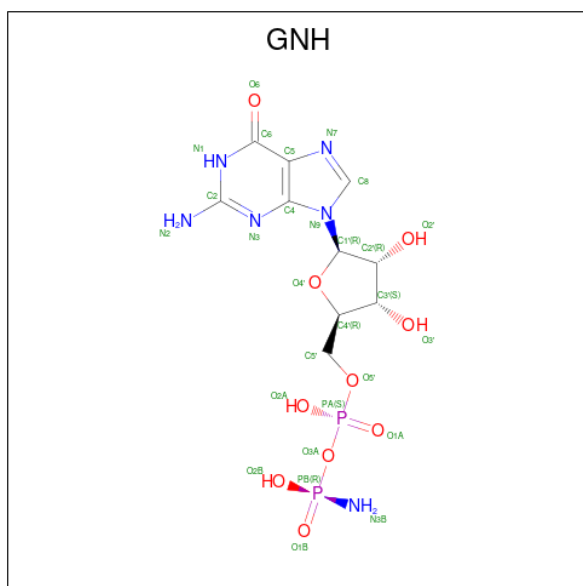
There are 4 unique types of molecules in this entry. The entry contains 9436 atoms, of which 2330 are hydrogens and 0 are deuteriums.

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 ADENYLOSUCCINATE SYNTHETASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	431	Total	C	H	N	O	S	0	0	0
			4069	2092	748	576	640	13			
1	B	431	Total	C	H	N	O	S	0	0	0
			4069	2092	748	576	640	13			

- Molecule 2 is AMINOPHOSPHONIC ACID-GUANYLATE ESTER (CCD ID: GNH) (formula: $C_{10}H_{16}N_6O_{10}P_2$).





Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	P	0	1
			44	10	12	6	13	3		

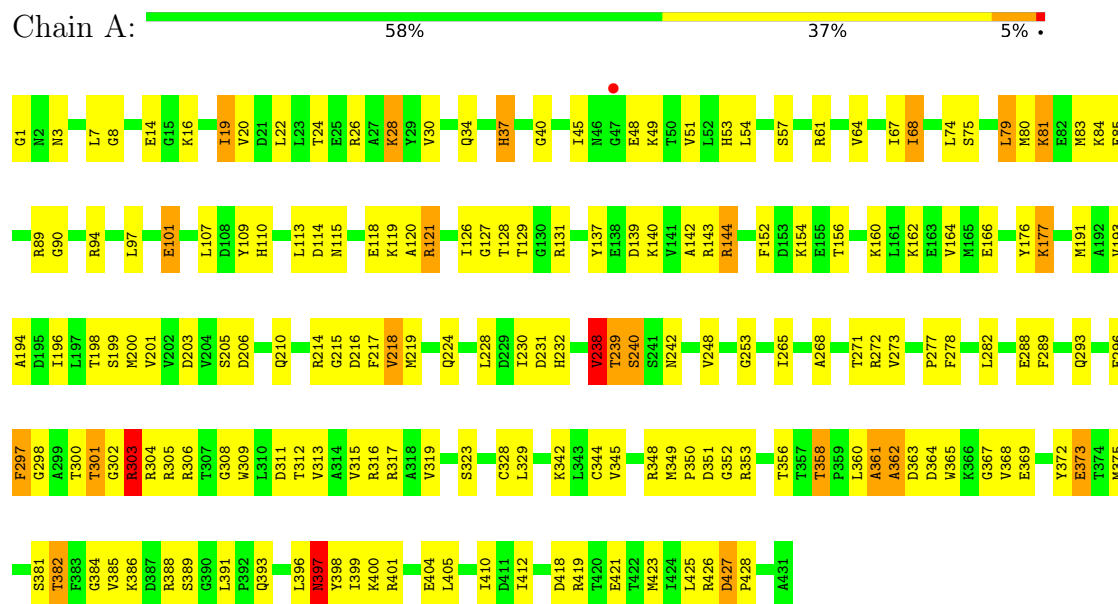
- Molecule 4 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	184	Total	H	O	0	0
			552	368	184		
4	B	220	Total	H	O	0	0
			660	440	220		

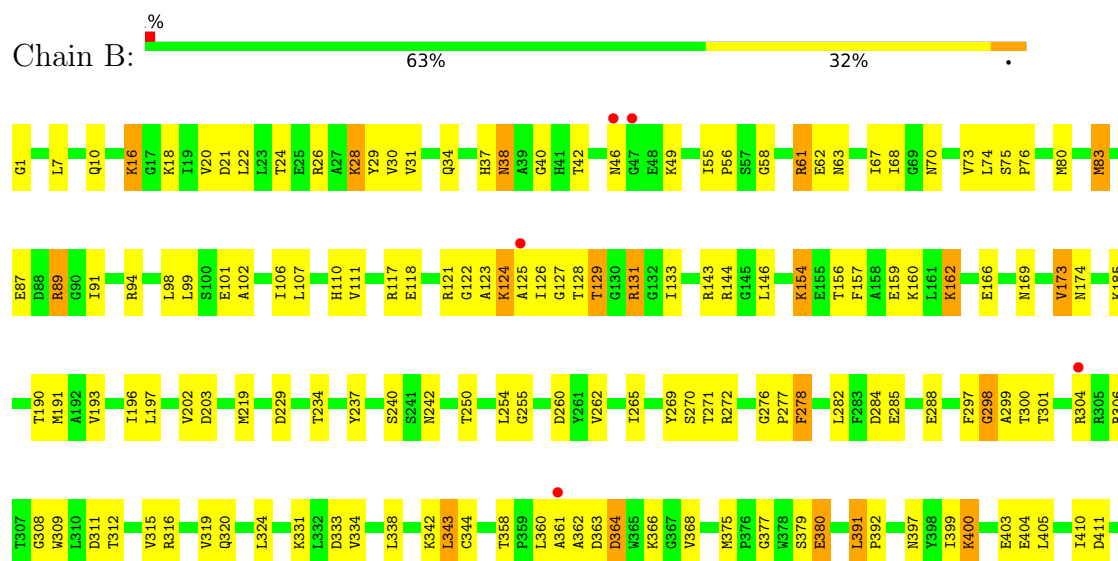
3 Residue-property plots

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: ADENYLOSUCCINATE SYNTHETASE



• Molecule 1: ADENYLOSUCCINATE SYNTHETASE



T415					
R419					
T420					
R426					
P427					
P428					
P429					
D430					
A431					

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.60Å 92.91Å 119.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	5.00 – 2.30 5.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	89.0 (5.00-2.30) 81.1 (5.00-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 2.29Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.185 , 0.210 0.186 , 0.195	Depositor DCC
R_{free} test set	2964 reflections (10.02%)	wwPDB-VP
Wilson B-factor (Å ²)	20.4	Xtriage
Anisotropy	0.280	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.18 , 50.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9436	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows:

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: GNH, GNP

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	0.94	3/3379 (0.1%)	1.18	13/4577 (0.3%)
1	B	0.64	0/3379	1.02	9/4577 (0.2%)
All	All	0.80	3/6758 (0.0%)	1.10	22/9154 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	399	ILE	CA-CB	8.09	1.63	1.54
1	A	248	VAL	CA-C	5.66	1.59	1.52
1	A	218	VAL	CA-CB	5.28	1.60	1.54

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	278	PHE	CA-C-N	7.99	128.42	119.32
1	A	278	PHE	C-N-CA	7.99	128.42	119.32
1	B	278	PHE	CA-C-N	7.31	127.65	119.32
1	B	278	PHE	C-N-CA	7.31	127.65	119.32
1	A	382	THR	N-CA-C	-6.58	105.79	113.88
1	A	398	TYR	N-CA-C	-6.35	104.28	111.07
1	A	131	ARG	N-CA-C	6.34	119.00	111.71
1	B	298	GLY	N-CA-C	6.31	120.30	112.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	358	THR	CA-C-N	6.23	126.25	119.90
1	B	358	THR	C-N-CA	6.23	126.25	119.90
1	A	215	GLY	N-CA-C	6.07	124.25	114.90
1	B	427	ASP	CA-C-N	5.97	127.30	119.84
1	B	427	ASP	C-N-CA	5.97	127.30	119.84
1	A	389	SER	N-CA-C	-5.94	105.12	112.90
1	A	319	VAL	CA-C-N	5.71	127.86	120.44
1	A	319	VAL	C-N-CA	5.71	127.86	120.44
1	B	74	LEU	N-CA-C	5.53	117.53	108.52
1	A	28	LYS	N-CA-C	-5.26	107.00	113.41
1	B	154	LYS	N-CA-C	5.25	116.69	111.07
1	A	67	ILE	CB-CA-C	-5.11	102.87	110.33
1	A	90	GLY	N-CA-C	5.07	122.89	114.48
1	A	238	VAL	N-CA-CB	-5.05	103.96	112.44

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	372	TYR	Sidechain

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	3321	748	3326	161	0
1	B	3321	748	3326	137	0
2	A	28	14	12	3	0
3	A	32	12	11	4	0
4	A	184	368	0	12	0
4	B	220	440	0	14	1
All	All	7106	2330	6675	287	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 21.

All (287) 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:418:ASP:HB3	1:A:421:GLU:HG3	1.43	1.01
1:A:20:VAL:HG13	1:A:219:MET:HE1	1.49	0.93
1:B:28:LYS:O	1:B:28:LYS:HD3	1.71	0.91
1:A:115:ASN:HA	1:A:129:THR:HA	1.52	0.88
1:A:118:GLU:HB2	1:A:128:THR:H	1.39	0.87
1:B:107:LEU:H	1:B:110:HIS:HD2	1.23	0.86
1:B:297:PHE:CD1	1:B:304:ARG:HG2	2.12	0.84
1:B:312:THR:HB	1:B:344:CYS:O	1.80	0.82
1:B:42:THR:HG21	1:B:49:LYS:HE2	1.63	0.81
1:A:118:GLU:HB3	1:A:129:THR:OG1	1.82	0.79
1:A:107:LEU:H	1:A:110:HIS:HD2	1.29	0.77
1:A:80:MET:O	1:A:84:LYS:HE2	1.85	0.77
1:A:301:THR:H	1:A:305:ARG:NH2	1.84	0.76
1:A:301:THR:H	1:A:305:ARG:HH22	1.34	0.76
1:B:1:GLY:HA3	1:B:260:ASP:HB3	1.68	0.75
1:A:203:ASP:HB2	1:B:360:LEU:HD13	1.71	0.73
1:B:28:LYS:HD3	1:B:28:LYS:C	2.13	0.72
1:B:297:PHE:HD1	1:B:304:ARG:HG2	1.54	0.70
1:A:54:LEU:HD21	1:A:74:LEU:HD12	1.75	0.69
1:B:20:VAL:HG13	1:B:219:MET:HE1	1.75	0.69
1:B:129:THR:O	1:B:131:ARG:HG2	1.92	0.68
1:A:94:ARG:HA	1:A:200:MET:CE	2.23	0.68
1:A:385:VAL:HG12	1:A:423:MET:HE2	1.75	0.67
1:B:107:LEU:H	1:B:110:HIS:CD2	2.11	0.67
1:A:301:THR:C	1:A:305:ARG:HH21	2.02	0.67
1:B:297:PHE:HE1	1:B:304:ARG:NE	1.93	0.67
1:B:68:ILE:HB	1:B:99:LEU:HD23	1.77	0.66
1:B:21:ASP:HB3	1:B:419:ARG:NH2	2.11	0.66
1:A:101:GLU:HG2	1:B:361:ALA:HA	1.77	0.65
1:A:16:LYS:HE3	3:A:432(B)[B]:GNP:O3G	1.97	0.65
1:B:143:ARG:HG2	1:B:143:ARG:HH11	1.59	0.65
1:A:26:ARG:O	1:A:217:PHE:HD2	1.80	0.65
1:B:118:GLU:HG2	1:B:127:GLY:O	1.97	0.65
1:A:311:ASP:O	1:A:315:VAL:HG23	1.96	0.65
1:B:154:LYS:HB3	1:B:191:MET:HE1	1.79	0.64
1:B:297:PHE:CE1	1:B:304:ARG:NE	2.66	0.64
1:B:117:ARG:HB3	1:B:121:ARG:HD2	1.78	0.63
1:B:129:THR:HA	1:B:131:ARG:NH2	2.13	0.63
1:A:144:ARG:O	1:A:164:VAL:HG21	1.98	0.63
1:B:129:THR:HA	1:B:131:ARG:CZ	2.28	0.63
1:B:360:LEU:HD22	4:B:433:HOH:O	1.97	0.63
1:A:400:LYS:O	1:A:404:GLU:HG3	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:317:ARG:HE	1:A:358:THR:HB	1.66	0.61
1:A:107:LEU:H	1:A:110:HIS:CD2	2.16	0.61
1:A:364:ASP:O	1:A:368:VAL:HG13	2.01	0.61
1:A:230:ILE:HD12	1:A:240:SER:O	2.01	0.60
1:B:24:THR:HG21	1:B:58:GLY:HA2	1.82	0.60
1:B:26:ARG:HH22	1:B:430:ASP:CG	2.08	0.60
1:A:120:ALA:O	1:A:121:ARG:C	2.44	0.60
1:A:94:ARG:HA	1:A:200:MET:HE3	1.84	0.60
1:B:400:LYS:HE2	1:B:404:GLU:OE2	2.02	0.59
1:A:19:ILE:HD12	1:A:19:ILE:N	2.17	0.59
1:A:206:ASP:O	1:A:210:GLN:HG3	2.02	0.59
1:B:21:ASP:HB3	1:B:419:ARG:HH21	1.68	0.59
1:B:430:ASP:O	1:B:431:ALA:HB2	2.03	0.59
1:A:121:ARG:HB3	4:A:530:HOH:O	2.02	0.59
1:B:237:TYR:CD2	1:B:276:GLY:HA2	2.38	0.59
1:A:397:ASN:HA	1:A:400:LYS:HE2	1.85	0.58
3:A:432(B)[B]:GNP:H8	3:A:432(B)[B]:GNP:H5'2	1.84	0.58
1:B:297:PHE:CE1	1:B:304:ARG:CZ	2.86	0.58
1:A:224:GLN:HE21	1:A:239:THR:HG21	1.68	0.58
1:A:375:MET:CE	1:A:401:ARG:HG3	2.33	0.58
1:B:399:ILE:HG23	1:B:410:ILE:HD12	1.86	0.58
1:A:289:PHE:CZ	1:A:293:GLN:HG3	2.37	0.58
1:B:397:ASN:O	1:B:400:LYS:HB3	2.03	0.58
1:A:19:ILE:N	1:A:19:ILE:CD1	2.66	0.57
1:A:296:GLU:HB3	1:A:305:ARG:HD2	1.86	0.57
1:B:419:ARG:HD3	4:B:532:HOH:O	2.04	0.57
1:B:254:LEU:HA	4:B:503:HOH:O	2.04	0.57
1:A:156:THR:HG22	1:A:160:LYS:HE3	1.86	0.57
1:A:177:LYS:HD3	1:A:177:LYS:N	2.19	0.57
1:A:193:VAL:HG12	1:A:196:ILE:HD12	1.87	0.57
1:B:297:PHE:CZ	1:B:304:ARG:NH2	2.73	0.57
1:A:80:MET:O	1:A:84:LYS:HG3	2.05	0.57
1:A:80:MET:C	1:A:84:LYS:HE2	2.30	0.56
1:A:205:SER:HB3	1:A:253:GLY:HA3	1.86	0.56
1:A:360:LEU:C	1:A:362:ALA:H	2.14	0.56
1:B:285:GLU:CD	1:B:285:GLU:H	2.13	0.56
1:B:426:ARG:O	1:B:427:ASP:HB3	2.05	0.56
1:B:31:VAL:HG22	1:B:67:ILE:HB	1.88	0.56
1:B:22:LEU:HD13	1:B:419:ARG:HB2	1.87	0.56
1:A:79:LEU:CD1	1:A:83:MET:HE1	2.36	0.56
1:A:118:GLU:HB2	1:A:128:THR:N	2.15	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:410:ILE:O	1:A:425:LEU:HD12	2.06	0.56
1:B:73:VAL:HG23	1:B:106:ILE:HD11	1.87	0.55
1:A:351:ASP:OD1	1:A:353:ARG:HG2	2.05	0.55
1:A:242:ASN:ND2	1:B:250:THR:HG23	2.21	0.55
1:B:18:LYS:HB3	4:B:465:HOH:O	2.05	0.55
1:B:117:ARG:HG2	1:B:121:ARG:NH1	2.21	0.55
1:A:37:HIS:H	1:A:37:HIS:CD2	2.24	0.55
1:A:238:VAL:HG12	1:B:143:ARG:NH1	2.22	0.55
1:A:391:LEU:HB2	1:A:396:LEU:HD21	1.87	0.55
1:A:152:PHE:CZ	1:B:361:ALA:HB1	2.42	0.55
1:A:418:ASP:HB3	1:A:421:GLU:CG	2.29	0.54
1:A:79:LEU:HD13	1:A:83:MET:HE1	1.89	0.54
1:B:362:ALA:O	1:B:363:ASP:C	2.49	0.54
1:A:1:GLY:HA2	4:A:529:HOH:O	2.08	0.54
1:A:79:LEU:HD13	1:A:83:MET:CE	2.37	0.54
1:A:345:VAL:HG12	1:A:405:LEU:HD13	1.89	0.53
1:A:384:GLY:O	1:A:386:LYS:HG3	2.08	0.53
1:A:393:GLN:HG2	4:A:598:HOH:O	2.08	0.53
2:A:432(A)[A]:GNH:H5'2	2:A:432(A)[A]:GNH:H8	1.89	0.53
1:B:146:LEU:HD21	1:B:160:LYS:HB3	1.90	0.53
1:A:129:THR:HG22	1:A:129:THR:O	2.08	0.53
1:A:22:LEU:HD13	1:A:419:ARG:HG3	1.90	0.53
1:B:297:PHE:CE1	1:B:304:ARG:HG2	2.43	0.53
1:A:426:ARG:HA	4:A:455:HOH:O	2.08	0.53
1:A:277:PRO:HB3	1:A:365:TRP:HB3	1.90	0.53
1:A:30:VAL:HG13	1:A:219:MET:HE2	1.90	0.53
1:B:124:LYS:NZ	1:B:124:LYS:HB3	2.24	0.53
1:A:81:LYS:HA	1:A:84:LYS:HE2	1.92	0.52
1:A:426:ARG:O	1:A:427:ASP:CB	2.57	0.52
1:A:268:ALA:HA	1:A:329:LEU:HD11	1.90	0.52
1:A:351:ASP:CG	1:A:352:GLY:N	2.68	0.52
1:B:62:GLU:HG2	1:B:63:ASN:N	2.24	0.52
1:B:111:VAL:HG22	1:B:131:ARG:O	2.10	0.52
1:B:343:LEU:O	1:B:343:LEU:HD12	2.10	0.52
1:A:85:GLU:O	1:A:89:ARG:HG2	2.09	0.52
1:A:16:LYS:HE3	3:A:432(B)[B]:GNP:PG	2.50	0.51
1:A:289:PHE:CE2	1:A:293:GLN:HG3	2.44	0.51
1:B:312:THR:HG22	1:B:316:ARG:HH21	1.75	0.51
1:B:73:VAL:HG21	1:B:133:ILE:HG23	1.93	0.51
1:B:24:THR:HB	1:B:61:ARG:HH11	1.75	0.51
1:A:24:THR:HG22	1:A:61:ARG:HD3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:ARG:NH2	1:B:240:SER:HB3	2.25	0.51
1:A:214:ARG:HD2	1:A:216:ASP:OD2	2.11	0.50
1:B:338:LEU:O	1:B:377:GLY:HA3	2.10	0.50
1:A:19:ILE:CD1	1:A:19:ILE:H	2.24	0.50
1:B:331:LYS:HD3	1:B:334:VAL:HG21	1.93	0.50
1:A:375:MET:HE3	1:A:401:ARG:HG3	1.92	0.50
1:A:297:PHE:O	1:A:298:GLY:C	2.54	0.50
1:A:119:LYS:HG2	1:A:129:THR:CG2	2.42	0.50
1:A:115:ASN:O	1:A:119:LYS:HB2	2.12	0.49
1:A:110:HIS:HE1	1:A:139:ASP:OD2	1.95	0.49
1:A:201:VAL:HG11	1:B:361:ALA:HB2	1.94	0.49
1:B:76:PRO:O	1:B:80:MET:HG2	2.12	0.49
1:A:137:TYR:O	1:A:140:LYS:HB3	2.13	0.49
1:A:154:LYS:CD	1:A:191:MET:HE1	2.43	0.49
1:B:271:THR:HA	1:B:306:ARG:O	2.13	0.49
1:B:380:GLU:HG3	1:B:392:PRO:HB3	1.94	0.49
1:B:411:ASP:CG	1:B:426:ARG:HB3	2.37	0.49
1:A:323:SER:HB3	1:B:255:GLY:HA3	1.95	0.49
1:A:385:VAL:HG12	1:A:423:MET:CE	2.42	0.49
1:B:16:LYS:HE2	4:B:508:HOH:O	2.12	0.49
1:A:194:ALA:HB3	4:A:456:HOH:O	2.11	0.49
3:A:432(B)[B]:GNP:H5'2	3:A:432(B)[B]:GNP:C8	2.42	0.49
1:B:73:VAL:CG2	1:B:106:ILE:HD11	2.42	0.49
1:B:193:VAL:HG12	1:B:193:VAL:O	2.13	0.49
1:B:91:ILE:HD12	1:B:91:ILE:H	1.78	0.49
1:B:30:VAL:HG11	4:B:524:HOH:O	2.14	0.48
1:A:316:ARG:NH1	1:A:356:THR:HB	2.28	0.48
1:B:262:VAL:HG12	1:B:324:LEU:HD23	1.95	0.48
1:A:304:ARG:HG3	1:A:304:ARG:O	2.14	0.48
1:B:156:THR:O	1:B:159:GLU:HG2	2.13	0.48
1:B:70:ASN:HB3	1:B:102:ALA:HB3	1.95	0.48
1:A:89:ARG:HA	4:A:556:HOH:O	2.12	0.48
1:A:176:TYR:HD1	4:A:548:HOH:O	1.96	0.48
1:A:8:GLY:HA3	1:A:14:GLU:CD	2.39	0.48
1:A:426:ARG:O	1:A:427:ASP:HB3	2.13	0.48
1:A:348:ARG:HG2	1:A:348:ARG:HH11	1.79	0.48
1:B:159:GLU:HG3	1:B:160:LYS:N	2.29	0.48
1:B:298:GLY:O	1:B:299:ALA:HB3	2.13	0.48
1:B:272:ARG:HD3	1:B:306:ARG:NH2	2.29	0.47
1:B:83:MET:O	1:B:87:GLU:HG3	2.14	0.47
1:A:154:LYS:HE3	1:A:191:MET:HE1	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:GLY:O	1:B:429:PHE:HA	2.14	0.47
1:B:202:VAL:HB	4:B:576:HOH:O	2.14	0.47
1:B:426:ARG:O	1:B:427:ASP:CB	2.63	0.47
1:A:232:HIS:HD2	4:B:442:HOH:O	1.97	0.46
1:B:360:LEU:HG	1:B:361:ALA:N	2.27	0.46
1:A:61:ARG:HB3	1:A:64:VAL:HG23	1.96	0.46
1:A:309:TRP:NE1	1:A:342:LYS:HD2	2.30	0.46
1:B:80:MET:SD	1:B:193:VAL:HG11	2.55	0.46
1:A:228:LEU:HB3	1:A:239:THR:OG1	2.15	0.46
1:B:282:LEU:HG	1:B:308:GLY:HA2	1.98	0.46
1:A:37:HIS:CD2	1:A:37:HIS:N	2.84	0.46
1:A:68:ILE:HD12	1:A:97:LEU:HD11	1.98	0.46
1:A:388:ARG:HD3	4:A:555:HOH:O	2.16	0.46
1:A:40:GLY:O	2:A:432(A)[A]:GNH:N3B	2.49	0.45
1:A:350:PRO:HG3	1:A:364:ASP:CB	2.46	0.45
1:B:185:LYS:HD3	4:B:631:HOH:O	2.16	0.45
1:A:51:VAL:HG12	1:A:53:HIS:CE1	2.51	0.45
1:A:303:ARG:HG2	1:A:304:ARG:N	2.32	0.45
1:B:311:ASP:O	1:B:315:VAL:HG23	2.15	0.45
1:A:48:GLU:HG2	1:A:49:LYS:N	2.30	0.45
1:A:375:MET:HE1	1:A:401:ARG:HG3	1.97	0.45
1:A:391:LEU:HD12	1:A:396:LEU:HD21	1.97	0.45
1:B:309:TRP:CD2	1:B:342:LYS:HB2	2.52	0.45
1:A:365:TRP:O	1:A:368:VAL:HG22	2.16	0.45
2:A:432(A)[A]:GNH:H5'2	2:A:432(A)[A]:GNH:C8	2.46	0.45
1:B:169:ASN:O	1:B:173:VAL:HG13	2.16	0.45
1:A:271:THR:HA	1:A:306:ARG:O	2.17	0.45
1:B:282:LEU:HD21	1:B:309:TRP:NE1	2.32	0.45
1:A:162:LYS:O	1:A:166:GLU:HG3	2.17	0.45
1:A:119:LYS:HG2	1:A:129:THR:HG21	1.99	0.45
1:A:206:ASP:HA	1:B:320:GLN:HE22	1.82	0.44
1:A:154:LYS:HE3	1:A:191:MET:CE	2.47	0.44
1:A:110:HIS:CE1	1:A:139:ASP:OD2	2.71	0.44
1:A:152:PHE:CD2	1:B:362:ALA:HB2	2.52	0.44
1:A:364:ASP:O	1:A:367:GLY:N	2.46	0.44
1:B:107:LEU:O	1:B:110:HIS:HB2	2.17	0.44
1:A:81:LYS:HE2	1:A:84:LYS:HZ1	1.82	0.44
1:A:101:GLU:HG2	1:B:361:ALA:CA	2.48	0.44
1:A:152:PHE:CE1	1:A:198:THR:HB	2.53	0.44
1:B:229:ASP:OD1	1:B:242:ASN:HA	2.17	0.44
1:A:142:ALA:O	1:A:143:ARG:HB2	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:143:ARG:HG2	1:B:143:ARG:NH1	2.31	0.43
1:B:300:THR:O	1:B:301:THR:C	2.61	0.43
1:B:312:THR:HG22	1:B:316:ARG:NH2	2.33	0.43
1:A:16:LYS:CB	1:A:265:ILE:HG13	2.47	0.43
1:A:152:PHE:CD2	1:B:362:ALA:CB	3.02	0.43
1:A:311:ASP:HA	1:A:344:CYS:HB3	2.00	0.43
1:A:313:VAL:HG13	1:A:356:THR:O	2.19	0.43
1:B:26:ARG:NH2	4:B:590:HOH:O	2.51	0.43
1:B:28:LYS:HD2	1:B:29:TYR:CZ	2.53	0.43
1:A:232:HIS:HE1	1:B:250:THR:O	2.01	0.43
1:A:109:TYR:O	1:A:113:LEU:HD23	2.19	0.43
1:A:361:ALA:HB2	4:A:586:HOH:O	2.17	0.43
1:B:1:GLY:HA3	1:B:260:ASP:CB	2.44	0.43
1:B:16:LYS:HB2	1:B:265:ILE:HG13	1.99	0.43
1:B:68:ILE:HD12	1:B:99:LEU:HD21	2.01	0.43
1:B:98:LEU:HD22	1:B:202:VAL:HB	2.00	0.43
1:A:363:ASP:OD1	1:A:364:ASP:N	2.50	0.43
1:B:55:ILE:HA	1:B:56:PRO:HD3	1.83	0.43
1:B:360:LEU:HD23	1:B:360:LEU:H	1.84	0.43
1:A:348:ARG:HG2	1:A:369:GLU:OE2	2.19	0.43
1:B:28:LYS:HE2	4:B:571:HOH:O	2.17	0.43
1:B:162:LYS:HE3	4:B:455:HOH:O	2.19	0.43
1:B:430:ASP:O	1:B:431:ALA:CB	2.65	0.43
1:A:144:ARG:HH11	1:A:144:ARG:HD3	1.70	0.43
1:A:119:LYS:HA	1:A:119:LYS:HD3	1.64	0.43
1:A:37:HIS:H	1:A:37:HIS:HD2	1.67	0.43
1:A:201:VAL:CG1	1:B:361:ALA:HB2	2.49	0.43
1:B:400:LYS:O	1:B:404:GLU:HG3	2.19	0.43
1:A:81:LYS:NZ	1:A:84:LYS:NZ	2.67	0.42
1:A:101:GLU:C	1:A:101:GLU:CD	2.87	0.42
1:A:1:GLY:N	1:A:428:PRO:O	2.52	0.42
1:B:38:ASN:C	1:B:38:ASN:HD22	2.27	0.42
1:B:173:VAL:CG2	1:B:174:ASN:N	2.82	0.42
1:A:61:ARG:HB3	1:A:64:VAL:CG2	2.49	0.42
1:B:117:ARG:O	1:B:121:ARG:HB2	2.20	0.42
1:B:202:VAL:HG22	1:B:203:ASP:N	2.34	0.42
1:A:3:ASN:OD1	1:A:218:VAL:HB	2.19	0.42
1:B:91:ILE:HD12	1:B:91:ILE:N	2.33	0.42
1:B:272:ARG:NH1	1:B:277:PRO:O	2.51	0.42
1:A:349:MET:O	1:A:352:GLY:N	2.52	0.42
1:A:373:GLU:OE2	1:A:401:ARG:NH1	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:70:ASN:CB	1:B:102:ALA:HB3	2.49	0.42
1:A:101:GLU:HG3	4:A:484:HOH:O	2.20	0.42
1:B:125:ALA:O	1:B:127:GLY:N	2.53	0.42
1:B:399:ILE:O	1:B:403:GLU:HG3	2.19	0.42
1:B:391:LEU:HA	1:B:392:PRO:HD2	1.96	0.42
1:A:302:GLY:O	1:A:303:ARG:HB3	2.19	0.42
1:A:397:ASN:ND2	1:A:400:LYS:NZ	2.68	0.42
1:B:101:GLU:HG2	4:B:485:HOH:O	2.20	0.42
1:B:269:TYR:CD1	1:B:269:TYR:N	2.88	0.42
1:B:319:VAL:HG22	1:B:324:LEU:HD12	2.02	0.42
1:A:16:LYS:HB2	1:A:265:ILE:HG13	2.02	0.41
1:A:54:LEU:HD21	1:A:74:LEU:CD1	2.47	0.41
1:B:128:THR:O	1:B:128:THR:HG23	2.20	0.41
1:B:312:THR:HG21	1:B:405:LEU:HB3	2.02	0.41
1:A:282:LEU:HG	1:A:308:GLY:HA2	2.02	0.41
1:A:273:VAL:HG22	4:A:505:HOH:O	2.20	0.41
1:A:297:PHE:O	1:A:297:PHE:HD1	2.03	0.41
1:A:201:VAL:HG11	1:B:361:ALA:CB	2.50	0.41
1:B:37:HIS:CD2	1:B:37:HIS:H	2.38	0.41
1:B:89:ARG:O	1:B:91:ILE:HD12	2.20	0.41
1:B:157:PHE:CZ	1:B:190:THR:HG21	2.56	0.41
1:B:270:SER:HB2	1:B:278:PHE:CZ	2.55	0.41
1:A:419:ARG:HH11	1:A:419:ARG:HG2	1.86	0.41
1:B:193:VAL:CG1	1:B:196:ILE:HB	2.51	0.41
1:A:301:THR:O	1:A:305:ARG:NH2	2.52	0.41
1:B:284:ASP:O	1:B:288:GLU:HG3	2.21	0.41
1:A:328:CYS:HA	1:A:412:ILE:O	2.20	0.41
1:B:122:GLY:O	1:B:123:ALA:HB2	2.20	0.41
1:B:300:THR:O	1:B:300:THR:HG23	2.21	0.41
1:A:238:VAL:HG13	1:A:239:THR:O	2.21	0.41
1:A:312:THR:OG1	1:A:316:ARG:NH2	2.54	0.41
1:A:348:ARG:HB3	1:A:369:GLU:CG	2.51	0.40
1:B:10:GLN:HG2	4:B:491:HOH:O	2.21	0.40
1:A:277:PRO:HG3	4:A:442:HOH:O	2.20	0.40
1:B:427:ASP:HA	1:B:428:PRO:HD3	1.84	0.40
1:A:24:THR:O	1:A:24:THR:CG2	2.70	0.40
1:A:156:THR:CG2	1:A:160:LYS:HE3	2.51	0.40
1:A:230:ILE:HG23	1:A:231:ASP:N	2.36	0.40
1:A:230:ILE:HG23	1:A:231:ASP:OD1	2.22	0.40
1:A:360:LEU:C	1:A:362:ALA:N	2.80	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:518:HOH:O	4:B:592:HOH:H2[4_455]	1.57	0.03

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	429/431 (100%)	383 (89%)	37 (9%)	9 (2%)	5	4
1	B	429/431 (100%)	389 (91%)	34 (8%)	6 (1%)	9	9
All	All	858/862 (100%)	772 (90%)	71 (8%)	15 (2%)	7	7

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	303	ARG
1	A	361	ALA
1	A	362	ALA
1	B	364	ASP
1	A	397	ASN
1	A	121	ARG
1	A	127	GLY
1	A	427	ASP
1	B	427	ASP
1	B	94	ARG
1	B	131	ARG
1	A	301	THR
1	B	40	GLY
1	B	126	ILE
1	A	45	ILE

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	353/353 (100%)	324 (92%)	29 (8%)	10	14
1	B	353/353 (100%)	323 (92%)	30 (8%)	10	13
All	All	706/706 (100%)	647 (92%)	59 (8%)	10	14

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	A	19	ILE
1	A	28	LYS
1	A	34	GLN
1	A	37	HIS
1	A	57	SER
1	A	68	ILE
1	A	75	SER
1	A	79	LEU
1	A	81	LYS
1	A	101	GLU
1	A	114	ASP
1	A	126	ILE
1	A	144	ARG
1	A	177	LYS
1	A	199	SER
1	A	238	VAL
1	A	239	THR
1	A	240	SER
1	A	272	ARG
1	A	288	GLU
1	A	297	PHE
1	A	300	THR
1	A	303	ARG
1	A	358	THR
1	A	373	GLU
1	A	381	SER

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Mol	Chain	Res	Type
1	A	382	THR
1	A	397	ASN
1	B	7	LEU
1	B	16	LYS
1	B	28	LYS
1	B	34	GLN
1	B	38	ASN
1	B	46	ASN
1	B	61	ARG
1	B	75	SER
1	B	83	MET
1	B	89	ARG
1	B	124	LYS
1	B	129	THR
1	B	144	ARG
1	B	162	LYS
1	B	166	GLU
1	B	173	VAL
1	B	197	LEU
1	B	234	THR
1	B	333	ASP
1	B	343	LEU
1	B	364	ASP
1	B	366	LYS
1	B	368	VAL
1	B	375	MET
1	B	379	SER
1	B	380	GLU
1	B	391	LEU
1	B	400	LYS
1	B	415	THR
1	B	420	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	HIS
1	A	110	HIS
1	A	210	GLN
1	A	242	ASN
1	A	397	ASN
1	B	2	ASN

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Mol	Chain	Res	Type
1	B	37	HIS
1	B	38	ASN
1	B	110	HIS
1	B	393	GLN
1	B	397	ASN

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](#)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	GNP	A	432(B)[B]	-	34,34,34	1.66	9 (26%)	47,54,54	1.77	9 (19%)
2	GNH	A	432(A)[A]	-	29,30,30	1.32	4 (13%)	42,47,47	1.52	7 (16%)

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	GNP	A	432(B)[B]	-	-	7/18/38/38	0/3/3/3
2	GNH	A	432(A)[A]	-	-	2/13/32/32	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	432(B)[B]	GNP	PA-O3A	3.63	1.63	1.59
3	A	432(B)[B]	GNP	PB-O2B	-3.46	1.47	1.56
2	A	432(A)[A]	GNH	PB-O2B	-3.43	1.47	1.56
3	A	432(B)[B]	GNP	PG-O2G	-3.04	1.48	1.56
3	A	432(B)[B]	GNP	PB-O3A	2.98	1.62	1.59
2	A	432(A)[A]	GNH	PB-O1B	2.89	1.50	1.46
3	A	432(B)[B]	GNP	C1'-N9	-2.75	1.39	1.47
2	A	432(A)[A]	GNH	C1'-N9	-2.73	1.39	1.47
3	A	432(B)[B]	GNP	PG-O1G	2.61	1.50	1.46
3	A	432(B)[B]	GNP	PG-O3G	-2.52	1.50	1.56
3	A	432(B)[B]	GNP	PG-N3B	2.47	1.69	1.63
2	A	432(A)[A]	GNH	PA-O3A	2.18	1.61	1.59
3	A	432(B)[B]	GNP	PB-N3B	2.17	1.69	1.63

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	432(B)[B]	GNP	O1G-PG-N3B	-6.28	102.52	111.77
3	A	432(B)[B]	GNP	O2B-PB-O1B	4.06	118.57	109.87
2	A	432(A)[A]	GNH	O2B-PB-O1B	3.63	119.65	110.26
3	A	432(B)[B]	GNP	O3A-PB-N3B	-3.61	96.58	106.59
3	A	432(B)[B]	GNP	N1-C2-N3	2.89	128.61	123.32
2	A	432(A)[A]	GNH	PB-O3A-PA	-2.87	122.92	132.10
2	A	432(A)[A]	GNH	N1-C2-N3	2.74	128.33	123.32
3	A	432(B)[B]	GNP	O2G-PG-O3G	2.54	114.41	107.59
3	A	432(B)[B]	GNP	N9-C8-N7	2.47	117.99	113.40
2	A	432(A)[A]	GNH	N9-C8-N7	2.47	117.98	113.40
2	A	432(A)[A]	GNH	O3'-C3'-C2'	2.40	119.52	111.82
2	A	432(A)[A]	GNH	C1'-N9-C8	2.22	133.04	126.73
2	A	432(A)[A]	GNH	C2-N3-C4	-2.10	108.68	112.30
3	A	432(B)[B]	GNP	C2-N3-C4	-2.05	108.77	112.30
3	A	432(B)[B]	GNP	C1'-N9-C8	2.05	132.55	126.73
3	A	432(B)[B]	GNP	O3'-C3'-C2'	2.00	118.23	111.82

There are no chirality outliers.

All (9) torsion outliers are listed below:

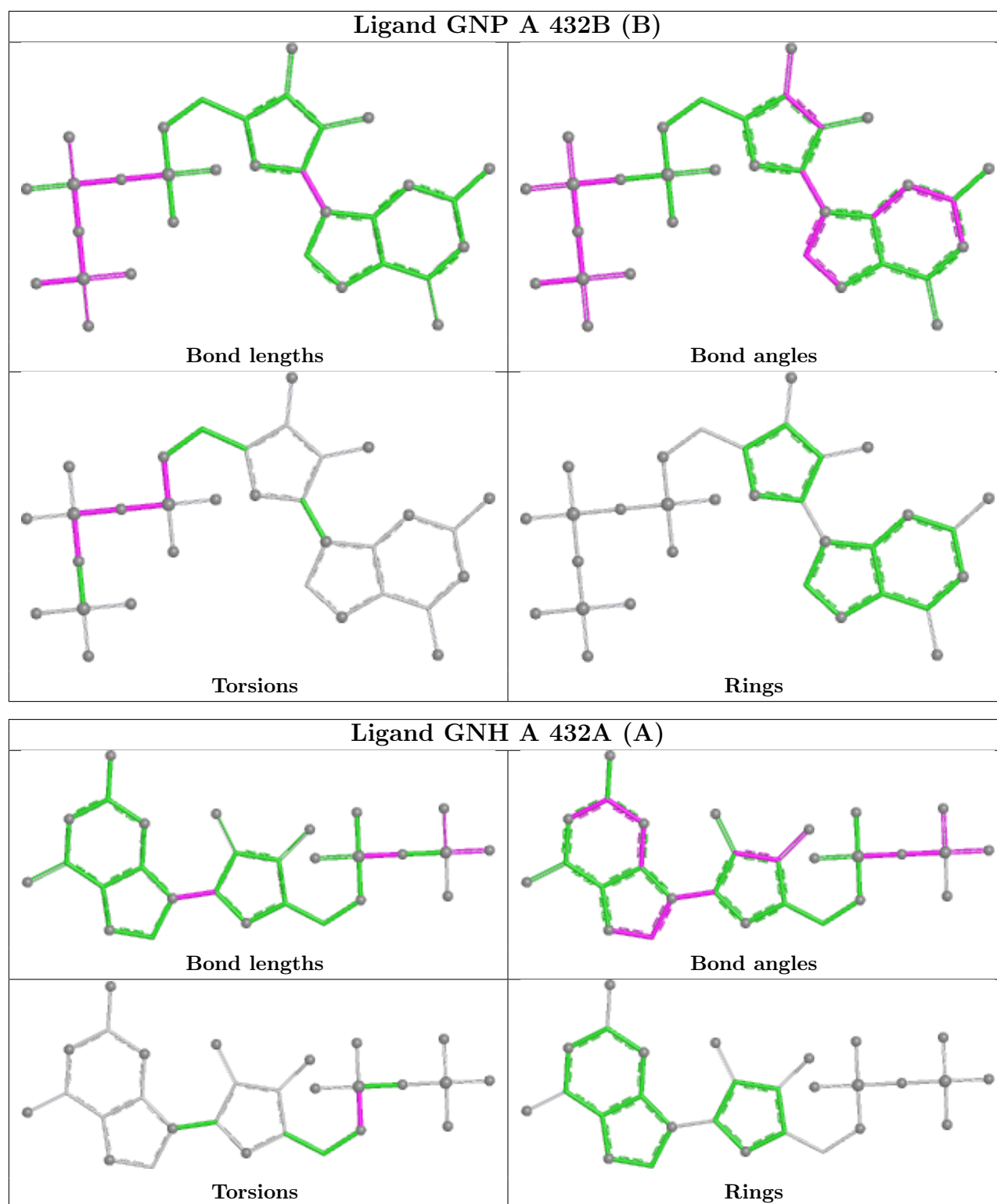
Mol	Chain	Res	Type	Atoms
2	A	432(A)[A]	GNH	C5'-O5'-PA-O3A
3	A	432(B)[B]	GNP	PG-N3B-PB-O1B
3	A	432(B)[B]	GNP	PG-N3B-PB-O3A
3	A	432(B)[B]	GNP	PB-O3A-PA-O5'
3	A	432(B)[B]	GNP	C5'-O5'-PA-O3A
3	A	432(B)[B]	GNP	C5'-O5'-PA-O1A
2	A	432(A)[A]	GNH	C5'-O5'-PA-O1A
3	A	432(B)[B]	GNP	C5'-O5'-PA-O2A
3	A	432(B)[B]	GNP	PA-O3A-PB-O2B

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	432(B)[B]	GNP	4	0
2	A	432(A)[A]	GNH	3	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	431/431 (100%)	-0.51	1 (0%) 91 91	7, 22, 46, 70	0
1	B	431/431 (100%)	-0.54	5 (1%) 76 77	8, 18, 46, 68	0
All	All	862/862 (100%)	-0.52	6 (0%) 84 85	7, 20, 46, 70	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	304	ARG	3.1
1	B	361	ALA	2.6
1	A	47	GLY	2.3
1	B	46	ASN	2.2
1	B	47	GLY	2.2
1	B	125	ALA	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](#)

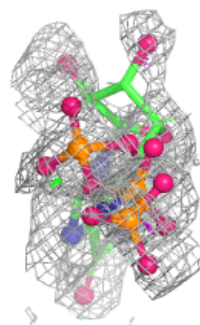
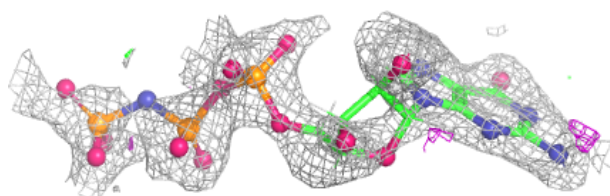
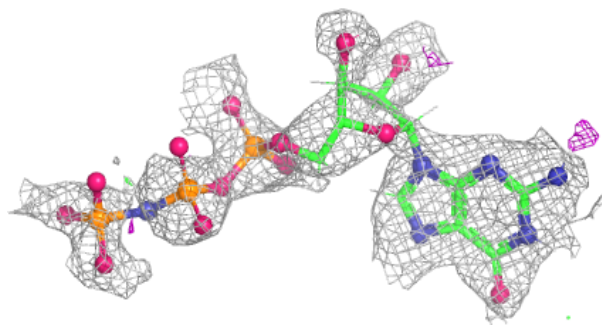
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GNP	A	432(B)[B]	32/32	0.86	0.09	0,31,39,41	44
2	GNH	A	432(A)[A]	28/28	0.87	0.08	0,33,43,44	42

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.

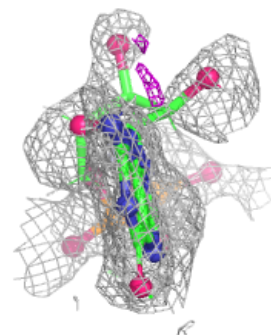
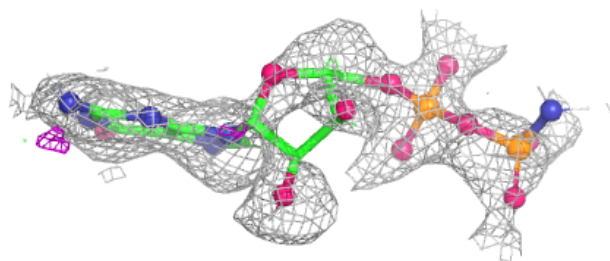
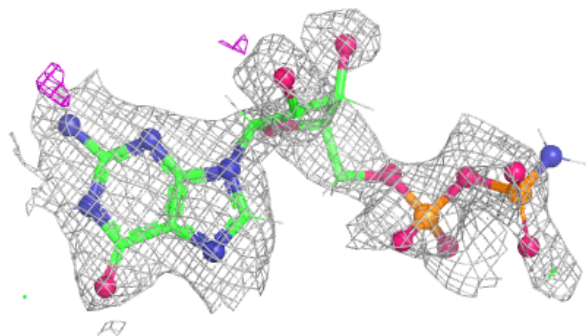
Electron density around GNP A 432B (B):

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 GNH A 432A (A):

$2mF_o-DF_c$ (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.