



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 09:17 AM UTC

PDB ID : 1IHI / pdb_00001ihi
Title : Crystal Structure of Human Type III 3-alpha-Hydroxysteroid Dehydrogenase/Bile Acid Binding Protein (AKR1C2) Complexed with NADP+ and Ursodeoxycholate
Authors : Jin, Y.; Stayrook, S.E.; Albert, R.H.; Palackal, N.T.; Penning, T.M.; Lewis, M.
Deposited on : 2001-04-19
Resolution : 3.00 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](#) ①) 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

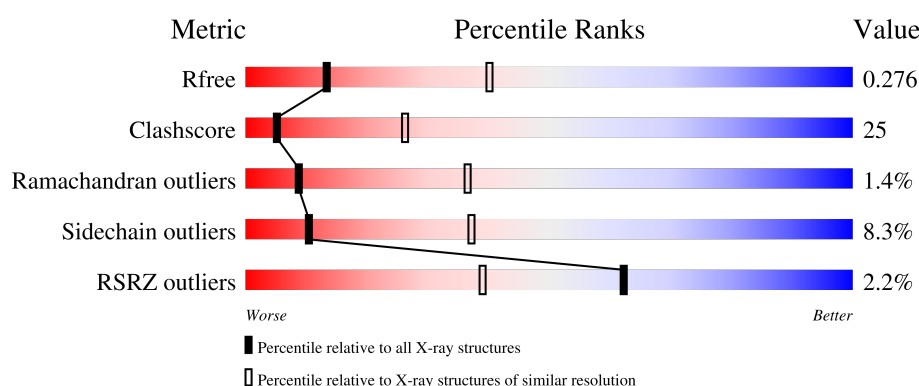
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	323	2% 49% 41% 9%
1	B	323	3% 53% 37% 8%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	IU5	A	326	X	-	-	-
3	IU5	B	327	X	-	-	-

2 Entry composition ⓘ

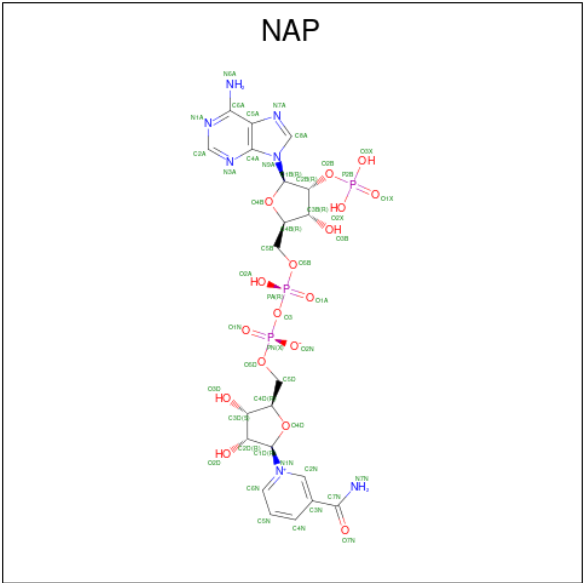
There are 3 unique types of molecules in this entry. The entry contains 5272 atoms, of which 0 are hydrogens and 0 are deuteriums.

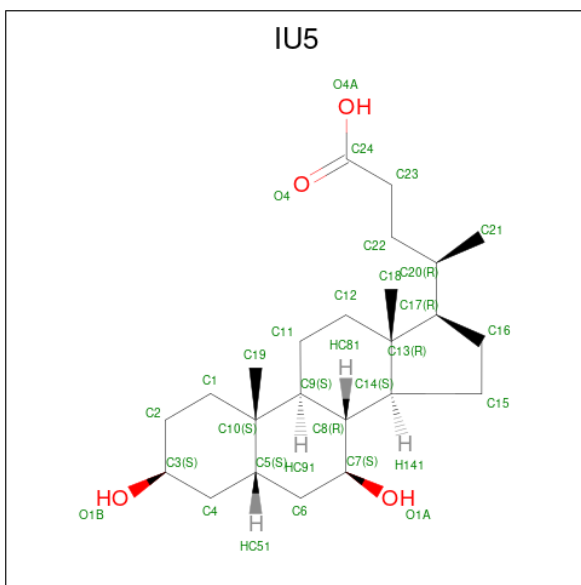
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called 3-ALPHA-HYDROXYSTEROID DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	320	Total	C	N	O	S	0	0	0
			2560	1640	442	467	11			
1	B	320	Total	C	N	O	S	0	0	0
			2560	1640	442	467	11			

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



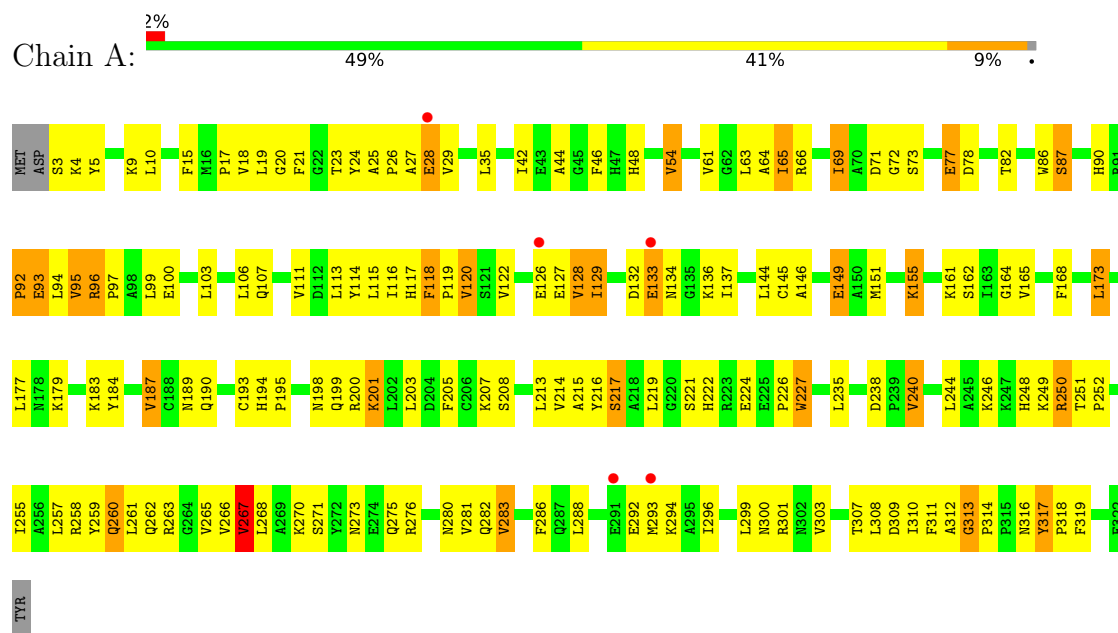


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 28 24 4	0	0
3	B	1	Total C O 28 24 4	0	0

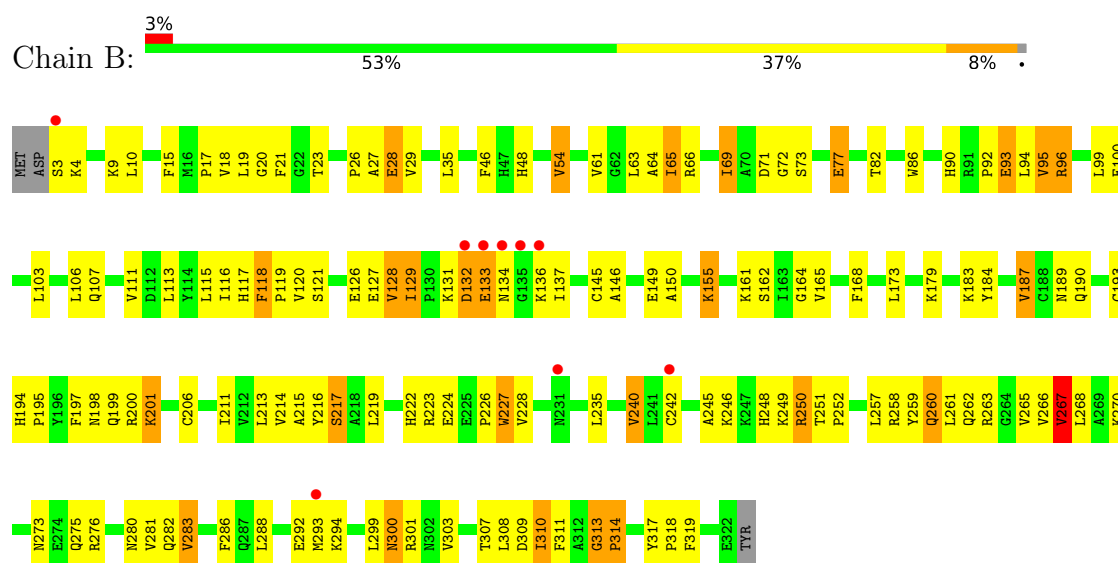
3 Residue-property plots

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: 3-ALPHA-HYDROXYSTEROID DEHYDROGENASE



• Molecule 1: 3-ALPHA-HYDROXYSTEROID DEHYDROGENASE



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	144.09Å 144.09Å 202.77Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	8.00 – 3.00 8.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	98.0 (8.00-3.00) 94.3 (8.00-3.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	15.17 (at 2.98Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.223 , 0.278 0.227 , 0.276	Depositor DCC
R_{free} test set	1620 reflections (9.89%)	wwPDB-VP
Wilson B-factor (Å ²)	29.8	Xtriage
Anisotropy	0.343	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.45 , 46.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	5272	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.89% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: IU5, NAP

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	0.48	1/2620 (0.0%)	1.06	22/3549 (0.6%)
1	B	0.49	1/2620 (0.0%)	1.06	24/3549 (0.7%)
All	All	0.48	2/5240 (0.0%)	1.06	46/7098 (0.6%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	293	MET	CG-SD	5.07	1.93	1.80
1	B	293	MET	CG-SD	5.07	1.93	1.80

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	319	PHE	N-CA-C	9.06	124.86	111.81
1	B	319	PHE	N-CA-C	9.00	124.78	111.81
1	A	313	GLY	CA-C-N	8.37	129.00	120.38
1	A	313	GLY	C-N-CA	8.37	129.00	120.38
1	A	129	ILE	CA-C-N	8.23	128.28	119.89
1	A	129	ILE	C-N-CA	8.23	128.28	119.89
1	A	72	GLY	N-CA-C	-8.19	103.43	115.63
1	B	72	GLY	N-CA-C	-7.44	104.78	115.64
1	B	313	GLY	CA-C-N	7.16	127.76	120.38
1	B	313	GLY	C-N-CA	7.16	127.76	120.38
1	B	69	ILE	N-CA-C	-6.85	104.31	111.58
1	B	227	TRP	N-CA-C	6.81	118.78	111.36
1	B	217	SER	N-CA-C	-6.74	103.62	112.24
1	B	317	TYR	CA-C-N	6.61	126.20	119.19
1	B	317	TYR	C-N-CA	6.61	126.20	119.19
1	B	118	PHE	CA-C-N	6.57	126.50	119.28
1	B	118	PHE	C-N-CA	6.57	126.50	119.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	69	ILE	N-CA-C	-6.49	104.70	111.58
1	A	116	ILE	N-CA-C	-6.43	97.10	106.88
1	B	116	ILE	N-CA-C	-6.40	97.15	106.88
1	A	217	SER	N-CA-C	-6.37	104.08	112.24
1	B	117	HIS	N-CA-C	6.29	118.66	111.11
1	B	318	PRO	N-CA-C	6.27	121.78	113.57
1	A	227	TRP	N-CA-C	6.25	119.06	111.82
1	A	317	TYR	CA-C-N	6.23	125.64	118.85
1	A	317	TYR	C-N-CA	6.23	125.64	118.85
1	B	129	ILE	CA-C-N	6.12	126.08	119.78
1	B	129	ILE	C-N-CA	6.12	126.08	119.78
1	B	293	MET	N-CA-C	-6.06	103.84	111.11
1	A	118	PHE	CA-C-N	6.04	125.93	119.28
1	A	118	PHE	C-N-CA	6.04	125.93	119.28
1	A	267	VAL	N-CA-C	6.00	118.01	108.89
1	A	318	PRO	N-CA-C	5.77	121.13	113.40
1	B	267	VAL	N-CA-C	5.69	117.79	108.97
1	A	303	VAL	N-CA-C	5.65	114.74	106.55
1	A	117	HIS	N-CA-C	5.61	117.84	111.11
1	A	312	ALA	N-CA-C	5.57	118.12	111.71
1	B	313	GLY	N-CA-C	5.50	123.56	112.34
1	A	293	MET	N-CA-C	-5.50	104.52	111.11
1	B	303	VAL	N-CA-C	5.45	114.46	106.55
1	B	310	ILE	N-CA-C	-5.35	106.04	113.00
1	A	313	GLY	N-CA-C	5.29	123.14	112.34
1	B	146	ALA	N-CA-C	-5.24	105.65	111.36
1	A	146	ALA	N-CA-C	-5.21	105.28	111.69
1	B	314	PRO	N-CA-C	5.13	116.95	110.70
1	B	317	TYR	N-CA-C	-5.06	100.72	109.58

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2560	0	2574	133	0
1	B	2560	0	2574	126	0
2	A	48	0	23	4	0
2	B	48	0	24	3	0
3	A	28	0	39	1	0
3	B	28	0	39	1	0
All	All	5272	0	5273	259	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 25.

All (259) 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:B:258:ARG:HE	1:B:262:GLN:HE21	1.03	1.03
1:B:246:LYS:HE2	1:B:246:LYS:HA	1.40	1.02
1:A:246:LYS:HE2	1:A:246:LYS:HA	1.40	1.01
1:B:273:ASN:HD22	1:B:276:ARG:H	1.12	0.93
1:A:258:ARG:HE	1:A:262:GLN:HE21	1.12	0.93
1:B:273:ASN:HD21	1:B:275:GLN:HB3	1.32	0.91
1:B:222:HIS:HD2	1:B:224:GLU:HB2	1.35	0.89
1:A:222:HIS:HD2	1:A:224:GLU:HB2	1.36	0.88
1:A:273:ASN:HD22	1:A:276:ARG:H	1.20	0.84
1:B:262:GLN:HE22	1:B:288:LEU:H	1.25	0.84
1:A:219:LEU:HD21	1:A:257:LEU:HD13	1.61	0.82
1:A:262:GLN:HE22	1:A:288:LEU:H	1.29	0.81
1:B:77:GLU:H	1:B:77:GLU:CD	1.87	0.81
1:B:10:LEU:HB3	1:B:187:VAL:CG2	2.10	0.81
1:A:273:ASN:HD21	1:A:275:GLN:HB3	1.47	0.80
1:A:10:LEU:HB3	1:A:187:VAL:CG2	2.12	0.79
1:A:132:ASP:HB2	1:A:136:LYS:HB2	1.62	0.79
1:B:10:LEU:HB3	1:B:187:VAL:HG22	1.65	0.79
1:B:136:LYS:HA	1:B:136:LYS:HE2	1.63	0.79
1:B:219:LEU:HD21	1:B:257:LEU:HD13	1.65	0.78
1:A:77:GLU:CD	1:A:77:GLU:H	1.90	0.78
1:B:132:ASP:HB2	1:B:136:LYS:HB2	1.67	0.76
1:B:118:PHE:CE1	1:B:308:LEU:HD12	2.21	0.75
1:A:136:LYS:HE2	1:A:136:LYS:HA	1.68	0.74
1:B:273:ASN:ND2	1:B:276:ARG:H	1.85	0.74
1:B:222:HIS:CD2	1:B:224:GLU:HB2	2.22	0.74
1:B:195:PRO:HG2	1:B:235:LEU:HD11	1.70	0.72
1:A:257:LEU:HB3	1:A:283:VAL:HG11	1.73	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:HIS:CD2	1:A:224:GLU:HB2	2.24	0.71
1:A:249:LYS:O	1:A:250:ARG:HD2	1.90	0.71
1:A:258:ARG:HE	1:A:262:GLN:NE2	1.86	0.71
1:A:10:LEU:HB3	1:A:187:VAL:HG22	1.73	0.70
1:B:259:TYR:CZ	1:B:263:ARG:HD2	2.27	0.70
1:B:133:GLU:C	1:B:134:ASN:HD22	2.00	0.70
1:B:198:ASN:O	1:B:200:ARG:N	2.25	0.69
1:B:258:ARG:HE	1:B:262:GLN:NE2	1.85	0.69
1:A:118:PHE:CE1	1:A:308:LEU:HD12	2.28	0.69
1:A:259:TYR:CZ	1:A:263:ARG:HD2	2.27	0.69
1:A:198:ASN:O	1:A:200:ARG:N	2.26	0.68
1:B:257:LEU:HB3	1:B:283:VAL:HG11	1.76	0.68
1:B:28:GLU:CD	1:B:28:GLU:H	2.00	0.67
1:A:270:LYS:HD3	1:A:270:LYS:C	2.20	0.67
1:B:155:LYS:HE3	1:B:161:LYS:O	1.95	0.67
1:A:128:VAL:HG12	1:A:129:ILE:HG13	1.77	0.67
1:B:257:LEU:HB3	1:B:283:VAL:CG1	2.25	0.67
1:B:198:ASN:HB2	1:B:263:ARG:NH2	2.11	0.66
1:A:273:ASN:ND2	1:A:276:ARG:H	1.92	0.66
1:A:226:PRO:HG2	1:A:227:TRP:CD1	2.31	0.65
1:B:249:LYS:O	1:B:250:ARG:HD2	1.97	0.65
1:A:257:LEU:HB3	1:A:283:VAL:CG1	2.26	0.65
1:B:26:PRO:HB2	1:B:28:GLU:HG2	1.78	0.65
1:A:28:GLU:H	1:A:28:GLU:CD	2.05	0.65
1:B:270:LYS:C	1:B:270:LYS:HD3	2.22	0.64
1:A:26:PRO:HB2	1:A:28:GLU:HG2	1.79	0.64
1:B:61:VAL:O	1:B:65:ILE:HG23	1.96	0.64
1:B:273:ASN:ND2	1:B:275:GLN:HB3	2.10	0.64
1:A:61:VAL:O	1:A:65:ILE:HG23	1.96	0.64
1:B:260:GLN:NE2	1:B:263:ARG:HE	1.95	0.63
1:A:195:PRO:HG2	1:A:235:LEU:HD11	1.81	0.63
1:B:128:VAL:HG12	1:B:129:ILE:HG13	1.80	0.62
1:B:260:GLN:HG2	1:B:267:VAL:HG13	1.80	0.62
1:A:200:ARG:NH2	1:A:201:LYS:HD2	2.15	0.62
1:A:35:LEU:HD12	1:A:64:ALA:HB2	1.81	0.62
1:A:260:GLN:HG2	1:A:267:VAL:HG13	1.80	0.61
1:A:155:LYS:HE3	1:A:161:LYS:O	2.01	0.61
1:A:198:ASN:HB2	1:A:263:ARG:NH2	2.16	0.61
1:B:35:LEU:HD12	1:B:64:ALA:HB2	1.81	0.61
1:A:258:ARG:NE	1:A:262:GLN:HE21	1.93	0.61
1:B:198:ASN:HB2	1:B:263:ARG:CZ	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:20:GLY:O	1:B:268:LEU:HA	2.02	0.60
1:B:119:PRO:HG3	1:B:168:PHE:CD1	2.37	0.60
1:A:19:LEU:HG	1:A:46:PHE:CE1	2.37	0.59
1:B:299:LEU:O	1:B:301:ARG:HG3	2.01	0.59
1:B:19:LEU:HG	1:B:46:PHE:CE1	2.37	0.59
1:B:240:VAL:CG2	1:B:299:LEU:HD11	2.33	0.59
1:A:198:ASN:HB2	1:A:263:ARG:CZ	2.33	0.59
1:A:93:GLU:CD	1:A:93:GLU:H	2.10	0.58
1:B:132:ASP:C	1:B:134:ASN:H	2.11	0.58
1:A:240:VAL:CG2	1:A:299:LEU:HD11	2.33	0.58
1:A:260:GLN:NE2	1:A:263:ARG:HE	2.01	0.58
1:A:71:ASP:HB3	1:A:73:SER:OG	2.03	0.58
1:A:20:GLY:O	1:A:268:LEU:HA	2.04	0.57
1:A:132:ASP:C	1:A:134:ASN:H	2.13	0.57
1:A:65:ILE:HG13	1:A:66:ARG:N	2.18	0.57
1:A:307:THR:O	1:A:308:LEU:C	2.48	0.57
1:B:226:PRO:HG2	1:B:227:TRP:CD1	2.39	0.57
1:B:65:ILE:O	1:B:69:ILE:HG13	2.04	0.56
1:A:54:VAL:HB	3:A:326:IU5:H183	1.88	0.56
1:A:54:VAL:HG12	1:A:128:VAL:HG11	1.87	0.56
1:A:119:PRO:HG3	1:A:168:PHE:CD1	2.41	0.55
1:B:99:LEU:O	1:B:103:LEU:HG	2.06	0.55
1:A:77:GLU:CD	1:A:77:GLU:N	2.61	0.55
1:A:183:LYS:HE2	1:A:184:TYR:OH	2.06	0.55
1:B:77:GLU:CD	1:B:77:GLU:N	2.60	0.55
1:A:240:VAL:CG2	1:A:299:LEU:HD21	2.36	0.55
1:A:3:SER:O	1:A:4:LYS:HB2	2.06	0.55
1:A:96:ARG:HD3	1:A:100:GLU:CD	2.32	0.55
1:A:299:LEU:O	1:A:301:ARG:HG3	2.06	0.55
1:B:200:ARG:NH2	1:B:201:LYS:HD2	2.21	0.55
1:B:276:ARG:HG2	2:B:325:NAP:C6A	2.37	0.55
1:A:248:HIS:HE2	1:A:292:GLU:CD	2.15	0.55
1:A:106:LEU:O	1:A:107:GLN:HB2	2.07	0.54
1:A:96:ARG:HD3	1:A:100:GLU:OE1	2.08	0.54
1:B:213:LEU:C	1:B:213:LEU:HD23	2.33	0.53
1:B:82:THR:OG1	1:B:113:LEU:HD23	2.08	0.53
1:A:198:ASN:O	1:A:198:ASN:ND2	2.41	0.53
1:B:18:VAL:HA	1:B:266:VAL:HG22	1.91	0.53
1:B:71:ASP:HB3	1:B:73:SER:OG	2.06	0.53
1:B:250:ARG:HG3	1:B:250:ARG:HH11	1.73	0.53
1:B:307:THR:O	1:B:308:LEU:C	2.51	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200:ARG:HH22	1:A:201:LYS:HD2	1.74	0.53
1:B:54:VAL:HG12	1:B:128:VAL:HG11	1.91	0.53
1:A:165:VAL:O	1:A:189:ASN:HA	2.09	0.52
1:B:183:LYS:HE2	1:B:184:TYR:OH	2.09	0.52
1:B:65:ILE:HG13	1:B:66:ARG:N	2.24	0.52
1:A:137:ILE:HD12	1:A:310:ILE:O	2.10	0.52
1:B:106:LEU:O	1:B:107:GLN:HB2	2.10	0.52
1:A:276:ARG:HG2	2:A:324:NAP:C6A	2.40	0.51
1:B:132:ASP:C	1:B:134:ASN:N	2.68	0.51
1:B:165:VAL:O	1:B:189:ASN:HA	2.10	0.51
1:A:82:THR:OG1	1:A:113:LEU:HD23	2.10	0.51
1:B:96:ARG:HD3	1:B:100:GLU:CD	2.36	0.51
1:B:3:SER:O	1:B:4:LYS:HB2	2.10	0.51
1:A:240:VAL:HG22	1:A:299:LEU:HD21	1.94	0.50
1:B:145:CYS:HB3	1:B:179:LYS:HD2	1.94	0.50
1:A:9:LYS:HG3	1:A:15:PHE:CZ	2.47	0.50
1:B:28:GLU:CD	1:B:28:GLU:N	2.70	0.50
1:A:248:HIS:O	1:A:250:ARG:HD3	2.12	0.49
1:A:86:TRP:NE1	1:A:128:VAL:HG13	2.26	0.49
1:B:54:VAL:HB	3:B:327:IU5:H183	1.93	0.49
1:A:99:LEU:O	1:A:103:LEU:HG	2.13	0.49
1:B:19:LEU:HG	1:B:46:PHE:HE1	1.77	0.49
1:A:19:LEU:HG	1:A:46:PHE:HE1	1.77	0.49
1:B:103:LEU:HD21	1:B:111:VAL:HG13	1.95	0.49
1:B:136:LYS:HA	1:B:136:LYS:CE	2.40	0.49
1:A:115:LEU:HA	1:A:164:GLY:O	2.13	0.49
1:A:133:GLU:C	1:A:134:ASN:HD22	2.20	0.49
1:A:213:LEU:C	1:A:213:LEU:HD23	2.38	0.49
1:A:10:LEU:HB3	1:A:187:VAL:HG21	1.94	0.48
1:A:132:ASP:C	1:A:134:ASN:N	2.70	0.48
1:B:194:HIS:HB2	1:B:195:PRO:HD2	1.95	0.48
1:A:213:LEU:O	1:A:265:VAL:HG13	2.14	0.48
1:B:137:ILE:HD12	1:B:310:ILE:O	2.14	0.48
1:A:219:LEU:HD21	1:A:257:LEU:CD1	2.40	0.48
1:B:93:GLU:CD	1:B:93:GLU:H	2.22	0.48
1:A:262:GLN:HE22	1:A:288:LEU:N	2.06	0.48
1:A:271:SER:HA	2:A:324:NAP:O1X	2.14	0.48
1:B:248:HIS:HE2	1:B:292:GLU:CD	2.22	0.48
1:B:246:LYS:HA	1:B:246:LYS:CE	2.26	0.48
1:A:65:ILE:O	1:A:69:ILE:HG13	2.14	0.47
1:B:240:VAL:CG2	1:B:299:LEU:HD21	2.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:250:ARG:HG2	1:A:286:PHE:CE2	2.50	0.47
1:B:119:PRO:HG3	1:B:168:PHE:CE1	2.49	0.47
1:B:265:VAL:HG12	1:B:266:VAL:N	2.29	0.47
1:A:28:GLU:CD	1:A:28:GLU:N	2.72	0.47
1:A:113:LEU:HA	1:A:162:SER:O	2.15	0.47
1:A:251:THR:HB	1:A:252:PRO:HD2	1.96	0.47
1:B:9:LYS:HG3	1:B:15:PHE:CZ	2.50	0.47
1:B:118:PHE:CZ	1:B:308:LEU:HD12	2.50	0.47
1:B:193:CYS:HB3	1:B:215:ALA:CB	2.45	0.47
1:B:265:VAL:CG1	1:B:266:VAL:N	2.78	0.47
1:B:198:ASN:O	1:B:198:ASN:ND2	2.48	0.46
1:A:119:PRO:HG3	1:A:168:PHE:CE1	2.50	0.46
1:A:194:HIS:HB2	1:A:195:PRO:HD2	1.96	0.46
1:A:126:GLU:O	1:A:127:GLU:C	2.58	0.46
1:B:115:LEU:HA	1:B:164:GLY:O	2.16	0.46
1:B:213:LEU:O	1:B:265:VAL:HG13	2.15	0.46
1:A:18:VAL:HA	1:A:266:VAL:HG22	1.96	0.46
1:A:103:LEU:HD21	1:A:111:VAL:HG13	1.98	0.46
1:B:260:GLN:HG3	1:B:265:VAL:HG11	1.98	0.46
1:A:21:PHE:HE1	1:A:23:THR:HG22	1.81	0.46
1:A:149:GLU:OE2	1:A:179:LYS:NZ	2.41	0.46
1:B:251:THR:HB	1:B:252:PRO:HD2	1.98	0.46
1:B:281:VAL:HG12	1:B:281:VAL:O	2.14	0.46
1:A:281:VAL:HG12	1:A:281:VAL:O	2.16	0.45
1:A:270:LYS:HD3	1:A:270:LYS:O	2.16	0.45
1:A:250:ARG:HG3	1:A:250:ARG:HH11	1.79	0.45
1:B:248:HIS:O	1:B:250:ARG:HD3	2.16	0.45
1:B:240:VAL:HG22	1:B:299:LEU:HD21	1.99	0.45
1:A:205:PHE:O	1:A:208:SER:OG	2.31	0.44
1:B:90:HIS:HD2	1:B:121:SER:OG	2.00	0.44
1:A:183:LYS:HE2	1:A:184:TYR:CZ	2.52	0.44
1:B:262:GLN:HE22	1:B:288:LEU:N	2.04	0.44
1:A:173:LEU:HD12	1:A:173:LEU:HA	1.85	0.44
1:A:193:CYS:HB3	1:A:215:ALA:CB	2.46	0.44
1:B:86:TRP:NE1	1:B:128:VAL:HG13	2.33	0.44
1:B:95:VAL:HG21	1:B:150:ALA:HB3	2.00	0.44
1:B:280:ASN:C	1:B:282:GLN:H	2.25	0.44
1:B:17:PRO:HB2	1:B:48:HIS:HB2	2.00	0.44
1:A:249:LYS:C	1:A:250:ARG:HD2	2.42	0.44
1:B:131:LYS:HA	1:B:136:LYS:O	2.17	0.44
1:A:17:PRO:HB2	1:A:48:HIS:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:SER:HA	1:A:90:HIS:NE2	2.33	0.44
1:A:244:LEU:HB3	1:A:255:ILE:HD13	1.99	0.44
1:A:145:CYS:HB3	1:A:179:LYS:HD2	1.99	0.43
1:B:133:GLU:C	1:B:134:ASN:ND2	2.73	0.43
1:A:24:TYR:CG	1:A:25:ALA:N	2.86	0.43
1:A:221:SER:HA	2:A:324:NAP:O3	2.18	0.43
1:B:219:LEU:HD21	1:B:257:LEU:CD1	2.43	0.43
1:B:270:LYS:O	2:B:325:NAP:H8A	2.18	0.43
1:A:177:LEU:HD23	1:A:177:LEU:HA	1.87	0.43
1:A:219:LEU:HB2	2:A:324:NAP:O1A	2.19	0.43
1:B:300:ASN:C	1:B:301:ARG:HG3	2.44	0.43
1:B:113:LEU:HA	1:B:162:SER:O	2.19	0.43
1:A:198:ASN:O	1:A:198:ASN:CG	2.62	0.43
1:B:27:ALA:C	1:B:29:VAL:N	2.76	0.43
1:A:21:PHE:CE1	1:A:23:THR:HG22	2.53	0.43
1:A:63:LEU:HD23	1:A:66:ARG:HH22	1.82	0.43
1:A:114:TYR:CD2	1:A:151:MET:HE1	2.53	0.43
1:B:250:ARG:HG2	1:B:286:PHE:CE2	2.53	0.43
1:A:273:ASN:ND2	1:A:275:GLN:HB3	2.23	0.43
1:A:190:GLN:HA	1:A:214:VAL:O	2.18	0.42
1:B:260:GLN:HG3	1:B:265:VAL:CG1	2.49	0.42
1:B:165:VAL:HG11	1:B:168:PHE:CE2	2.53	0.42
1:B:309:ASP:C	1:B:311:PHE:N	2.77	0.42
1:A:95:VAL:O	1:A:96:ARG:C	2.61	0.42
1:B:242:CYS:O	1:B:245:ALA:HB3	2.18	0.42
1:A:92:PRO:C	1:A:94:LEU:H	2.27	0.42
1:B:10:LEU:HB3	1:B:187:VAL:HG21	1.97	0.42
1:B:200:ARG:HH22	1:B:201:LYS:HD2	1.85	0.42
1:B:21:PHE:HE1	1:B:23:THR:HG22	1.84	0.42
1:A:132:ASP:CB	1:A:136:LYS:HB2	2.41	0.42
1:A:195:PRO:HB3	1:A:296:ILE:CG2	2.50	0.42
1:B:94:LEU:HD23	1:B:94:LEU:HA	1.77	0.42
1:B:126:GLU:O	1:B:127:GLU:C	2.62	0.42
1:B:113:LEU:HD11	1:B:115:LEU:CD2	2.50	0.42
1:A:120:VAL:HG11	1:A:311:PHE:HB3	2.00	0.42
1:B:96:ARG:HD3	1:B:100:GLU:OE1	2.19	0.42
1:B:206:CYS:HB3	1:B:211:ILE:O	2.20	0.42
1:A:87:SER:C	1:A:122:VAL:HG22	2.45	0.41
1:B:26:PRO:HG3	1:B:224:GLU:OE1	2.19	0.41
1:B:96:ARG:O	1:B:100:GLU:HG3	2.20	0.41
1:B:219:LEU:HB2	2:B:325:NAP:O1A	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:223:ARG:HA	1:B:228:VAL:HB	2.02	0.41
1:A:216:TYR:O	1:A:217:SER:HB2	2.19	0.41
1:A:240:VAL:HG21	1:A:299:LEU:HD21	2.01	0.41
1:A:316:ASN:O	1:A:317:TYR:C	2.63	0.41
1:B:63:LEU:HD23	1:B:66:ARG:HH22	1.84	0.41
1:B:190:GLN:HA	1:B:214:VAL:O	2.20	0.41
1:A:27:ALA:C	1:A:29:VAL:N	2.78	0.41
1:B:249:LYS:C	1:B:250:ARG:HD2	2.45	0.41
1:A:65:ILE:HD12	1:A:69:ILE:HD11	2.02	0.41
1:A:118:PHE:CZ	1:A:308:LEU:HD12	2.56	0.41
1:A:144:LEU:HD23	1:A:144:LEU:HA	1.88	0.41
1:A:203:LEU:HD11	1:A:207:LYS:HE3	2.03	0.41
1:A:260:GLN:HG3	1:A:265:VAL:HG11	2.02	0.41
1:A:309:ASP:C	1:A:311:PHE:H	2.29	0.41
1:B:194:HIS:CE1	1:B:197:PHE:HB2	2.55	0.41
1:B:222:HIS:CD2	1:B:224:GLU:H	2.39	0.41
1:B:260:GLN:CG	1:B:267:VAL:HG13	2.51	0.41
1:A:44:ALA:HB1	1:A:281:VAL:HG21	2.04	0.40
1:A:309:ASP:C	1:A:311:PHE:N	2.77	0.40
1:A:280:ASN:C	1:A:282:GLN:H	2.30	0.40
1:B:216:TYR:O	1:B:217:SER:HB2	2.21	0.40
1:A:5:TYR:OH	1:A:78:ASP:HA	2.22	0.40
1:B:21:PHE:CE1	1:B:23:THR:HG22	2.56	0.40
1:B:92:PRO:C	1:B:94:LEU:H	2.29	0.40
1:A:238:ASP:OD2	1:A:240:VAL:HG22	2.21	0.40
1:A:94:LEU:O	1:A:97:PRO:HG2	2.22	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	318/323 (98%)	289 (91%)	24 (8%)	5 (2%)	7	34
1	B	318/323 (98%)	288 (91%)	26 (8%)	4 (1%)	9	38
All	All	636/646 (98%)	577 (91%)	50 (8%)	9 (1%)	9	36

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	199	GLN
1	B	199	GLN
1	B	314	PRO
1	A	87	SER
1	A	314	PRO
1	B	313	GLY
1	A	95	VAL
1	A	313	GLY
1	B	95	VAL

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	282/285 (99%)	258 (92%)	24 (8%)	10	36
1	B	282/285 (99%)	259 (92%)	23 (8%)	10	37
All	All	564/570 (99%)	517 (92%)	47 (8%)	10	37

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	GLU
1	A	42	ILE
1	A	54	VAL
1	A	65	ILE
1	A	77	GLU
1	A	92	PRO
1	A	93	GLU

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Mol	Chain	Res	Type
1	A	96	ARG
1	A	120	VAL
1	A	128	VAL
1	A	133	GLU
1	A	149	GLU
1	A	155	LYS
1	A	173	LEU
1	A	187	VAL
1	A	201	LYS
1	A	240	VAL
1	A	250	ARG
1	A	260	GLN
1	A	261	LEU
1	A	267	VAL
1	A	283	VAL
1	A	294	LYS
1	A	300	ASN
1	B	28	GLU
1	B	54	VAL
1	B	65	ILE
1	B	77	GLU
1	B	93	GLU
1	B	96	ARG
1	B	120	VAL
1	B	128	VAL
1	B	132	ASP
1	B	133	GLU
1	B	149	GLU
1	B	155	LYS
1	B	173	LEU
1	B	187	VAL
1	B	201	LYS
1	B	240	VAL
1	B	250	ARG
1	B	260	GLN
1	B	261	LEU
1	B	267	VAL
1	B	283	VAL
1	B	294	LYS
1	B	300	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	14	HIS
1	A	56	ASN
1	A	90	HIS
1	A	105	ASN
1	A	134	ASN
1	A	167	ASN
1	A	178	ASN
1	A	190	GLN
1	A	198	ASN
1	A	260	GLN
1	A	262	GLN
1	A	273	ASN
1	A	282	GLN
1	A	287	GLN
1	A	300	ASN
1	B	6	GLN
1	B	14	HIS
1	B	48	HIS
1	B	56	ASN
1	B	90	HIS
1	B	105	ASN
1	B	134	ASN
1	B	167	ASN
1	B	178	ASN
1	B	198	ASN
1	B	260	GLN
1	B	262	GLN
1	B	273	ASN
1	B	282	GLN
1	B	287	GLN
1	B	300	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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](#)

4 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	IU5	A	326	-	31,31,31	2.06	12 (38%)	49,49,49	1.49	7 (14%)
2	NAP	A	324	-	50,52,52	2.68	16 (32%)	71,80,80	1.61	12 (16%)
3	IU5	B	327	-	31,31,31	2.08	13 (41%)	49,49,49	1.49	8 (16%)
2	NAP	B	325	-	50,52,52	2.73	17 (34%)	71,80,80	1.61	13 (18%)

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	IU5	A	326	-	1/1/11/11	1/9/70/70	0/4/4/4
3	IU5	B	327	-	1/1/11/11	1/9/70/70	0/4/4/4
2	NAP	A	324	-	-	4/35/67/67	0/5/5/5
2	NAP	B	325	-	-	4/35/67/67	0/5/5/5

All (58) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	324	NAP	C3B-C4B	-9.73	1.28	1.53
2	B	325	NAP	C3B-C4B	-9.49	1.28	1.53
2	A	324	NAP	C2N-N1N	7.76	1.43	1.35
2	B	325	NAP	C2N-N1N	7.61	1.43	1.35
2	B	325	NAP	PN-O3	5.33	1.65	1.59
2	B	325	NAP	O4B-C4B	5.01	1.56	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	324	NAP	PN-O3	4.92	1.64	1.59
2	A	324	NAP	O4B-C4B	4.82	1.55	1.45
2	B	325	NAP	C2N-C3N	4.68	1.46	1.39
2	A	324	NAP	C2N-C3N	4.34	1.45	1.39
2	B	325	NAP	O2B-C2B	-4.03	1.30	1.44
2	A	324	NAP	O2B-C2B	-3.82	1.31	1.44
2	B	325	NAP	C6N-N1N	3.76	1.43	1.35
2	A	324	NAP	C6N-N1N	3.75	1.43	1.35
3	A	326	IU5	C8-C14	3.71	1.61	1.53
3	A	326	IU5	C23-C24	3.70	1.59	1.50
3	B	327	IU5	C8-C14	3.56	1.60	1.53
3	B	327	IU5	C23-C24	3.53	1.58	1.50
2	B	325	NAP	PA-O3	3.49	1.63	1.59
2	B	325	NAP	C4A-N3A	3.41	1.40	1.34
2	A	324	NAP	C4A-N3A	3.33	1.40	1.34
3	B	327	IU5	C20-C17	3.10	1.59	1.54
3	A	326	IU5	C20-C17	3.09	1.59	1.54
2	B	325	NAP	C3N-C7N	-3.04	1.46	1.50
3	A	326	IU5	C18-C13	3.04	1.59	1.54
3	B	327	IU5	C18-C13	3.03	1.59	1.54
3	A	326	IU5	C10-C5	2.96	1.60	1.55
3	B	327	IU5	O4A-C24	-2.94	1.21	1.30
3	A	326	IU5	O4A-C24	-2.93	1.21	1.30
3	B	327	IU5	C10-C5	2.92	1.60	1.55
2	A	324	NAP	P2B-O2B	2.91	1.64	1.59
2	A	324	NAP	C4N-C3N	2.85	1.43	1.39
2	B	325	NAP	C4N-C3N	2.83	1.43	1.39
2	B	325	NAP	O5B-C5B	2.80	1.55	1.44
2	A	324	NAP	C3N-C7N	-2.80	1.46	1.50
2	B	325	NAP	P2B-O2B	2.78	1.64	1.59
3	A	326	IU5	C10-C9	2.73	1.60	1.56
3	B	327	IU5	C10-C9	2.69	1.60	1.56
2	A	324	NAP	O5B-C5B	2.62	1.54	1.44
2	A	324	NAP	P2B-O2X	-2.55	1.45	1.54
2	B	325	NAP	P2B-O2X	-2.54	1.45	1.54
3	B	327	IU5	C8-C7	2.50	1.58	1.53
2	A	324	NAP	PA-O2A	-2.45	1.44	1.55
2	B	325	NAP	C2B-C1B	-2.40	1.47	1.53
2	A	324	NAP	C2B-C1B	-2.32	1.47	1.53
3	A	326	IU5	C12-C11	2.28	1.57	1.53
2	B	325	NAP	PA-O2A	-2.27	1.44	1.55
3	B	327	IU5	C16-C17	2.25	1.59	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	326	IU5	C2-C3	2.22	1.56	1.51
3	A	326	IU5	C1-C10	2.19	1.58	1.54
3	B	327	IU5	C12-C11	2.16	1.57	1.53
2	B	325	NAP	C8A-N9A	2.16	1.41	1.37
3	A	326	IU5	C16-C17	2.16	1.58	1.54
3	B	327	IU5	C8-C9	2.12	1.57	1.53
3	B	327	IU5	C2-C3	2.11	1.56	1.51
2	A	324	NAP	P2B-O3X	-2.05	1.47	1.54
3	A	326	IU5	C8-C7	2.03	1.57	1.53
3	B	327	IU5	C1-C10	2.02	1.57	1.54

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	325	NAP	O3-PA-O1A	-4.95	95.82	110.70
2	A	324	NAP	O3-PA-O1A	-4.88	96.03	110.70
3	A	326	IU5	C6-C5-C10	-4.26	108.12	112.66
2	B	325	NAP	C1B-N9A-C8A	4.23	136.48	127.09
2	A	324	NAP	C5N-C4N-C3N	4.23	124.52	120.36
3	B	327	IU5	C6-C5-C10	-4.22	108.17	112.66
2	A	324	NAP	C1B-N9A-C8A	4.13	136.27	127.09
2	B	325	NAP	C5N-C4N-C3N	4.02	124.31	120.36
2	A	324	NAP	O7N-C7N-N7N	-3.79	117.14	122.62
2	A	324	NAP	C2B-C3B-C4B	3.73	110.02	101.99
2	B	325	NAP	O7N-C7N-N7N	-3.64	117.36	122.62
2	B	325	NAP	C2B-C3B-C4B	3.61	109.76	101.99
3	A	326	IU5	C9-C8-C7	-3.33	107.67	111.86
3	B	327	IU5	C9-C8-C7	-3.32	107.68	111.86
2	B	325	NAP	C4A-N9A-C1B	-3.13	119.31	126.63
2	A	324	NAP	C4A-N9A-C1B	-3.04	119.52	126.63
2	A	324	NAP	N3A-C2A-N1A	-2.98	124.07	128.58
2	B	325	NAP	N3A-C2A-N1A	-2.89	124.20	128.58
2	B	325	NAP	P2B-O2B-C2B	-2.89	115.70	123.43
3	B	327	IU5	C17-C13-C14	-2.89	96.79	100.10
2	A	324	NAP	P2B-O2B-C2B	-2.79	115.99	123.43
3	A	326	IU5	C17-C13-C14	-2.78	96.91	100.10
3	A	326	IU5	C1-C2-C3	-2.75	106.84	110.48
3	B	327	IU5	C1-C2-C3	-2.74	106.86	110.48
2	A	324	NAP	C3N-C7N-N7N	2.71	121.07	117.74
3	B	327	IU5	C5-C4-C3	-2.66	108.70	112.71
3	A	326	IU5	C5-C4-C3	-2.59	108.81	112.71
2	B	325	NAP	C3N-C7N-N7N	2.51	120.83	117.74

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	325	NAP	O5B-C5B-C4B	-2.44	100.67	108.99
3	B	327	IU5	C5-C6-C7	-2.34	111.62	114.40
3	A	326	IU5	C5-C6-C7	-2.30	111.66	114.40
2	A	324	NAP	O5B-C5B-C4B	-2.28	101.25	108.99
2	B	325	NAP	O2A-PA-O3	2.24	113.32	107.27
2	B	325	NAP	O2A-PA-O1A	2.18	122.59	112.44
3	A	326	IU5	C1-C10-C5	-2.13	104.69	107.75
3	B	327	IU5	C1-C10-C5	-2.09	104.74	107.75
2	A	324	NAP	C4B-O4B-C1B	-2.09	104.85	109.47
2	B	325	NAP	C4B-O4B-C1B	-2.09	104.85	109.47
3	B	327	IU5	C16-C17-C13	-2.09	101.38	103.84
2	A	324	NAP	O2A-PA-O1A	2.07	122.08	112.44

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	326	IU5	C3
3	B	327	IU5	C3

All (10) torsion outliers are listed below:

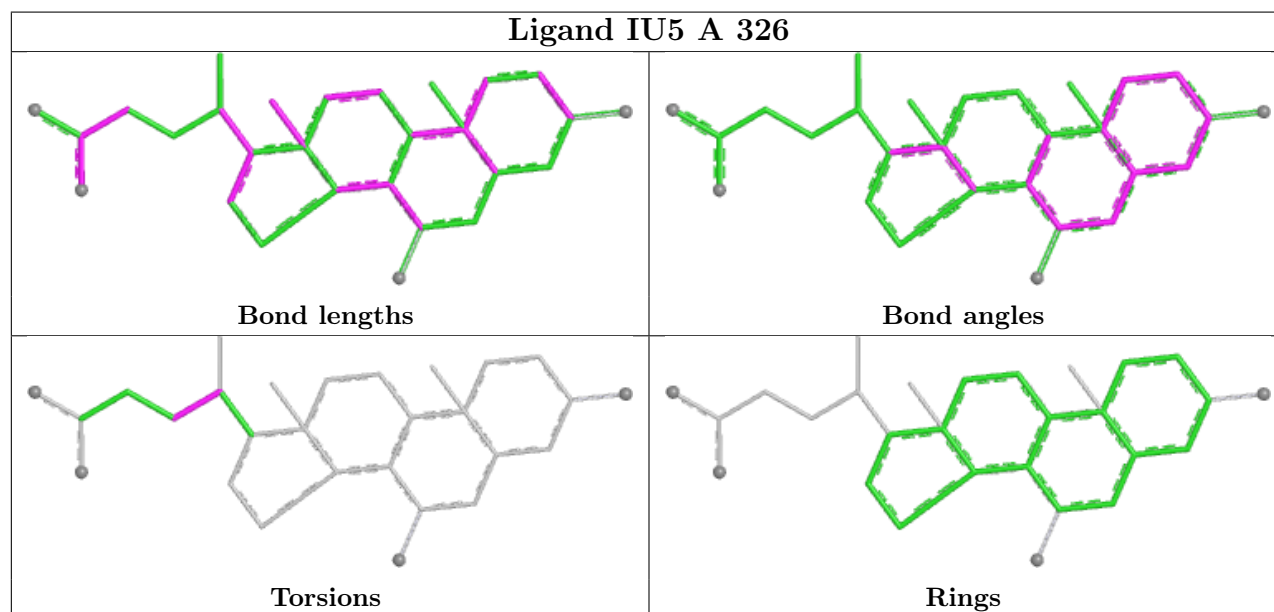
Mol	Chain	Res	Type	Atoms
2	A	324	NAP	C3D-C4D-C5D-O5D
2	B	325	NAP	C3D-C4D-C5D-O5D
2	A	324	NAP	O4D-C4D-C5D-O5D
2	B	325	NAP	O4D-C4D-C5D-O5D
2	A	324	NAP	PA-O3-PN-O5D
2	B	325	NAP	PA-O3-PN-O5D
3	A	326	IU5	C17-C20-C22-C23
3	B	327	IU5	C17-C20-C22-C23
2	A	324	NAP	PA-O3-PN-O2N
2	B	325	NAP	PA-O3-PN-O2N

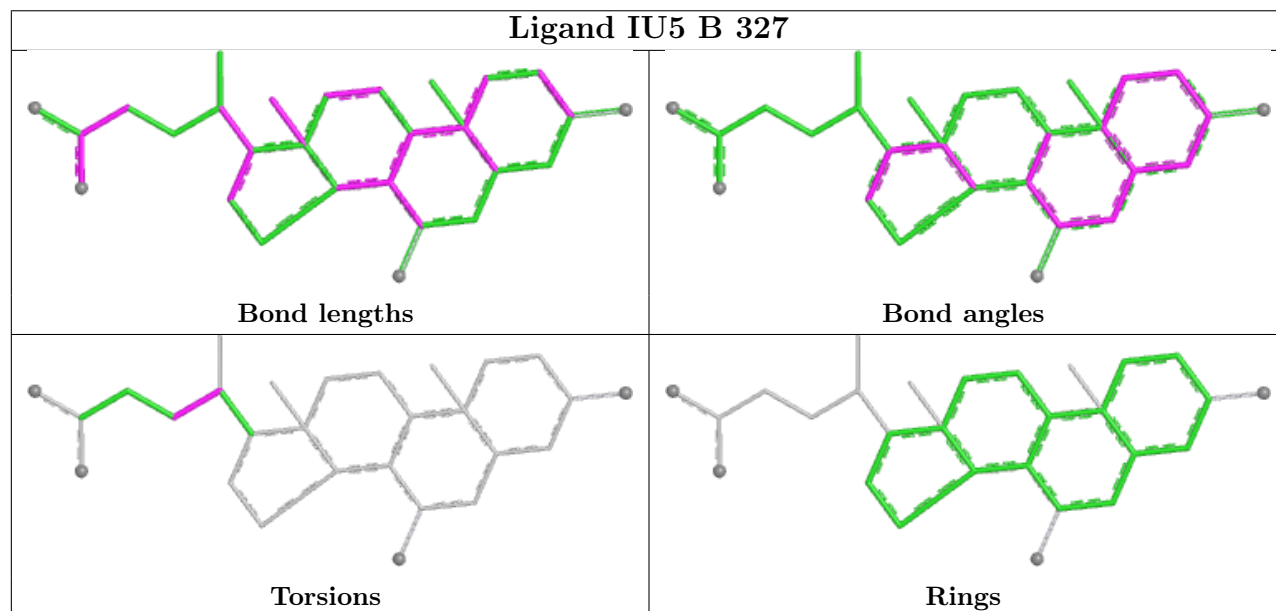
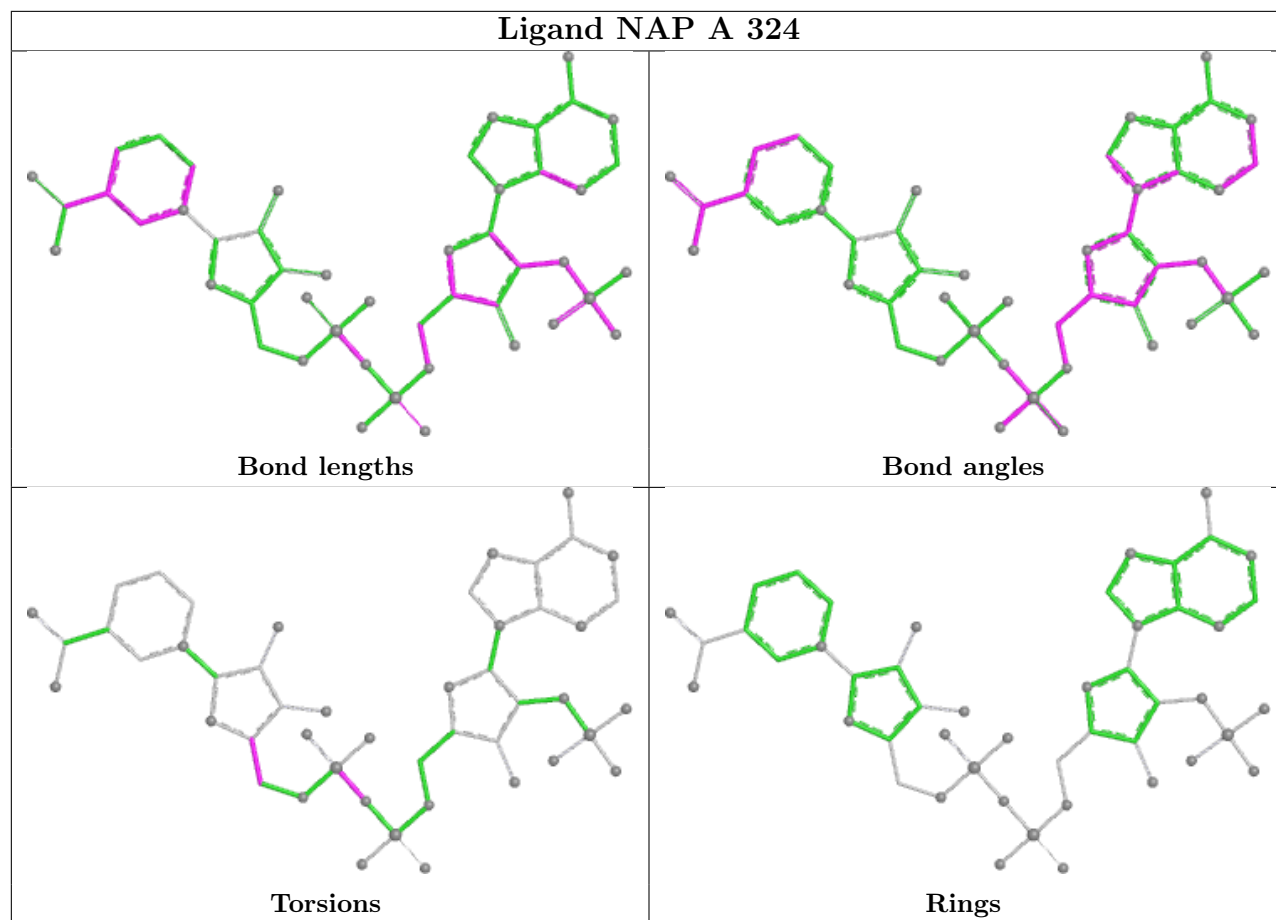
There are no ring outliers.

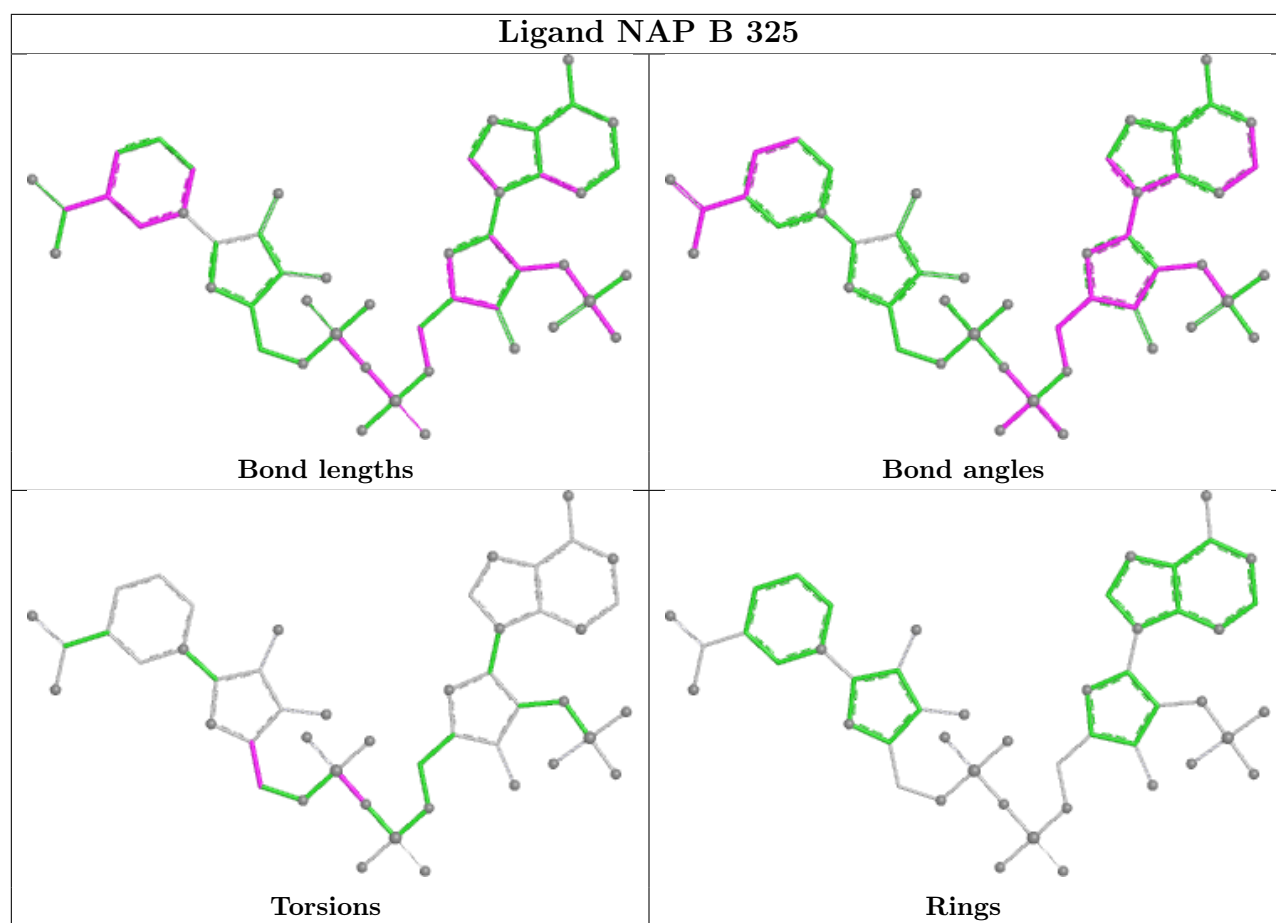
4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	326	IU5	1	0
2	A	324	NAP	4	0
3	B	327	IU5	1	0
2	B	325	NAP	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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	320/323 (99%)	-0.41	5 (1%) 70 47	3, 14, 26, 34	0
1	B	320/323 (99%)	-0.32	9 (2%) 55 32	3, 15, 34, 41	0
All	All	640/646 (99%)	-0.36	14 (2%) 62 39	3, 14, 30, 41	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	133	GLU	4.9
1	B	132	ASP	4.6
1	B	134	ASN	4.0
1	B	242	CYS	3.8
1	A	293	MET	3.6
1	B	136	LYS	3.1
1	B	293	MET	2.6
1	B	135	GLY	2.6
1	A	28	GLU	2.5
1	A	126	GLU	2.5
1	B	231	ASN	2.3
1	A	133	GLU	2.2
1	B	3	SER	2.1
1	A	291	GLU	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](#)

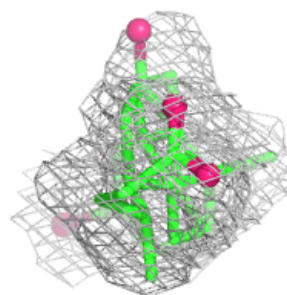
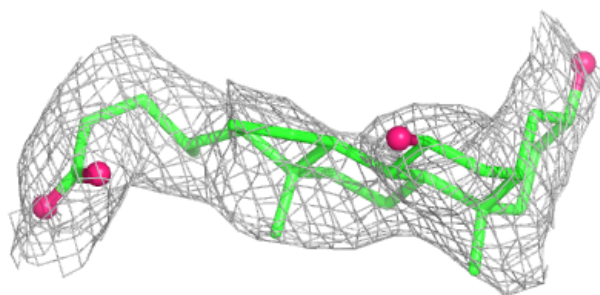
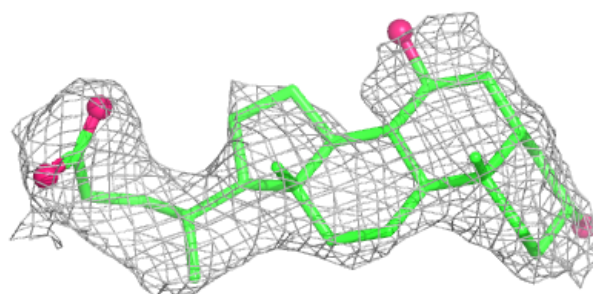
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	IU5	B	327	28/28	0.85	0.11	13,29,31,33	0
3	IU5	A	326	28/28	0.87	0.10	8,18,21,22	0
2	NAP	B	325	48/48	0.91	0.10	18,32,37,37	0
2	NAP	A	324	48/48	0.96	0.07	13,18,23,25	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.

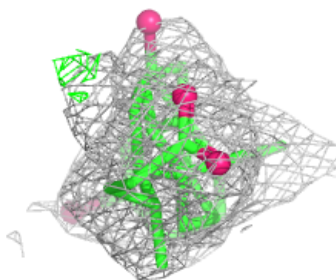
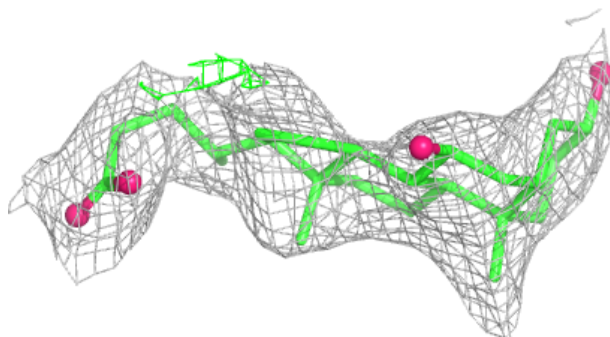
Electron density around IU5 B 327:

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

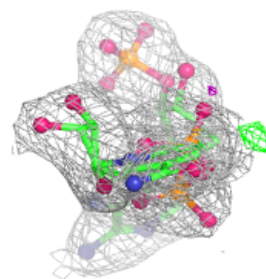
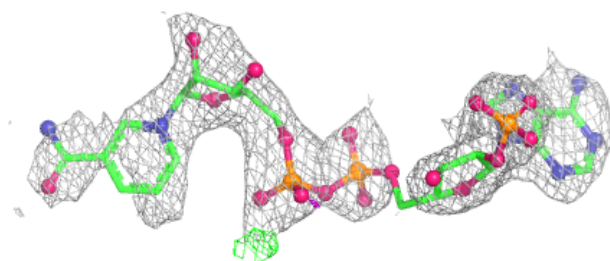
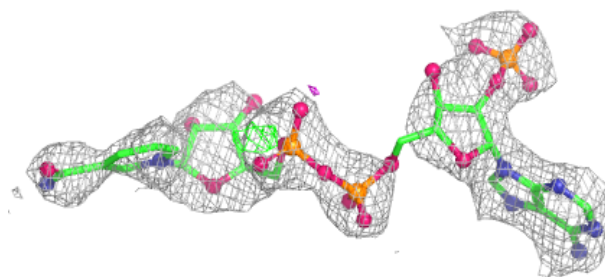


Electron density around IU5 A 326:

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

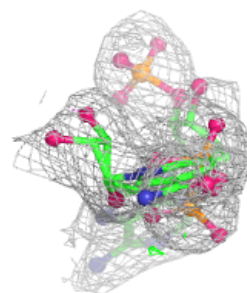
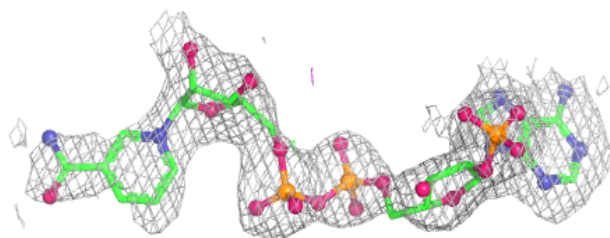
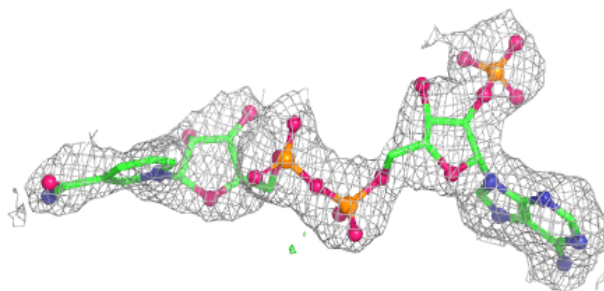
**Electron density around NAP B 325:**

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



Electron density around NAP A 324:

$2mF_o-DF_c$ (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.