



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:53 PM GMT

PDB ID : 4GI2
Title : Crotonyl-CoA Carboxylase/Reductase
Authors : Weidenweber, S.; Erb, T.J.; Ermler, U.
Deposited on : 2012-08-08
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

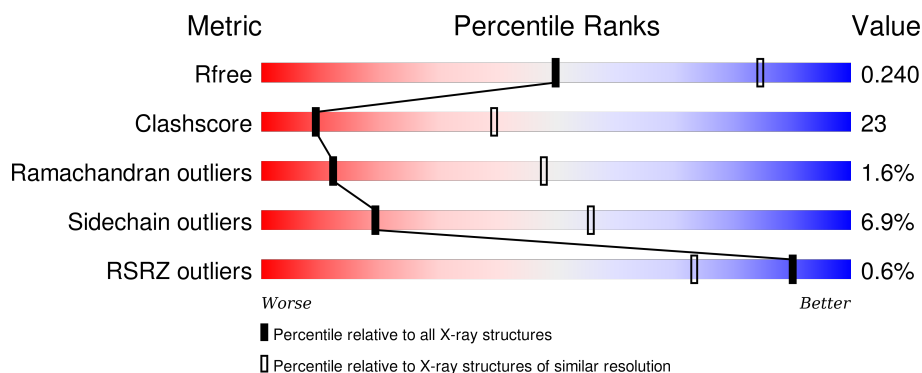
1 Overall quality at a glance ⓘ

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}	91344	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	443	<div> <div>59%</div> <div>33%</div> <div>• •</div> </div>
1	B	443	<div> <div>%</div> <div>61%</div> <div>30%</div> <div>5%</div> <div>•</div> </div>

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
2	MG	A	501	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6730 atoms, of which 1 is hydrogen 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 Crotonyl-CoA carboxylase/reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	428	Total	C	N	O	S	0	0	0
			3322	2110	587	604	21			
1	B	426	Total	C	N	O	S	0	0	0
			3310	2102	585	602	21			

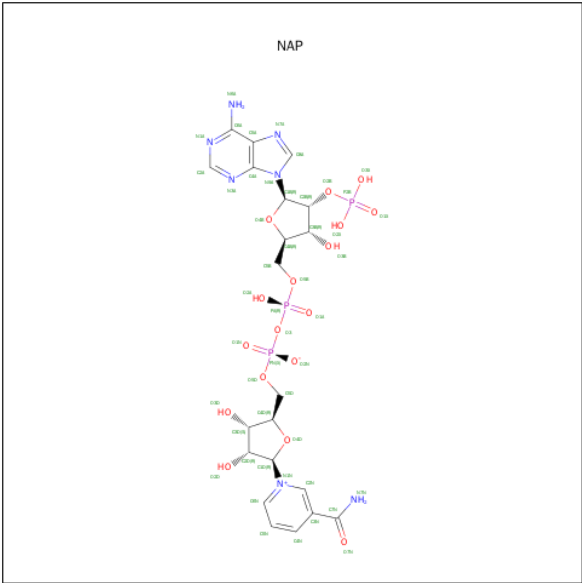
There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	433	ALA	-	EXPRESSION TAG	UNP C5AP81
A	434	ALA	-	EXPRESSION TAG	UNP C5AP81
A	435	ALA	-	EXPRESSION TAG	UNP C5AP81
A	436	LEU	-	EXPRESSION TAG	UNP C5AP81
A	437	GLU	-	EXPRESSION TAG	UNP C5AP81
A	438	HIS	-	EXPRESSION TAG	UNP C5AP81
A	439	HIS	-	EXPRESSION TAG	UNP C5AP81
A	440	HIS	-	EXPRESSION TAG	UNP C5AP81
A	441	HIS	-	EXPRESSION TAG	UNP C5AP81
A	442	HIS	-	EXPRESSION TAG	UNP C5AP81
A	443	HIS	-	EXPRESSION TAG	UNP C5AP81
B	433	ALA	-	EXPRESSION TAG	UNP C5AP81
B	434	ALA	-	EXPRESSION TAG	UNP C5AP81
B	435	ALA	-	EXPRESSION TAG	UNP C5AP81
B	436	LEU	-	EXPRESSION TAG	UNP C5AP81
B	437	GLU	-	EXPRESSION TAG	UNP C5AP81
B	438	HIS	-	EXPRESSION TAG	UNP C5AP81
B	439	HIS	-	EXPRESSION TAG	UNP C5AP81
B	440	HIS	-	EXPRESSION TAG	UNP C5AP81
B	441	HIS	-	EXPRESSION TAG	UNP C5AP81
B	442	HIS	-	EXPRESSION TAG	UNP C5AP81
B	443	HIS	-	EXPRESSION TAG	UNP C5AP81

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).

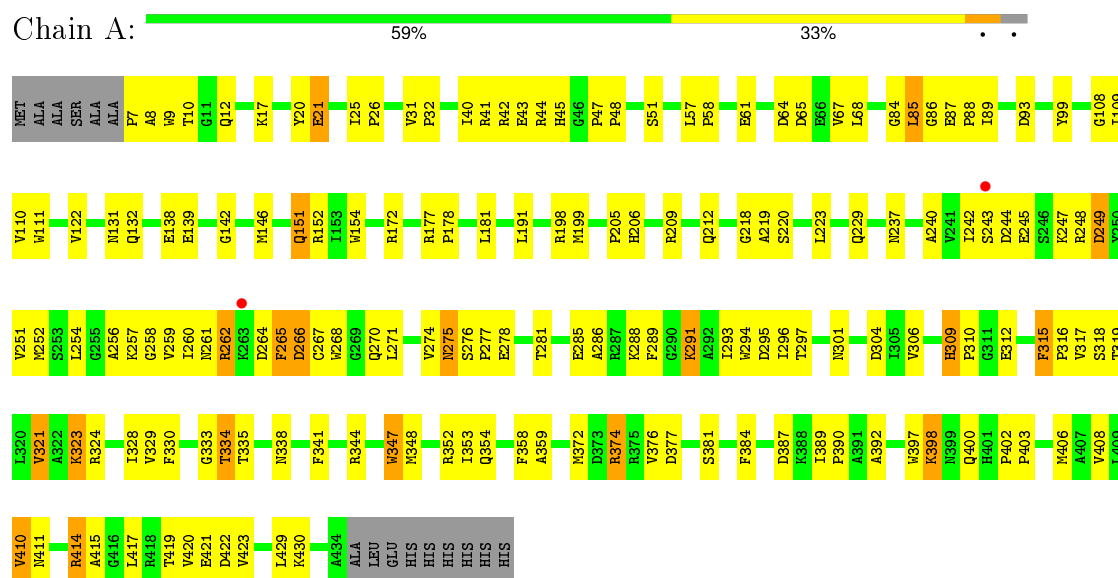


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
3	A	1	Total 48	C 21	N 7	O 17	P 3	0	0	
3	B	1	Total 49	C 21	H 1	N 7	O 17	P 3	0	0

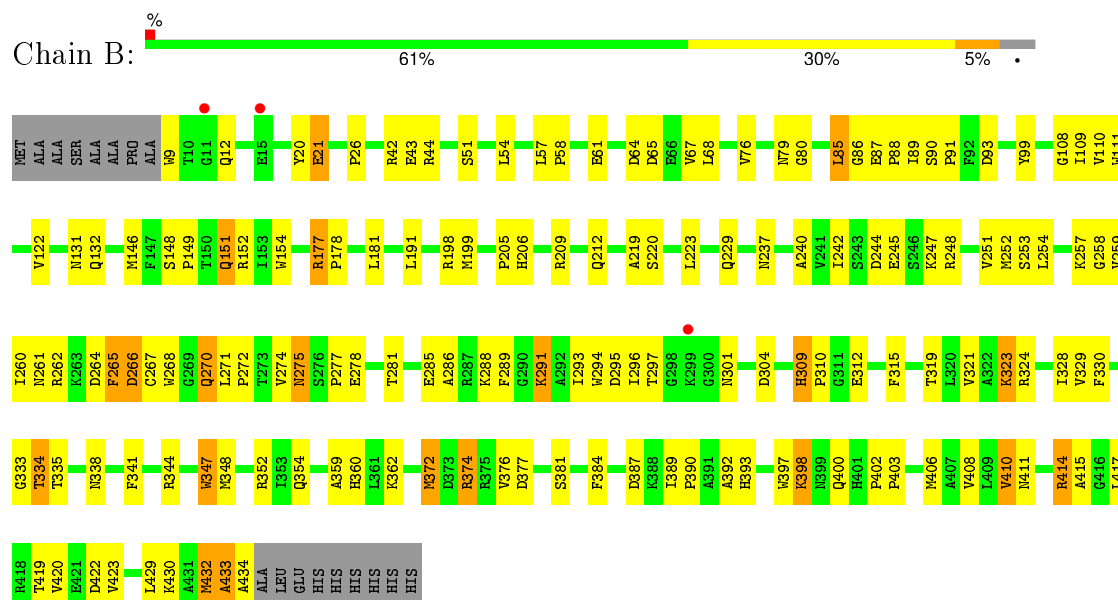
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Crotonyl-CoA carboxylase/reductase



• Molecule 1: Crotonyl-CoA carboxylase/reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	91.40 Å 91.40 Å 206.10 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.18 – 3.00 43.18 – 2.52	Depositor EDS
% Data completeness (in resolution range)	99.6 (43.18-3.00) 95.1 (43.18-2.52)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.25 (at 2.51 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.182 , 0.241 0.181 , 0.240	Depositor DCC
R_{free} test set	1039 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	73.2	Xtriage
Anisotropy	0.245	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 67.4	EDS
Estimated twinning fraction	0.027 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 32697 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6730	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.83% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: NAP, 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	0.45	0/3416	0.60	0/4642
1	B	0.45	0/3403	0.60	0/4624
All	All	0.45	0/6819	0.60	0/9266

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	3322	0	3218	160	1
1	B	3310	0	3205	148	1
2	A	1	0	0	0	0
3	A	48	0	25	2	0
3	B	48	1	25	3	0
All	All	6729	1	6473	302	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 23.

All (302) 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:374:ARG:HH11	1:A:374:ARG:HG3	1.13	1.13
1:B:374:ARG:HH11	1:B:374:ARG:HG3	1.11	1.08
1:A:374:ARG:CG	1:A:374:ARG:HH11	1.79	0.96
1:A:229:GLN:HE22	1:A:377:ASP:H	1.09	0.95
1:B:352:ARG:HE	1:B:354:GLN:HE21	1.13	0.94
1:A:352:ARG:HE	1:A:354:GLN:HE21	1.12	0.94
1:B:374:ARG:HH11	1:B:374:ARG:CG	1.79	0.94
1:A:309:HIS:H	1:A:310:PRO:HD2	1.33	0.92
1:B:229:GLN:HE22	1:B:377:ASP:H	1.16	0.92
1:A:110:VAL:HG23	1:A:122:VAL:HA	1.53	0.91
1:B:297:THR:CG2	1:B:301:ASN:HB3	2.02	0.90
1:B:110:VAL:HG23	1:B:122:VAL:HA	1.56	0.86
1:B:309:HIS:H	1:B:310:PRO:HD2	1.40	0.86
1:A:278:GLU:N	1:A:278:GLU:OE1	2.09	0.85
1:B:278:GLU:N	1:B:278:GLU:OE1	2.07	0.85
1:A:297:THR:CG2	1:A:301:ASN:HB3	2.06	0.85
1:A:392:ALA:HB1	1:A:406:MET:HE1	1.59	0.84
1:A:9:TRP:NE1	1:A:420:VAL:HG11	1.93	0.84
1:A:258:GLY:HA3	1:A:296:ILE:HD13	1.60	0.83
1:B:297:THR:HG22	1:B:301:ASN:HB3	1.58	0.82
1:B:245:GLU:OE1	1:B:248:ARG:NH1	2.12	0.82
1:B:392:ALA:HB1	1:B:406:MET:HE1	1.61	0.82
1:B:352:ARG:HE	1:B:354:GLN:NE2	1.77	0.82
1:A:352:ARG:HE	1:A:354:GLN:NE2	1.78	0.81
1:A:352:ARG:NE	1:A:354:GLN:HE21	1.78	0.80
1:B:374:ARG:NH1	1:B:374:ARG:HG3	1.87	0.80
1:A:277:PRO:O	1:A:281:THR:HG23	1.80	0.80
1:B:344:ARG:O	1:B:348:MET:HG3	1.81	0.80
1:B:352:ARG:NE	1:B:354:GLN:HE21	1.79	0.80
1:B:277:PRO:O	1:B:281:THR:HG23	1.82	0.79
1:A:245:GLU:OE1	1:A:248:ARG:NH1	2.16	0.79
1:A:43:GLU:N	1:A:43:GLU:OE1	2.17	0.77
1:A:344:ARG:O	1:A:348:MET:HG3	1.84	0.77
1:B:252:MET:CE	1:B:259:VAL:HG23	2.14	0.77
1:A:229:GLN:HE22	1:A:377:ASP:N	1.82	0.77
1:B:258:GLY:HA3	1:B:296:ILE:HD13	1.66	0.77
1:A:297:THR:HG22	1:A:301:ASN:HB3	1.65	0.77
1:B:43:GLU:OE1	1:B:43:GLU:N	2.18	0.76
1:A:374:ARG:NH1	1:A:374:ARG:HG3	1.89	0.76
1:B:109:ILE:N	1:B:109:ILE:HD12	2.02	0.75
1:A:392:ALA:HB1	1:A:406:MET:CE	2.16	0.75
1:B:398:LYS:H	1:B:398:LYS:HD2	1.51	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:398:LYS:HD2	1:A:398:LYS:H	1.52	0.74
1:A:252:MET:CE	1:A:259:VAL:HG23	2.18	0.74
1:A:381:SER:HB2	1:A:402:PRO:HG2	1.71	0.73
1:B:297:THR:HG22	1:B:301:ASN:CB	2.18	0.73
1:B:392:ALA:HB1	1:B:406:MET:CE	2.18	0.72
1:B:178:PRO:HA	1:B:372:MET:CE	2.20	0.72
1:B:9:TRP:NE1	1:B:420:VAL:HG11	2.05	0.72
1:A:57:LEU:HB3	1:A:58:PRO:HD2	1.70	0.71
1:B:257:LYS:O	1:B:257:LYS:HG2	1.91	0.71
1:A:309:HIS:N	1:A:310:PRO:HD2	2.06	0.71
1:B:381:SER:HB2	1:B:402:PRO:HG2	1.72	0.71
1:A:64:ASP:HB2	1:B:64:ASP:HB2	1.72	0.70
1:B:229:GLN:HE22	1:B:377:ASP:N	1.90	0.70
1:A:419:THR:HG23	1:A:422:ASP:H	1.57	0.69
1:B:178:PRO:HB2	1:B:181:LEU:HD12	1.75	0.69
1:A:109:ILE:HD12	1:A:109:ILE:N	2.06	0.69
1:A:110:VAL:CG2	1:A:122:VAL:HA	2.23	0.69
1:A:297:THR:HG22	1:A:301:ASN:CB	2.23	0.68
1:A:257:LYS:HG2	1:A:257:LYS:O	1.92	0.68
1:B:419:THR:HG23	1:B:422:ASP:H	1.58	0.68
1:A:319:THR:HA	1:A:328:ILE:CD1	2.24	0.68
1:A:178:PRO:HA	1:A:372:MET:CE	2.24	0.68
1:A:291:LYS:HE3	1:A:295:ASP:OD2	1.92	0.68
1:A:9:TRP:CD1	1:A:420:VAL:HG11	2.29	0.67
1:B:252:MET:HE2	1:B:259:VAL:HG23	1.76	0.67
1:B:297:THR:HG21	1:B:301:ASN:HB3	1.77	0.67
1:B:319:THR:HA	1:B:328:ILE:CD1	2.26	0.66
1:A:7:PRO:HA	1:A:10:THR:CG2	2.26	0.66
1:B:398:LYS:O	1:B:398:LYS:HG2	1.95	0.66
1:A:297:THR:HG21	1:A:301:ASN:HB3	1.77	0.65
1:B:309:HIS:N	1:B:310:PRO:HD2	2.10	0.65
1:B:54:LEU:HD23	1:B:432:MET:HG3	1.78	0.65
1:B:110:VAL:CG2	1:B:122:VAL:HA	2.26	0.65
1:A:352:ARG:HH21	1:A:354:GLN:NE2	1.94	0.64
1:B:252:MET:HE2	1:B:259:VAL:CG2	2.27	0.64
1:B:54:LEU:HD23	1:B:432:MET:HE2	1.80	0.64
1:B:291:LYS:HE3	1:B:295:ASP:OD2	1.98	0.64
1:B:374:ARG:NH1	1:B:374:ARG:CG	2.50	0.64
1:A:297:THR:HG22	1:A:297:THR:O	1.98	0.64
1:A:178:PRO:HB2	1:A:181:LEU:HD12	1.80	0.64
1:B:108:GLY:C	1:B:109:ILE:HD12	2.18	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:GLY:HA3	1:B:360:HIS:CE1	2.33	0.63
1:B:252:MET:HE1	1:B:259:VAL:HG23	1.80	0.63
1:A:252:MET:HE2	1:A:259:VAL:HG23	1.81	0.63
1:A:252:MET:HE1	1:A:259:VAL:HG23	1.80	0.63
1:A:43:GLU:CD	1:A:43:GLU:H	2.01	0.62
1:A:87:GLU:HA	1:A:88:PRO:C	2.20	0.62
1:A:265:PHE:N	1:A:265:PHE:CD2	2.68	0.62
1:B:54:LEU:HD23	1:B:432:MET:CE	2.29	0.61
1:B:26:PRO:HG2	1:B:111:TRP:CZ2	2.34	0.61
1:B:352:ARG:HH21	1:B:354:GLN:NE2	1.99	0.61
1:A:9:TRP:CD1	1:A:420:VAL:CG1	2.83	0.61
1:A:21:GLU:HG2	1:A:430:LYS:HA	1.82	0.61
1:A:251:VAL:HA	1:A:254:LEU:HD12	1.82	0.61
1:A:319:THR:HA	1:A:328:ILE:HD11	1.83	0.61
1:B:432:MET:O	1:B:434:ALA:N	2.34	0.60
1:A:264:ASP:HB2	1:A:265:PHE:CE2	2.37	0.60
1:A:26:PRO:HG2	1:A:111:TRP:CZ2	2.37	0.60
1:B:398:LYS:HG2	1:B:400:GLN:HG3	1.84	0.60
1:B:21:GLU:HG2	1:B:430:LYS:HA	1.83	0.60
1:A:108:GLY:C	1:A:109:ILE:HD12	2.21	0.60
1:A:398:LYS:HG2	1:A:398:LYS:O	2.01	0.60
1:B:178:PRO:HA	1:B:372:MET:HE1	1.83	0.60
1:B:9:TRP:CD1	1:B:420:VAL:HG11	2.37	0.59
1:B:191:LEU:C	1:B:191:LEU:HD23	2.21	0.59
1:B:347:TRP:C	1:B:347:TRP:CD1	2.76	0.59
1:B:57:LEU:HB3	1:B:58:PRO:HD2	1.83	0.59
1:A:304:ASP:OD1	1:A:323:LYS:HE2	2.03	0.59
1:A:266:ASP:O	1:A:268:TRP:N	2.36	0.59
1:B:265:PHE:CD2	1:B:265:PHE:N	2.70	0.58
1:A:429:LEU:HD12	1:A:430:LYS:N	2.17	0.58
1:B:87:GLU:HA	1:B:88:PRO:C	2.21	0.58
1:A:266:ASP:C	1:A:268:TRP:H	2.06	0.58
1:B:43:GLU:CD	1:B:43:GLU:H	2.05	0.58
1:A:229:GLN:NE2	1:A:377:ASP:H	1.89	0.57
1:A:252:MET:HE2	1:A:259:VAL:CG2	2.34	0.57
1:B:251:VAL:HA	1:B:254:LEU:HD12	1.85	0.57
1:B:319:THR:HA	1:B:328:ILE:HD11	1.85	0.57
1:B:304:ASP:OD1	1:B:323:LYS:HE2	2.05	0.57
1:B:266:ASP:C	1:B:268:TRP:H	2.07	0.57
1:A:44:ARG:HB3	1:A:51:SER:HA	1.86	0.57
1:B:67:VAL:HG12	1:B:68:LEU:N	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:LEU:HD23	1:A:191:LEU:C	2.26	0.56
1:B:392:ALA:CB	1:B:406:MET:HE1	2.35	0.56
1:B:242:ILE:HD11	1:B:259:VAL:HG11	1.88	0.56
1:B:264:ASP:HB2	1:B:265:PHE:CE2	2.41	0.56
1:A:7:PRO:HA	1:A:10:THR:HG23	1.89	0.55
1:A:410:VAL:HG12	1:A:411:ASN:N	2.22	0.55
1:A:398:LYS:HG2	1:A:400:GLN:HG3	1.88	0.55
1:B:109:ILE:CD1	1:B:109:ILE:N	2.67	0.55
1:B:54:LEU:CD2	1:B:432:MET:HG3	2.37	0.55
1:A:309:HIS:H	1:A:310:PRO:CD	2.09	0.55
1:A:242:ILE:HD11	1:A:259:VAL:HG11	1.88	0.55
1:A:132:GLN:NE2	1:A:154:TRP:HB2	2.21	0.55
1:B:87:GLU:HA	1:B:89:ILE:N	2.21	0.55
1:B:334:THR:HG22	3:B:501:NAP:O3D	2.08	0.54
1:B:304:ASP:HA	1:B:323:LYS:HE3	1.89	0.54
1:A:392:ALA:CB	1:A:406:MET:HE1	2.34	0.54
1:B:429:LEU:HD12	1:B:430:LYS:N	2.21	0.54
1:A:304:ASP:HA	1:A:323:LYS:HE3	1.90	0.54
1:B:199:MET:HE3	1:B:330:PHE:HA	1.90	0.54
1:A:178:PRO:HA	1:A:372:MET:HE3	1.88	0.53
1:A:142:GLY:CA	1:B:360:HIS:CE1	2.91	0.53
1:B:44:ARG:HB3	1:B:51:SER:HA	1.91	0.53
1:A:347:TRP:C	1:A:347:TRP:CD1	2.82	0.53
1:A:334:THR:HG22	3:A:502:NAP:O3D	2.09	0.53
1:B:229:GLN:NE2	1:B:377:ASP:H	1.97	0.53
1:B:132:GLN:NE2	1:B:154:TRP:HB2	2.24	0.52
1:B:42:ARG:HH22	1:B:93:ASP:CG	2.13	0.52
1:B:252:MET:CE	1:B:259:VAL:CG2	2.85	0.52
1:A:297:THR:CG2	1:A:297:THR:O	2.57	0.52
1:B:266:ASP:O	1:B:268:TRP:N	2.39	0.52
1:A:218:GLY:HA3	3:A:502:NAP:H1B	1.91	0.52
1:A:333:GLY:O	1:A:335:THR:N	2.43	0.52
1:A:309:HIS:CD2	1:A:310:PRO:HD3	2.45	0.52
1:B:309:HIS:CD2	1:B:310:PRO:HD3	2.45	0.52
1:A:220:SER:HB2	1:A:403:PRO:HB3	1.92	0.52
1:B:410:VAL:HG12	1:B:411:ASN:N	2.25	0.52
1:A:323:LYS:HD2	1:A:324:ARG:O	2.09	0.51
1:B:9:TRP:CD1	1:B:420:VAL:CG1	2.94	0.51
1:B:309:HIS:H	1:B:310:PRO:CD	2.15	0.51
1:A:109:ILE:CD1	1:A:109:ILE:N	2.74	0.51
1:A:309:HIS:N	1:A:310:PRO:CD	2.72	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:GLU:HA	1:A:61:GLU:OE1	2.11	0.51
1:A:87:GLU:HA	1:A:89:ILE:N	2.25	0.51
1:B:259:VAL:CG1	1:B:260:ILE:N	2.74	0.50
1:A:259:VAL:CG1	1:A:260:ILE:N	2.75	0.50
1:B:259:VAL:HG12	1:B:260:ILE:N	2.26	0.50
1:A:131:ASN:OD1	1:A:151:GLN:NE2	2.45	0.50
1:A:286:ALA:O	1:A:289:PHE:HB3	2.11	0.50
1:A:293:ILE:HG22	1:A:294:TRP:N	2.24	0.50
1:A:252:MET:CE	1:A:259:VAL:CG2	2.88	0.50
1:A:199:MET:HE3	1:A:330:PHE:HA	1.93	0.50
1:A:199:MET:CE	1:A:330:PHE:HA	2.40	0.50
1:B:323:LYS:HD2	1:B:324:ARG:O	2.11	0.50
1:B:297:THR:HG22	1:B:297:THR:O	2.11	0.50
1:B:199:MET:CE	1:B:330:PHE:HA	2.42	0.49
1:A:285:GLU:O	1:A:288:LYS:HB2	2.12	0.49
1:B:178:PRO:HA	1:B:372:MET:HE3	1.94	0.49
1:B:209:ARG:O	1:B:212:GLN:HG3	2.13	0.49
1:B:274:VAL:O	1:B:275:ASN:HB2	2.12	0.49
1:B:285:GLU:O	1:B:288:LYS:HB2	2.12	0.49
1:A:67:VAL:HG12	1:A:68:LEU:N	2.27	0.49
1:B:329:VAL:HA	1:B:354:GLN:O	2.12	0.49
1:B:21:GLU:CG	1:B:430:LYS:HA	2.43	0.49
1:B:414:ARG:HD2	1:B:415:ALA:H	1.78	0.49
1:A:259:VAL:HG12	1:A:260:ILE:N	2.28	0.49
1:A:57:LEU:HD13	1:A:99:TYR:CE1	2.49	0.48
1:A:8:ALA:HB3	1:A:421:GLU:OE2	2.14	0.48
1:B:61:GLU:HA	1:B:61:GLU:OE1	2.14	0.48
1:B:57:LEU:HD13	1:B:99:TYR:CE1	2.49	0.48
1:B:86:GLY:O	1:B:87:GLU:HG2	2.13	0.48
1:A:86:GLY:O	1:A:87:GLU:HG2	2.13	0.48
1:A:244:ASP:HB3	1:A:247:LYS:HG3	1.96	0.48
1:A:198:ARG:HD2	1:A:359:ALA:HA	1.96	0.48
1:A:219:ALA:CB	1:A:240:ALA:HB1	2.44	0.47
1:A:7:PRO:CA	1:A:10:THR:HG23	2.44	0.47
1:A:209:ARG:O	1:A:212:GLN:HG3	2.14	0.47
1:A:374:ARG:CG	1:A:374:ARG:NH1	2.50	0.47
1:A:229:GLN:NE2	1:A:376:VAL:HA	2.29	0.47
1:B:384:PHE:HB2	1:B:408:VAL:HG12	1.97	0.47
1:B:333:GLY:O	1:B:335:THR:N	2.47	0.47
1:B:309:HIS:N	1:B:310:PRO:CD	2.76	0.47
1:B:319:THR:HA	1:B:328:ILE:HD13	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85:LEU:HA	1:B:85:LEU:HD12	1.69	0.46
1:B:286:ALA:O	1:B:289:PHE:HB3	2.16	0.46
1:A:12:GLN:N	1:A:12:GLN:OE1	2.49	0.46
1:B:20:TYR:CE2	1:B:26:PRO:HB3	2.51	0.46
1:A:274:VAL:O	1:A:275:ASN:HB2	2.15	0.45
1:A:9:TRP:CE2	1:A:420:VAL:HG11	2.52	0.45
1:A:85:LEU:HD12	1:A:85:LEU:HA	1.70	0.45
1:A:306:VAL:HG11	1:A:318:SER:HB3	1.97	0.45
1:A:329:VAL:HA	1:A:354:GLN:O	2.16	0.45
1:A:41:ARG:HD2	1:A:44:ARG:HH12	1.81	0.45
1:A:205:PRO:HD2	1:A:206:HIS:ND1	2.32	0.45
1:B:76:VAL:HG21	1:B:406:MET:HE2	1.99	0.45
1:B:244:ASP:HB3	1:B:247:LYS:HG3	1.98	0.45
1:A:20:TYR:CE2	1:A:26:PRO:HB3	2.51	0.45
1:A:384:PHE:HB2	1:A:408:VAL:HG12	1.98	0.45
1:B:242:ILE:CD1	1:B:259:VAL:HG11	2.46	0.45
1:A:7:PRO:CA	1:A:10:THR:CG2	2.93	0.45
1:A:42:ARG:HH22	1:A:93:ASP:CG	2.20	0.45
1:A:352:ARG:NH2	1:A:354:GLN:NE2	2.62	0.45
1:A:319:THR:HA	1:A:328:ILE:HD13	1.96	0.45
1:B:198:ARG:HD2	1:B:359:ALA:HA	1.98	0.45
1:A:312:GLU:HG3	1:A:338:ASN:O	2.17	0.45
1:B:199:MET:HE3	1:B:330:PHE:CA	2.47	0.44
1:A:199:MET:HE3	1:A:330:PHE:CA	2.48	0.44
1:A:389:ILE:HB	1:A:390:PRO:HD3	1.99	0.44
1:A:260:ILE:HG22	1:A:261:ASN:N	2.32	0.44
1:A:242:ILE:HD11	1:A:259:VAL:CG1	2.48	0.44
1:B:262:ARG:NH2	3:B:501:NAP:O1X	2.47	0.44
1:B:260:ILE:HG22	1:B:261:ASN:N	2.33	0.44
1:A:9:TRP:HE3	1:A:9:TRP:O	2.00	0.44
1:A:25:ILE:HG22	1:A:26:PRO:O	2.18	0.44
1:B:131:ASN:OD1	1:B:151:GLN:NE2	2.51	0.43
1:B:352:ARG:NH2	1:B:354:GLN:NE2	2.65	0.43
1:A:7:PRO:C	1:A:10:THR:HG23	2.39	0.43
1:B:229:GLN:NE2	1:B:376:VAL:HA	2.34	0.43
1:B:177:ARG:HD3	1:B:178:PRO:O	2.19	0.43
1:A:317:VAL:O	1:A:321:VAL:HG23	2.18	0.43
1:B:417:LEU:HD23	1:B:423:VAL:HA	2.01	0.43
1:A:278:GLU:CD	1:A:278:GLU:H	2.08	0.43
1:B:297:THR:O	1:B:297:THR:CG2	2.66	0.43
1:B:205:PRO:HD2	1:B:206:HIS:ND1	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:242:ILE:CD1	1:A:259:VAL:HG11	2.48	0.43
1:A:237:ASN:OD1	1:A:257:LYS:HD2	2.18	0.43
1:B:242:ILE:HD11	1:B:259:VAL:CG1	2.49	0.43
1:A:249:ASP:N	1:A:249:ASP:OD1	2.51	0.43
1:A:252:MET:HA	1:A:256:ALA:HB3	2.01	0.42
1:A:20:TYR:CZ	1:A:26:PRO:HB3	2.53	0.42
1:B:9:TRP:O	1:B:9:TRP:HE3	2.03	0.42
1:A:243:SER:HB3	1:A:262:ARG:NE	2.33	0.42
1:A:31:VAL:HA	1:A:32:PRO:HD2	1.85	0.42
1:B:220:SER:HB2	1:B:403:PRO:HB3	2.02	0.42
1:B:389:ILE:O	1:B:393:HIS:HD2	2.02	0.42
1:B:270:GLN:O	1:B:272:PRO:HD3	2.20	0.42
1:A:7:PRO:O	1:A:10:THR:HG23	2.20	0.42
1:A:21:GLU:CG	1:A:430:LYS:HA	2.48	0.42
1:B:293:ILE:HG22	1:B:294:TRP:N	2.34	0.42
1:B:79:ASN:OD1	1:B:80:GLY:N	2.53	0.42
1:B:352:ARG:NE	1:B:354:GLN:NE2	2.51	0.41
1:A:142:GLY:HA3	1:B:360:HIS:ND1	2.34	0.41
1:B:148:SER:HA	1:B:149:PRO:HD3	1.96	0.41
1:B:67:VAL:CG1	1:B:68:LEU:N	2.82	0.41
1:A:151:GLN:O	1:A:151:GLN:HG3	2.19	0.41
1:A:414:ARG:HD2	1:A:415:ALA:H	1.85	0.41
1:A:315:PHE:N	1:A:316:PRO:HD2	2.34	0.41
1:A:429:LEU:HD12	1:A:430:LYS:H	1.83	0.41
1:A:44:ARG:O	1:A:45:HIS:C	2.59	0.41
1:A:237:ASN:HB3	1:A:257:LYS:HD3	2.03	0.41
1:B:219:ALA:CB	1:B:240:ALA:HB1	2.51	0.41
1:A:9:TRP:CE3	1:A:9:TRP:O	2.73	0.41
1:B:398:LYS:CG	1:B:398:LYS:O	2.67	0.41
1:A:17:LYS:HG3	1:A:20:TYR:CZ	2.56	0.41
1:A:353:ILE:HG22	1:A:353:ILE:O	2.20	0.41
1:B:312:GLU:HG3	1:B:338:ASN:O	2.21	0.41
1:A:341:PHE:CD1	1:A:341:PHE:C	2.94	0.41
1:A:398:LYS:HD2	1:A:398:LYS:N	2.28	0.41
1:B:266:ASP:C	1:B:268:TRP:N	2.73	0.41
1:A:47:PRO:HA	1:A:48:PRO:HD3	1.96	0.41
1:B:237:ASN:OD1	1:B:257:LYS:HD2	2.21	0.40
1:A:304:ASP:HA	1:A:323:LYS:HG2	2.03	0.40
1:B:151:GLN:HE21	1:B:151:GLN:HB2	1.73	0.40
1:B:341:PHE:CD1	1:B:341:PHE:C	2.94	0.40
1:B:252:MET:HE2	1:B:259:VAL:HG21	2.01	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:262:ARG:HH21	3:B:501:NAP:P2B	2.43	0.40
1:B:12:GLN:N	1:B:12:GLN:OE1	2.53	0.40
1:A:139:GLU:OE1	1:B:362:LYS:NZ	2.54	0.40
1:A:417:LEU:HD23	1:A:423:VAL:HA	2.03	0.40
1:A:178:PRO:HA	1:A:372:MET:HE1	2.01	0.40
1:B:389:ILE:HB	1:B:390:PRO:HD3	2.04	0.40
1:B:90:SER:HA	1:B:91:PRO:HD3	1.85	0.40
1:A:276:SER:HB3	1:A:278:GLU:CD	2.42	0.40
1:B:398:LYS:HD2	1:B:398:LYS:N	2.28	0.40
1:B:9:TRP:O	1:B:9:TRP:CE3	2.75	0.40
1:A:417:LEU:HD23	1:A:423:VAL:HG22	2.03	0.40
1:A:172:ARG:HD3	1:A:172:ARG:HH11	1.77	0.40
1:A:138:GLU:HB3	1:B:362:LYS:HG3	2.04	0.40
1:A:40:ILE:HD13	1:A:84:GLY:HA2	2.03	0.40

All (1) 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:93:ASP:O	1:B:433:ALA:O[4_455]	1.83	0.37

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	426/443 (96%)	397 (93%)	22 (5%)	7 (2%)	12	48
1	B	424/443 (96%)	392 (92%)	25 (6%)	7 (2%)	11	46
All	All	850/886 (96%)	789 (93%)	47 (6%)	14 (2%)	12	48

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	267	CYS
1	A	309	HIS
1	A	397	TRP
1	B	267	CYS
1	B	309	HIS
1	B	397	TRP
1	A	334	THR
1	B	433	ALA
1	B	334	THR
1	A	270	GLN
1	B	270	GLN
1	A	358	PHE
1	A	410	VAL
1	B	410	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	341/351 (97%)	318 (93%)	23 (7%)	20	57
1	B	340/351 (97%)	316 (93%)	24 (7%)	18	54
All	All	681/702 (97%)	634 (93%)	47 (7%)	19	56

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

Mol	Chain	Res	Type
1	A	21	GLU
1	A	65	ASP
1	A	85	LEU
1	A	146	MET
1	A	151	GLN
1	A	152	ARG
1	A	177	ARG
1	A	223	LEU
1	A	249	ASP
1	A	262	ARG

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Mol	Chain	Res	Type
1	A	265	PHE
1	A	266	ASP
1	A	271	LEU
1	A	275	ASN
1	A	291	LYS
1	A	315	PHE
1	A	321	VAL
1	A	323	LYS
1	A	347	TRP
1	A	374	ARG
1	A	387	ASP
1	A	398	LYS
1	A	414	ARG
1	B	21	GLU
1	B	65	ASP
1	B	85	LEU
1	B	146	MET
1	B	151	GLN
1	B	152	ARG
1	B	177	ARG
1	B	223	LEU
1	B	253	SER
1	B	265	PHE
1	B	266	ASP
1	B	271	LEU
1	B	275	ASN
1	B	291	LYS
1	B	315	PHE
1	B	321	VAL
1	B	323	LYS
1	B	347	TRP
1	B	372	MET
1	B	374	ARG
1	B	387	ASP
1	B	398	LYS
1	B	414	ARG
1	B	432	MET

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

Mol	Chain	Res	Type
1	A	131	ASN

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Mol	Chain	Res	Type
1	A	151	GLN
1	A	229	GLN
1	A	275	ASN
1	A	309	HIS
1	A	354	GLN
1	A	400	GLN
1	B	52	HIS
1	B	131	ASN
1	B	151	GLN
1	B	229	GLN
1	B	275	ASN
1	B	309	HIS
1	B	354	GLN
1	B	400	GLN

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 carbohydrates 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	NAP	A	502	-	42,52,52	1.85	11 (26%)	54,80,80	2.68	8 (14%)
3	NAP	B	501	-	42,52,52	1.82	10 (23%)	54,80,80	4.08	15 (27%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAP	A	502	-	-	0/27/67/67	0/5/5/5
3	NAP	B	501	-	-	0/27/67/67	0/5/5/5

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	NAP	C5A-C4A	-3.35	1.32	1.40
3	A	502	NAP	C3B-C2B	-3.34	1.45	1.53
3	A	502	NAP	C2D-C3D	-2.96	1.45	1.53
3	B	501	NAP	C5A-N7A	-2.88	1.29	1.39
3	B	501	NAP	C3B-C2B	-2.82	1.46	1.53
3	A	502	NAP	O2D-C2D	-2.54	1.36	1.43
3	B	501	NAP	C2D-C3D	-2.53	1.46	1.53
3	A	502	NAP	C5A-N7A	-2.43	1.31	1.39
3	A	502	NAP	O4B-C1B	-2.30	1.38	1.41
3	B	501	NAP	O2B-C2B	-2.13	1.37	1.44
3	A	502	NAP	O2B-C2B	-2.10	1.37	1.44
3	B	501	NAP	O2D-C2D	-2.08	1.38	1.43
3	A	502	NAP	C3D-C4D	-2.02	1.47	1.53
3	A	502	NAP	C4N-C3N	2.04	1.42	1.39
3	B	501	NAP	C6A-N6A	2.29	1.42	1.34
3	A	502	NAP	C6A-N6A	2.44	1.42	1.34
3	B	501	NAP	C6N-C5N	2.84	1.45	1.38
3	B	501	NAP	C4N-C3N	3.31	1.45	1.39
3	B	501	NAP	C5N-C4N	4.15	1.47	1.38
3	B	501	NAP	C7N-N7N	5.69	1.44	1.33
3	A	502	NAP	C7N-N7N	6.09	1.45	1.33

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	NAP	N3A-C2A-N1A	-12.22	119.54	128.89
3	B	501	NAP	C2N-C3N-C4N	-10.32	106.80	118.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	NAP	C4B-O4B-C1B	-7.75	101.20	109.72
3	B	501	NAP	C4B-O4B-C1B	-7.53	101.45	109.72
3	A	502	NAP	C4D-O4D-C1D	-7.04	101.99	109.72
3	B	501	NAP	C2A-N1A-C6A	-5.76	108.48	118.77
3	B	501	NAP	C4D-O4D-C1D	-5.23	103.97	109.72
3	A	502	NAP	PN-O3-PA	-4.91	118.95	132.73
3	B	501	NAP	PN-O3-PA	-4.52	120.03	132.73
3	B	501	NAP	O7N-C7N-N7N	-2.15	119.57	122.59
3	B	501	NAP	P2B-O2B-C2B	2.09	126.58	121.56
3	B	501	NAP	C2B-C3B-C4B	2.50	107.77	101.85
3	A	502	NAP	O4D-C1D-N1N	2.66	111.05	108.13
3	B	501	NAP	O3-PN-O5D	2.81	110.40	102.94
3	A	502	NAP	C4A-C5A-N7A	3.01	112.24	109.48
3	A	502	NAP	O4B-C4B-C3B	3.16	111.51	105.15
3	B	501	NAP	C4A-C5A-N7A	3.22	112.44	109.48
3	B	501	NAP	C4N-C3N-C7N	3.66	130.75	121.09
3	B	501	NAP	O3-PA-O5B	4.23	114.15	102.94
3	B	501	NAP	O4B-C1B-C2B	5.29	116.17	106.60
3	A	502	NAP	O4B-C1B-C2B	6.22	117.85	106.60
3	B	501	NAP	N3A-C2A-N1A	15.38	140.67	128.89
3	B	501	NAP	C3N-C2N-N1N	17.11	140.08	120.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	NAP	2	0
3	B	501	NAP	3	0

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	428/443 (96%)	-0.52	2 (0%) 91 76	55, 84, 132, 156	0
1	B	426/443 (96%)	-0.38	3 (0%) 89 70	53, 83, 128, 157	0
All	All	854/886 (96%)	-0.45	5 (0%) 90 73	53, 84, 130, 157	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	11	GLY	3.7
1	B	15	GLU	2.8
1	A	243	SER	2.8
1	B	299	LYS	2.6
1	A	263	LYS	2.4

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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	MG	A	501	1/1	0.86	0.28	5.33	85,85,85,85	0
3	NAP	B	501	48/48	0.92	0.15	-0.45	75,101,130,187	0
3	NAP	A	502	48/48	0.94	0.13	-0.61	54,106,146,267	0

6.5 Other polymers ⓘ

There are no such residues in this entry.