Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 07:37 PM GMT

PDB ID	:	1GIQ
Title	:	Crystal Structure of the Enzymatic Componet of Iota-Toxin from Clostridium Perfringens with NADH
Authors	:	Tsuge, H.; Nagahama, M.; Nishimura, H.; Hisatsune, J.; Sakaguchi, Y.; Itogawa, Y.; Katunuma, N.; Sakurai, J.
Deposited on	:	2001-03-12
Resolution	:	1.80 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 (i) symbol.

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

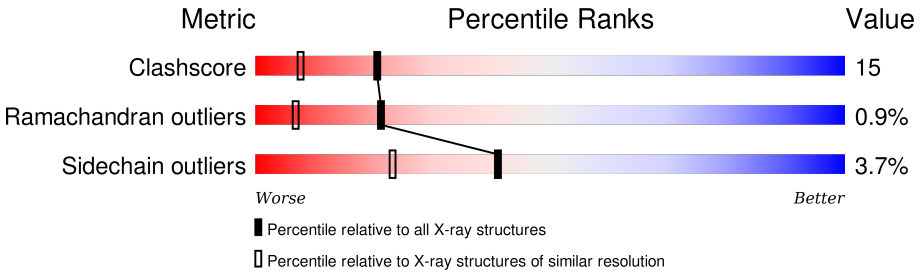
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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 1.80 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))
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)

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 <=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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	413	70% 27% .
1	B	413	74% 24% .

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	NAI	A	500	X	-	-	-
2	NAI	B	1500	X	-	-	-

2 Entry composition

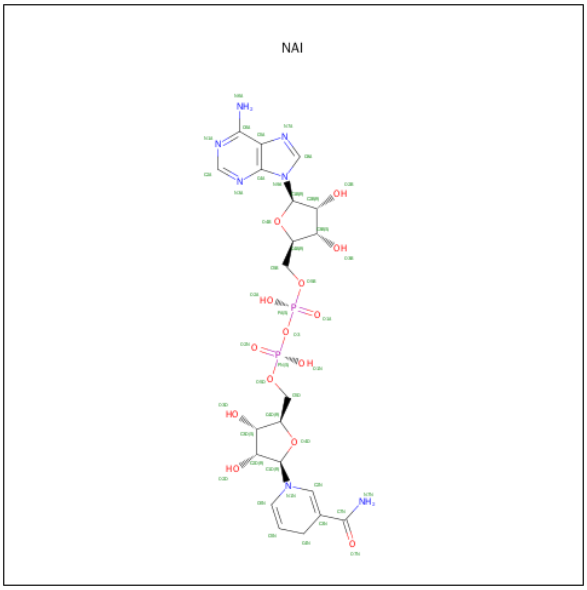
There are 3 unique types of molecules in this entry. The entry contains 7068 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 IOTA TOXIN COMPONENT IA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	411	Total	C	N	O	S	0	0	0
			3347	2134	551	659	3			
1	B	411	Total	C	N	O	S	0	0	0
			3347	2134	551	659	3			

- Molecule 2 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (three-letter code: NAI) (formula: $C_{21}H_{29}N_7O_{14}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	154	Total O 154 154	0	0
3	B	132	Total O 132 132	0	0

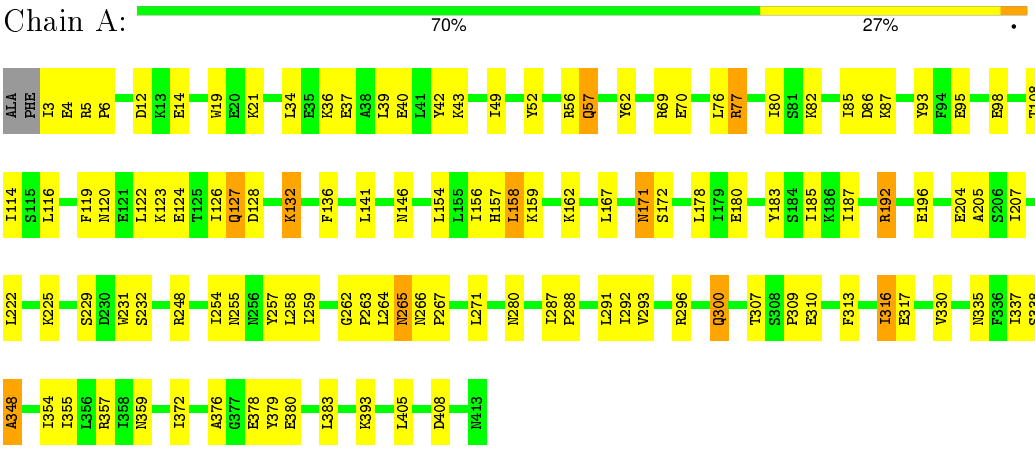
3 Residue-property plots (i)

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.

Note EDS was not executed.

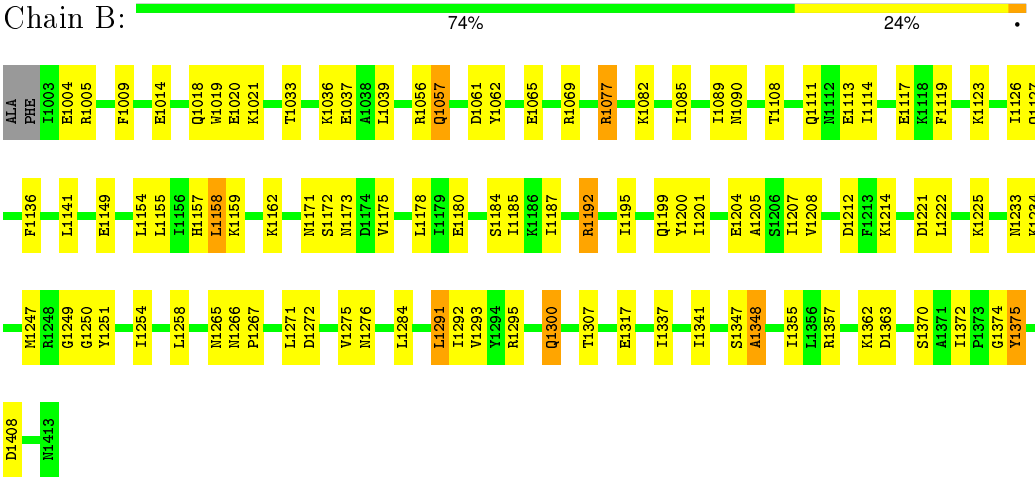
• Molecule 1: IOTA TOXIN COMPONENT IA

Chain A:



• Molecule 1: IOTA TOXIN COMPONENT IA

Chain B:



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	48.20Å 54.44Å 103.38Å 99.20° 93.10° 106.90°	Depositor
Resolution (Å)	14.97 – 1.80	Depositor
% Data completeness (in resolution range)	91.0 (14.97-1.80)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.224 , 0.250	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7068	wwPDB-VP
Average B, all atoms (Å^2)	30.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.30	0/3415	0.59	0/4614
1	B	0.31	0/3415	0.59	0/4614
All	All	0.30	0/6830	0.59	0/9228

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	3347	0	3329	108	0
1	B	3347	0	3329	89	0
2	A	44	0	26	0	0
2	B	44	0	26	0	0
3	A	154	0	0	11	0
3	B	132	0	0	6	0
All	All	7068	0	6710	196	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 15.

All (196) 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:338:SER:HB2	3:A:4246:HOH:O	1.66	0.95
1:A:56:ARG:HD2	1:A:70:GLU:OE1	1.74	0.87
1:B:1108:THR:HG22	1:B:1111:GLN:HG3	1.58	0.85
1:A:316:ILE:H	1:A:316:ILE:HD13	1.44	0.80
1:B:1132:LYS:HD3	1:B:1357:ARG:NH1	1.96	0.80
1:A:57:GLN:HG2	1:A:214:LYS:HD3	1.68	0.75
1:A:108:THR:HG22	1:A:111:GLN:HG3	1.68	0.75
1:A:14:GLU:CD	1:A:14:GLU:H	1.90	0.74
1:B:1085:ILE:HG13	1:B:1162:LYS:O	1.88	0.74
1:B:1300:GLN:H	1:B:1300:GLN:NE2	1.87	0.73
1:A:380:GLU:HG3	3:A:4246:HOH:O	1.88	0.72
1:B:1184:SER:HB2	1:B:1208:VAL:HG13	1.71	0.71
1:A:300:GLN:H	1:A:300:GLN:NE2	1.87	0.71
1:B:1212:ASP:HB2	1:B:1292:ILE:HD12	1.73	0.71
1:B:1337:ILE:CG2	1:B:1383:LEU:HB2	2.22	0.70
1:B:1250:GLY:O	1:B:1254:ILE:HD13	1.93	0.69
1:B:1126:ILE:HD12	1:B:1187:ILE:HD13	1.73	0.69
1:A:291:LEU:HD13	1:A:292:ILE:N	2.07	0.68
1:A:291:LEU:CD1	1:A:293:VAL:HG13	2.22	0.68
1:A:347:SER:O	1:A:348:ALA:HB3	1.93	0.68
1:B:1119:PHE:HB2	1:B:1192:ARG:NH2	2.07	0.68
1:B:1347:SER:O	1:B:1348:ALA:HB3	1.91	0.68
1:B:1005:ARG:NH1	1:B:1089:ILE:HD12	2.09	0.68
1:B:1014:GLU:H	1:B:1014:GLU:CD	1.97	0.68
1:A:254:ILE:O	1:A:258:LEU:HD23	1.95	0.67
1:B:1057:GLN:HG2	1:B:1214:LYS:HD3	1.77	0.67
1:A:162:LYS:HG3	3:A:4127:HOH:O	1.95	0.66
1:A:291:LEU:HD12	1:A:293:VAL:HG13	1.77	0.66
1:A:36:LYS:HB3	1:A:36:LYS:NZ	2.09	0.66
1:A:85:ILE:HG13	1:A:162:LYS:O	1.94	0.66
1:B:1291:LEU:HD23	1:B:1292:ILE:N	2.10	0.66
1:B:1234:LYS:HB2	1:B:1284:LEU:HD21	1.78	0.66
1:A:5:ARG:HD3	1:A:19:TRP:CZ2	2.31	0.66
1:B:1337:ILE:HG22	1:B:1383:LEU:HB2	1.78	0.65
1:B:1133:GLN:HG3	1:B:1185:ILE:HD12	1.76	0.65
1:A:132:LYS:HD3	1:A:357:ARG:NH1	2.12	0.65
1:B:1347:SER:O	1:B:1348:ALA:CB	2.45	0.65
1:A:52:TYR:CZ	1:A:56:ARG:HG3	2.32	0.64
1:B:1005:ARG:HD3	1:B:1019:TRP:CZ2	2.33	0.64
1:A:254:ILE:HD13	1:A:271:LEU:HD22	1.80	0.64
1:A:132:LYS:HG3	1:A:408:ASP:OD1	1.98	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:248:ARG:O	1:A:248:ARG:HG2	1.98	0.64
1:B:1077:ARG:HG2	1:B:1180:GLU:OE1	1.98	0.63
1:B:1108:THR:CG2	1:B:1111:GLN:HG3	2.29	0.62
1:A:157:HIS:HD2	1:A:204:GLU:OE2	1.82	0.61
1:B:1341:ILE:HG22	1:B:1379:TYR:HB3	1.82	0.61
1:A:337:ILE:HG23	1:A:383:LEU:HB2	1.83	0.61
1:B:1108:THR:HG22	1:B:1111:GLN:CG	2.30	0.61
1:B:1295:ARG:NE	1:B:1337:ILE:HD11	2.16	0.61
1:B:1127:GLN:HG2	1:B:1128:ASP:OD1	2.01	0.60
1:A:108:THR:CG2	1:A:111:GLN:HG3	2.32	0.60
1:A:36:LYS:O	1:A:40:GLU:HG2	2.02	0.60
1:A:231:TRP:HZ3	1:A:372:ILE:HD11	1.67	0.60
1:B:1126:ILE:O	1:B:1126:ILE:HG13	2.01	0.60
1:B:1039:LEU:O	1:B:1039:LEU:HD23	2.02	0.59
1:A:183:TYR:CD2	1:A:207:ILE:HD12	2.37	0.59
1:A:185:ILE:CG2	1:A:205:ALA:HB1	2.33	0.58
1:B:1033:THR:O	1:B:1037:GLU:HG2	2.03	0.58
1:B:1085:ILE:O	1:B:1162:LYS:O	2.21	0.58
1:B:1184:SER:HB2	1:B:1208:VAL:CG1	2.32	0.58
1:B:1221:ASP:O	1:B:1225:LYS:HG2	2.03	0.58
1:A:77:ARG:HG2	1:A:180:GLU:OE1	2.03	0.58
1:B:1157:HIS:HD2	1:B:1204:GLU:OE2	1.87	0.58
1:A:52:TYR:OH	1:A:56:ARG:HG3	2.04	0.57
1:A:93:TYR:CE1	1:A:156:ILE:HD12	2.40	0.57
1:A:330:VAL:HG22	3:A:4241:HOH:O	2.04	0.57
1:B:1005:ARG:HG3	1:B:1009:PHE:CZ	2.40	0.57
1:A:162:LYS:HB3	3:A:4236:HOH:O	2.05	0.57
1:A:376:ALA:O	1:A:378:GLU:HG3	2.06	0.56
1:B:1355:ILE:N	1:B:1355:ILE:HD12	2.20	0.56
1:B:1158:LEU:HD21	1:B:1207:ILE:HG13	1.87	0.55
1:A:80:ILE:HD11	1:A:178:LEU:HD23	1.89	0.55
1:A:337:ILE:CG2	1:A:383:LEU:HB2	2.36	0.55
1:B:1033:THR:HG21	3:B:4185:HOH:O	2.07	0.55
1:A:316:ILE:HG12	1:A:317:GLU:OE2	2.07	0.55
1:A:76:LEU:O	1:A:80:ILE:HG12	2.07	0.55
1:A:347:SER:O	1:A:348:ALA:CB	2.55	0.55
1:A:127:GLN:HG2	1:A:128:ASP:OD1	2.07	0.54
1:A:85:ILE:O	1:A:162:LYS:O	2.25	0.54
1:B:1295:ARG:CD	1:B:1337:ILE:HD11	2.37	0.53
1:A:132:LYS:HE3	1:A:408:ASP:HB2	1.89	0.53
1:A:108:THR:HG22	1:A:111:GLN:CG	2.39	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1005:ARG:HD3	1:B:1019:TRP:CE2	2.43	0.53
1:B:1132:LYS:HG3	1:B:1408:ASP:OD1	2.08	0.53
1:A:259:ILE:HD11	1:A:335:ASN:OD1	2.10	0.52
1:B:1266:ASN:N	1:B:1267:PRO:HD3	2.25	0.52
1:B:1126:ILE:HB	1:B:1187:ILE:HD12	1.91	0.52
1:B:1132:LYS:HD2	3:B:4073:HOH:O	2.10	0.51
1:B:1272:ASP:O	1:B:1275:VAL:HG22	2.10	0.51
1:A:316:ILE:N	1:A:316:ILE:HD13	2.19	0.51
1:A:126:ILE:HD12	1:A:187:ILE:HD13	1.91	0.51
1:A:355:ILE:N	1:A:355:ILE:HD12	2.25	0.51
1:B:1291:LEU:HD22	1:B:1293:VAL:HG13	1.93	0.51
1:B:1119:PHE:HB2	1:B:1192:ARG:HH22	1.74	0.51
1:A:86:ASP:C	1:A:87:LYS:HD2	2.31	0.50
1:A:221:ASP:O	1:A:225:LYS:HG2	2.11	0.50
1:B:1020:GLU:HG2	1:B:1175:VAL:HG11	1.94	0.50
1:B:1014:GLU:O	1:B:1018:GLN:HG3	2.12	0.50
1:A:220:GLY:C	1:A:341:ILE:HD12	2.33	0.49
1:B:1275:VAL:HG23	1:B:1276:ASN:N	2.28	0.49
1:A:232:SER:HB2	3:A:4221:HOH:O	2.12	0.49
1:A:292:ILE:HD13	1:A:359:ASN:HA	1.95	0.49
1:A:95:GLU:HG3	1:A:156:ILE:HD12	1.95	0.49
1:A:119:PHE:CD2	1:A:192:ARG:HD2	2.48	0.49
1:A:108:THR:HG23	1:A:111:GLN:H	1.78	0.48
1:A:264:LEU:O	1:A:265:ASN:HB2	2.12	0.48
1:A:291:LEU:HD13	1:A:291:LEU:C	2.33	0.48
1:A:316:ILE:H	1:A:316:ILE:CD1	2.19	0.48
1:B:1132:LYS:HE3	1:B:1408:ASP:HB2	1.94	0.48
1:A:21:LYS:HB3	1:A:21:LYS:NZ	2.29	0.48
1:A:296:ARG:HG3	1:A:296:ARG:HH11	1.79	0.47
1:B:1089:ILE:HD11	1:B:1141:LEU:HG	1.96	0.47
1:B:1062:TYR:HA	1:B:1065:GLU:HG2	1.97	0.47
1:A:158:LEU:HD21	1:A:207:ILE:CD1	2.45	0.47
1:B:1271:LEU:O	1:B:1275:VAL:HG13	2.15	0.47
1:B:1247:MET:SD	1:B:1372:ILE:HD12	2.55	0.47
1:A:126:ILE:HB	1:A:187:ILE:HD12	1.96	0.47
1:A:5:ARG:HG2	1:A:6:PRO:O	2.15	0.46
1:A:116:LEU:C	1:A:116:LEU:HD13	2.36	0.46
1:A:122:LEU:O	1:A:126:ILE:HG12	2.15	0.46
1:A:86:ASP:OD2	1:A:87:LYS:HD3	2.16	0.46
1:A:120:ASN:O	1:A:124:GLU:HG3	2.15	0.46
1:A:40:GLU:HA	1:A:43:LYS:HD3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1275:VAL:HG21	3:B:4242:HOH:O	2.15	0.46
1:B:1136:PHE:HB3	1:B:1178:LEU:HD11	1.97	0.46
1:A:171:ASN:HD22	1:A:171:ASN:C	2.19	0.46
1:B:1236:THR:OG1	1:B:1239:GLU:HG3	2.15	0.46
1:B:1254:ILE:O	1:B:1258:LEU:CD2	2.64	0.46
1:B:1113:GLU:HG2	1:B:1199:GLN:NE2	2.30	0.46
1:B:1254:ILE:O	1:B:1258:LEU:HD23	2.15	0.46
1:A:266:ASN:N	1:A:267:PRO:HD3	2.30	0.46
1:A:159:LYS:HG3	3:A:4034:HOH:O	2.16	0.45
1:B:1337:ILE:O	1:B:1337:ILE:HG23	2.16	0.45
1:A:291:LEU:HD13	1:A:292:ILE:C	2.36	0.45
1:A:136:PHE:HB3	1:A:178:LEU:HD11	1.98	0.45
1:A:123:LYS:HE3	3:A:4233:HOH:O	2.16	0.45
1:A:39:LEU:O	1:A:39:LEU:HD23	2.16	0.45
1:A:114:ILE:O	1:A:192:ARG:NH2	2.43	0.45
1:A:4:GLU:HG3	1:A:4:GLU:O	2.17	0.45
1:A:300:GLN:H	1:A:300:GLN:CD	2.20	0.45
1:B:1362:LYS:O	1:B:1363:ASP:HB2	2.17	0.45
1:B:1249:GLY:HA2	1:B:1251:TYR:CE1	2.52	0.45
1:B:1222:LEU:HD13	1:B:1222:LEU:C	2.37	0.45
1:B:1225:LYS:HE2	1:B:1370:SER:O	2.16	0.45
1:B:1185:ILE:CG2	1:B:1205:ALA:HB1	2.46	0.45
1:B:1208:VAL:HG13	1:B:1208:VAL:O	2.17	0.44
1:B:1192:ARG:HE	1:B:1192:ARG:HB3	1.62	0.44
1:A:36:LYS:HB3	1:A:36:LYS:HZ3	1.82	0.44
1:A:36:LYS:HB3	1:A:36:LYS:HZ2	1.81	0.44
1:A:158:LEU:HD21	1:A:207:ILE:HG12	1.99	0.44
1:A:313:PHE:CE2	1:A:354:ILE:HD11	2.52	0.44
1:B:1172:SER:OG	1:B:1173:ASN:N	2.47	0.44
1:B:1123:LYS:O	1:B:1127:GLN:HB2	2.18	0.44
1:A:225:LYS:O	1:A:229:SER:HB3	2.18	0.44
1:A:34:LEU:O	1:A:37:GLU:HB2	2.17	0.44
1:A:82:LYS:HB3	3:A:4164:HOH:O	2.16	0.44
1:B:1082:LYS:HB3	3:B:4091:HOH:O	2.18	0.44
1:A:317:GLU:CD	1:A:317:GLU:H	2.21	0.43
1:A:255:ASN:O	1:A:259:ILE:HG12	2.18	0.43
1:B:1061:ASP:OD1	1:B:1062:TYR:N	2.50	0.43
1:B:1090:ASN:ND2	1:B:1159:LYS:HB3	2.33	0.43
1:A:132:LYS:HD2	3:A:4084:HOH:O	2.18	0.43
1:A:86:ASP:O	1:A:87:LYS:HD2	2.19	0.43
1:B:1114:ILE:CG2	1:B:1192:ARG:NH2	2.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:GLN:NE2	1:B:1233:ASN:HD21	2.16	0.43
1:A:341:ILE:HG12	1:A:379:TYR:HB3	2.00	0.43
1:B:1254:ILE:N	1:B:1254:ILE:HD12	2.34	0.43
1:A:172:SER:N	3:A:4108:HOH:O	2.52	0.43
1:B:1195:ILE:HB	1:B:1200:TYR:CE2	2.54	0.42
1:B:1291:LEU:CD2	1:B:1292:ILE:N	2.81	0.42
1:A:146:ASN:HA	1:A:196:GLU:OE2	2.19	0.42
1:A:307:THR:O	1:A:307:THR:HG22	2.19	0.42
1:B:1192:ARG:HH21	1:B:1201:ILE:HD11	1.84	0.42
1:B:1275:VAL:CG2	1:B:1276:ASN:N	2.83	0.42
1:A:132:LYS:HG3	1:A:408:ASP:CG	2.40	0.42
1:A:287:ILE:HA	1:A:288:PRO:HD3	1.80	0.42
1:B:1374:GLY:O	1:B:1375:TYR:HB2	2.20	0.42
1:A:12:ASP:OD2	1:A:14:GLU:HG2	2.20	0.42
1:A:119:PHE:CE2	1:A:192:ARG:HD2	2.55	0.41
1:A:257:TYR:CE2	1:A:262:GLY:HA3	2.55	0.41
1:A:3:ILE:HD13	1:A:87:LYS:HE3	2.00	0.41
1:A:262:GLY:N	1:A:263:PRO:CD	2.84	0.41
1:B:1375:TYR:HA	3:B:4113:HOH:O	2.21	0.41
1:B:1036:LYS:HD3	3:B:4244:HOH:O	2.21	0.41
1:A:42:TYR:HA	1:A:49:ILE:HD12	2.01	0.41
1:A:80:ILE:HD13	1:A:167:LEU:HB2	2.02	0.41
1:B:1119:PHE:CD2	1:B:1192:ARG:NH2	2.89	0.41
1:B:1222:LEU:HD13	1:B:1222:LEU:O	2.21	0.41
1:A:108:THR:HG23	1:A:110:ASN:H	1.86	0.41
1:B:1307:THR:HG22	1:B:1307:THR:O	2.20	0.41
1:B:1317:GLU:H	1:B:1317:GLU:CD	2.24	0.41
1:A:280:ASN:HA	1:A:280:ASN:HD22	1.73	0.40
1:B:1114:ILE:CG2	1:B:1192:ARG:HH22	2.34	0.40
1:A:309:PRO:HD2	1:A:310:GLU:OE1	2.21	0.40
1:A:98:GLU:O	1:A:393:LYS:HE2	2.20	0.40
1:B:1021:LYS:NZ	1:B:1021:LYS:HB3	2.36	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	409/413 (99%)	390 (95%)	16 (4%)	3 (1%)	26 11
1	B	409/413 (99%)	385 (94%)	20 (5%)	4 (1%)	19 5
All	All	818/826 (99%)	775 (95%)	36 (4%)	7 (1%)	21 7

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	265	ASN
1	B	1375	TYR
1	B	1004	GLU
1	B	1348	ALA
1	B	1265	ASN
1	A	127	GLN
1	A	348	ALA

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	375/376 (100%)	361 (96%)	14 (4%)	41 23
1	B	375/376 (100%)	361 (96%)	14 (4%)	41 23
All	All	750/752 (100%)	722 (96%)	28 (4%)	41 23

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	GLN
1	A	62	TYR
1	A	69	ARG
1	A	77	ARG
1	A	132	LYS

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Mol	Chain	Res	Type
1	A	141	LEU
1	A	154	LEU
1	A	158	LEU
1	A	171	ASN
1	A	192	ARG
1	A	222	LEU
1	A	300	GLN
1	A	316	ILE
1	A	405	LEU
1	B	1056	ARG
1	B	1057	GLN
1	B	1069	ARG
1	B	1077	ARG
1	B	1117	GLU
1	B	1132	LYS
1	B	1149	GLU
1	B	1154	LEU
1	B	1155	LEU
1	B	1158	LEU
1	B	1171	ASN
1	B	1192	ARG
1	B	1291	LEU
1	B	1300	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	ASN
1	A	48	GLN
1	A	51	ASN
1	A	54	GLN
1	A	78	ASN
1	A	90	ASN
1	A	120	ASN
1	A	127	GLN
1	A	157	HIS
1	A	171	ASN
1	A	227	ASN
1	A	244	ASN
1	A	266	ASN
1	A	280	ASN
1	A	300	GLN

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Mol	Chain	Res	Type
1	A	413	ASN
1	B	1015	ASN
1	B	1018	GLN
1	B	1051	ASN
1	B	1054	GLN
1	B	1078	ASN
1	B	1090	ASN
1	B	1120	ASN
1	B	1127	GLN
1	B	1157	HIS
1	B	1171	ASN
1	B	1233	ASN
1	B	1244	ASN
1	B	1266	ASN
1	B	1280	ASN
1	B	1300	GLN

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

There are no carbohydrates in this entry.

5.6 Ligand geometry

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	$\#\|Z\|>2$
2	NAI	A	500	-	38,48,48	1.57	6 (15%)	48,73,73	3.02	15 (31%)
2	NAI	B	1500	-	38,48,48	1.56	6 (15%)	48,73,73	3.03	15 (31%)

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
2	NAI	A	500	-	1/1/13/16	0/25/72/72	0/5/5/5
2	NAI	B	1500	-	1/1/13/16	0/25/72/72	0/5/5/5

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	NAI	C4N-C5N	-2.82	1.43	1.49
2	B	1500	NAI	C4N-C5N	-2.68	1.43	1.49
2	A	500	NAI	C5A-N7A	-2.33	1.31	1.39
2	B	1500	NAI	C5A-N7A	-2.24	1.31	1.39
2	B	1500	NAI	C2N-C3N	3.01	1.42	1.34
2	A	500	NAI	C2N-C3N	3.05	1.42	1.34
2	A	500	NAI	C2A-N3A	3.29	1.38	1.32
2	B	1500	NAI	C2A-N3A	3.33	1.38	1.32
2	A	500	NAI	C6N-N1N	3.68	1.48	1.37
2	B	1500	NAI	C6N-N1N	3.73	1.48	1.37
2	B	1500	NAI	C6N-C5N	4.93	1.42	1.33
2	A	500	NAI	C6N-C5N	4.98	1.43	1.33

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	500	NAI	N3A-C2A-N1A	-12.97	118.96	128.89
2	B	1500	NAI	N3A-C2A-N1A	-12.95	118.98	128.89
2	A	500	NAI	C4N-C5N-C6N	-3.82	116.28	122.58
2	B	1500	NAI	C4N-C5N-C6N	-3.77	116.37	122.58
2	A	500	NAI	C3N-C2N-N1N	-2.54	119.50	123.14
2	B	1500	NAI	C3N-C2N-N1N	-2.27	119.88	123.14
2	B	1500	NAI	O3-PN-O5D	2.01	108.28	102.94
2	A	500	NAI	O3-PN-O5D	2.06	108.40	102.94
2	A	500	NAI	O4D-C4D-C3D	2.09	109.37	105.15
2	B	1500	NAI	O4D-C4D-C3D	2.16	109.49	105.15
2	B	1500	NAI	O4B-C4B-C3B	2.19	109.56	105.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	500	NAI	O5D-C5D-C4D	2.21	117.25	109.12
2	A	500	NAI	O4B-C4B-C3B	2.23	109.64	105.15
2	B	1500	NAI	N6A-C6A-N1A	2.33	124.21	119.20
2	A	500	NAI	N6A-C6A-N1A	2.34	124.22	119.20
2	B	1500	NAI	O5D-C5D-C4D	2.34	117.74	109.12
2	B	1500	NAI	O4D-C4D-C5D	2.46	118.14	109.32
2	A	500	NAI	O4D-C4D-C5D	2.50	118.27	109.32
2	A	500	NAI	PN-O3-PA	2.76	140.49	132.73
2	B	1500	NAI	PN-O3-PA	2.80	140.59	132.73
2	B	1500	NAI	C5N-C4N-C3N	4.63	125.28	112.52
2	A	500	NAI	C5N-C4N-C3N	4.67	125.40	112.52
2	A	500	NAI	C4A-C5A-N7A	4.93	114.01	109.48
2	B	1500	NAI	C4A-C5A-N7A	4.95	114.03	109.48
2	A	500	NAI	C2B-C1B-N9A	6.06	123.55	114.29
2	A	500	NAI	C1B-N9A-C4A	6.27	136.40	126.94
2	B	1500	NAI	C2B-C1B-N9A	6.31	123.94	114.29
2	B	1500	NAI	C1B-N9A-C4A	6.60	136.89	126.94
2	B	1500	NAI	O4B-C1B-N9A	7.24	123.26	108.10
2	A	500	NAI	O4B-C1B-N9A	7.40	123.58	108.10

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	500	NAI	C1B
2	B	1500	NAI	C1B

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.