



Full wwPDB X-ray Structure Validation Report i

Feb 28, 2014 – 01:16 PM GMT

PDB ID : 1S3N
Title : Structural and Functional Characterization of a Novel Archaeal Phosphodiesterase
Authors : Chen, S.; Busso, D.; Yakunin, A.F.; Kuznetsova, E.; Proudfoot, M.; Jancrick, J.; Kim, R.; Kim, S.-H.; Berkeley Structural Genomics Center (BSGC)
Deposited on : 2004-01-13
Resolution : 2.50 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

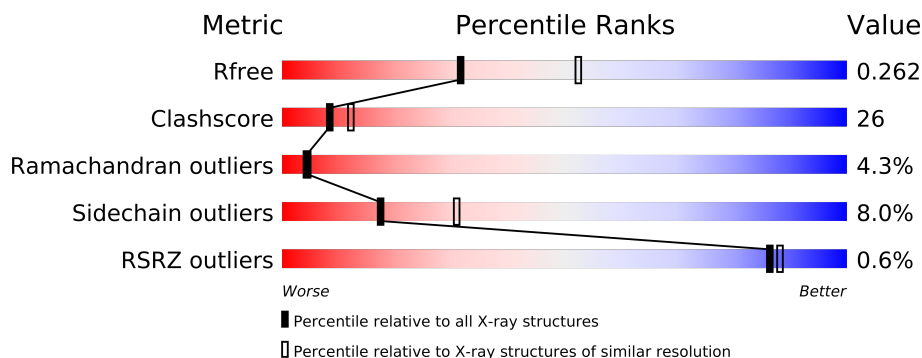
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.50 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}	66092	2784 (2.50-2.50)
Clashscore	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)
RSRZ outliers	66119	2785 (2.50-2.50)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	190	
1	B	190	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	MN	A	503	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2724 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Hypothetical protein MJ0936.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	165	Total	C	N	O	S	0	0	0
			1330	847	216	260	7			
1	B	165	Total	C	N	O	S	0	0	0
			1330	847	216	260	7			

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-24	MET	-	EXPRESSION TAG	UNP Q58346
A	-23	GLY	-	EXPRESSION TAG	UNP Q58346
A	-22	SER	-	EXPRESSION TAG	UNP Q58346
A	-21	SER	-	EXPRESSION TAG	UNP Q58346
A	-20	HIS	-	EXPRESSION TAG	UNP Q58346
A	-19	HIS	-	EXPRESSION TAG	UNP Q58346
A	-18	HIS	-	EXPRESSION TAG	UNP Q58346
A	-17	HIS	-	EXPRESSION TAG	UNP Q58346
A	-16	HIS	-	EXPRESSION TAG	UNP Q58346
A	-15	HIS	-	EXPRESSION TAG	UNP Q58346
A	-14	ASP	-	EXPRESSION TAG	UNP Q58346
A	-13	TYR	-	EXPRESSION TAG	UNP Q58346
A	-12	ASP	-	EXPRESSION TAG	UNP Q58346
A	-11	ILE	-	EXPRESSION TAG	UNP Q58346
A	-10	PRO	-	EXPRESSION TAG	UNP Q58346
A	-9	THR	-	EXPRESSION TAG	UNP Q58346
A	-8	THR	-	EXPRESSION TAG	UNP Q58346
A	-7	GLU	-	EXPRESSION TAG	UNP Q58346
A	-6	ASN	-	EXPRESSION TAG	UNP Q58346
A	-5	LEU	-	EXPRESSION TAG	UNP Q58346
A	-4	TYR	-	EXPRESSION TAG	UNP Q58346
A	-3	PHE	-	EXPRESSION TAG	UNP Q58346
A	-2	GLN	-	EXPRESSION TAG	UNP Q58346
A	-1	GLY	-	EXPRESSION TAG	UNP Q58346
A	0	HIS	-	EXPRESSION TAG	UNP Q58346

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Chain	Residue	Modelled	Actual	Comment	Reference
B	176	MET	-	EXPRESSION TAG	UNP Q58346
B	177	GLY	-	EXPRESSION TAG	UNP Q58346
B	178	SER	-	EXPRESSION TAG	UNP Q58346
B	179	SER	-	EXPRESSION TAG	UNP Q58346
B	180	HIS	-	EXPRESSION TAG	UNP Q58346
B	181	HIS	-	EXPRESSION TAG	UNP Q58346
B	182	HIS	-	EXPRESSION TAG	UNP Q58346
B	183	HIS	-	EXPRESSION TAG	UNP Q58346
B	184	HIS	-	EXPRESSION TAG	UNP Q58346
B	185	HIS	-	EXPRESSION TAG	UNP Q58346
B	186	ASP	-	EXPRESSION TAG	UNP Q58346
B	187	TYR	-	EXPRESSION TAG	UNP Q58346
B	188	ASP	-	EXPRESSION TAG	UNP Q58346
B	189	ILE	-	EXPRESSION TAG	UNP Q58346
B	190	PRO	-	EXPRESSION TAG	UNP Q58346
B	191	THR	-	EXPRESSION TAG	UNP Q58346
B	192	THR	-	EXPRESSION TAG	UNP Q58346
B	193	GLU	-	EXPRESSION TAG	UNP Q58346
B	194	ASN	-	EXPRESSION TAG	UNP Q58346
B	195	LEU	-	EXPRESSION TAG	UNP Q58346
B	196	TYR	-	EXPRESSION TAG	UNP Q58346
B	197	PHE	-	EXPRESSION TAG	UNP Q58346
B	198	GLN	-	EXPRESSION TAG	UNP Q58346
B	199	GLY	-	EXPRESSION TAG	UNP Q58346
B	200	HIS	-	EXPRESSION TAG	UNP Q58346

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	2	Total Mn 2 2	0	0
2	A	2	Total Mn 2 2	0	0

- Molecule 3 is water.

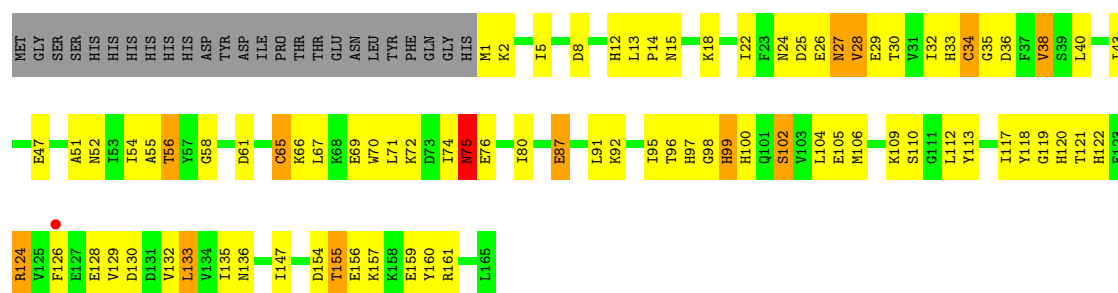
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	34	Total O 34 34	0	0
3	B	26	Total O 26 26	0	0

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.

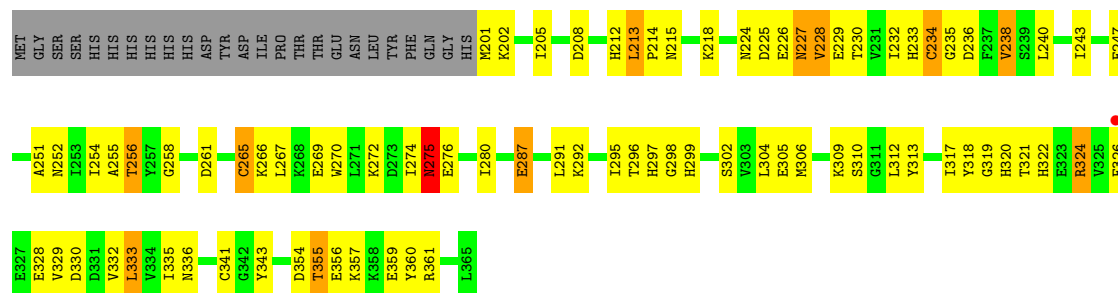
• Molecule 1: Hypothetical protein MJ0936

Chain A: 



• Molecule 1: Hypothetical protein MJ0936

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	I 41	Depositor
Cell constants a, b, c, α , β , γ	70.40Å 70.40Å 196.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.20 – 2.50 35.20 – 1.90	Depositor EDS
% Data completeness (in resolution range)	89.7 (35.20-2.50) 80.6 (35.20-1.90)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.12 (at 1.91Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.220 , 0.253 0.230 , 0.262	Depositor DCC
R_{free} test set	1560 reflections (9.88%)	DCC
Wilson B-factor (Å ²)	38.9	Xtriage
Anisotropy	0.336	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.6	EDS
Estimated twinning fraction	0.499 for -h,k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 36521 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2724	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.78% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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

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	0.65	0/1355	0.84	0/1832
1	B	0.65	0/1355	0.85	0/1832
All	All	0.65	0/2710	0.85	0/3664

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1330	0	1298	71	1
1	B	1330	0	1295	66	1
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	34	0	0	2	0
3	B	26	0	0	1	0
All	All	2724	0	2593	137	2

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 26.

All (137) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:208:ASP:H	1:B:235:GLY:HA3	1.32	0.95
1:A:8:ASP:H	1:A:35:GLY:HA3	1.32	0.95
1:B:324:ARG:HD2	1:B:324:ARG:O	1.77	0.84
1:A:124:ARG:O	1:A:124:ARG:HD2	1.78	0.83
1:B:230:THR:HG22	1:B:252:ASN:OD1	1.80	0.81
1:A:30:THR:HG22	1:A:52:ASN:OD1	1.80	0.80
1:A:36:ASP:HB3	1:A:38:VAL:HG23	1.64	0.78
1:B:236:ASP:HB3	1:B:238:VAL:HG23	1.64	0.77
1:A:2:LYS:HB2	1:A:28:VAL:H	1.53	0.73
1:B:202:LYS:HB2	1:B:228:VAL:H	1.52	0.73
1:A:106:MET:O	1:A:110:SER:HB3	1.90	0.71
1:B:306:MET:O	1:B:310:SER:HB3	1.90	0.70
1:B:202:LYS:O	1:B:228:VAL:HA	1.96	0.65
1:A:80:ILE:O	1:A:80:ILE:HG23	1.96	0.65
1:B:224:ASN:O	1:B:226:GLU:N	2.30	0.65
1:B:280:ILE:O	1:B:280:ILE:HG23	1.97	0.65
1:B:234:CYS:O	1:B:256:THR:HG22	1.97	0.64
1:A:154:ASP:OD1	1:A:157:LYS:N	2.31	0.64
1:A:1:MET:HA	1:A:29:GLU:OE1	1.98	0.64
1:A:24:ASN:O	1:A:26:GLU:N	2.31	0.64
1:A:70:TRP:O	1:A:74:ILE:HG12	1.99	0.63
1:B:354:ASP:OD1	1:B:357:LYS:N	2.31	0.63
1:A:34:CYS:O	1:A:56:THR:HG22	1.98	0.62
1:B:270:TRP:O	1:B:274:ILE:HG12	1.99	0.62
1:A:124:ARG:C	1:A:124:ARG:HD2	2.20	0.62
1:B:324:ARG:HD2	1:B:324:ARG:C	2.20	0.61
1:B:208:ASP:H	1:B:235:GLY:CA	2.11	0.61
1:A:36:ASP:OD2	1:A:58:GLY:HA3	2.00	0.60
1:A:2:LYS:O	1:A:28:VAL:HA	2.01	0.60
1:B:236:ASP:OD2	1:B:258:GLY:HA3	2.01	0.60
1:A:98:GLY:HA3	1:A:120:HIS:HB3	1.83	0.59
1:B:256:THR:HG21	1:B:297:HIS:N	2.17	0.59
1:A:56:THR:HG21	1:A:97:HIS:N	2.18	0.59
1:B:298:GLY:HA3	1:B:320:HIS:HB3	1.84	0.59
1:A:8:ASP:H	1:A:35:GLY:CA	2.11	0.58
1:B:321:THR:O	1:B:322:HIS:HB2	2.04	0.58
1:B:305:GLU:HG3	1:B:309:LYS:HE3	1.85	0.57
1:A:105:GLU:HG3	1:A:109:LYS:HE3	1.86	0.57
1:A:43:ILE:HD11	1:A:67:LEU:HD23	1.87	0.57
1:B:243:ILE:HD11	1:B:267:LEU:HD23	1.87	0.56
1:A:121:THR:O	1:A:122:HIS:HB2	2.05	0.56
1:B:240:LEU:HD21	1:B:266:LYS:HB3	1.88	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:12:HIS:CD2	1:A:15:ASN:H	2.24	0.55
1:B:212:HIS:CD2	1:B:215:ASN:H	2.24	0.55
1:A:40:LEU:HD21	1:A:66:LYS:HB3	1.89	0.55
1:B:224:ASN:C	1:B:226:GLU:H	2.10	0.54
1:A:54:ILE:N	1:A:54:ILE:HD12	2.23	0.54
1:B:274:ILE:O	1:B:275:ASN:HB2	2.08	0.54
1:B:254:ILE:HD12	1:B:254:ILE:N	2.23	0.54
1:A:24:ASN:C	1:A:26:GLU:H	2.11	0.53
1:A:74:ILE:O	1:A:75:ASN:HB2	2.08	0.52
1:B:243:ILE:HG21	1:B:270:TRP:HB3	1.92	0.51
1:B:272:LYS:O	1:B:275:ASN:N	2.40	0.51
1:B:296:THR:O	1:B:318:TYR:HA	2.11	0.51
1:A:8:ASP:N	1:A:35:GLY:HA3	2.12	0.51
1:B:202:LYS:HB2	1:B:227:ASN:HB3	1.93	0.51
1:B:208:ASP:N	1:B:235:GLY:HA3	2.12	0.51
1:A:96:THR:O	1:A:118:TYR:HA	2.10	0.50
1:A:43:ILE:HG21	1:A:70:TRP:HB3	1.93	0.50
1:A:12:HIS:CD2	1:A:14:PRO:HD2	2.47	0.50
1:B:265:CYS:O	1:B:269:GLU:HG3	2.11	0.50
1:A:2:LYS:HB2	1:A:27:ASN:HB3	1.94	0.49
1:A:72:LYS:HA	1:A:75:ASN:O	2.13	0.49
1:A:1:MET:O	1:A:155:THR:OG1	2.31	0.48
1:A:72:LYS:O	1:A:75:ASN:N	2.41	0.48
1:B:258:GLY:HA2	1:B:297:HIS:CG	2.48	0.48
1:B:212:HIS:CD2	1:B:214:PRO:HD2	2.48	0.48
1:A:65:CYS:O	1:A:69:GLU:HG3	2.12	0.48
1:A:1:MET:O	1:A:156:GLU:OE2	2.32	0.48
1:B:272:LYS:HA	1:B:275:ASN:O	2.13	0.48
1:A:58:GLY:HA2	1:A:97:HIS:CG	2.49	0.47
1:B:201:MET:O	1:B:355:THR:OG1	2.31	0.47
1:A:87:GLU:HA	1:A:91:LEU:O	2.14	0.47
1:B:330:ASP:C	1:B:332:VAL:H	2.18	0.47
1:A:12:HIS:HD2	1:A:15:ASN:H	1.63	0.47
1:B:232:ILE:HD13	1:B:295:ILE:CD1	2.44	0.47
1:B:287:GLU:HA	1:B:291:LEU:O	2.14	0.47
1:B:212:HIS:HD2	1:B:215:ASN:H	1.64	0.46
1:A:32:ILE:HD13	1:A:95:ILE:CD1	2.46	0.46
1:A:55:ALA:O	1:A:80:ILE:HA	2.16	0.46
1:A:75:ASN:HD22	1:A:76:GLU:N	2.14	0.46
1:A:130:ASP:C	1:A:132:VAL:H	2.18	0.46
1:B:224:ASN:C	1:B:226:GLU:N	2.70	0.45
1:B:218:LYS:NZ	3:B:601:HOH:O	2.39	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:275:ASN:HD22	1:B:276:GLU:N	2.15	0.45
1:B:229:GLU:O	1:B:251:ALA:HB1	2.16	0.45
1:A:24:ASN:C	1:A:26:GLU:N	2.70	0.45
1:B:201:MET:HA	1:B:229:GLU:OE1	2.16	0.45
1:B:255:ALA:O	1:B:280:ILE:HA	2.17	0.45
1:B:247:GLU:HA	1:B:274:ILE:HG23	1.98	0.45
1:A:110:SER:OG	1:A:112:LEU:HG	2.17	0.45
1:A:29:GLU:O	1:A:51:ALA:HB1	2.16	0.44
1:B:319:GLY:O	1:B:320:HIS:HB3	2.17	0.44
1:B:280:ILE:O	1:B:280:ILE:CG2	2.64	0.44
1:A:47:GLU:HA	1:A:74:ILE:HG23	1.98	0.44
1:B:354:ASP:HB3	1:B:359:GLU:O	2.17	0.44
1:B:232:ILE:HD13	1:B:295:ILE:HD13	1.99	0.44
1:A:154:ASP:HB3	1:A:159:GLU:O	2.17	0.44
1:A:1:MET:HB2	3:A:628:HOH:O	2.17	0.44
1:A:32:ILE:HD13	1:A:95:ILE:HD13	2.00	0.44
1:A:80:ILE:CG2	1:A:80:ILE:O	2.64	0.43
1:A:129:VAL:O	1:A:129:VAL:HG22	2.18	0.43
1:B:310:SER:OG	1:B:312:LEU:HG	2.18	0.43
1:B:326:PHE:CE1	1:B:360:TYR:HB3	2.54	0.43
1:B:329:VAL:O	1:B:329:VAL:HG22	2.18	0.43
1:B:304:LEU:HD11	1:B:318:TYR:CD2	2.54	0.43
1:A:126:PHE:CE1	1:A:160:TYR:HB3	2.54	0.43
1:A:104:LEU:HD11	1:A:118:TYR:CD2	2.54	0.43
1:A:119:GLY:O	1:A:120:HIS:HB3	2.20	0.42
1:B:326:PHE:CZ	1:B:360:TYR:HB3	2.54	0.42
1:A:126:PHE:CZ	1:A:160:TYR:HB3	2.54	0.42
1:A:5:ILE:HD13	1:A:117:ILE:HD13	2.02	0.42
1:B:233:HIS:O	1:B:255:ALA:HA	2.19	0.42
1:B:205:ILE:HD13	1:B:317:ILE:HD13	2.02	0.42
1:A:126:PHE:HD1	1:A:135:ILE:HG12	1.85	0.42
1:B:292:LYS:HB2	1:B:313:TYR:HA	2.02	0.42
1:A:154:ASP:HB3	1:A:159:GLU:H	1.85	0.42
1:A:33:HIS:O	1:A:55:ALA:HA	2.19	0.41
1:A:119:GLY:HA2	1:A:136:ASN:ND2	2.35	0.41
1:A:102:SER:OG	3:A:642:HOH:O	2.22	0.41
1:A:92:LYS:HB2	1:A:113:TYR:HA	2.02	0.41
1:B:213:LEU:HD12	1:B:213:LEU:HA	1.92	0.41
1:A:124:ARG:HH12	1:A:147:ILE:HG21	1.86	0.41
1:A:126:PHE:CE1	1:A:133:LEU:HD11	2.56	0.41
1:B:319:GLY:HA2	1:B:336:ASN:ND2	2.36	0.41
1:B:201:MET:O	1:B:356:GLU:OE2	2.39	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:326:PHE:HD1	1:B:335:ILE:HG12	1.86	0.41
1:A:99:HIS:HB3	1:A:100:HIS:H	1.78	0.41
1:A:43:ILE:HD12	1:A:71:LEU:HG	2.03	0.41
1:A:118:TYR:O	1:A:136:ASN:HA	2.21	0.41
1:A:98:GLY:N	1:A:120:HIS:HB2	2.36	0.40
1:B:354:ASP:HB3	1:B:359:GLU:H	1.86	0.40
1:A:18:LYS:O	1:A:22:ILE:HG13	2.21	0.40
1:B:298:GLY:N	1:B:320:HIS:HB2	2.36	0.40
1:B:326:PHE:CE1	1:B:333:LEU:HD11	2.56	0.40
1:A:129:VAL:O	1:A:130:ASP:HB2	2.22	0.40
1:B:341:CYS:HG	1:B:343:TYR:HD2	1.61	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:38:VAL:CG1	1:A:38:VAL:CG1[6_765]	1.88	0.32
1:B:238:VAL:CG1	1:B:238:VAL:CG1[6_765]	1.89	0.31

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	163/190 (86%)	142 (87%)	14 (9%)	7 (4%)	4	4
1	B	163/190 (86%)	141 (86%)	15 (9%)	7 (4%)	4	4
All	All	326/380 (86%)	283 (87%)	29 (9%)	14 (4%)	4	4

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	25	ASP
1	A	99	HIS
1	B	225	ASP
1	B	299	HIS

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Mol	Chain	Res	Type
1	A	28	VAL
1	B	228	VAL
1	A	27	ASN
1	B	227	ASN
1	A	38	VAL
1	B	261	ASP
1	A	61	ASP
1	A	75	ASN
1	B	238	VAL
1	B	275	ASN

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	150/173 (87%)	138 (92%)	12 (8%)	17	31
1	B	150/173 (87%)	138 (92%)	12 (8%)	17	31
All	All	300/346 (87%)	276 (92%)	24 (8%)	17	31

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LEU
1	A	34	CYS
1	A	56	THR
1	A	65	CYS
1	A	75	ASN
1	A	87	GLU
1	A	102	SER
1	A	124	ARG
1	A	128	GLU
1	A	133	LEU
1	A	155	THR
1	A	161	ARG
1	B	213	LEU
1	B	234	CYS
1	B	256	THR

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Mol	Chain	Res	Type
1	B	265	CYS
1	B	275	ASN
1	B	287	GLU
1	B	302	SER
1	B	324	ARG
1	B	328	GLU
1	B	333	LEU
1	B	355	THR
1	B	361	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	HIS
1	A	27	ASN
1	A	50	ASN
1	A	75	ASN
1	B	212	HIS
1	B	250	ASN
1	B	275	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	165/190 (86%)	-0.12	1 (0%) 86 88	35, 61, 85, 91	0
1	B	165/190 (86%)	-0.10	1 (0%) 86 88	34, 61, 86, 91	0
All	All	330/380 (86%)	-0.11	2 (0%) 86 88	34, 61, 86, 91	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	326	PHE	2.9
1	A	126	PHE	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	MN	A	503	1/1	0.18	6.09	52,52,52,52	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	MN	B	501	1/1	0.15	1.77	54,54,54,54	0
2	MN	A	504	1/1	0.14	1.17	58,58,58,58	0
2	MN	B	502	1/1	0.11	-0.44	59,59,59,59	0

6.5 Other polymers ⓘ

There are no such residues in this entry.