



Full wwPDB X-ray Structure Validation Report ⓘ

May 28, 2020 – 08:28 pm BST

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

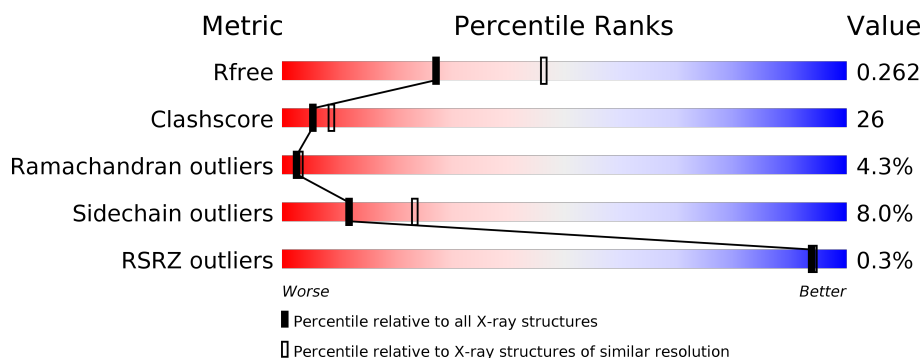
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (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 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	190	
1	B	190	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2724 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 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

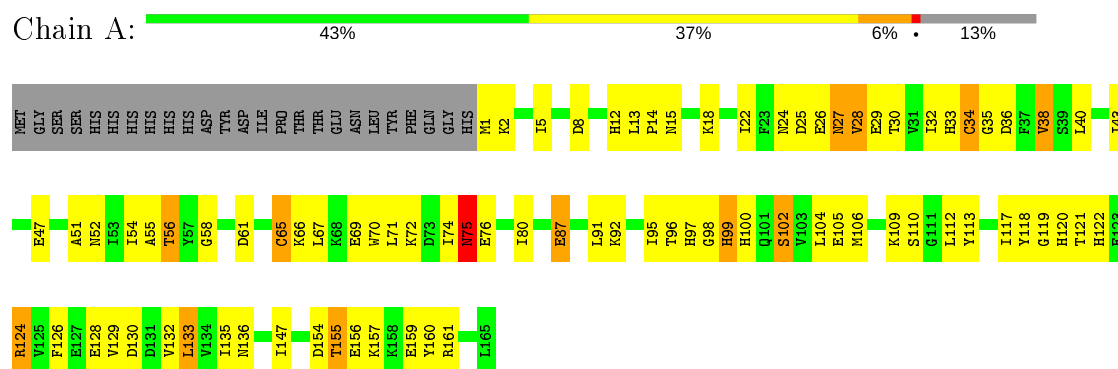
- Molecule 3 is water.

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

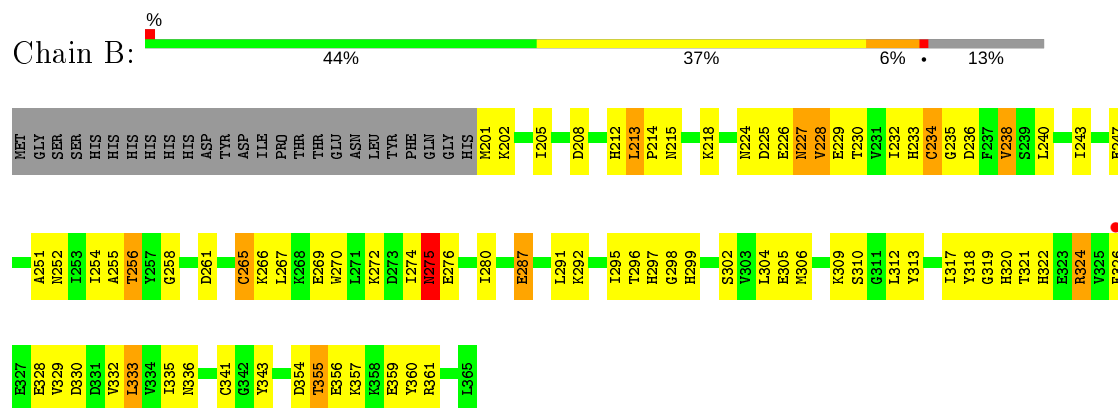
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Hypothetical protein MJ0936



• Molecule 1: Hypothetical protein MJ0936



4 Data and refinement statistics

Property	Value	Source
Space group	I 41	Depositor
Cell constants a, b, c, α , β , γ	70.40 Å 70.40 Å 196.71 Å 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 1.91 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.220 , 0.253 0.230 , 0.262	Depositor DCC
R_{free} test set	3643 reflections (9.98%)	wwPDB-VP
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 , 43.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.499 for -h,k,-l	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.

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: 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 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	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

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 26.

All (137) 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: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:A:154:ASP:OD1	1:A:157:LYS:N	2.31	0.64
1:B:234:CYS:O	1:B:256:THR:HG22	1.97	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:A:121:THR:O	1:A:122:HIS:HB2	2.05	0.56
1:B:243:ILE:HD11	1:B:267:LEU:HD23	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:240:LEU:HD21	1:B:266:LYS:HB3	1.88	0.55
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:218:LYS:NZ	3:B:601:HOH:O	2.39	0.45
1:B:224:ASN:C	1:B:226:GLU:N	2.70	0.45
1:B:229:GLU:O	1:B:251:ALA:HB1	2.16	0.45
1:B:275:ASN:HD22	1:B:276:GLU:N	2.15	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:247:GLU:HA	1:B:274:ILE:HG23	1.98	0.45
1:B:255:ALA:O	1:B:280:ILE:HA	2.17	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:A:47:GLU:HA	1:A:74:ILE:HG23	1.98	0.44
1:B:280:ILE:O	1:B:280:ILE:CG2	2.64	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:329:VAL:O	1:B:329:VAL:HG22	2.18	0.43
1:B:326:PHE:CE1	1:B:360:TYR:HB3	2.54	0.43
1:A:126:PHE:CE1	1:A:160:TYR:HB3	2.54	0.43
1:B:304:LEU:HD11	1:B:318:TYR:CD2	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:119:GLY:HA2	1:A:136:ASN:ND2	2.35	0.41
1:A:33:HIS:O	1:A:55:ALA:HA	2.19	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:PHE:CE1	1:A:133:LEU:HD11	2.56	0.41
1:A:99:HIS:HB3	1:A:100:HIS:H	1.78	0.41
1:B:319:GLY:HA2	1:B:336:ASN:ND2	2.36	0.41
1:B:326:PHE:HD1	1:B:335:ILE:HG12	1.86	0.41
1:B:201:MET:O	1:B:356:GLU:OE2	2.39	0.41
1:A:118:TYR:O	1:A:136:ASN:HA	2.21	0.41
1:A:43:ILE:HD12	1:A:71:LEU:HG	2.03	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	Interatomic distance (Å)	Clash overlap (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 [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	163/190 (86%)	142 (87%)	14 (9%)	7 (4%)	2	3
1	B	163/190 (86%)	141 (86%)	15 (9%)	7 (4%)	2	3
All	All	326/380 (86%)	283 (87%)	29 (9%)	14 (4%)	2	3

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
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%)	12	23
1	B	150/173 (87%)	138 (92%)	12 (8%)	12	23
All	All	300/346 (87%)	276 (92%)	24 (8%)	12	23

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

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Mol	Chain	Res	Type
1	A	155	THR
1	A	161	ARG
1	B	213	LEU
1	B	234	CYS
1	B	256	THR
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 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

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.

No monomer is involved in short contacts.

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	165/190 (86%)	-0.22	0 100 100	35, 61, 85, 91	0
1	B	165/190 (86%)	-0.20	1 (0%) 89 90	34, 61, 86, 91	0
All	All	330/380 (86%)	-0.21	1 (0%) 94 94	34, 61, 86, 91	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	326	PHE	2.9

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. 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(Å ²)	Q<0.9
2	MN	A	503	1/1	0.95	0.19	52,52,52,52	0
2	MN	B	501	1/1	0.96	0.15	54,54,54,54	0
2	MN	B	502	1/1	0.96	0.11	59,59,59,59	0
2	MN	A	504	1/1	0.97	0.15	58,58,58,58	0

6.5 Other polymers

There are no such residues in this entry.