



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:39 PM UTC

PDB ID : 4D1D / pdb_00004d1d
Title : STRUCTURE OF MHP1, A NUCLEOBASE-CATION-SYMPORT-1 FAMILY TRANSPORTER with the inhibitor 5-(2-naphthylmethyl)-L-hydantoin
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Deposited on : 2014-05-01
Resolution : 3.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

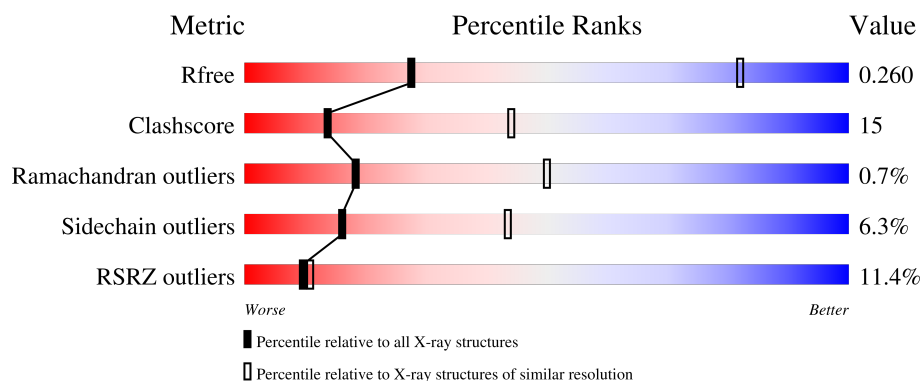
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1131 (3.80-3.60)
Clashscore	190562	1171 (3.80-3.60)
Ramachandran outliers	187476	1129 (3.80-3.60)
Sidechain outliers	187428	1126 (3.80-3.60)
RSRZ outliers	180081	1130 (3.80-3.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	495	11% 57% 33% 8%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3550 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.

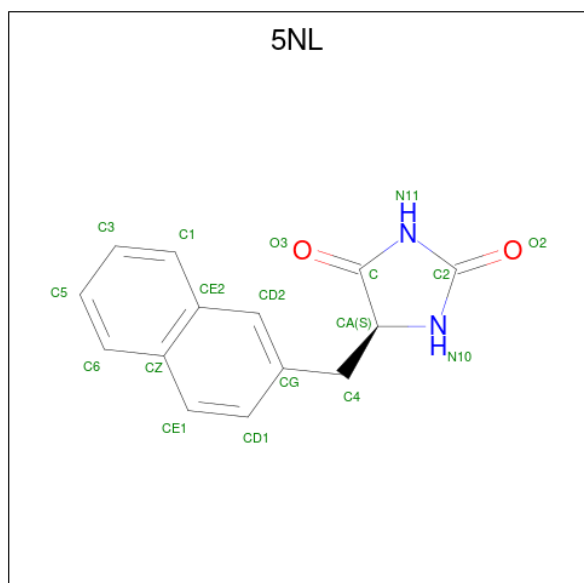
- Molecule 1 is a protein called HYDANTOIN TRANSPORT PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	456	3513	2336	563	592	22	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

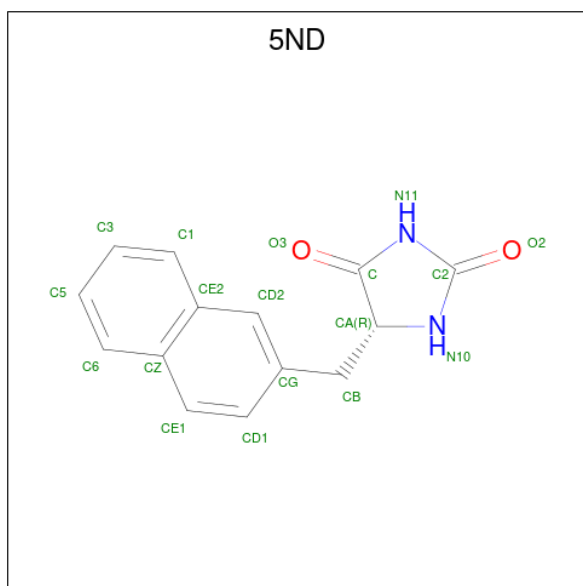
Chain	Residue	Modelled	Actual	Comment	Reference
A	488	GLY	-	expression tag	UNP D6R8X8
A	489	SER	-	expression tag	UNP D6R8X8
A	490	GLU	-	expression tag	UNP D6R8X8
A	491	ASN	-	expression tag	UNP D6R8X8
A	492	LEU	-	expression tag	UNP D6R8X8
A	493	TYR	-	expression tag	UNP D6R8X8
A	494	PHE	-	expression tag	UNP D6R8X8
A	495	GLN	-	expression tag	UNP D6R8X8

- Molecule 2 is 5-(2-NAPHTHYLMETHYL)-L-HYDANTOIN (CCD ID: 5NL) (formula: C₁₄H₁₂N₂O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			18	14	2	2		

- Molecule 3 is 5-(2-NAPHTHYLMETHYL)-D-HYDANTOIN (CCD ID: 5ND) (formula: $C_{14}H_{12}N_2O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			18	14	2	2		

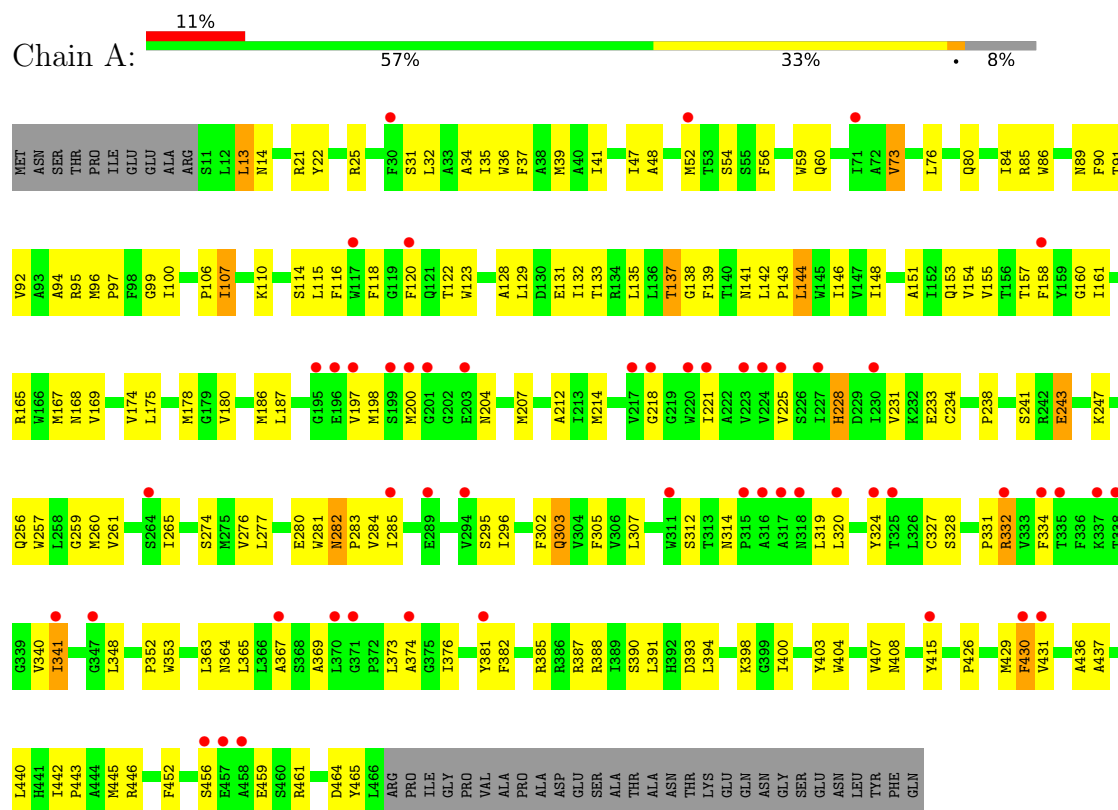
- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Na	0	0
			1	1		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: HYDANTOIN TRANSPORT PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	87.60Å 106.90Å 106.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.78 – 3.70 37.78 – 3.70	Depositor EDS
% Data completeness (in resolution range)	88.9 (37.78-3.70) 95.6 (37.78-3.70)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.13 (at 3.66Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.256 , 0.308 (Not available) , 0.260	Depositor DCC
R_{free} test set	1074 reflections (10.02%)	wwPDB-VP
Wilson B-factor (Å ²)	130.0	Xtriage
Anisotropy	0.714	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 73.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.21$	Xtriage
Estimated twinning fraction	0.048 for -h,l,k	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	3550	wwPDB-VP
Average B, all atoms (Å ²)	194.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.46% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 5NL, NA, 5ND

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.48	0/3610	0.96	7/4936 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	382	PHE	N-CA-C	5.42	118.01	111.40
1	A	73	VAL	N-CA-C	5.36	115.56	110.42
1	A	285	ILE	N-CA-C	5.30	115.51	110.42
1	A	430	PHE	N-CA-C	5.04	117.16	111.11
1	A	282	ASN	CA-C-N	5.03	124.82	119.28
1	A	282	ASN	C-N-CA	5.03	124.82	119.28
1	A	398	LYS	N-CA-C	5.03	116.83	108.02

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3513	0	3604	104	0
2	A	18	0	3	0	0
3	A	18	0	2	0	0
4	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3550	0	3609	104	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 (104) 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:186:MET:HE2	1:A:277:LEU:HD12	1.65	0.78
1:A:52:MET:HE1	1:A:212:ALA:HB1	1.71	0.71
1:A:59:TRP:HB3	1:A:198:MET:HG3	1.73	0.71
1:A:32:LEU:HA	1:A:35:ILE:HD12	1.70	0.71
1:A:436:ALA:O	1:A:440:LEU:HG	1.97	0.63
1:A:13:LEU:HD12	1:A:14:ASN:N	2.13	0.63
1:A:324:TYR:O	1:A:328:SER:HB3	2.00	0.62
1:A:131:GLU:HB3	1:A:303:GLN:HG2	1.82	0.61
1:A:155:VAL:HA	1:A:158:PHE:CD2	2.37	0.59
1:A:48:ALA:O	1:A:52:MET:HG2	2.04	0.57
1:A:110:LYS:HD3	1:A:110:LYS:C	2.29	0.57
1:A:283:PRO:HG2	1:A:302:PHE:HZ	1.67	0.57
1:A:154:VAL:HG13	1:A:340:VAL:HG12	1.86	0.57
1:A:96:MET:HB2	1:A:97:PRO:HD3	1.87	0.56
1:A:369:ALA:O	1:A:415:TYR:OH	2.14	0.56
1:A:186:MET:HE1	1:A:274:SER:HA	1.88	0.55
1:A:442:ILE:HB	1:A:443:PRO:HD3	1.87	0.55
1:A:385:ARG:NH2	1:A:403:TYR:OH	2.40	0.55
1:A:257:TRP:CE2	1:A:261:VAL:HG21	2.42	0.54
1:A:221:ILE:HG21	1:A:367:ALA:HB1	1.90	0.54
1:A:95:ARG:HG2	1:A:100:ILE:HA	1.89	0.54
1:A:131:GLU:O	1:A:135:LEU:HG	2.08	0.54
1:A:96:MET:HE2	1:A:394:LEU:HD11	1.89	0.54
1:A:214:MET:HE1	1:A:430:PHE:H	1.74	0.53
1:A:165:ARG:O	1:A:169:VAL:HG23	2.08	0.53
1:A:426:PRO:HA	1:A:429:MET:HG3	1.91	0.53
1:A:97:PRO:O	1:A:408:ASN:ND2	2.41	0.53
1:A:302:PHE:O	1:A:305:PHE:HB3	2.09	0.52
1:A:207:MET:HE3	1:A:212:ALA:HA	1.92	0.52
1:A:461:ARG:O	1:A:464:ASP:HB2	2.10	0.52
1:A:302:PHE:O	1:A:302:PHE:HD1	1.92	0.52
1:A:327:CYS:HA	1:A:334:PHE:HB2	1.92	0.52
1:A:22:TYR:OH	1:A:85:ARG:NH1	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:ASN:ND2	1:A:331:PRO:O	2.34	0.51
1:A:21:ARG:NH1	1:A:464:ASP:O	2.43	0.51
1:A:31:SER:O	1:A:35:ILE:HG13	2.10	0.51
1:A:34:ALA:O	1:A:168:ASN:HB3	2.11	0.50
1:A:157:THR:HG21	1:A:319:LEU:HD23	1.93	0.50
1:A:25:ARG:HD2	1:A:234:CYS:O	2.12	0.50
1:A:284:VAL:HA	1:A:302:PHE:CE2	2.46	0.50
1:A:36:TRP:CD2	1:A:39:MET:HE3	2.47	0.50
1:A:388:ARG:NH1	1:A:459:GLU:OE2	2.45	0.49
1:A:256:GLN:O	1:A:260:MET:HG2	2.12	0.49
1:A:174:VAL:O	1:A:178:MET:HG3	2.13	0.48
1:A:178:MET:CE	1:A:305:PHE:HB2	2.43	0.48
1:A:408:ASN:HD21	1:A:445:MET:HE1	1.77	0.48
1:A:13:LEU:O	1:A:332:ARG:NH2	2.46	0.48
1:A:128:ALA:HA	1:A:303:GLN:HG3	1.96	0.48
1:A:76:LEU:O	1:A:80:GLN:HG2	2.14	0.48
1:A:37:PHE:CE1	1:A:41:ILE:HD11	2.49	0.47
1:A:142:LEU:HB3	1:A:143:PRO:HD3	1.95	0.47
1:A:37:PHE:HA	1:A:259:GLY:O	2.15	0.47
1:A:274:SER:HB3	1:A:280:GLU:O	2.15	0.47
1:A:106:PRO:HB2	1:A:373:LEU:HD13	1.97	0.47
1:A:54:SER:O	1:A:204:ASN:HB3	2.16	0.46
1:A:122:THR:HG21	1:A:348:LEU:HD21	1.97	0.46
1:A:352:PRO:HD2	1:A:353:TRP:CZ3	2.50	0.46
1:A:89:ASN:HB3	1:A:92:VAL:H	1.81	0.45
1:A:129:LEU:HG	1:A:307:LEU:HD13	1.98	0.45
1:A:137:THR:C	1:A:139:PHE:H	2.24	0.45
1:A:85:ARG:HD2	1:A:86:TRP:CE2	2.51	0.45
1:A:25:ARG:NH1	1:A:84:ILE:O	2.44	0.45
1:A:99:GLY:HA3	1:A:407:VAL:HA	1.98	0.45
1:A:129:LEU:O	1:A:133:THR:HG23	2.17	0.45
1:A:218:GLY:HA2	1:A:221:ILE:HB	1.99	0.45
1:A:225:VAL:HG23	1:A:374:ALA:HB2	1.99	0.45
1:A:151:ALA:O	1:A:155:VAL:HG23	2.17	0.44
1:A:118:PHE:HA	1:A:314:ASN:OD1	2.18	0.44
1:A:391:LEU:HA	1:A:394:LEU:HD12	1.99	0.44
1:A:153:GLN:O	1:A:157:THR:N	2.48	0.44
1:A:56:PHE:HE1	1:A:207:MET:HE2	1.82	0.44
1:A:400:ILE:HD13	1:A:456:SER:HB2	2.00	0.43
1:A:404:TRP:CE2	1:A:452:PHE:HE2	2.36	0.43
1:A:73:VAL:HA	1:A:76:LEU:HD12	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:LYS:HD3	1:A:110:LYS:O	2.18	0.43
1:A:47:ILE:HG12	1:A:281:TRP:CE2	2.53	0.43
1:A:89:ASN:HB2	1:A:92:VAL:HB	2.00	0.43
1:A:22:TYR:CZ	1:A:238:PRO:HB3	2.53	0.43
1:A:132:ILE:CG1	1:A:303:GLN:HB3	2.49	0.43
1:A:144:LEU:HG	1:A:148:ILE:HD12	2.01	0.42
1:A:158:PHE:HZ	1:A:341:ILE:HD11	1.84	0.42
1:A:376:ILE:HG13	1:A:437:ALA:HB1	2.00	0.42
1:A:115:LEU:HD23	1:A:115:LEU:HA	1.84	0.42
1:A:282:ASN:HA	1:A:283:PRO:HD2	1.90	0.42
1:A:243:GLU:O	1:A:247:LYS:HG2	2.19	0.42
1:A:231:VAL:C	1:A:233:GLU:H	2.27	0.42
1:A:141:ASN:OD1	1:A:143:PRO:HD2	2.19	0.42
1:A:391:LEU:HD12	1:A:465:TYR:HB2	2.01	0.42
1:A:120:PHE:O	1:A:123:TRP:HB3	2.20	0.41
1:A:381:TYR:CE2	1:A:387:ARG:HG2	2.54	0.41
1:A:94:ALA:O	1:A:97:PRO:HD2	2.19	0.41
1:A:319:LEU:HD21	1:A:340:VAL:HG13	2.02	0.41
1:A:56:PHE:HB3	1:A:60:GLN:OE1	2.20	0.41
1:A:320:LEU:HD23	1:A:320:LEU:HA	1.82	0.41
1:A:91:THR:HG23	1:A:107:ILE:HD13	2.03	0.41
1:A:116:PHE:CE2	1:A:363:LEU:HD11	2.55	0.41
1:A:128:ALA:O	1:A:132:ILE:HG13	2.20	0.41
1:A:41:ILE:HD13	1:A:175:LEU:HB3	2.02	0.41
1:A:90:PHE:HB2	1:A:228:HIS:HB2	2.03	0.41
1:A:197:VAL:HA	1:A:200:MET:HG3	2.02	0.41
1:A:390:SER:OG	1:A:393:ASP:OD1	2.38	0.41
1:A:158:PHE:C	1:A:160:GLY:H	2.29	0.41
1:A:178:MET:HE1	1:A:305:PHE:HB2	2.02	0.40
1:A:440:LEU:O	1:A:443:PRO:HD2	2.22	0.40

There are no symmetry-related clashes.

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	454/495 (92%)	403 (89%)	48 (11%)	3 (1%)	18	50

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	228	HIS
1	A	137	THR
1	A	138	GLY

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	367/399 (92%)	344 (94%)	23 (6%)	16	43

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LEU
1	A	107	ILE
1	A	114	SER
1	A	144	LEU
1	A	146	ILE
1	A	161	ILE
1	A	167	MET
1	A	180	VAL
1	A	187	LEU
1	A	241	SER
1	A	243	GLU
1	A	265	ILE
1	A	276	VAL
1	A	295	SER
1	A	296	ILE
1	A	303	GLN

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Mol	Chain	Res	Type
1	A	312	SER
1	A	332	ARG
1	A	341	ILE
1	A	364	ASN
1	A	365	LEU
1	A	431	VAL
1	A	446	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	GLN
1	A	168	ASN

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 oligosaccharides 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	5ND	A	1468	-	20,20,20	2.51	4 (20%)	26,28,28	2.59	3 (11%)
2	5NL	A	1467	-	20,20,20	2.50	4 (20%)	26,28,28	2.47	3 (11%)

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	5ND	A	1468	-	-	3/4/16/16	0/3/3/3
2	5NL	A	1467	-	-	0/4/16/16	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1468	5ND	CA-C	-7.24	1.47	1.52
2	A	1467	5NL	CA-C	-6.85	1.47	1.52
2	A	1467	5NL	C2-N10	5.62	1.45	1.35
3	A	1468	5ND	C2-N10	5.57	1.45	1.35
2	A	1467	5NL	C-N11	4.15	1.43	1.37
3	A	1468	5ND	C-N11	3.74	1.42	1.37
2	A	1467	5NL	C2-N11	3.08	1.46	1.38
3	A	1468	5ND	C2-N11	2.95	1.45	1.38

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1468	5ND	C-CA-N10	11.53	108.04	100.99
2	A	1467	5NL	C-CA-N10	11.31	107.90	100.99
3	A	1468	5ND	C-N11-C2	-4.24	108.03	111.61
2	A	1467	5NL	C-N11-C2	-4.11	108.14	111.61
3	A	1468	5ND	CB-CA-N10	-3.63	109.92	113.84
2	A	1467	5NL	C4-CA-N10	-2.05	111.63	113.84

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1468	5ND	CA-CB-CG-CD1
3	A	1468	5ND	CA-CB-CG-CD2
3	A	1468	5ND	N10-CA-CB-CG

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

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	456/495 (92%)	0.57	52 (11%)	10 11	102, 185, 281, 399	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	200	MET	13.3
1	A	203	GLU	13.1
1	A	430	PHE	5.7
1	A	197	VAL	5.4
1	A	201	GLY	5.4
1	A	431	VAL	5.1
1	A	335	THR	5.1
1	A	218	GLY	4.8
1	A	196	GLU	4.6
1	A	458	ALA	4.5
1	A	341	ILE	4.3
1	A	225	VAL	4.1
1	A	337	LYS	3.9
1	A	381	TYR	3.8
1	A	217	VAL	3.7
1	A	338	THR	3.5
1	A	30	PHE	3.2
1	A	223	VAL	3.0
1	A	224	VAL	3.0
1	A	317	ALA	3.0
1	A	227	ILE	3.0
1	A	371	GLY	2.9
1	A	230	ILE	2.8
1	A	289	GLU	2.8
1	A	316	ALA	2.8
1	A	264	SER	2.8
1	A	456	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	220	TRP	2.7
1	A	117	TRP	2.6
1	A	195	GLY	2.6
1	A	120	PHE	2.6
1	A	374	ALA	2.6
1	A	318	ASN	2.6
1	A	199	SER	2.5
1	A	332	ARG	2.4
1	A	158	PHE	2.4
1	A	221	ILE	2.4
1	A	324	TYR	2.4
1	A	334	PHE	2.3
1	A	315	PRO	2.3
1	A	285	ILE	2.3
1	A	320	LEU	2.2
1	A	347	GLY	2.2
1	A	457	GLU	2.2
1	A	311	TRP	2.2
1	A	71	ILE	2.1
1	A	415	TYR	2.1
1	A	367	ALA	2.1
1	A	294	VAL	2.1
1	A	370	LEU	2.0
1	A	52	MET	2.0
1	A	325	THR	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 oligosaccharides 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. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	5NL	A	1467	18/18	0.83	0.20	142,154,170,171	18
3	5ND	A	1468	18/18	0.84	0.23	140,151,167,169	18
4	NA	A	1471	1/1	0.98	0.03	115,115,115,115	0

6.5 Other polymers [i](#)

There are no such residues in this entry.