



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 3, 2017 – 12:49 PM EDT

PDB ID : 5HP3
Title : Human Adenosine Deaminase Acting on dsRNA (ADAR2) bound to dsRNA sequence derived from *S. cerevisiae* BDF2 gene with AC mismatch at reaction site
Authors : Matthews, M.M.; Fisher, A.J.; Beal, P.A.
Deposited on : unknown
Resolution : 3.09 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029824
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : rb-20029824

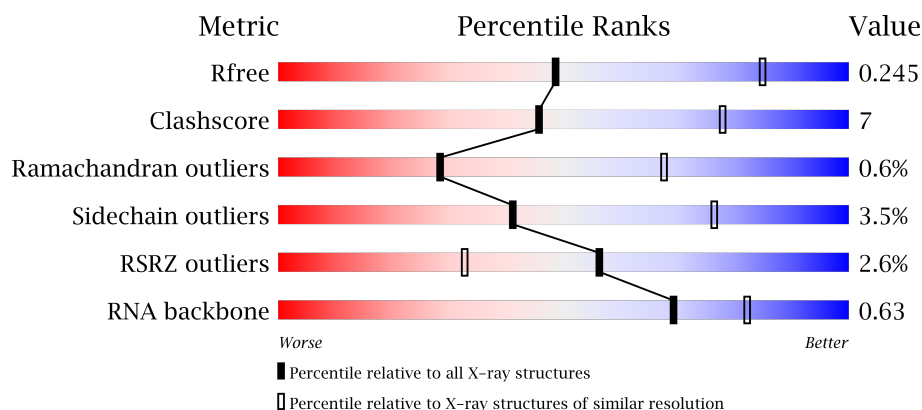
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.09 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}	100719	1001 (3.12-3.08)
Clashscore	112137	1099 (3.12-3.08)
Ramachandran outliers	110173	1057 (3.12-3.08)
Sidechain outliers	110143	1057 (3.12-3.08)
RSRZ outliers	101464	1006 (3.12-3.08)
RNA backbone	2435	1112 (3.50-2.70)

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	403	<div> <div>81%</div> <div>16%</div> <div>..</div> </div>
1	D	403	<div> <div>5%</div> <div>77%</div> <div>21%</div> <div>..</div> </div>
2	B	23	<div> <div>65%</div> <div>30%</div> <div>.</div> </div>
3	C	23	<div> <div>43%</div> <div>57%</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7205 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 Double-stranded RNA-specific editase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	394	Total	C	N	O	S	0	0	0
			3085	1946	560	568	11			
1	D	397	Total	C	N	O	S	0	0	0
			3072	1940	553	568	11			

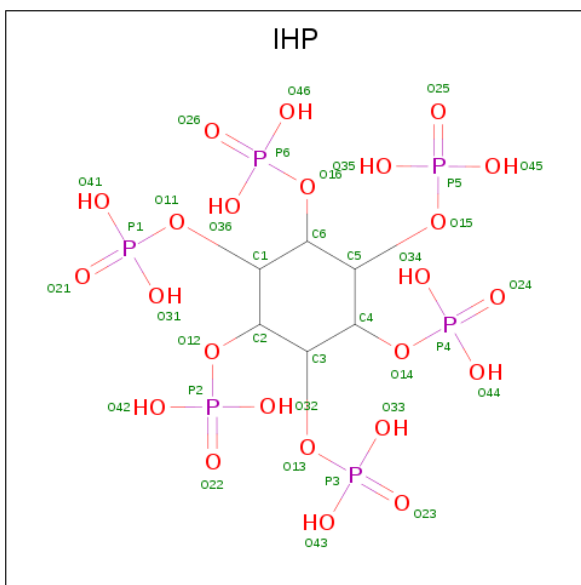
- Molecule 2 is a RNA chain called RNA (5'-R(*UP*UP*CP*CP*CP*CP*AP*CP*AP*UP*UP*(8AZ)P*GP*AP*CP*GP*UP*UP*CP*AP*GP*UP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	23	Total	C	N	O	P	0	0	0
			476	214	78	162	22			

- Molecule 3 is a RNA chain called RNA (5'-R(*GP*AP*CP*UP*GP*AP*AP*CP*GP*AP*CP*CP*AP*AP*UP*GP*UP*GP*GP*GP*GP*AP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	23	Total	C	N	O	P	0	0	0
			497	223	98	154	22			

- Molecule 4 is INOSITOL HEXAKISPHOSPHATE (three-letter code: IHP) (formula: C₆H₁₈O₂₄P₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	P	0	0
			36	6	24	6		
4	D	1	Total	C	O	P	0	0
			36	6	24	6		

- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Zn	0	0
			1	1		
5	D	1	Total	Zn	0	0
			1	1		

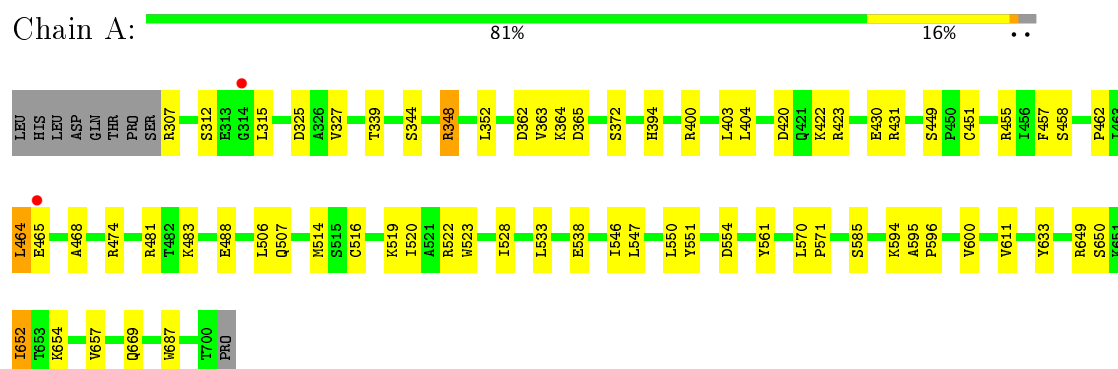
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	1	Total	O	0	0
			1	1		

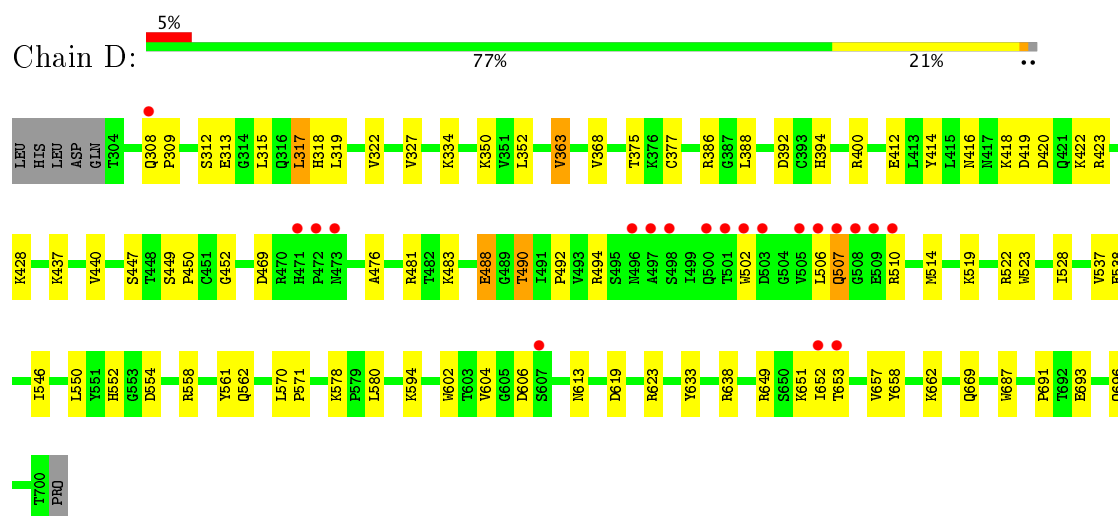
3 Residue-property plots [i](#)

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.

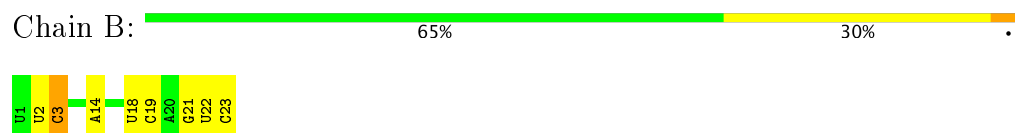
• Molecule 1: Double-stranded RNA-specific editase 1



• Molecule 1: Double-stranded RNA-specific editase 1



• Molecule 2: RNA (5'-R(*UP*UP*CP*CP*CP*CP*AP*CP*AP*UP*UP*(8AZ)P*GP*AP*CP*GP*UP*UP*CP*AP*GP*UP*C)-3')

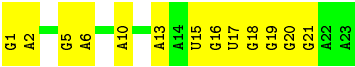


● Molecule 3: RNA (5'-R(*GP*AP*CP*UP*GP*AP*AP*CP*GP*AP*CP*CP*AP*AP*UP*GP*UP*GP*GP*GP*GP*AP*A)-3')

Chain C:

43%

57%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.51Å 107.21Å 120.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.33 – 3.09 36.33 – 3.09	Depositor EDS
% Data completeness (in resolution range)	96.9 (36.33-3.09) 96.9 (36.33-3.09)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.163 , 0.238 0.174 , 0.245	Depositor DCC
R_{free} test set	989 reflections (5.39%)	DCC
Wilson B-factor (Å ²)	65.4	Xtriage
Anisotropy	0.601	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 64.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7205	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.00% 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: ZN, IHP, 8AZ

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	0/3149	0.68	0/4259
1	D	0.44	0/3135	0.65	1/4245 (0.0%)
2	B	0.56	0/503	1.04	1/777 (0.1%)
3	C	0.53	0/558	1.02	0/870
All	All	0.47	0/7345	0.74	2/10151 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	317	LEU	CA-CB-CG	5.64	128.27	115.30
2	B	3	C	C6-N1-C2	-5.20	118.22	120.30

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	3085	0	3112	38	2
1	D	3072	0	3071	53	2
2	B	476	0	237	7	0
3	C	497	0	252	12	0
4	A	36	0	6	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	36	0	6	0	0
5	A	1	0	0	0	0
5	D	1	0	0	0	0
6	D	1	0	0	0	0
All	All	7205	0	6684	102	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 7.

All (102) 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:348:ARG:NH1	3:C:5:G:OP1	2.21	0.73
3:C:18:G:H2'	3:C:19:G:C8	2.24	0.73
1:D:392:ASP:OD2	1:D:483:LYS:NZ	2.22	0.71
1:D:420:ASP:OD1	1:D:423:ARG:NH1	2.26	0.69
1:D:481:ARG:HE	1:D:490:THR:CG2	2.09	0.66
1:D:537:VAL:O	1:D:649:ARG:NH1	2.28	0.66
1:D:400:ARG:HD3	1:D:523:TRP:CE2	2.32	0.65
1:D:514:MET:HG3	1:D:687:TRP:CE2	2.32	0.65
1:D:400:ARG:NH1	1:D:528:ILE:O	2.31	0.64
1:D:327:VAL:HG21	1:D:546:ILE:HG13	1.80	0.64
1:D:377:CYS:SG	1:D:483:LYS:HD3	2.39	0.63
1:A:451:CYS:HA	1:A:455:ARG:HD2	1.82	0.62
1:D:313:GLU:HG3	1:D:578:LYS:HB2	1.83	0.60
1:D:394:HIS:CE1	1:D:483:LYS:HE2	2.36	0.60
1:D:437:LYS:O	1:D:440:VAL:HG12	2.03	0.59
1:D:350:LYS:HB3	1:D:375:THR:HG21	1.85	0.59
1:D:313:GLU:CG	1:D:578:LYS:HB2	2.34	0.58
1:D:481:ARG:HB3	1:D:490:THR:HG23	1.84	0.58
1:A:519:LYS:NZ	4:A:801:IHP:O31	2.26	0.58
1:D:400:ARG:HD3	1:D:523:TRP:CZ2	2.39	0.58
1:D:552:HIS:CE1	1:D:554:ASP:HB2	2.39	0.58
1:A:523:TRP:HB3	1:A:528:ILE:HG22	1.86	0.57
1:A:400:ARG:HD3	1:A:523:TRP:CE2	2.39	0.57
1:A:394:HIS:CE1	1:A:483:LYS:HE2	2.40	0.56
2:B:14:A:N1	3:C:10:A:H2	2.04	0.56
1:D:602:TRP:CZ2	1:D:604:VAL:HA	2.40	0.56
1:A:594:LYS:HB2	3:C:6:A:OP2	2.06	0.56
1:D:693:GLU:HA	1:D:696:GLN:HG3	1.88	0.56
1:D:386:ARG:HE	1:D:388:LEU:HD11	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:514:MET:HG3	1:A:687:TRP:CE2	2.42	0.55
1:D:558:ARG:HA	1:D:562:GLN:HB3	1.90	0.54
1:D:452:GLY:HA2	1:D:490:THR:HG21	1.88	0.54
1:A:474:ARG:NH1	3:C:16:G:OP1	2.41	0.54
1:D:318:HIS:HD2	1:D:363:VAL:HG22	1.72	0.54
1:A:652:ILE:HD11	1:A:654:LYS:HE2	1.90	0.53
1:D:481:ARG:HE	1:D:490:THR:HG22	1.73	0.53
1:A:554:ASP:N	1:A:554:ASP:OD1	2.42	0.53
3:C:1:G:H2'	3:C:2:A:C8	2.45	0.52
1:D:318:HIS:CD2	1:D:363:VAL:HG13	2.45	0.52
1:A:519:LYS:O	1:A:522:ARG:HB3	2.10	0.51
1:D:315:LEU:HD13	1:D:317:LEU:HD21	1.92	0.51
1:A:404:LEU:HD22	1:A:533:LEU:HD12	1.92	0.51
3:C:18:G:H2'	3:C:19:G:H8	1.75	0.50
1:A:420:ASP:OD1	1:A:423:ARG:NH1	2.45	0.50
3:C:16:G:H2'	3:C:17:U:C6	2.47	0.49
1:D:494:ARG:HD2	1:D:510:ARG:HH22	1.77	0.49
2:B:18:U:H2'	2:B:19:C:C6	2.48	0.49
1:A:464:LEU:HD13	1:A:465:GLU:H	1.79	0.48
1:A:307:ARG:HA	1:A:551:TYR:CD1	2.49	0.48
1:A:400:ARG:O	1:A:403:LEU:HB3	2.14	0.48
1:D:412:GLU:OE1	1:D:638:ARG:HD2	2.14	0.47
1:A:538:GLU:HA	1:A:649:ARG:NH1	2.29	0.47
1:A:352:LEU:HD12	1:A:372:SER:O	2.15	0.47
2:B:21:G:H2'	2:B:22:U:H6	1.78	0.47
1:A:339:THR:HB	1:A:344:SER:HB3	1.96	0.47
1:A:362:ASP:OD1	1:A:364:LYS:HG2	2.14	0.46
1:D:602:TRP:CD1	1:D:606:ASP:HB2	2.50	0.46
1:A:600:VAL:HG22	1:A:611:VAL:HG22	1.98	0.46
2:B:18:U:H2'	2:B:19:C:H6	1.81	0.45
1:D:394:HIS:ND1	1:D:483:LYS:HE2	2.31	0.45
3:C:19:G:H2'	3:C:20:G:O4'	2.16	0.45
1:D:449:SER:OG	1:D:450:PRO:HD2	2.17	0.45
2:B:2:U:H2'	2:B:3:C:C6	2.52	0.45
1:A:474:ARG:HG2	2:B:3:C:OP2	2.17	0.45
1:A:325:ASP:OD2	1:A:585:SER:OG	2.31	0.44
1:D:481:ARG:HG2	1:D:492:PRO:HA	2.00	0.44
2:B:21:G:H2'	2:B:22:U:C6	2.53	0.44
1:D:319:LEU:HB2	1:D:322:VAL:HG12	1.99	0.44
1:D:352:LEU:O	1:D:447:SER:N	2.49	0.44
1:A:422:LYS:HD3	1:D:507:GLN:NE2	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:538:GLU:HA	1:D:649:ARG:NH1	2.33	0.43
1:A:516:CYS:O	1:A:520:ILE:HG13	2.18	0.43
1:D:312:SER:HB3	1:D:580:LEU:N	2.34	0.43
1:D:502:TRP:HE3	1:D:691:PRO:HG2	1.83	0.43
1:D:334:LYS:HG2	1:D:368:VAL:HG11	2.00	0.43
3:C:5:G:H2'	3:C:6:A:C8	2.53	0.43
1:A:449:SER:HA	1:A:547:LEU:HD13	1.99	0.43
1:D:519:LYS:O	1:D:522:ARG:HB3	2.19	0.43
1:A:430:GLU:OE2	1:D:613:ASN:ND2	2.44	0.42
1:D:422:LYS:HE2	1:D:422:LYS:HB3	1.84	0.42
1:D:651:LYS:HB2	1:D:653:THR:HG23	2.00	0.42
1:A:650:SER:OG	1:A:652:ILE:HG23	2.19	0.42
3:C:5:G:H2'	3:C:6:A:H8	1.83	0.42
1:D:658:TYR:CE2	1:D:662:LYS:HE3	2.54	0.42
1:A:327:VAL:HG21	1:A:546:ILE:HG13	2.01	0.42
1:A:430:GLU:HG3	1:A:431:ARG:N	2.34	0.42
1:A:538:GLU:HG2	1:A:649:ARG:CZ	2.50	0.41
1:A:570:LEU:HA	1:A:571:PRO:HD3	1.93	0.41
1:A:362:ASP:O	1:A:365:ASP:HB2	2.20	0.41
1:A:462:PRO:C	1:A:464:LEU:H	2.24	0.41
1:A:481:ARG:NH2	3:C:15:U:OP1	2.54	0.41
1:D:308:GLN:HA	1:D:309:PRO:HD2	1.84	0.41
1:D:552:HIS:HE1	1:D:554:ASP:HB2	1.86	0.41
1:D:502:TRP:O	1:D:506:LEU:HG	2.21	0.41
1:D:623:ARG:HE	1:D:623:ARG:HB2	1.60	0.41
1:D:414:TYR:OH	1:D:428:LYS:HG3	2.21	0.41
1:D:594:LYS:H	1:D:594:LYS:HG2	1.68	0.41
1:A:595:ALA:HA	1:A:596:PRO:HD2	1.84	0.41
1:D:412:GLU:O	1:D:416:ASN:HB2	2.21	0.40
1:D:619:ASP:OD2	1:D:623:ARG:NH2	2.54	0.40
1:A:506:LEU:HA	1:A:506:LEU:HD23	1.87	0.40
1:D:570:LEU:HA	1:D:571:PRO:HD3	1.94	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:344:SER:OG	1:D:469:ASP:OD2[2_445]	2.11	0.09
1:A:507:GLN:NE2	1:D:488:GLU:OE2[3_555]	2.17	0.03

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	392/403 (97%)	373 (95%)	16 (4%)	3 (1%)	22	62
1	D	395/403 (98%)	379 (96%)	14 (4%)	2 (0%)	32	71
All	All	787/806 (98%)	752 (96%)	30 (4%)	5 (1%)	28	67

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	468	ALA
1	D	476	ALA
1	A	312	SER
1	A	561	TYR
1	D	561	TYR

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	336/347 (97%)	324 (96%)	12 (4%)	40	75
1	D	330/347 (95%)	319 (97%)	11 (3%)	43	77
All	All	666/694 (96%)	643 (96%)	23 (4%)	41	76

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

Mol	Chain	Res	Type
1	A	315	LEU

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Mol	Chain	Res	Type
1	A	348	ARG
1	A	363	VAL
1	A	457	PHE
1	A	458	SER
1	A	464	LEU
1	A	488	GLU
1	A	550	LEU
1	A	633	TYR
1	A	652	ILE
1	A	657	VAL
1	A	669	GLN
1	D	363	VAL
1	D	418	LYS
1	D	419	ASP
1	D	488	GLU
1	D	490	THR
1	D	507	GLN
1	D	550	LEU
1	D	633	TYR
1	D	652	ILE
1	D	657	VAL
1	D	669	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	316	GLN
1	A	318	HIS
1	A	630	HIS
1	D	318	HIS
1	D	507	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	21/23 (91%)	1 (4%)	0
3	C	22/23 (95%)	2 (9%)	0
All	All	43/46 (93%)	3 (6%)	0

All (3) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	23	C
3	C	13	A
3	C	21	G

There are no RNA pucker outliers to report.

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

1 non-standard protein/DNA/RNA residue is 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	8AZ	B	12	2,5	16,24,25	0.95	1 (6%)	15,35,38	0.59	0

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	8AZ	B	12	2,5	-	0/3/35/36	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	12	8AZ	O5'-C5'	-2.30	1.41	1.44

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.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are 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	$\# Z > 2$
4	IHP	A	801	-	36,36,36	0.89	1 (2%)	54,60,60	1.10	2 (3%)
4	IHP	D	801	-	36,36,36	0.84	0	54,60,60	1.13	4 (7%)

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
4	IHP	A	801	-	-	0/30/54/54	0/1/1/1
4	IHP	D	801	-	-	0/30/54/54	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	801	IHP	P4-O14	2.34	1.63	1.59

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	801	IHP	O16-P6-O26	-2.35	100.05	109.26
4	D	801	IHP	O15-C5-C6	-2.21	103.47	108.68
4	D	801	IHP	O16-P6-O26	-2.09	101.08	109.26
4	A	801	IHP	O34-P4-O24	2.03	118.44	110.50
4	D	801	IHP	O14-C4-C5	2.03	113.48	108.68
4	D	801	IHP	C5-C6-C1	2.48	115.95	110.46

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	801	IHP	1	0

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 ⓘ

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	394/403 (97%)	-0.41	2 (0%) 90 80	31, 54, 100, 160	0
1	D	397/403 (98%)	-0.02	20 (5%) 30 13	43, 70, 126, 165	0
2	B	22/23 (95%)	-0.04	0 100 100	51, 75, 99, 105	0
3	C	23/23 (100%)	-0.12	0 100 100	52, 80, 104, 108	0
All	All	836/852 (98%)	-0.21	22 (2%) 56 33	31, 62, 117, 165	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	473	ASN	6.7
1	D	472	PRO	6.0
1	D	505	VAL	4.7
1	D	501	THR	4.0
1	A	314	GLY	3.6
1	D	498	SER	3.5
1	D	653	THR	3.3
1	D	497	ALA	3.3
1	D	507	GLN	3.1
1	D	503	ASP	3.1
1	D	506	LEU	2.9
1	D	471	HIS	2.9
1	D	510	ARG	2.8
1	D	502	TRP	2.7
1	D	509	GLU	2.7
1	D	508	GLY	2.6
1	D	496	ASN	2.5
1	A	465	GLU	2.3
1	D	652	ILE	2.2
1	D	308	GLN	2.2
1	D	607	SER	2.1
1	D	500	GLN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
2	8AZ	B	12	22/23	0.98	0.20	-	39,54,63,64	0

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(Å ²)	Q<0.9
4	IHP	A	801	36/36	0.98	0.23	0.88	18,36,54,67	0
4	IHP	D	801	36/36	0.98	0.24	0.53	23,48,66,74	0
5	ZN	A	802	1/1	1.00	0.17	-1.42	47,47,47,47	0
5	ZN	D	802	1/1	0.99	0.17	-1.66	50,50,50,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.