



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 10:57 AM UTC

PDB ID : 5VDU / pdb_00005vdu
Title : Human cyclic GMP-AMP synthase (cGAS) in complex with Compound F2
Authors : Byrnes, L.J.; Hall, J.D.
Deposited on : 2017-04-03
Resolution : 2.73 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

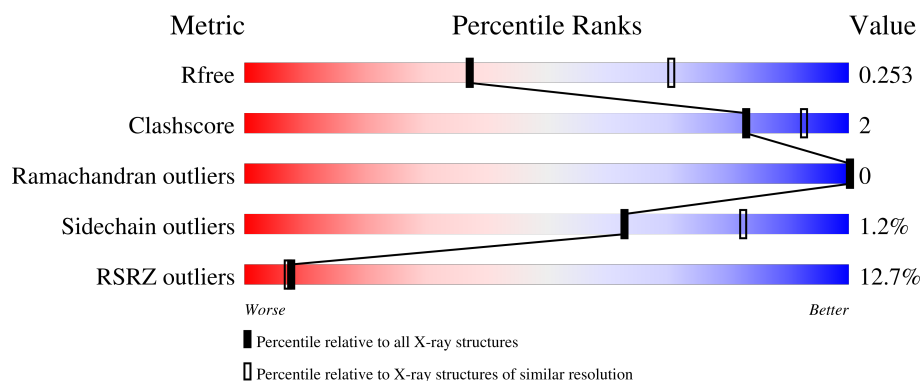
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.73 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	4348 (2.74-2.70)
Clashscore	190562	4665 (2.74-2.70)
Ramachandran outliers	187476	4584 (2.74-2.70)
Sidechain outliers	187428	4585 (2.74-2.70)
RSRZ outliers	180081	4348 (2.74-2.70)

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 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	363	
1	B	363	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5734 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 Cyclic GMP-AMP synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	361	Total	C	N	O	S	0	0	0
			2785	1785	471	514	15			
1	B	361	Total	C	N	O	S	0	2	0
			2799	1797	474	514	14			

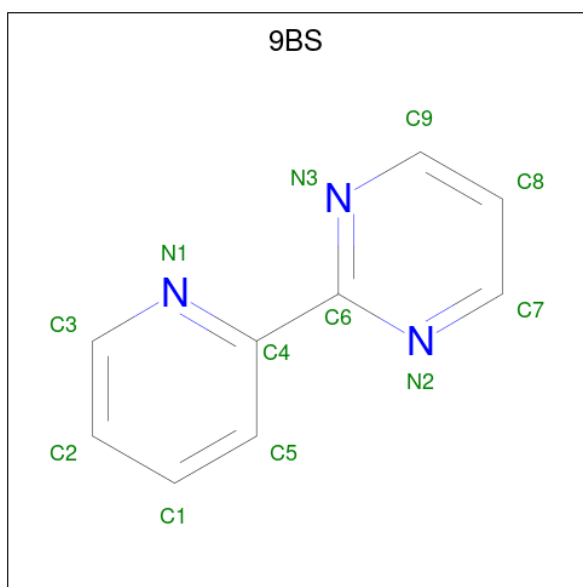
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	160	MET	-	expression tag	UNP Q8N884
B	160	MET	-	expression tag	UNP Q8N884

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

- Molecule 3 is 2-(pyridin-2-yl)pyrimidine (CCD ID: 9BS) (formula: C₉H₇N₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	1
			36	27	9		
3	B	1	Total	C	N	0	1
			36	27	9		
3	B	1	Total	C	N	0	0
			12	9	3		
3	B	1	Total	C	N	0	0
			12	9	3		

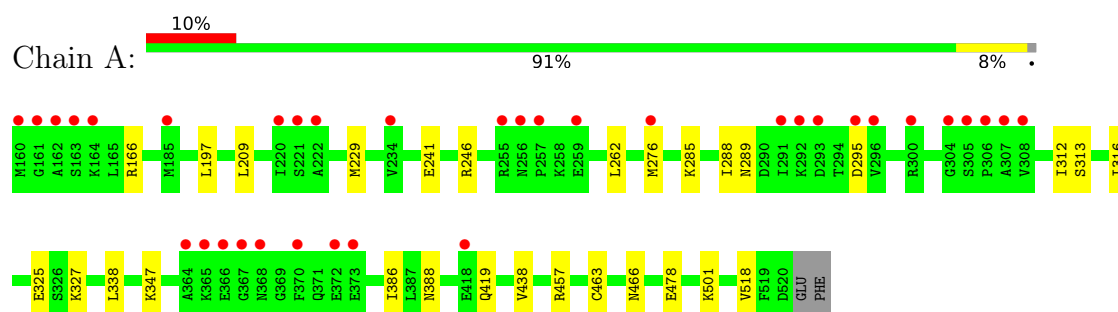
- Molecule 4 is water.

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

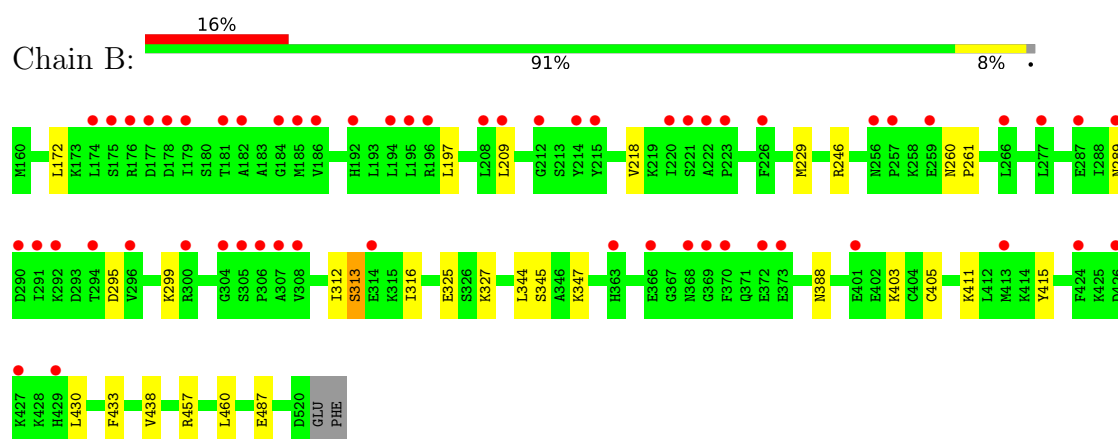
3 Residue-property plots [i](#)

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: Cyclic GMP-AMP synthase



• Molecule 1: Cyclic GMP-AMP synthase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	216.41Å 48.02Å 86.47Å 90.00° 104.94° 90.00°	Depositor
Resolution (Å)	52.27 – 2.73 52.27 – 2.73	Depositor EDS
% Data completeness (in resolution range)	98.1 (52.27-2.73) 98.1 (52.27-2.73)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 2.73Å)	Xtriage
Refinement program	PHENIX dev_1999	Depositor
R, R_{free}	0.222 , 0.252 0.226 , 0.253	Depositor DCC
R_{free} test set	1062 reflections (4.55%)	wwPDB-VP
Wilson B-factor (Å ²)	47.8	Xtriage
Anisotropy	0.127	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 55.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5734	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.14% 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, 9BS

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	0.26	0/2842	0.70	0/3843
1	B	0.26	0/2864	0.71	1/3874 (0.0%)
All	All	0.26	0/5706	0.70	1/7717 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	313	SER	CB-CA-C	-5.04	110.74	116.54

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	2785	0	2661	14	0
1	B	2799	0	2671	14	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	36	0	0	0	0
3	B	60	0	0	0	0
4	A	26	0	0	0	0
4	B	26	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5734	0	5332	26	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 2.

All (26) 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:241:GLU:OE2	1:A:246:ARG:NH1	2.32	0.62
1:A:463:CYS:O	1:A:466:ASN:HB2	2.05	0.57
1:A:347:LYS:NZ	1:B:388:ASN:O	2.38	0.55
1:A:209:LEU:HD23	1:A:229:MET:HE3	1.89	0.55
1:B:209:LEU:HD23	1:B:229:MET:HE3	1.88	0.55
1:A:262:LEU:HD13	1:A:276:MET:HE1	1.90	0.54
1:A:312:ILE:HB	1:A:316:ILE:HB	1.90	0.53
1:A:388:ASN:O	1:B:347:LYS:NZ	2.38	0.52
1:B:312:ILE:HB	1:B:316:ILE:HB	1.93	0.51
1:B:246:ARG:NH1	1:B:487:GLU:OE2	2.44	0.50
1:B:403:LYS:HB3	1:B:457:ARG:HD2	1.94	0.49
1:A:478:GLU:OE2	1:A:501:LYS:NZ	2.48	0.44
1:A:166:ARG:NH2	1:A:457:ARG:O	2.48	0.44
1:A:419:GLN:HB3	1:A:518:VAL:HG22	2.00	0.44
1:A:295:ASP:HB3	1:A:313:SER:HA	2.00	0.43
1:B:295:ASP:HB3	1:B:313:SER:HA	1.99	0.43
1:B:430:LEU:HA	1:B:433:PHE:HD2	1.84	0.42
1:B:344:LEU:O	1:B:345:SER:OG	2.34	0.42
1:B:325:GLU:OE1	1:B:327:LYS:NZ	2.53	0.42
1:A:285:LYS:O	1:A:288:ILE:HG13	2.20	0.42
1:B:218:VAL:HA	1:B:411:LYS:HE3	2.01	0.42
1:B:172:LEU:HD21	1:B:415:TYR:CE1	2.55	0.41
1:B:405:CYS:HB2	1:B:460:LEU:HD13	2.01	0.41
1:A:338:LEU:HD11	1:A:386:ILE:HD11	2.02	0.41
1:A:325:GLU:OE1	1:A:327:LYS:NZ	2.54	0.41
1:B:260:ASN:HA	1:B:261:PRO:HD3	1.89	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	359/363 (99%)	346 (96%)	13 (4%)	0	100	100
1	B	361/363 (99%)	346 (96%)	15 (4%)	0	100	100
All	All	720/726 (99%)	692 (96%)	28 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	288/335 (86%)	285 (99%)	3 (1%)	68	85
1	B	289/335 (86%)	285 (99%)	4 (1%)	59	80
All	All	577/670 (86%)	570 (99%)	7 (1%)	63	82

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	197	LEU
1	A	289	ASN
1	A	438	VAL
1	B	197	LEU
1	B	289	ASN
1	B	299	LYS
1	B	438	VAL

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	210	ASN
1	A	419	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 10 ligands modelled in this entry, 2 are monoatomic - leaving 8 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	9BS	B	604	-	13,13,13	0.12	0	16,16,16	0.32	0
3	9BS	B	602[B]	-	13,13,13	0.10	0	16,16,16	0.31	0
3	9BS	A	602[B]	-	13,13,13	0.10	0	16,16,16	0.30	0
3	9BS	B	603	-	13,13,13	0.10	0	16,16,16	0.29	0
3	9BS	B	602[C]	-	13,13,13	0.11	0	16,16,16	0.30	0
3	9BS	A	602[C]	-	13,13,13	0.11	0	16,16,16	0.31	0
3	9BS	B	602[A]	-	13,13,13	0.11	0	16,16,16	0.31	0
3	9BS	A	602[A]	-	13,13,13	0.10	0	16,16,16	0.31	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
3	9BS	B	604	-	-	0/4/4/4	0/2/2/2
3	9BS	B	602[B]	-	-	0/4/4/4	0/2/2/2
3	9BS	A	602[B]	-	-	0/4/4/4	0/2/2/2
3	9BS	B	603	-	-	0/4/4/4	0/2/2/2
3	9BS	B	602[C]	-	-	0/4/4/4	0/2/2/2
3	9BS	A	602[C]	-	-	0/4/4/4	0/2/2/2
3	9BS	B	602[A]	-	-	0/4/4/4	0/2/2/2
3	9BS	A	602[A]	-	-	0/4/4/4	0/2/2/2

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 [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	361/363 (99%)	0.52	35 (9%) 13 11	21, 56, 138, 224	0
1	B	361/363 (99%)	0.71	57 (15%) 5 4	19, 64, 172, 207	2 (0%)
All	All	722/726 (99%)	0.62	92 (12%) 8 7	19, 59, 161, 224	2 (0%)

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	305	SER	5.3
1	B	305	SER	5.2
1	B	174	LEU	4.2
1	B	304	GLY	4.1
1	A	291	ILE	3.9
1	B	179	ILE	3.9
1	A	257	PRO	3.9
1	A	304	GLY	3.8
1	A	373	GLU	3.8
1	A	161	GLY	3.8
1	B	426	ASP	3.7
1	A	306	PRO	3.6
1	B	427	LYS	3.6
1	B	294	THR	3.5
1	A	367	GLY	3.5
1	B	424	PHE	3.5
1	A	160	MET	3.5
1	A	368	ASN	3.4
1	A	256	ASN	3.4
1	B	366	GLU	3.3
1	B	178	ASP	3.2
1	A	366	GLU	3.1
1	B	257	PRO	3.1
1	B	307	ALA	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	296	VAL	3.1
1	A	293	ASP	3.1
1	B	182	ALA	3.1
1	B	256	ASN	3.0
1	B	212	GLY	3.0
1	B	429	HIS	2.9
1	A	221	SER	2.9
1	B	181	THR	2.9
1	B	176	ARG	2.9
1	B	214	TYR	2.8
1	B	369	GLY	2.8
1	B	175	SER	2.8
1	B	373	GLU	2.8
1	A	222	ALA	2.7
1	A	300	ARG	2.7
1	B	186	VAL	2.7
1	B	223	PRO	2.7
1	A	372	GLU	2.7
1	B	222	ALA	2.6
1	A	259	GLU	2.6
1	A	163	SER	2.5
1	B	221	SER	2.5
1	B	208	LEU	2.5
1	B	226	PHE	2.5
1	A	307	ALA	2.5
1	B	296	VAL	2.5
1	B	289	ASN	2.4
1	A	292	LYS	2.4
1	A	364	ALA	2.4
1	A	276	MET	2.4
1	B	292	LYS	2.4
1	B	368	ASN	2.3
1	B	372	GLU	2.3
1	B	308	VAL	2.3
1	B	370	PHE	2.3
1	A	255	ARG	2.3
1	B	300	ARG	2.3
1	B	194	LEU	2.3
1	B	220	ILE	2.3
1	B	185	MET	2.3
1	A	295	ASP	2.3
1	A	162	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	365	LYS	2.2
1	B	306	PRO	2.2
1	B	192	HIS	2.2
1	A	418	GLU	2.2
1	B	314	GLU	2.2
1	A	308	VAL	2.2
1	B	291	ILE	2.2
1	B	184	GLY	2.2
1	A	164	LYS	2.2
1	B	195	LEU	2.2
1	B	177	ASP	2.2
1	B	413	MET	2.2
1	A	234	VAL	2.2
1	B	287	GLU	2.1
1	B	401	GLU	2.1
1	B	266	LEU	2.1
1	B	215	TYR	2.1
1	B	290	ASP	2.1
1	B	196	ARG	2.1
1	B	259	GLU	2.1
1	B	363	HIS	2.1
1	A	185	MET	2.0
1	B	277	LEU	2.0
1	A	220	ILE	2.0
1	A	370	PHE	2.0
1	B	209	LEU	2.0

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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
3	9BS	B	604	12/12	0.73	0.26	84,85,86,87	12
3	9BS	B	603	12/12	0.76	0.20	84,85,86,86	12
3	9BS	A	602[C]	12/12	0.78	0.23	73,75,75,75	12
3	9BS	A	602[A]	12/12	0.78	0.23	74,75,75,76	12
3	9BS	A	602[B]	12/12	0.78	0.23	74,75,75,78	12
3	9BS	B	602[C]	12/12	0.89	0.32	82,85,87,88	12
3	9BS	B	602[A]	12/12	0.89	0.32	84,85,87,87	12
3	9BS	B	602[B]	12/12	0.89	0.32	84,85,86,87	12
2	ZN	B	601	1/1	0.99	0.04	38,38,38,38	0
2	ZN	A	601	1/1	1.00	0.01	33,33,33,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.