



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:55 PM UTC

PDB ID : 5R9S / pdb_00005r9s
Title : PanDDA analysis group deposition Form1 MAP kinase p38-alpha – Fragment N13470a in complex with MAP kinase p38-alpha
Authors : De Nicola, G.F.; Nichols, C.E.
Deposited on : 2020-03-04
Resolution : 1.70 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

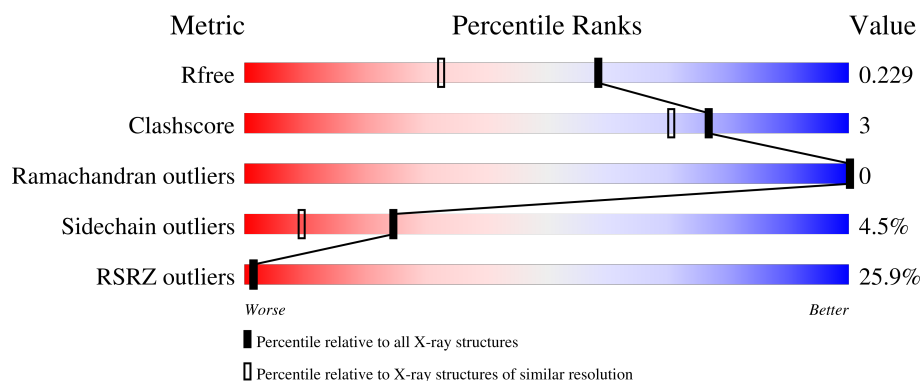
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	360	25% 87% 9% ..

2 Entry composition [i](#)

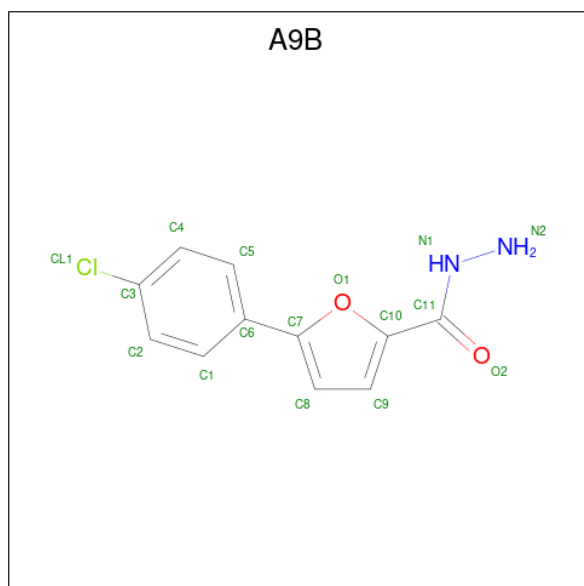
There are 7 unique types of molecules in this entry. The entry contains 3050 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Mitogen-activated protein kinase 14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	348	Total	C	N	O	S	0	4	0
			2781	1781	469	516	15			

- Molecule 2 is 5-(4-chlorophenyl)furan-2-carbohydrazide (CCD ID: A9B) (formula: $C_{11}H_9ClN_2O_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	O	0	0
			16	11	1	2	2		

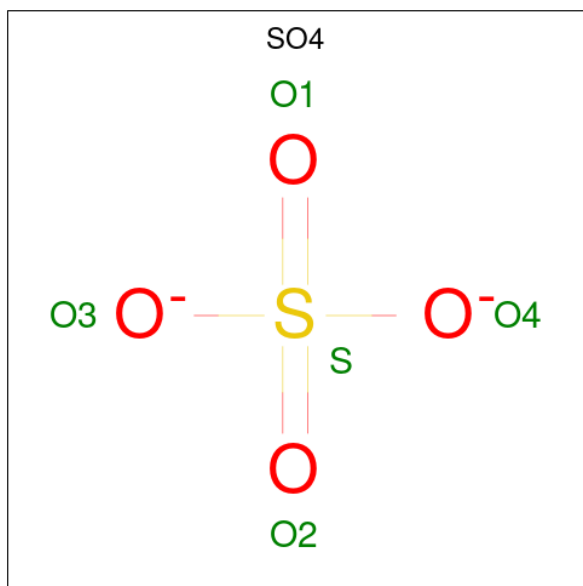
- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	5	Total	Cl	0	0
			5	5		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

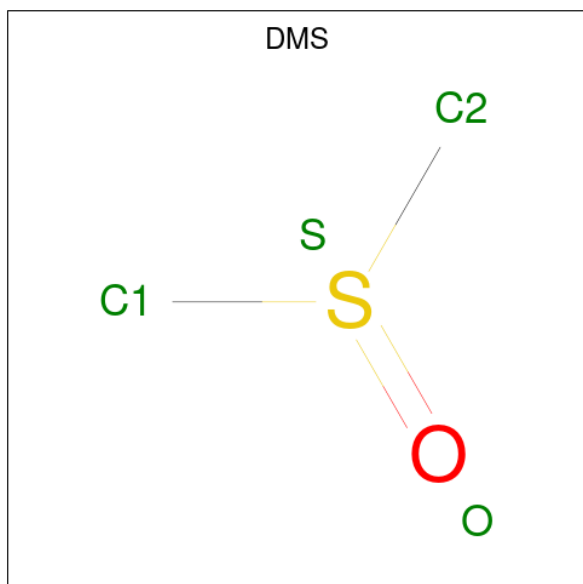
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C₂H₆OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	O	S	0	0
			4	2	1	1		

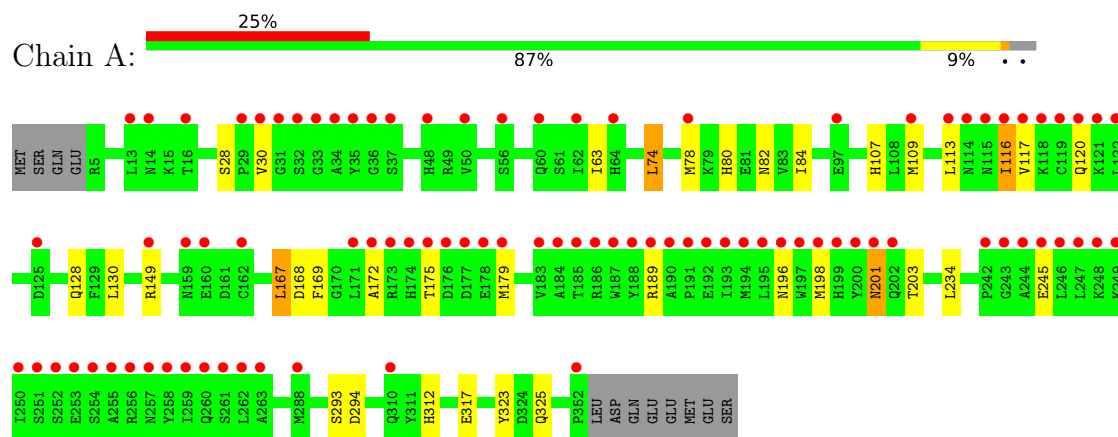
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	238	Total	O	0	0
			238	238		

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: Mitogen-activated protein kinase 14



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	45.78Å 86.04Å 127.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.30 – 1.70 29.30 – 1.70	Depositor EDS
% Data completeness (in resolution range)	98.0 (29.30-1.70) 98.2 (29.30-1.70)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.91 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.180 , 0.199 0.223 , 0.229	Depositor DCC
R_{free} test set	2805 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	23.4	Xtriage
Anisotropy	0.029	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 36.6	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	0.93	EDS
Total number of atoms	3050	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.00% 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: MG, DMS, SO4, CL, A9B

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	1.32	4/2847 (0.1%)	1.25	2/3873 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	116	ILE	C-N	-12.80	1.17	1.33
1	A	128	GLN	C-O	5.31	1.30	1.24
1	A	234	LEU	C-O	5.18	1.30	1.24
1	A	130	LEU	C-O	5.04	1.29	1.24

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	294	ASP	CA-CB-CG	5.54	118.14	112.60
1	A	116	ILE	O-C-N	-5.49	114.86	122.05

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	2781	0	2702	16	0
2	A	16	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	5	0	0	0	0
4	A	1	0	0	0	0
5	A	5	0	0	0	0
6	A	4	0	6	0	0
7	A	238	0	0	3	0
All	All	3050	0	2708	16	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 3.

All (16) 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:80:HIS:HD2	1:A:82:ASN:H	1.31	0.76
1:A:80:HIS:CD2	1:A:82:ASN:H	2.05	0.74
1:A:84:ILE:HD13	1:A:109:MET:HE3	1.81	0.61
1:A:312:HIS:HD2	1:A:317:GLU:OE2	1.87	0.58
1:A:201:ASN:ND2	7:A:504:HOH:O	2.38	0.56
1:A:78[B]:MET:HE3	1:A:169:PHE:CZ	2.46	0.51
1:A:117:VAL:CB	7:A:686:HOH:O	2.58	0.51
1:A:107:HIS:HD2	7:A:576:HOH:O	1.93	0.50
1:A:179:MET:HE1	1:A:189:ARG:NH1	2.29	0.48
1:A:167:LEU:HD12	1:A:168:ASP:CG	2.40	0.46
1:A:84:ILE:CD1	1:A:109:MET:HE3	2.46	0.43
1:A:74:LEU:HD13	1:A:78[B]:MET:HE2	2.01	0.43
1:A:323:TYR:CE2	1:A:325:GLN:HG2	2.55	0.41
1:A:149[B]:ARG:NH2	1:A:172:ALA:HA	2.35	0.41
1:A:84:ILE:HD12	1:A:84:ILE:HA	1.92	0.41
1:A:203:THR:OG1	1:A:293:SER:HB2	2.22	0.40

There are no symmetry-related clashes.

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	350/360 (97%)	337 (96%)	13 (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	295/319 (92%)	282 (96%)	13 (4%)	25	10

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	SER
1	A	30	VAL
1	A	63	ILE
1	A	74	LEU
1	A	113	LEU
1	A	116	ILE
1	A	120	GLN
1	A	167	LEU
1	A	175	THR
1	A	196	ASN
1	A	198	MET
1	A	201	ASN
1	A	245	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	HIS
1	A	80	HIS
1	A	120	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	128	GLN
1	A	201	ASN
1	A	272	ASN
1	A	310	GLN
1	A	312	HIS

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 9 ligands modelled in this entry, 6 are monoatomic - leaving 3 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	$\# Z > 2$
6	DMS	A	409	-	3,3,3	0.21	0	3,3,3	0.08	0
5	SO4	A	408	-	4,4,4	0.22	0	6,6,6	0.24	0
2	A9B	A	401	-	17,17,17	1.98	3 (17%)	22,23,23	2.34	9 (40%)

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	A9B	A	401	-	-	0/10/10/10	0/2/2/2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	A9B	C11-N1	6.47	1.39	1.33
2	A	401	A9B	C5-C4	2.87	1.43	1.38
2	A	401	A9B	C8-C7	2.21	1.39	1.35

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	A9B	O1-C7-C6	4.92	122.56	116.73
2	A	401	A9B	C1-C6-C7	-4.66	113.36	120.77
2	A	401	A9B	C5-C6-C7	4.31	127.61	120.77
2	A	401	A9B	O1-C10-C11	3.52	123.23	117.52
2	A	401	A9B	C9-C10-C11	-3.25	124.56	132.04
2	A	401	A9B	C2-C3-CL1	-2.28	115.99	119.36
2	A	401	A9B	C4-C3-CL1	2.24	122.66	119.36
2	A	401	A9B	O1-C10-C9	2.16	112.18	109.68
2	A	401	A9B	C6-C7-C8	-2.01	127.21	132.99

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	116:ILE	C	117:VAL	N	1.17

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	348/360 (96%)	1.87	90 (25%) ⓘ ⓘ	4, 21, 43, 62	66 (18%)

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	30	VAL	14.1
1	A	35	TYR	14.0
1	A	34	ALA	13.3
1	A	116	ILE	12.8
1	A	187	TRP	12.4
1	A	117	VAL	11.9
1	A	50	VAL	11.7
1	A	113	LEU	11.0
1	A	258	TYR	10.8
1	A	183	VAL	10.7
1	A	259	ILE	10.5
1	A	263	ALA	10.5
1	A	29	PRO	10.4
1	A	122	LEU	9.8
1	A	262	LEU	9.8
1	A	33	GLY	9.7
1	A	188	TYR	9.5
1	A	184	ALA	9.3
1	A	119	CYS	9.2
1	A	31	GLY	9.0
1	A	193	ILE	9.0
1	A	255	ALA	8.9
1	A	185	THR	8.8
1	A	37	SER	8.7
1	A	121	LYS	8.5
1	A	120	GLN	8.3
1	A	32	SER	8.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	190	ALA	8.0
1	A	250	ILE	7.8
1	A	260	GLN	7.7
1	A	118	LYS	7.7
1	A	242	PRO	7.7
1	A	149[A]	ARG	7.6
1	A	125[A]	ASP	7.6
1	A	36	GLY	7.6
1	A	115	ASN	7.4
1	A	194	MET	7.2
1	A	191	PRO	7.1
1	A	78[A]	MET	7.1
1	A	261	SER	7.0
1	A	256	ARG	7.0
1	A	244	ALA	7.0
1	A	195	LEU	6.9
1	A	114	ASN	6.9
1	A	257	ASN	6.8
1	A	254	SER	6.8
1	A	198	MET	6.8
1	A	288[A]	MET	6.6
1	A	186	ARG	6.6
1	A	201	ASN	6.6
1	A	247	LEU	6.5
1	A	243	GLY	6.3
1	A	253	GLU	6.2
1	A	197	TRP	6.2
1	A	189	ARG	6.2
1	A	174	HIS	6.0
1	A	246	LEU	6.0
1	A	199	HIS	5.8
1	A	175	THR	5.6
1	A	249	LYS	5.5
1	A	245	GLU	5.5
1	A	196	ASN	5.4
1	A	200	TYR	5.3
1	A	202	GLN	5.2
1	A	192	GLU	5.1
1	A	248	LYS	4.9
1	A	252	SER	4.8
1	A	173	ARG	4.8
1	A	172	ALA	4.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	178	GLU	4.2
1	A	251	SER	4.2
1	A	177	ASP	4.2
1	A	56	SER	4.0
1	A	97	GLU	3.9
1	A	352	PRO	3.8
1	A	13	LEU	3.5
1	A	14	ASN	3.3
1	A	109	MET	3.0
1	A	162	CYS	2.8
1	A	171	LEU	2.6
1	A	48	HIS	2.5
1	A	16	THR	2.5
1	A	160	GLU	2.5
1	A	310	GLN	2.4
1	A	60	GLN	2.4
1	A	62	ILE	2.3
1	A	64	HIS	2.2
1	A	159	ASN	2.1
1	A	176	ASP	2.1
1	A	179	MET	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	A9B	A	401	16/16	0.69	0.27	31,36,51,52	16
6	DMS	A	409	4/4	0.73	0.29	72,72,77,84	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	SO4	A	408	5/5	0.77	0.19	51,65,73,75	0
3	CL	A	404	1/1	0.94	0.15	45,45,45,45	0
3	CL	A	403	1/1	0.97	0.09	39,39,39,39	0
4	MG	A	407	1/1	0.97	0.21	29,29,29,29	0
3	CL	A	402	1/1	0.98	0.32	43,43,43,43	0
3	CL	A	405	1/1	0.98	0.27	37,37,37,37	0
3	CL	A	406	1/1	0.98	0.13	33,33,33,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.