



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:07 AM UTC

PDB ID : 2WI2 / pdb_00002wi2
Title : Orally Active 2-Amino Thienopyrimidine Inhibitors of the Hsp90 Chaperone
Authors : Brough, P.A.; Barril, X.; Borgognoni, J.; Chene, P.; Davies, N.G.M.; Davis, B.; Drysdale, M.J.; Dymock, B.; Eccles, S.A.; Garcia-Echeverria, C.; Fromont, C.; Hayes, A.; Hubbard, R.E.; Jordan, A.M.; Rugaard-Jensen, M.; Massey, A.; Merret, A.; Padfield, A.; Parsons, R.; Radimerski, T.; Raynaud, F.I.; Robertson, A.; Roughley, S.D.; Schoepfer, J.; Simmonite, H.; Surgenor, A.; Valenti, M.; Walls, S.; Webb, P.; Wood, M.; Workman, P.; Wright, L.M.
Deposited on : 2009-05-08
Resolution : 2.09 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

<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)

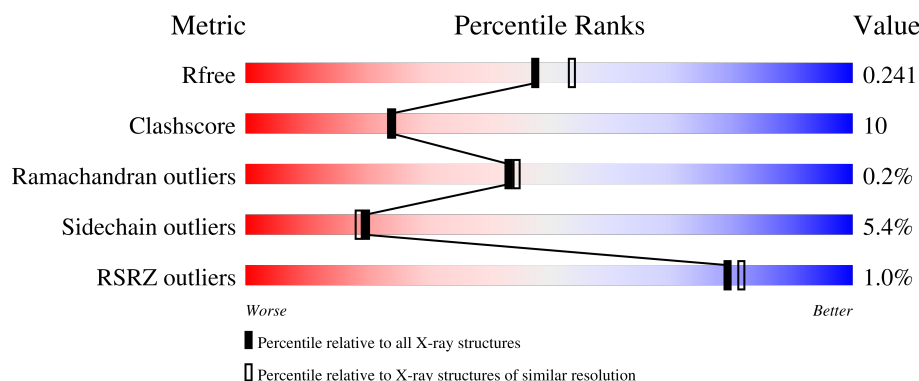
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

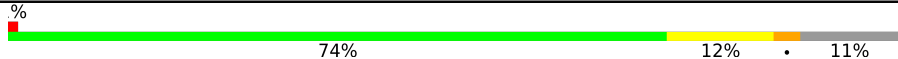
The reported resolution of this entry is 2.09 Å.

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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	236	 74% 12% • 11%
1	B	236	 68% 19% • 11%

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3528 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 HEAT SHOCK PROTEIN HSP 90-ALPHA.

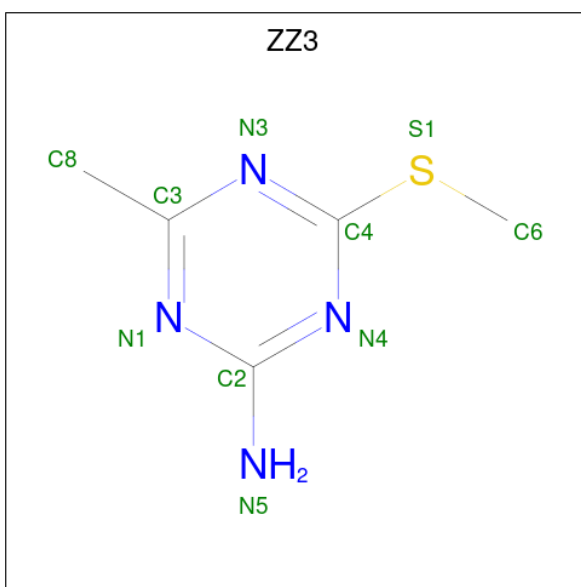
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	209	Total	C	N	O	S	0	1	1
			1634	1038	269	322	5			
1	B	209	Total	C	N	O	S	0	0	1
			1632	1036	270	321	5			

- Molecule 2 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C₂H₆OS).



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

- Molecule 3 is 4-METHYL-6-(METHYLSULFANYL)-1,3,5-TRIAZIN-2-AMINE (CCD ID: ZZ3) (formula: C₅H₈N₄S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	S	0	0
			10	5	4	1		
3	B	1	Total	C	N	S	0	0
			10	5	4	1		

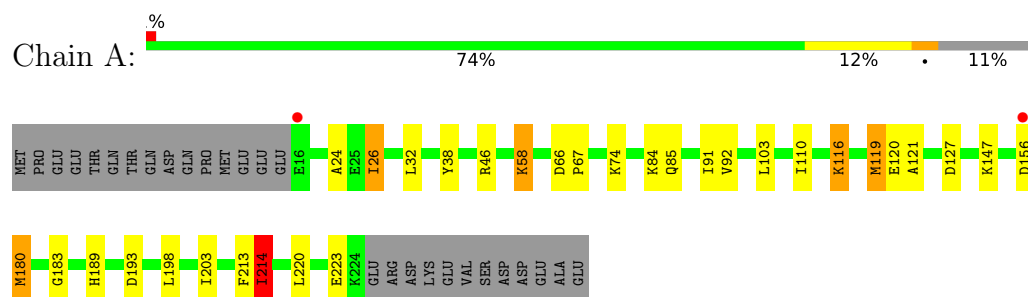
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	119	Total	O	0	0
			119	119		
4	B	119	Total	O	0	0
			119	119		

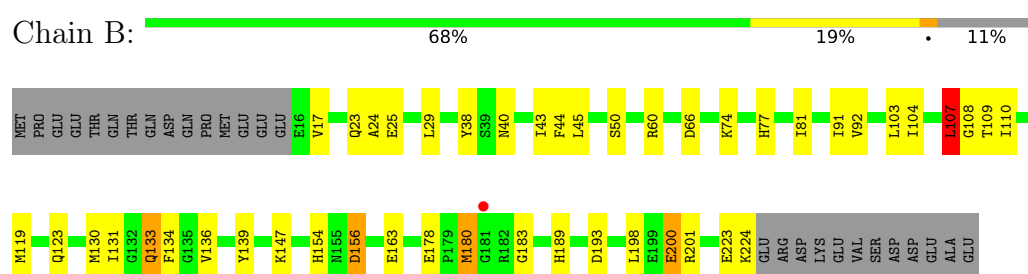
3 Residue-property plots [i](#)

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: HEAT SHOCK PROTEIN HSP 90-ALPHA



- Molecule 1: HEAT SHOCK PROTEIN HSP 90-ALPHA



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	64.27Å 88.44Å 98.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	100.00 – 2.09 98.59 – 2.09	Depositor EDS
% Data completeness (in resolution range)	92.3 (100.00-2.09) 92.3 (98.59-2.09)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 2.08Å)	Xtrriage
Refinement program	REFMAC 5.5.0066	Depositor
R, R_{free}	0.237 , 0.289 0.231 , 0.241	Depositor DCC
R_{free} test set	1591 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	31.8	Xtrriage
Anisotropy	0.385	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 38.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3528	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 40.09 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.8787e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: DMS, ZZ3

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	1.23	4/1663 (0.2%)	1.18	5/2244 (0.2%)
1	B	1.26	2/1658 (0.1%)	1.30	13/2237 (0.6%)
All	All	1.24	6/3321 (0.2%)	1.24	18/4481 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	223	GLU	C-N	-7.75	1.22	1.33
1	B	223	GLU	C-N	-6.10	1.24	1.33
1	B	24	ALA	CA-CB	5.61	1.62	1.53
1	A	38	TYR	CA-C	-5.31	1.46	1.52
1	A	92	VAL	C-O	-5.28	1.18	1.24
1	A	24	ALA	CA-CB	5.04	1.61	1.53

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	107	LEU	N-CA-C	8.96	124.34	113.41
1	B	156	ASP	N-CA-C	6.79	120.54	112.93
1	B	66	ASP	CA-C-N	6.55	126.18	119.56
1	B	66	ASP	C-N-CA	6.55	126.18	119.56
1	B	180	MET	CA-C-N	-6.43	108.81	121.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	180	MET	C-N-CA	-6.43	108.81	121.41
1	B	107	LEU	CA-C-N	6.17	132.80	121.70
1	B	107	LEU	C-N-CA	6.17	132.80	121.70
1	B	81	ILE	N-CA-C	6.01	114.10	108.15
1	B	81	ILE	CA-C-N	5.99	126.45	119.99
1	B	81	ILE	C-N-CA	5.99	126.45	119.99
1	B	17	VAL	N-CA-C	5.96	116.45	107.75
1	B	92	VAL	CB-CA-C	-5.83	101.94	110.62
1	A	58	LYS	N-CA-C	5.61	117.07	111.07
1	A	214	ILE	N-CA-C	-5.56	99.80	108.81
1	A	121	ALA	N-CA-C	-5.22	105.67	111.36
1	A	156	ASP	N-CA-C	5.05	119.02	113.21
1	A	85	GLN	N-CA-C	5.04	116.47	111.07

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	107	LEU	Peptide

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	1634	0	1630	22	0
1	B	1632	0	1630	43	0
2	A	4	0	6	0	1
3	A	10	0	8	0	0
3	B	10	0	8	1	0
4	A	119	0	0	4	1
4	B	119	0	0	8	0
All	All	3528	0	3282	66	1

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 10.

All (66) 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:32:LEU:HD13	1:A:119:MET:CE	1.87	1.05
1:A:32:LEU:HD13	1:A:119:MET:HE3	1.39	1.04
1:A:32:LEU:CD1	1:A:119:MET:HE2	1.90	1.01
1:B:29:LEU:CA	1:B:119:MET:HE2	1.94	0.97
1:B:154:HIS:CD2	1:B:156:ASP:H	1.87	0.93
1:A:32:LEU:HD12	1:A:119:MET:HE2	1.50	0.93
1:B:154:HIS:HD2	1:B:156:ASP:H	0.96	0.91
1:B:25:GLU:O	1:B:119:MET:HE1	1.69	0.90
1:A:32:LEU:CD1	1:A:119:MET:CE	2.48	0.89
1:B:29:LEU:HB2	1:B:119:MET:CE	2.07	0.84
1:A:189:HIS:HB3	4:A:2043:HOH:O	1.77	0.83
1:B:114:GLY:HA3	1:B:134:PHE:O	1.83	0.79
1:B:23:GLN:NE2	1:B:110:ILE:H	1.81	0.79
1:B:29:LEU:CB	1:B:119:MET:HE2	2.16	0.76
1:B:154:HIS:HD2	1:B:156:ASP:N	1.80	0.75
1:B:50:SER:HB2	4:B:2027:HOH:O	1.84	0.75
1:B:224:LYS:N	4:B:2119:HOH:O	2.19	0.75
1:B:29:LEU:HA	1:B:119:MET:HE2	1.69	0.75
1:B:29:LEU:HB2	1:B:119:MET:HE2	1.69	0.73
1:B:178:GLU:HG3	4:B:2096:HOH:O	1.89	0.72
3:B:1225:ZZ3:H63C	4:B:2011:HOH:O	1.92	0.69
1:A:26:ILE:HD12	1:A:26:ILE:N	2.07	0.69
1:A:110:ILE:O	1:A:110:ILE:HG13	1.91	0.69
1:B:104:ILE:O	1:B:108:GLY:HA2	1.92	0.69
1:B:29:LEU:N	1:B:119:MET:HE2	2.12	0.64
1:B:25:GLU:O	1:B:119:MET:CE	2.44	0.62
1:A:26:ILE:N	1:A:26:ILE:CD1	2.62	0.61
1:B:180:MET:HE2	1:B:183:GLY:C	2.26	0.60
1:A:127:ASP:OD1	4:A:2063:HOH:O	2.15	0.59
1:B:29:LEU:HB2	1:B:119:MET:HE3	1.84	0.59
1:B:200:GLU:HG2	4:B:2116:HOH:O	2.02	0.59
1:B:91:ILE:HD12	1:B:91:ILE:N	2.19	0.58
1:B:29:LEU:N	1:B:119:MET:CE	2.67	0.58
1:B:113:SER:HB2	1:B:136:VAL:HG13	1.85	0.57
1:A:180:MET:HE2	1:A:183:GLY:C	2.30	0.57
1:B:107:LEU:O	1:B:139:TYR:OH	2.17	0.56
1:B:23:GLN:HE22	1:B:110:ILE:H	1.54	0.56
1:B:104:ILE:O	1:B:108:GLY:CA	2.54	0.56
1:B:189:HIS:HB3	4:B:2054:HOH:O	2.06	0.56
1:B:180:MET:HE3	4:B:2058:HOH:O	2.07	0.55
1:A:26:ILE:CD1	1:A:26:ILE:H	2.21	0.54
1:B:103:LEU:O	1:B:107:LEU:HB2	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:ILE:O	1:A:110:ILE:CG1	2.54	0.50
1:B:130:MET:CE	1:B:133:GLN:HE22	2.25	0.49
1:B:43:ILE:C	1:B:43:ILE:HD12	2.38	0.49
1:B:23:GLN:NE2	1:B:109:THR:HB	2.27	0.48
1:B:130:MET:HE3	1:B:133:GLN:HE22	1.79	0.47
1:A:84:LYS:HA	1:A:198:LEU:HD13	1.98	0.45
1:B:147:LYS:HE3	1:B:163:GLU:CD	2.42	0.45
1:B:44:PHE:CG	1:B:45:LEU:N	2.84	0.45
1:A:213:PHE:HA	4:A:2107:HOH:O	2.16	0.44
1:B:29:LEU:CB	1:B:119:MET:CE	2.80	0.44
1:A:203:ILE:HG21	1:A:220:LEU:HD13	2.00	0.43
1:B:23:GLN:HE21	1:B:109:THR:HB	1.83	0.43
1:A:91:ILE:N	1:A:91:ILE:HD12	2.33	0.43
1:B:23:GLN:NE2	1:B:110:ILE:N	2.61	0.42
1:A:213:PHE:O	1:A:214:ILE:HD13	2.19	0.42
1:B:198:LEU:HD23	1:B:198:LEU:HA	1.72	0.42
1:A:66:ASP:HA	1:A:67:PRO:HD2	1.81	0.41
1:B:131:ILE:CG1	1:B:136:VAL:HG23	2.51	0.41
1:B:193:ASP:C	4:B:2099:HOH:O	2.64	0.41
1:A:116:LYS:O	1:A:120:GLU:HG3	2.21	0.41
1:B:77:HIS:CD2	1:B:77:HIS:N	2.89	0.40
1:B:38:TYR:CZ	1:B:40:ASN:HB3	2.57	0.40
1:A:176:THR:C	1:A:177:GLY:O	2.63	0.40
1:A:193:ASP:CG	4:A:2087:HOH:O	2.65	0.40

All (1) 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)
2:A:1225:DMS:C2	4:A:2118:HOH:O[2_755]	2.00	0.20

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	208/236 (88%)	200 (96%)	7 (3%)	1 (0%)	24	22
1	B	207/236 (88%)	195 (94%)	12 (6%)	0	100	100
All	All	415/472 (88%)	395 (95%)	19 (5%)	1 (0%)	43	44

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	177	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	178/206 (86%)	167 (94%)	11 (6%)	16	14
1	B	178/206 (86%)	170 (96%)	8 (4%)	24	25
All	All	356/412 (86%)	337 (95%)	19 (5%)	20	19

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	ILE
1	A	46	ARG
1	A	58	LYS
1	A	74	LYS
1	A	103	LEU
1	A	116	LYS
1	A	119	MET
1	A	147	LYS
1	A	178	GLU
1	A	180	MET
1	A	214	ILE
1	B	60	ARG
1	B	74	LYS

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Mol	Chain	Res	Type
1	B	107	LEU
1	B	116	LYS
1	B	123	GLN
1	B	133	GLN
1	B	200	GLU
1	B	201	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	ASN
1	A	40	ASN
1	A	79	ASN
1	A	105	ASN
1	A	106	ASN
1	B	23	GLN
1	B	79	ASN
1	B	105	ASN
1	B	106	ASN
1	B	123	GLN
1	B	133	GLN
1	B	154	HIS

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 ⓘ

3 ligands are 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$
3	ZZ3	B	1225	-	10,10,10	2.31	1 (10%)	13,13,13	4.48	7 (53%)
3	ZZ3	A	1226	-	10,10,10	2.34	1 (10%)	13,13,13	4.48	7 (53%)
2	DMS	A	1225	-	3,3,3	2.30	1 (33%)	3,3,3	0.23	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	ZZ3	B	1225	-	-	2/2/2/2	0/1/1/1
3	ZZ3	A	1226	-	-	0/2/2/2	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1226	ZZ3	C4-S1	-7.07	1.69	1.75
3	B	1225	ZZ3	C4-S1	-6.97	1.70	1.75
2	A	1225	DMS	O-S	3.90	1.76	1.50

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1225	ZZ3	C6-S1-C4	9.62	109.47	102.25
3	A	1226	ZZ3	C6-S1-C4	9.60	109.45	102.25
3	A	1226	ZZ3	C4-N4-C2	9.47	119.99	113.46
3	B	1225	ZZ3	C4-N4-C2	9.46	119.99	113.46
3	A	1226	ZZ3	N3-C4-N4	-3.84	120.00	127.00
3	B	1225	ZZ3	N3-C4-N4	-3.83	120.01	127.00
3	A	1226	ZZ3	C2-N1-C3	3.70	120.02	115.25
3	B	1225	ZZ3	C2-N1-C3	3.69	120.01	115.25
3	A	1226	ZZ3	N1-C2-N4	-3.60	119.96	125.48
3	B	1225	ZZ3	N1-C2-N4	-3.57	120.01	125.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1225	ZZ3	N3-C3-N1	-3.28	119.99	125.77
3	A	1226	ZZ3	N3-C3-N1	-3.26	120.01	125.77
3	A	1226	ZZ3	C3-N3-C4	3.10	120.02	113.90
3	B	1225	ZZ3	C3-N3-C4	3.10	120.01	113.90

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1225	ZZ3	N4-C4-S1-C6
3	B	1225	ZZ3	N3-C4-S1-C6

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1225	ZZ3	1	0
2	A	1225	DMS	0	1

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 [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	209/236 (88%)	-0.22	3 (1%) 73 75	16, 24, 35, 44	1 (0%)
1	B	209/236 (88%)	-0.21	1 (0%) 87 88	16, 25, 38, 44	0
All	All	418/472 (88%)	-0.21	4 (0%) 79 81	16, 25, 37, 44	1 (0%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	177	GLY	3.6
1	A	16	GLU	2.4
1	B	181	GLY	2.1
1	A	156	ASP	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(Å ²)	Q<0.9
2	DMS	A	1225	4/4	0.78	0.29	55,55,57,58	4

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ZZ3	B	1225	10/10	0.92	0.10	18,24,34,35	0
3	ZZ3	A	1226	10/10	0.97	0.06	19,23,26,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.