



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:09 PM UTC

PDB ID : 5QYI / pdb_00005qyi
Title : PanDDA analysis group deposition – Aar2/RNaseH in complex with fragment F2X-Entry G12a
Authors : Weiss, M.S.; Wollenhaupt, J.; Metz, A.; Barthel, T.; Lima, G.M.A.; Heine, A.; Mueller, U.; Klebe, G.
Deposited on : 2020-02-12
Resolution : 1.58 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:

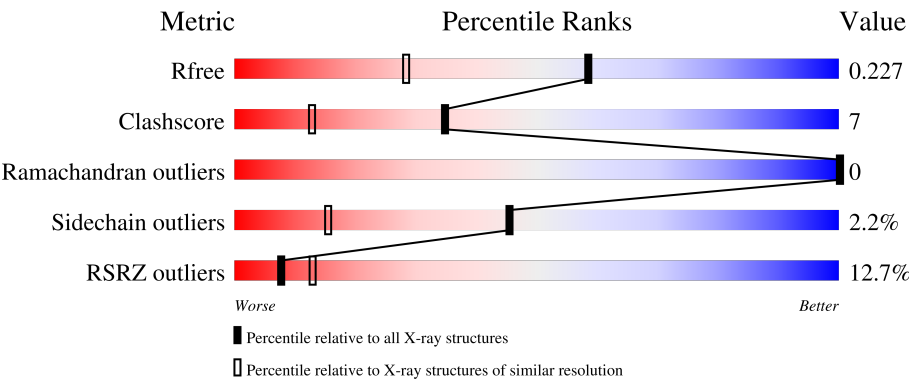
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal 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 i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.58 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	1094 (1.58-1.58)
Clashscore	190562	1105 (1.58-1.58)
Ramachandran outliers	187476	1082 (1.58-1.58)
Sidechain outliers	187428	1081 (1.58-1.58)
RSRZ outliers	180081	1094 (1.58-1.58)

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	258	10% 83%8%8%
2	B	308	14% 85%12%.

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6466 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 Pre-mRNA-splicing factor 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	237	Total	C	N	O	S	0	13	0
			2429	1526	421	463	19			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1833	GLY	-	expression tag	UNP P33334
A	1834	ALA	-	expression tag	UNP P33334
A	1835	MET	-	expression tag	UNP P33334

- Molecule 2 is a protein called A1 cistron-splicing factor AAR2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	300	Total	C	N	O	S	28	32	0
			3854	2488	633	700	33			

There are 9 discrepancies between the modelled and reference sequences:

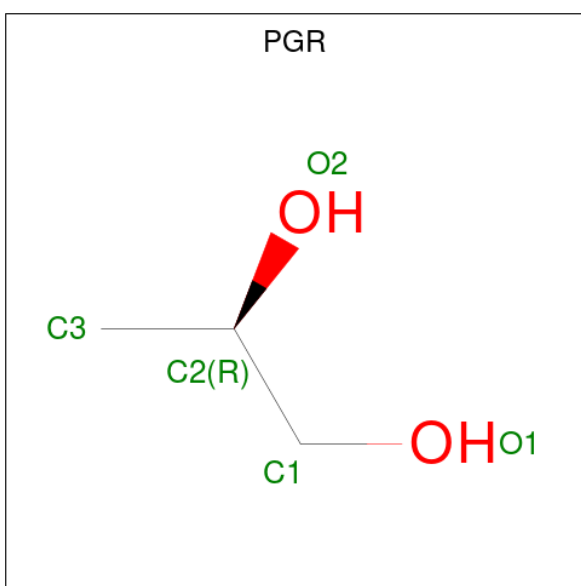
Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	GLY	-	expression tag	UNP P32357
B	-2	ALA	-	expression tag	UNP P32357
B	-1	MET	-	expression tag	UNP P32357
B	0	ALA	-	expression tag	UNP P32357
B	166	SER	-	linker	UNP P32357
B	167	SER	-	linker	UNP P32357
B	168	SER	-	linker	UNP P32357
B	169	SER	-	linker	UNP P32357
B	170	SER	-	linker	UNP P32357

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 4 is R-1,2-PROPANEDIOL (CCD ID: PGR) (formula: $C_3H_8O_2$).



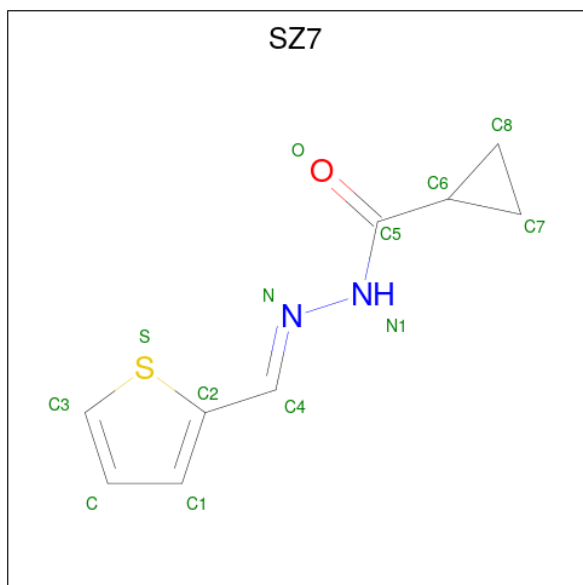
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			5	3	2		
4	A	1	Total	C	O	0	0
			5	3	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			5	3	2		

- Molecule 5 is {N}-[({E})-thiophen-2-ylmethylideneamino]cyclopropanecarboxamide (CCD ID: SZ7) (formula: C₉H₁₀N₂OS) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	S	0	1
			39	27	6	3	3		

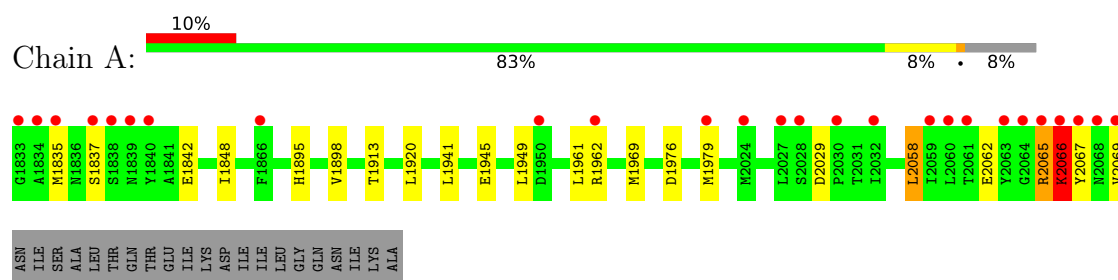
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	57	Total	O	0	0
			57	57		
6	B	55	Total	O	0	2
			62	62		

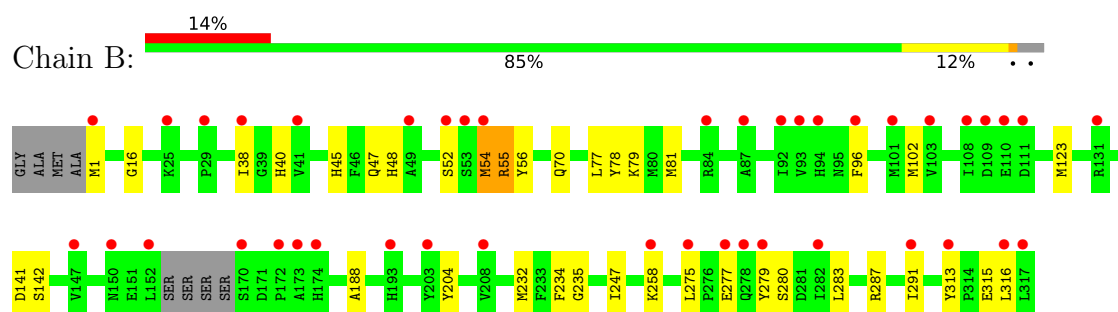
3 Residue-property plots [i](#)

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: Pre-mRNA-splicing factor 8



- Molecule 2: A1 cistron-splicing factor AAR2



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	89.20Å 82.17Å 93.29Å 90.00° 108.02° 90.00°	Depositor
Resolution (Å)	44.92 – 1.58 44.92 – 1.58	Depositor EDS
% Data completeness (in resolution range)	97.3 (44.92-1.58) 97.5 (44.92-1.58)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.95 (at 1.58Å)	Xtriage
Refinement program	REFMAC 5.8.0238, PHENIX 1.16.3549	Depositor
R, R_{free}	0.219 , 0.220 0.229 , 0.227	Depositor DCC
R_{free} test set	4353 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	34.8	Xtriage
Anisotropy	0.073	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 40.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6466	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.66% 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: SZ7, PGR, SO4

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	1.08	2/2470 (0.1%)	1.25	4/3332 (0.1%)
2	B	0.99	0/3965	1.13	3/5358 (0.1%)
All	All	1.02	2/6435 (0.0%)	1.18	7/8690 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1920	LEU	C-O	6.23	1.31	1.24
1	A	1913	THR	C-O	5.05	1.29	1.24

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	2066	LYS	CB-CA-C	-6.56	100.58	110.88
2	B	247	ILE	N-CA-C	-5.66	105.21	110.53
2	B	141	ASP	CA-C-N	5.47	127.87	120.38
2	B	141	ASP	C-N-CA	5.47	127.87	120.38
1	A	2069	VAL	CA-C-O	-5.28	111.83	120.80
1	A	1949	LEU	CA-C-N	5.02	128.63	120.60
1	A	1949	LEU	C-N-CA	5.02	128.63	120.60

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	2429	0	2462	45	0
2	B	3854	0	3646	45	0
3	A	10	0	0	0	0
4	A	15	0	24	1	0
5	B	39	0	0	0	0
6	A	57	0	0	1	0
6	B	62	0	0	1	0
All	All	6466	0	6132	90	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 7.

All (90) 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:1979[B]:MET:HA	1:A:1979[B]:MET:HE2	1.62	0.79
1:A:1979[E]:MET:HA	1:A:1979[E]:MET:HE2	1.63	0.79
1:A:2065[A]:ARG:CG	1:A:2065[A]:ARG:HH11	1.96	0.78
1:A:2065[B]:ARG:CG	1:A:2065[B]:ARG:HH11	1.96	0.78
1:A:2065[C]:ARG:HH11	1:A:2065[C]:ARG:CG	1.97	0.77
2:B:56[A]:TYR:HH	2:B:78:TYR:H	1.32	0.74
2:B:56[C]:TYR:HH	2:B:78:TYR:H	1.32	0.74
2:B:56[D]:TYR:HH	2:B:78:TYR:H	1.33	0.72
2:B:56[F]:TYR:HH	2:B:78:TYR:H	1.33	0.72
2:B:70:GLN:HB3	2:B:81:MET:CE	2.19	0.72
1:A:2066:LYS:CA	1:A:2066:LYS:CE	2.69	0.70
2:B:96:PHE:HB2	2:B:102:MET:HE3	1.72	0.70
2:B:56[B]:TYR:HH	2:B:78:TYR:H	1.35	0.70
1:A:2066:LYS:N	1:A:2066:LYS:HE3	2.08	0.68
2:B:70:GLN:HB3	2:B:81:MET:HE2	1.75	0.67
1:A:2062:GLU:O	1:A:2066:LYS:HE3	1.94	0.67
1:A:2065[C]:ARG:HH11	1:A:2065[C]:ARG:HG2	1.61	0.65
1:A:2058:LEU:C	1:A:2058:LEU:HD23	2.22	0.65
1:A:2065[A]:ARG:HH11	1:A:2065[A]:ARG:HG2	1.62	0.65
1:A:2065[B]:ARG:HH11	1:A:2065[B]:ARG:HG2	1.62	0.65
2:B:16:GLY:HA3	2:B:45[A]:HIS:CE1	2.32	0.64
1:A:1941:LEU:O	1:A:1945:GLU:HG3	1.99	0.63
2:B:280:SER:HB3	2:B:313:TYR:CE1	2.34	0.63
1:A:1835:MET:HE1	1:A:1961:LEU:HD12	1.81	0.62
1:A:2066:LYS:CA	1:A:2066:LYS:HE3	2.28	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1969:MET:HA	1:A:1969:MET:HE2	1.84	0.60
2:B:16:GLY:HA3	2:B:45[C]:HIS:CE1	2.36	0.60
2:B:47[A]:GLN:OE1	2:B:52[A]:SER:O	2.20	0.59
1:A:1842:GLU:OE1	6:A:2201:HOH:O	2.18	0.55
2:B:96:PHE:CB	2:B:102:MET:HE3	2.37	0.54
2:B:40:HIS:HD2	6:B:533:HOH:O	1.91	0.53
2:B:287:ARG:O	2:B:291:ILE:HD13	2.08	0.53
1:A:2065[A]:ARG:CG	1:A:2065[A]:ARG:NH1	2.62	0.52
1:A:2065[B]:ARG:CG	1:A:2065[B]:ARG:NH1	2.62	0.52
1:A:2066:LYS:CE	1:A:2066:LYS:N	2.72	0.52
2:B:16:GLY:HA3	2:B:45[B]:HIS:CE1	2.45	0.52
1:A:2065[C]:ARG:CG	1:A:2065[C]:ARG:NH1	2.63	0.52
2:B:48[B]:HIS:HE1	2:B:56[B]:TYR:OH	1.93	0.52
2:B:77:LEU:HD21	2:B:79:LYS:HE3	1.93	0.51
1:A:2066:LYS:CE	1:A:2066:LYS:HA	2.40	0.51
2:B:1:MET:N	2:B:38:ILE:HG13	2.27	0.50
2:B:48[A]:HIS:HE1	2:B:56[A]:TYR:OH	1.95	0.49
2:B:48[F]:HIS:HE1	2:B:56[F]:TYR:OH	1.96	0.49
2:B:48[D]:HIS:HE1	2:B:56[D]:TYR:OH	1.96	0.49
2:B:234[D]:PHE:O	2:B:235[D]:GLY:C	2.56	0.48
2:B:234[F]:PHE:O	2:B:235[F]:GLY:C	2.56	0.48
2:B:48[C]:HIS:HE1	2:B:56[C]:TYR:OH	1.96	0.48
2:B:47[D]:GLN:NE2	2:B:48[D]:HIS:O	2.46	0.48
2:B:47[F]:GLN:NE2	2:B:48[F]:HIS:O	2.46	0.48
1:A:2062:GLU:O	1:A:2066:LYS:CE	2.61	0.47
1:A:2066:LYS:HB3	1:A:2067:TYR:CD1	2.49	0.47
2:B:258:LYS:HD2	2:B:258:LYS:H	1.81	0.46
2:B:279[C]:TYR:HB3	2:B:283[C]:LEU:HD12	1.98	0.46
2:B:279[A]:TYR:HB3	2:B:283[A]:LEU:HD12	1.98	0.45
1:A:1848:ILE:HG12	4:A:2104:PGR:H11	1.98	0.45
1:A:2065[A]:ARG:HH11	1:A:2065[A]:ARG:HG3	1.79	0.44
1:A:2065[B]:ARG:HH11	1:A:2065[B]:ARG:HG3	1.79	0.44
1:A:2065[A]:ARG:NH1	1:A:2065[A]:ARG:HG3	2.31	0.44
1:A:2065[B]:ARG:NH1	1:A:2065[B]:ARG:HG3	2.31	0.44
2:B:54[B]:MET:HG2	2:B:55[B]:ARG:N	2.31	0.44
2:B:45[D]:HIS:N	2:B:45[D]:HIS:CD2	2.86	0.44
2:B:45[F]:HIS:N	2:B:45[F]:HIS:CD2	2.86	0.44
2:B:55[B]:ARG:CZ	2:B:55[B]:ARG:HB2	2.48	0.43
1:A:2065[C]:ARG:NH1	1:A:2065[C]:ARG:HG3	2.32	0.43
2:B:232[D]:MET:O	2:B:232[D]:MET:HG2	2.19	0.43
2:B:232[F]:MET:O	2:B:232[F]:MET:HG2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:54[A]:MET:HG2	2:B:55[A]:ARG:N	2.34	0.43
2:B:277:GLU:CD	2:B:277:GLU:H	2.27	0.43
1:A:2065[C]:ARG:HH11	1:A:2065[C]:ARG:HG3	1.81	0.42
2:B:77:LEU:HD21	2:B:79:LYS:CE	2.49	0.42
1:A:2029:ASP:OD1	1:A:2029:ASP:C	2.63	0.41
2:B:54[D]:MET:HG2	2:B:55[D]:ARG:N	2.35	0.41
2:B:54[F]:MET:HG2	2:B:55[F]:ARG:N	2.35	0.41
1:A:1979[B]:MET:HA	1:A:1979[B]:MET:CE	2.43	0.41
1:A:2058:LEU:C	1:A:2058:LEU:CD2	2.92	0.41
1:A:1895:HIS:O	1:A:1898[C]:VAL:HG22	2.20	0.41
1:A:1895:HIS:O	1:A:1898[D]:VAL:HG22	2.20	0.41
1:A:1895:HIS:O	1:A:1898[F]:VAL:HG22	2.20	0.41
1:A:1979[E]:MET:HA	1:A:1979[E]:MET:CE	2.43	0.41
1:A:1895:HIS:O	1:A:1898[A]:VAL:HG22	2.20	0.41
2:B:142:SER:HA	2:B:234[F]:PHE:O	2.20	0.41
2:B:188:ALA:HA	2:B:204:TYR:CD1	2.57	0.40
2:B:142:SER:HA	2:B:234[D]:PHE:O	2.20	0.40
2:B:275:LEU:HD22	2:B:283[A]:LEU:HD13	2.03	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	300/258 (116%)	297 (99%)	3 (1%)	0	100	100
2	B	456/308 (148%)	445 (98%)	11 (2%)	0	100	100
All	All	756/566 (134%)	742 (98%)	14 (2%)	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	280/233 (120%)	269 (96%)	11 (4%)	28	4
2	B	422/284 (149%)	407 (96%)	15 (4%)	31	5
All	All	702/517 (136%)	676 (96%)	26 (4%)	45	5

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1837	SER
1	A	1962	ARG
1	A	1976	ASP
1	A	2058	LEU
1	A	2065[A]	ARG
1	A	2065[B]	ARG
1	A	2065[C]	ARG
1	A	2065[D]	ARG
1	A	2065[E]	ARG
1	A	2065[F]	ARG
1	A	2066	LYS
2	B	54[A]	MET
2	B	54[B]	MET
2	B	54[C]	MET
2	B	54[D]	MET
2	B	54[E]	MET
2	B	54[F]	MET
2	B	55[A]	ARG
2	B	55[B]	ARG
2	B	55[C]	ARG
2	B	55[D]	ARG
2	B	55[E]	ARG
2	B	55[F]	ARG
2	B	123	MET
2	B	315	GLU
2	B	316	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	40	HIS
2	B	136	GLN
2	B	300	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](#)

8 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	SO4	A	2102	-	4,4,4	0.35	0	6,6,6	0.17	0
5	SZ7	B	401[F]	-	14,14,14	0.66	0	18,18,18	0.39	0
4	PGR	A	2103	-	4,4,4	0.60	0	4,4,4	0.38	0
4	PGR	A	2105	-	4,4,4	0.53	0	4,4,4	0.95	0
5	SZ7	B	401[E]	-	14,14,14	0.78	1 (7%)	18,18,18	0.47	0
4	PGR	A	2104	-	4,4,4	0.55	0	4,4,4	1.13	0
5	SZ7	B	401[D]	-	14,14,14	0.66	0	18,18,18	0.38	0
3	SO4	A	2101	-	4,4,4	0.27	0	6,6,6	0.13	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
5	SZ7	B	401[F]	-	-	1/10/12/12	0/2/2/2
4	PGR	A	2103	-	-	2/2/2/2	-
5	SZ7	B	401[E]	-	-	1/10/12/12	0/2/2/2
4	PGR	A	2105	-	-	1/2/2/2	-
4	PGR	A	2104	-	-	0/2/2/2	-
5	SZ7	B	401[D]	-	-	1/10/12/12	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	401[E]	SZ7	C3-S	2.32	1.80	1.70

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	2103	PGR	O1-C1-C2-C3
4	A	2103	PGR	O1-C1-C2-O2
4	A	2105	PGR	O1-C1-C2-O2
5	B	401[D]	SZ7	N1-C5-C6-C7
5	B	401[E]	SZ7	N1-C5-C6-C7
5	B	401[F]	SZ7	N1-C5-C6-C7

There are no ring outliers.

1 monomer is involved in 1 short contact:

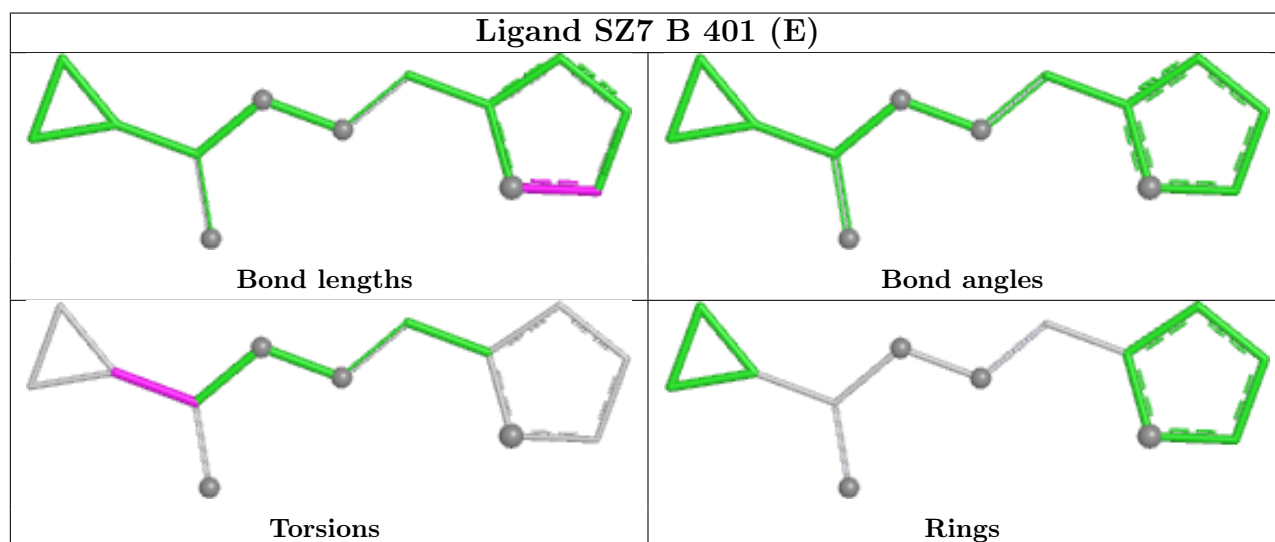
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	2104	PGR	1	0

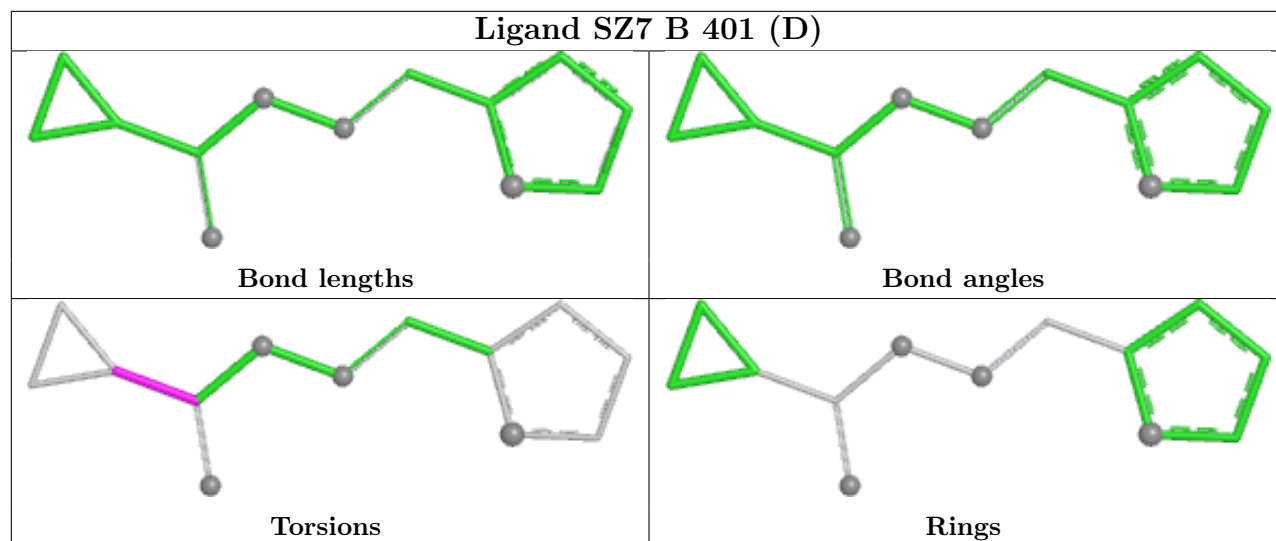
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring

Ligand SZ7 B 401 (F)

The figure displays four panels illustrating the structural analysis of Ligand SZ7 B 401 (F). The panels are arranged in a 2x2 grid, showing the molecule's geometry and highlighting specific features.

- Bond lengths:** The top-left panel shows the molecule with green sticks, highlighting the bond lengths.
- Bond angles:** The top-right panel shows the molecule with green sticks, highlighting the bond angles.
- Torsions:** The bottom-left panel shows the molecule with green sticks, highlighting the torsions. The cyclopropyl group is highlighted in magenta.
- Rings:** The bottom-right panel shows the molecule with green sticks, highlighting the rings. The cyclopropyl group is highlighted in magenta.





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	237/258 (91%)	0.76	26 (10%)	10 17	4, 32, 61, 76	13 (5%)
2	B	300/308 (97%)	0.90	42 (14%)	6 10	3, 34, 62, 101	32 (10%)
All	All	537/566 (94%)	0.84	68 (12%)	8 13	3, 34, 62, 101	45 (8%)

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	152	LEU	7.3
1	A	1979[A]	MET	5.6
2	B	316	LEU	5.4
2	B	54[A]	MET	5.3
2	B	174	HIS	4.6
1	A	2069	VAL	4.5
2	B	52[A]	SER	4.5
2	B	173	ALA	4.3
2	B	108	ILE	4.2
1	A	2065[A]	ARG	4.0
1	A	2067	TYR	3.8
2	B	203[A]	TYR	3.8
2	B	170	SER	3.7
1	A	1834	ALA	3.5
1	A	1866	PHE	3.5
1	A	1833	GLY	3.4
1	A	2060	LEU	3.4
2	B	1	MET	3.4
1	A	1837	SER	3.4
2	B	279[A]	TYR	3.4
1	A	2066	LYS	3.3
2	B	172	PRO	3.3
2	B	84	ARG	3.3
1	A	2063	TYR	3.2

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Mol	Chain	Res	Type	RSRZ
2	B	87	ALA	3.2
2	B	131	ARG	3.1
2	B	275	LEU	3.1
1	A	2024[A]	MET	3.1
2	B	317	LEU	3.1
1	A	2027	LEU	3.0
1	A	2064	GLY	3.0
2	B	96	PHE	3.0
2	B	101	MET	3.0
1	A	2068	ASN	2.9
2	B	282[A]	ILE	2.9
2	B	277	GLU	2.9
2	B	41[A]	VAL	2.9
2	B	313	TYR	2.8
1	A	1835	MET	2.8
2	B	150	ASN	2.8
1	A	2030	PRO	2.7
2	B	49[A]	ALA	2.7
2	B	147	VAL	2.6
1	A	2028	SER	2.6
2	B	93	VAL	2.6
2	B	38	ILE	2.5
2	B	29	PRO	2.5
2	B	53[A]	SER	2.5
2	B	109	ASP	2.5
1	A	2059	ILE	2.5
2	B	258	LYS	2.5
2	B	110	GLU	2.4
1	A	1962	ARG	2.4
1	A	2032	ILE	2.4
1	A	1838	SER	2.4
1	A	1950	ASP	2.3
1	A	1840	TYR	2.3
2	B	92	ILE	2.3
2	B	291	ILE	2.3
2	B	278	GLN	2.2
1	A	2061	THR	2.2
2	B	103	VAL	2.1
2	B	111	ASP	2.1
2	B	94	HIS	2.0
2	B	25	LYS	2.0
2	B	193	HIS	2.0

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Mol	Chain	Res	Type	RSRZ
2	B	208	VAL	2.0
1	A	1839	ASN	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](#)

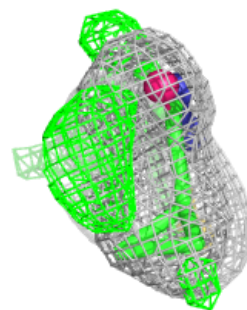
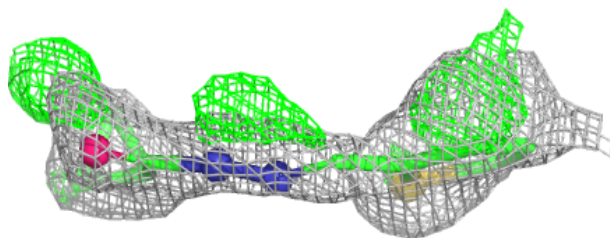
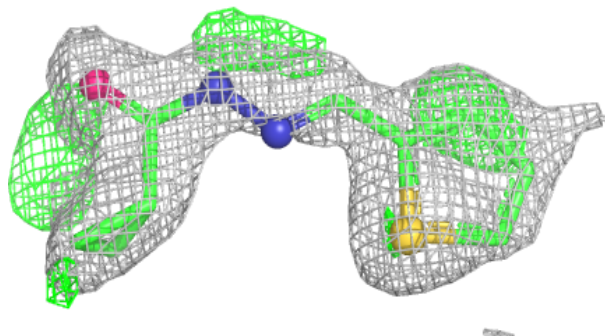
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	PGR	A	2104	5/5	0.78	0.24	34,45,52,59	0
4	PGR	A	2105	5/5	0.81	0.20	46,47,54,59	0
5	SZ7	B	401[D]	13/13	0.86	0.32	43,46,48,48	13
5	SZ7	B	401[E]	13/13	0.86	0.32	45,46,48,48	13
5	SZ7	B	401[F]	13/13	0.86	0.32	43,46,48,48	13
4	PGR	A	2103	5/5	0.90	0.15	39,44,46,48	0
3	SO4	A	2102	5/5	0.93	0.09	56,62,70,70	0
3	SO4	A	2101	5/5	0.98	0.05	37,46,50,52	0

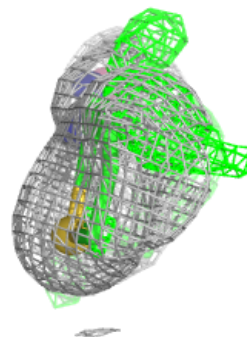
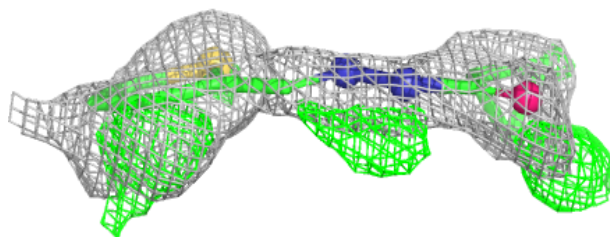
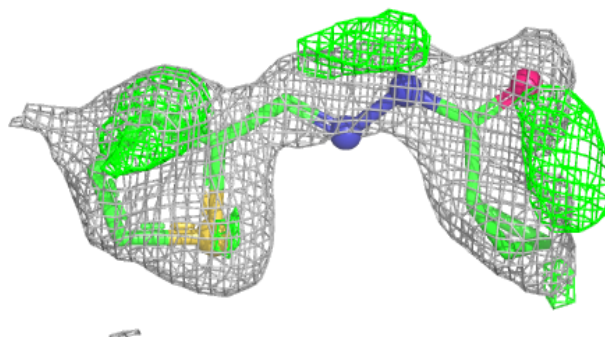
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around SZ7 B 401 (D):

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

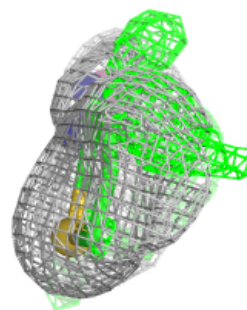
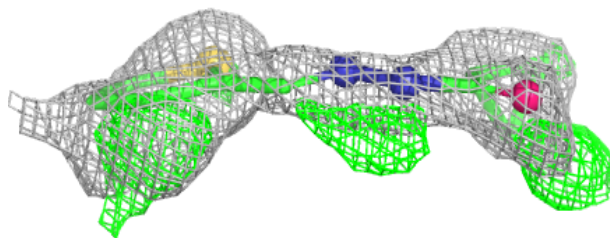
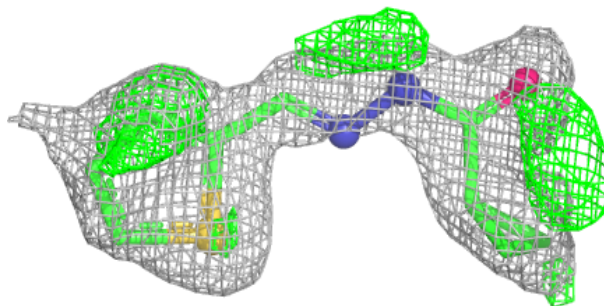
**Electron density around SZ7 B 401 (E):**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around SZ7 B 401 (F):

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.