



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 8, 2025 – 04:16 pm BST

PDB ID : 9S02 / pdb_00009s02
Title : PYCR1 in complex with 3-(2-thiazolyl)propionic acid
Authors : Ragin-Oh, W.; Czerwonka, D.; Ruszkowski, M.
Deposited on : 2025-07-16
Resolution : 1.65 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 : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 2.0
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
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) : 2.46

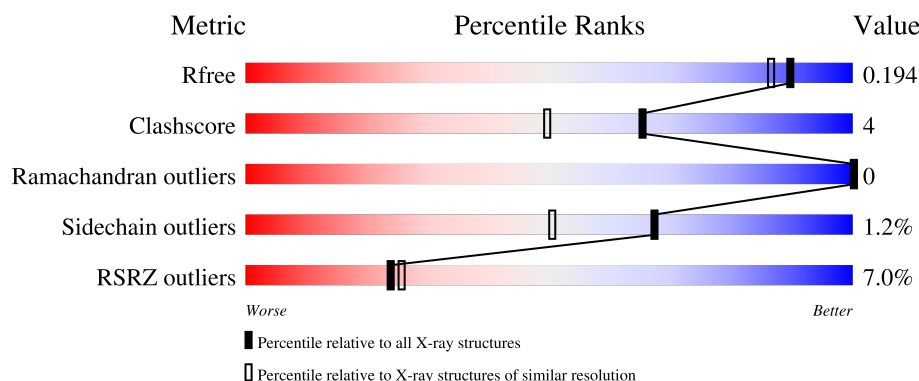
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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}	164625	2328 (1.66-1.66)
Clashscore	180529	2515 (1.66-1.66)
Ramachandran outliers	177936	2475 (1.66-1.66)
Sidechain outliers	177891	2475 (1.66-1.66)
RSRZ outliers	164620	2328 (1.66-1.66)

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	342	<div> <div>6%</div> <div> <div></div> <div>74%</div> <div>6%</div> <div>20%</div> </div> </div>
1	B	342	<div> <div>5%</div> <div> <div></div> <div>75%</div> <div>6%</div> <div>19%</div> </div> </div>
1	C	342	<div> <div>6%</div> <div> <div></div> <div>73%</div> <div>7%</div> <div>20%</div> </div> </div>
1	D	342	<div> <div>5%</div> <div> <div></div> <div>74%</div> <div>6%</div> <div>20%</div> </div> </div>
1	E	342	<div> <div>6%</div> <div> <div></div> <div>74%</div> <div>6%</div> <div>20%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	EDO	D	404	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11140 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 Isoform 3 of Pyrroline-5-carboxylate reductase 1, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	273	Total	C	N	O	S	0	4	0
			2027	1278	358	377	14			
1	B	278	Total	C	N	O	S	0	5	0
			2075	1310	365	386	14			
1	C	273	Total	C	N	O	S	0	3	0
			2023	1275	358	376	14			
1	D	275	Total	C	N	O	S	0	4	0
			2042	1286	361	381	14			
1	E	273	Total	C	N	O	S	0	4	0
			2027	1278	358	377	14			

There are 110 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	initiating methionine	UNP P32322
A	-21	HIS	-	expression tag	UNP P32322
A	-20	HIS	-	expression tag	UNP P32322
A	-19	HIS	-	expression tag	UNP P32322
A	-18	HIS	-	expression tag	UNP P32322
A	-17	HIS	-	expression tag	UNP P32322
A	-16	HIS	-	expression tag	UNP P32322
A	-15	SER	-	expression tag	UNP P32322
A	-14	SER	-	expression tag	UNP P32322
A	-13	GLY	-	expression tag	UNP P32322
A	-12	VAL	-	expression tag	UNP P32322
A	-11	ASP	-	expression tag	UNP P32322
A	-10	LEU	-	expression tag	UNP P32322
A	-9	GLY	-	expression tag	UNP P32322
A	-8	THR	-	expression tag	UNP P32322
A	-7	GLU	-	expression tag	UNP P32322
A	-6	ASN	-	expression tag	UNP P32322
A	-5	ASN	-	expression tag	UNP P32322
A	-4	LEU	-	expression tag	UNP P32322

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	TYR	-	expression tag	UNP P32322
A	-2	PHE	-	expression tag	UNP P32322
A	-1	GLN	-	expression tag	UNP P32322
B	-22	MET	-	initiating methionine	UNP P32322
B	-21	HIS	-	expression tag	UNP P32322
B	-20	HIS	-	expression tag	UNP P32322
B	-19	HIS	-	expression tag	UNP P32322
B	-18	HIS	-	expression tag	UNP P32322
B	-17	HIS	-	expression tag	UNP P32322
B	-16	HIS	-	expression tag	UNP P32322
B	-15	SER	-	expression tag	UNP P32322
B	-14	SER	-	expression tag	UNP P32322
B	-13	GLY	-	expression tag	UNP P32322
B	-12	VAL	-	expression tag	UNP P32322
B	-11	ASP	-	expression tag	UNP P32322
B	-10	LEU	-	expression tag	UNP P32322
B	-9	GLY	-	expression tag	UNP P32322
B	-8	THR	-	expression tag	UNP P32322
B	-7	GLU	-	expression tag	UNP P32322
B	-6	ASN	-	expression tag	UNP P32322
B	-5	ASN	-	expression tag	UNP P32322
B	-4	LEU	-	expression tag	UNP P32322
B	-3	TYR	-	expression tag	UNP P32322
B	-2	PHE	-	expression tag	UNP P32322
B	-1	GLN	-	expression tag	UNP P32322
C	-22	MET	-	initiating methionine	UNP P32322
C	-21	HIS	-	expression tag	UNP P32322
C	-20	HIS	-	expression tag	UNP P32322
C	-19	HIS	-	expression tag	UNP P32322
C	-18	HIS	-	expression tag	UNP P32322
C	-17	HIS	-	expression tag	UNP P32322
C	-16	HIS	-	expression tag	UNP P32322
C	-15	SER	-	expression tag	UNP P32322
C	-14	SER	-	expression tag	UNP P32322
C	-13	GLY	-	expression tag	UNP P32322
C	-12	VAL	-	expression tag	UNP P32322
C	-11	ASP	-	expression tag	UNP P32322
C	-10	LEU	-	expression tag	UNP P32322
C	-9	GLY	-	expression tag	UNP P32322
C	-8	THR	-	expression tag	UNP P32322
C	-7	GLU	-	expression tag	UNP P32322
C	-6	ASN	-	expression tag	UNP P32322

Continued on next page...

Continued from previous page...

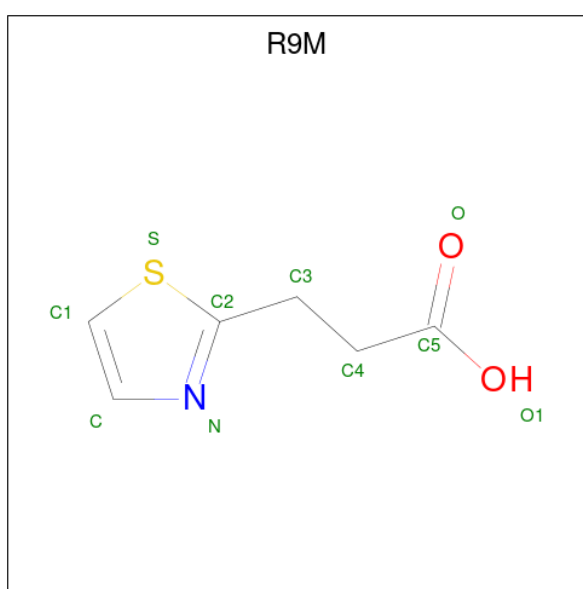
Chain	Residue	Modelled	Actual	Comment	Reference
C	-5	ASN	-	expression tag	UNP P32322
C	-4	LEU	-	expression tag	UNP P32322
C	-3	TYR	-	expression tag	UNP P32322
C	-2	PHE	-	expression tag	UNP P32322
C	-1	GLN	-	expression tag	UNP P32322
D	-22	MET	-	initiating methionine	UNP P32322
D	-21	HIS	-	expression tag	UNP P32322
D	-20	HIS	-	expression tag	UNP P32322
D	-19	HIS	-	expression tag	UNP P32322
D	-18	HIS	-	expression tag	UNP P32322
D	-17	HIS	-	expression tag	UNP P32322
D	-16	HIS	-	expression tag	UNP P32322
D	-15	SER	-	expression tag	UNP P32322
D	-14	SER	-	expression tag	UNP P32322
D	-13	GLY	-	expression tag	UNP P32322
D	-12	VAL	-	expression tag	UNP P32322
D	-11	ASP	-	expression tag	UNP P32322
D	-10	LEU	-	expression tag	UNP P32322
D	-9	GLY	-	expression tag	UNP P32322
D	-8	THR	-	expression tag	UNP P32322
D	-7	GLU	-	expression tag	UNP P32322
D	-6	ASN	-	expression tag	UNP P32322
D	-5	ASN	-	expression tag	UNP P32322
D	-4	LEU	-	expression tag	UNP P32322
D	-3	TYR	-	expression tag	UNP P32322
D	-2	PHE	-	expression tag	UNP P32322
D	-1	GLN	-	expression tag	UNP P32322
E	-22	MET	-	initiating methionine	UNP P32322
E	-21	HIS	-	expression tag	UNP P32322
E	-20	HIS	-	expression tag	UNP P32322
E	-19	HIS	-	expression tag	UNP P32322
E	-18	HIS	-	expression tag	UNP P32322
E	-17	HIS	-	expression tag	UNP P32322
E	-16	HIS	-	expression tag	UNP P32322
E	-15	SER	-	expression tag	UNP P32322
E	-14	SER	-	expression tag	UNP P32322
E	-13	GLY	-	expression tag	UNP P32322
E	-12	VAL	-	expression tag	UNP P32322
E	-11	ASP	-	expression tag	UNP P32322
E	-10	LEU	-	expression tag	UNP P32322
E	-9	GLY	-	expression tag	UNP P32322
E	-8	THR	-	expression tag	UNP P32322

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
E	-7	GLU	-	expression tag	UNP P32322
E	-6	ASN	-	expression tag	UNP P32322
E	-5	ASN	-	expression tag	UNP P32322
E	-4	LEU	-	expression tag	UNP P32322
E	-3	TYR	-	expression tag	UNP P32322
E	-2	PHE	-	expression tag	UNP P32322
E	-1	GLN	-	expression tag	UNP P32322

- Molecule 2 is 3-(1,3-thiazol-2-yl)propanoic acid (CCD ID: R9M) (formula: $C_6H_7NO_2S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			10	6	1	2	1		
2	B	1	Total	C	N	O	S	0	0
			10	6	1	2	1		
2	C	1	Total	C	N	O	S	0	0
			10	6	1	2	1		
2	D	1	Total	C	N	O	S	0	0
			10	6	1	2	1		
2	E	1	Total	C	N	O	S	0	0
			10	6	1	2	1		

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

Continued on next page...

Continued from previous page...

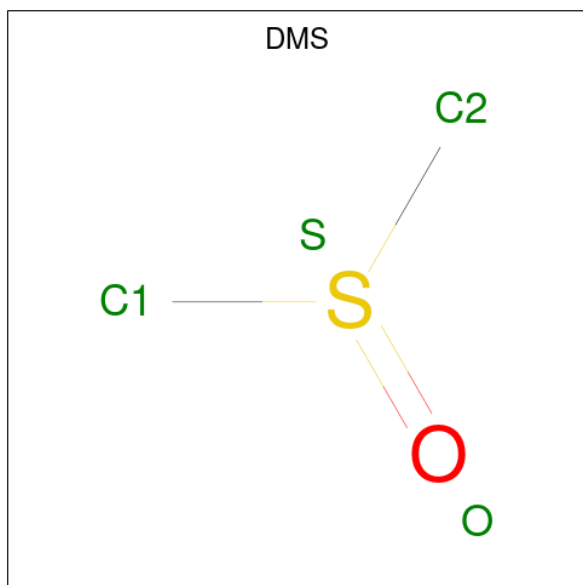
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	C	1	Total 4	C 2	O 2	0	0
3	C	1	Total 4	C 2	O 2	0	0
3	C	1	Total 4	C 2	O 2	0	0
3	C	1	Total 4	C 2	O 2	0	0
3	C	1	Total 4	C 2	O 2	0	0
3	C	1	Total 4	C 2	O 2	0	0
3	C	1	Total 4	C 2	O 2	0	0
3	C	1	Total 4	C 2	O 2	0	0
3	C	1	Total 4	C 2	O 2	0	0
3	C	1	Total 4	C 2	O 2	0	0
3	C	1	Total 4	C 2	O 2	0	0
3	D	1	Total 4	C 2	O 2	0	0
3	D	1	Total 4	C 2	O 2	0	0
3	D	1	Total 4	C 2	O 2	0	0
3	D	1	Total 4	C 2	O 2	0	0
3	D	1	Total 4	C 2	O 2	0	0
3	D	1	Total 4	C 2	O 2	0	0
3	E	1	Total 4	C 2	O 2	0	0
3	E	1	Total 4	C 2	O 2	0	0

Continued on next page...

Continued from previous page...

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

- Molecule 4 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C_2H_6OS).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	137	Total	O	0	0
			137	137		

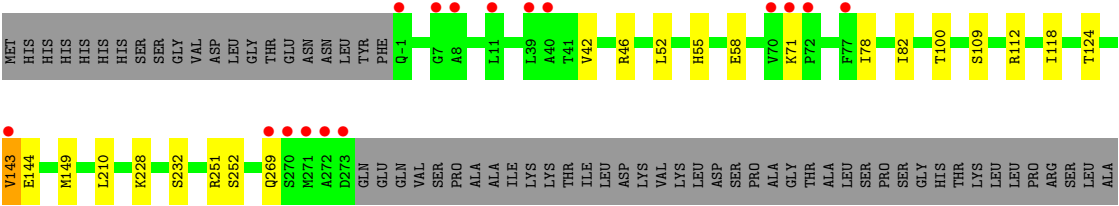
Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	147	Total 147	O 147	0	0
5	C	156	Total 156	O 156	0	0
5	D	160	Total 160	O 160	0	0
5	E	124	Total 124	O 124	0	0

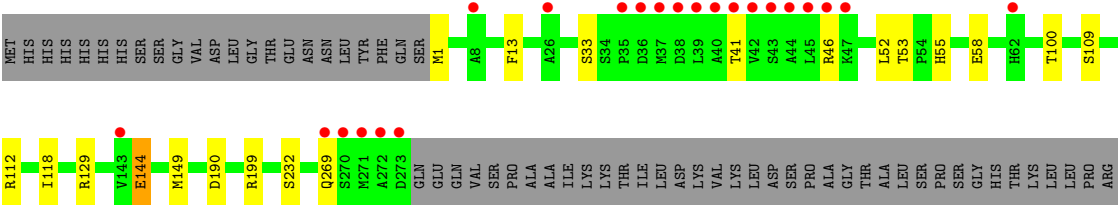
GLY
LYS
ASP

● Molecule 1: Isoform 3 of Pyrroline-5-carboxylate reductase 1, mitochondrial



PRO
ALA
GLY
LYS
ASP

● Molecule 1: Isoform 3 of Pyrroline-5-carboxylate reductase 1, mitochondrial



SER
LEU
ALA
PRO
ALA
GLY
LYS
ASP

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	163.67Å 88.04Å 116.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.11 – 1.65 37.11 – 1.65	Depositor EDS
% Data completeness (in resolution range)	99.9 (37.11-1.65) 99.9 (37.11-1.65)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 1.65Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.174 , 0.193 0.174 , 0.194	Depositor DCC
R_{free} test set	2000 reflections (0.99%)	wwPDB-VP
Wilson B-factor (Å ²)	28.1	Xtriage
Anisotropy	0.473	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	11140	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.84% 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: R9M, EDO, DMS

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	0.44	0/2070	0.58	0/2803
1	B	0.44	0/2126	0.59	0/2879
1	C	0.44	0/2063	0.59	0/2792
1	D	0.46	0/2085	0.59	0/2823
1	E	0.41	0/2070	0.56	0/2803
All	All	0.44	0/10414	0.58	0/14100

There are no bond length outliers.

There are no bond angle outliers.

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	2027	0	2096	14	0
1	B	2075	0	2131	16	0
1	C	2023	0	2089	19	0
1	D	2042	0	2109	18	0
1	E	2027	0	2096	13	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
2	C	10	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	10	0	0	0	0
2	E	10	0	0	0	0
3	A	32	0	48	4	0
3	B	36	0	54	3	0
3	C	40	0	60	3	0
3	D	24	0	36	5	0
3	E	20	0	30	1	0
4	A	4	0	6	3	0
4	B	4	0	6	2	0
4	C	8	0	12	4	0
4	D	4	0	6	1	0
5	A	137	0	0	1	0
5	B	147	0	0	0	0
5	C	156	0	0	2	0
5	D	160	0	0	2	0
5	E	124	0	0	2	0
All	All	11140	0	10779	79	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 4.

All (79) 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:E:13:PHE:HB2	1:E:41:THR:HG23	1.49	0.94
1:C:269:GLN:HE21	1:C:271:MET:H	1.34	0.74
1:A:269:GLN:NE2	1:A:273:ASP:OD1	2.25	0.70
1:B:42:VAL:HG13	1:B:52:LEU:HD13	1.75	0.68
4:B:406:DMS:H23	3:B:411:EDO:H22	1.76	0.67
1:E:190:ASP:OD1	1:E:199[A]:ARG:NH1	2.23	0.66
1:C:210[A]:LEU:HG	1:D:210[A]:LEU:HG	1.77	0.66
1:A:210[A]:LEU:HG	1:B:210[A]:LEU:HG	1.79	0.64
1:C:41:THR:HG21	5:C:521:HOH:O	2.02	0.59
1:C:233:SER:HB2	1:D:71:LYS:HE3	1.85	0.58
3:A:403:EDO:O2	4:A:405:DMS:H23	2.03	0.57
1:A:233:SER:HA	4:A:405:DMS:H13	1.87	0.57
1:C:28:HIS:O	1:C:51:LYS:NZ	2.36	0.57
1:C:138:GLY:HA2	3:C:411:EDO:H12	1.87	0.56
1:A:42:VAL:HG13	1:A:52:LEU:HD13	1.87	0.56
3:C:409:EDO:O1	4:D:406:DMS:O	2.23	0.55
1:C:269:GLN:NE2	1:C:271:MET:H	2.03	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:233:SER:HA	4:C:404:DMS:H13	1.89	0.54
1:A:266:ARG:HG2	3:A:404:EDO:H11	1.90	0.53
1:B:55:HIS:HB2	1:B:58:GLU:HG3	1.90	0.53
1:D:55:HIS:HB2	1:D:58:GLU:HG3	1.90	0.52
1:C:37:MET:HE1	1:C:52:LEU:HD22	1.92	0.52
1:E:13:PHE:HB2	1:E:41:THR:CG2	2.33	0.52
1:B:129:ARG:HG3	3:B:405:EDO:H21	1.92	0.51
1:B:144:GLU:HG2	1:B:144:GLU:O	2.11	0.51
3:C:403:EDO:O1	4:C:404:DMS:H23	2.11	0.51
4:C:404:DMS:H21	1:D:71:LYS:HZ2	1.76	0.51
1:C:233:SER:CB	1:D:71:LYS:HE3	2.41	0.51
1:B:46:ARG:CG	1:B:52:LEU:HD12	2.40	0.50
1:D:78:ILE:O	1:D:82:ILE:HG13	2.11	0.50
1:A:98:GLY:O	1:A:269:GLN:HG2	2.11	0.50
1:D:46:ARG:HG3	5:D:606:HOH:O	2.12	0.50
1:E:55:HIS:HB2	1:E:58:GLU:HG3	1.92	0.50
1:C:200:ARG:HG2	4:C:405:DMS:H21	1.93	0.49
1:D:251:ARG:HH22	3:D:404:EDO:C1	2.25	0.49
1:C:204:ARG:HH12	1:C:208:GLN:HB2	1.78	0.49
1:D:42:VAL:HG13	1:D:52:LEU:HD13	1.93	0.49
1:A:109:SER:HA	1:A:112:ARG:O	2.13	0.49
1:C:210[A]:LEU:HD11	1:D:210[A]:LEU:HD11	1.95	0.48
1:D:252:SER:OG	3:D:405:EDO:H21	2.14	0.48
3:A:404:EDO:H21	5:A:569:HOH:O	2.13	0.48
1:E:109:SER:HA	1:E:112:ARG:O	2.14	0.48
1:E:144:GLU:O	1:E:144:GLU:HG2	2.12	0.47
1:D:143:VAL:HG13	1:D:144:GLU:OE1	2.15	0.47
1:D:228:LYS:NZ	3:D:404:EDO:O2	2.29	0.46
1:E:1:MET:N	5:E:504:HOH:O	2.47	0.46
1:A:71:LYS:HE3	1:B:233:SER:HB2	1.95	0.46
1:C:109:SER:HA	1:C:112:ARG:O	2.16	0.46
1:A:103:SER:HB2	1:A:271:MET:HB2	1.98	0.46
1:A:55:HIS:HB2	1:A:58:GLU:HG3	1.97	0.46
1:A:210[B]:LEU:HD22	1:B:181:ALA:HB2	1.97	0.46
1:B:186:ASP:OD2	3:D:404:EDO:H22	2.17	0.45
1:B:112:ARG:HG2	1:B:113:PRO:HD2	1.98	0.45
1:E:129:ARG:HD3	3:E:406:EDO:H21	1.99	0.44
3:B:404:EDO:O1	4:B:406:DMS:H11	2.17	0.44
1:C:112:ARG:HG2	1:C:113:PRO:HD2	1.98	0.44
1:E:1:MET:N	5:E:506:HOH:O	2.50	0.44
1:E:46:ARG:HG3	1:E:52:LEU:HD12	1.97	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:109:SER:HA	1:D:112:ARG:O	2.17	0.44
1:D:118:ILE:HD12	1:D:149[B]:MET:HG3	2.00	0.44
1:E:100:THR:H	1:E:269:GLN:HE22	1.64	0.44
1:E:118:ILE:HD12	1:E:149[B]:MET:HG3	1.99	0.43
1:B:46:ARG:HG2	1:B:52:LEU:HD12	2.01	0.43
1:A:190:ASP:OD1	1:A:199[A]:ARG:NH1	2.46	0.43
1:E:33:SER:HA	1:E:53:THR:O	2.19	0.42
1:C:222:GLN:OE1	5:C:501:HOH:O	2.22	0.42
1:B:133:THR:O	1:B:159:CYS:HA	2.20	0.42
1:C:204:ARG:NH1	1:C:208:GLN:HB2	2.35	0.42
1:A:234:PRO:HD3	4:A:405:DMS:H13	2.01	0.42
1:D:124[A]:THR:HG21	5:D:584:HOH:O	2.19	0.42
1:B:46:ARG:HG3	1:B:52:LEU:HD12	2.00	0.42
1:A:46:ARG:CG	1:A:52:LEU:HD12	2.49	0.42
1:D:228:LYS:HZ3	3:D:404:EDO:C2	2.27	0.41
1:D:100:THR:H	1:D:269:GLN:NE2	2.18	0.41
1:C:33:SER:HA	1:C:53:THR:O	2.20	0.41
1:B:33:SER:HA	1:B:53:THR:O	2.20	0.41
3:A:409:EDO:H11	1:B:264:ARG:NH2	2.36	0.41
1:C:39:LEU:C	1:C:41:THR:H	2.29	0.41
1:B:4:GLY:HA3	1:B:63:SER:OG	2.21	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	275/342 (80%)	273 (99%)	2 (1%)	0	100	100
1	B	281/342 (82%)	275 (98%)	6 (2%)	0	100	100
1	C	274/342 (80%)	269 (98%)	5 (2%)	0	100	100
1	D	277/342 (81%)	274 (99%)	3 (1%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	275/342 (80%)	272 (99%)	3 (1%)	0	100	100
All	All	1382/1710 (81%)	1363 (99%)	19 (1%)	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	216/270 (80%)	213 (99%)	3 (1%)	62	43
1	B	222/270 (82%)	219 (99%)	3 (1%)	62	43
1	C	215/270 (80%)	212 (99%)	3 (1%)	62	43
1	D	218/270 (81%)	216 (99%)	2 (1%)	75	63
1	E	216/270 (80%)	214 (99%)	2 (1%)	75	63
All	All	1087/1350 (80%)	1074 (99%)	13 (1%)	67	50

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

Mol	Chain	Res	Type
1	A	1	MET
1	A	37	MET
1	A	269	GLN
1	B	1	MET
1	B	143	VAL
1	B	144	GLU
1	C	1	MET
1	C	42	VAL
1	C	144	GLU
1	D	143	VAL
1	D	232	SER
1	E	144	GLU
1	E	232	SER

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	55	HIS
1	A	219	HIS
1	A	240	HIS
1	B	219	HIS
1	C	56	ASN
1	C	240	HIS
1	E	55	HIS
1	E	226	GLN

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](#)

48 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	EDO	C	409	-	3,3,3	0.49	0	2,2,2	0.21	0
3	EDO	E	403	-	3,3,3	0.50	0	2,2,2	0.20	0
3	EDO	C	412	-	3,3,3	0.50	0	2,2,2	0.35	0
3	EDO	A	403	-	3,3,3	0.41	0	2,2,2	0.42	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EDO	D	402	-	3,3,3	0.46	0	2,2,2	0.40	0
3	EDO	B	411	-	3,3,3	0.45	0	2,2,2	0.26	0
3	EDO	D	408	-	3,3,3	0.51	0	2,2,2	0.27	0
3	EDO	C	402	-	3,3,3	0.50	0	2,2,2	0.17	0
3	EDO	C	413	-	3,3,3	0.40	0	2,2,2	0.81	0
3	EDO	C	406	-	3,3,3	0.27	0	2,2,2	0.74	0
3	EDO	E	402	-	3,3,3	0.47	0	2,2,2	0.31	0
3	EDO	D	405	-	3,3,3	0.53	0	2,2,2	0.20	0
4	DMS	C	404	-	3,3,3	0.62	0	3,3,3	0.70	0
3	EDO	A	410	-	3,3,3	0.44	0	2,2,2	0.27	0
3	EDO	E	404	-	3,3,3	0.43	0	2,2,2	0.42	0
3	EDO	C	403	-	3,3,3	0.37	0	2,2,2	0.51	0
3	EDO	D	403	-	3,3,3	0.45	0	2,2,2	0.50	0
4	DMS	C	405	-	3,3,3	0.70	0	3,3,3	0.44	0
3	EDO	B	408	-	3,3,3	0.44	0	2,2,2	0.45	0
3	EDO	A	407	-	3,3,3	0.47	0	2,2,2	0.43	0
3	EDO	E	405	-	3,3,3	0.50	0	2,2,2	0.30	0
3	EDO	A	402	-	3,3,3	0.48	0	2,2,2	0.41	0
2	R9M	A	401	-	9,10,10	1.92	1 (11%)	6,12,12	1.91	1 (16%)
2	R9M	E	401	-	9,10,10	1.84	2 (22%)	6,12,12	1.77	2 (33%)
3	EDO	C	410	-	3,3,3	0.50	0	2,2,2	0.36	0
2	R9M	B	402	-	9,10,10	2.02	1 (11%)	6,12,12	1.95	2 (33%)
3	EDO	C	411	-	3,3,3	0.49	0	2,2,2	0.34	0
3	EDO	C	407	-	3,3,3	0.44	0	2,2,2	0.37	0
3	EDO	A	406	-	3,3,3	0.27	0	2,2,2	1.14	0
3	EDO	B	409	-	3,3,3	0.52	0	2,2,2	0.35	0
4	DMS	D	406	-	3,3,3	0.63	0	3,3,3	0.42	0
3	EDO	C	408	-	3,3,3	0.43	0	2,2,2	0.58	0
3	EDO	B	403	-	3,3,3	0.48	0	2,2,2	0.34	0
3	EDO	D	407	-	3,3,3	0.47	0	2,2,2	0.16	0
3	EDO	A	404	-	3,3,3	0.57	0	2,2,2	0.27	0
3	EDO	D	404	-	3,3,3	0.43	0	2,2,2	0.33	0
2	R9M	D	401	-	9,10,10	1.90	2 (22%)	6,12,12	1.89	1 (16%)
3	EDO	B	404	-	3,3,3	0.44	0	2,2,2	0.21	0
3	EDO	A	409	-	3,3,3	0.45	0	2,2,2	0.38	0
2	R9M	C	401	-	9,10,10	1.92	1 (11%)	6,12,12	1.89	2 (33%)
3	EDO	B	401	-	3,3,3	0.46	0	2,2,2	0.41	0
4	DMS	B	406	-	3,3,3	0.62	0	3,3,3	0.59	0
4	DMS	A	405	-	3,3,3	0.54	0	3,3,3	0.46	0
3	EDO	B	407	-	3,3,3	0.49	0	2,2,2	0.34	0
3	EDO	E	406	-	3,3,3	0.46	0	2,2,2	0.20	0
3	EDO	B	405	-	3,3,3	0.41	0	2,2,2	0.24	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	EDO	A	408	-	3,3,3	0.54	0	2,2,2	0.25	0
3	EDO	B	410	-	3,3,3	0.48	0	2,2,2	0.25	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	EDO	C	409	-	-	1/1/1/1	-
3	EDO	E	403	-	-	0/1/1/1	-
3	EDO	C	412	-	-	0/1/1/1	-
3	EDO	A	403	-	-	0/1/1/1	-
3	EDO	D	402	-	-	0/1/1/1	-
3	EDO	B	411	-	-	1/1/1/1	-
3	EDO	D	408	-	-	1/1/1/1	-
3	EDO	C	402	-	-	0/1/1/1	-
3	EDO	C	413	-	-	1/1/1/1	-
3	EDO	C	406	-	-	1/1/1/1	-
3	EDO	E	402	-	-	0/1/1/1	-
3	EDO	D	405	-	-	0/1/1/1	-
3	EDO	A	410	-	-	0/1/1/1	-
3	EDO	E	404	-	-	1/1/1/1	-
3	EDO	C	403	-	-	0/1/1/1	-
3	EDO	D	403	-	-	1/1/1/1	-
3	EDO	B	408	-	-	0/1/1/1	-
3	EDO	A	407	-	-	0/1/1/1	-
3	EDO	E	405	-	-	1/1/1/1	-
3	EDO	A	402	-	-	0/1/1/1	-
2	R9M	A	401	-	-	2/4/5/5	0/1/1/1
2	R9M	E	401	-	-	2/4/5/5	0/1/1/1
3	EDO	C	410	-	-	0/1/1/1	-
2	R9M	B	402	-	-	2/4/5/5	0/1/1/1
3	EDO	C	411	-	-	1/1/1/1	-
3	EDO	C	407	-	-	0/1/1/1	-
3	EDO	A	406	-	-	0/1/1/1	-
3	EDO	B	409	-	-	0/1/1/1	-
3	EDO	C	408	-	-	1/1/1/1	-
3	EDO	B	403	-	-	0/1/1/1	-
3	EDO	D	407	-	-	1/1/1/1	-
3	EDO	A	404	-	-	0/1/1/1	-
3	EDO	D	404	-	-	0/1/1/1	-

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	R9M	D	401	-	-	2/4/5/5	0/1/1/1
3	EDO	B	404	-	-	0/1/1/1	-
3	EDO	A	409	-	-	1/1/1/1	-
2	R9M	C	401	-	-	2/4/5/5	0/1/1/1
3	EDO	B	401	-	-	1/1/1/1	-
3	EDO	B	407	-	-	0/1/1/1	-
3	EDO	E	406	-	-	0/1/1/1	-
3	EDO	B	405	-	-	0/1/1/1	-
3	EDO	A	408	-	-	0/1/1/1	-
3	EDO	B	410	-	-	1/1/1/1	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	402	R9M	C3-C2	5.40	1.52	1.49
2	A	401	R9M	C3-C2	5.37	1.52	1.49
2	C	401	R9M	C3-C2	5.18	1.52	1.49
2	D	401	R9M	C3-C2	4.94	1.52	1.49
2	E	401	R9M	C3-C2	4.83	1.52	1.49
2	D	401	R9M	O-C5	2.21	1.29	1.22
2	E	401	R9M	O-C5	2.05	1.28	1.22

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	402	R9M	C-N-C2	3.57	115.22	103.14
2	A	401	R9M	C-N-C2	3.46	114.85	103.14
2	E	401	R9M	C-N-C2	3.45	114.82	103.14
2	C	401	R9M	C-N-C2	3.45	114.80	103.14
2	D	401	R9M	C-N-C2	3.40	114.63	103.14
2	B	402	R9M	C4-C3-C2	-2.65	108.22	112.67
2	C	401	R9M	C4-C3-C2	-2.26	108.88	112.67
2	E	401	R9M	C3-C4-C5	-2.01	108.12	113.76

There are no chirality outliers.

All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	409	EDO	O1-C1-C2-O2
3	E	404	EDO	O1-C1-C2-O2
3	B	410	EDO	O1-C1-C2-O2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
3	B	411	EDO	O1-C1-C2-O2
3	C	411	EDO	O1-C1-C2-O2
2	D	401	R9M	C3-C4-C5-O1
2	B	402	R9M	C3-C4-C5-O1
2	E	401	R9M	C3-C4-C5-O1
3	C	409	EDO	O1-C1-C2-O2
2	C	401	R9M	C3-C4-C5-O1
2	C	401	R9M	C3-C4-C5-O
2	A	401	R9M	C3-C4-C5-O1
2	A	401	R9M	C3-C4-C5-O
2	D	401	R9M	C3-C4-C5-O
2	B	402	R9M	C3-C4-C5-O
3	D	403	EDO	O1-C1-C2-O2
3	D	407	EDO	O1-C1-C2-O2
2	E	401	R9M	C3-C4-C5-O
3	B	401	EDO	O1-C1-C2-O2
3	C	413	EDO	O1-C1-C2-O2
3	E	405	EDO	O1-C1-C2-O2
3	C	406	EDO	O1-C1-C2-O2
3	C	408	EDO	O1-C1-C2-O2
3	D	408	EDO	O1-C1-C2-O2

There are no ring outliers.

17 monomers are involved in 21 short contacts:

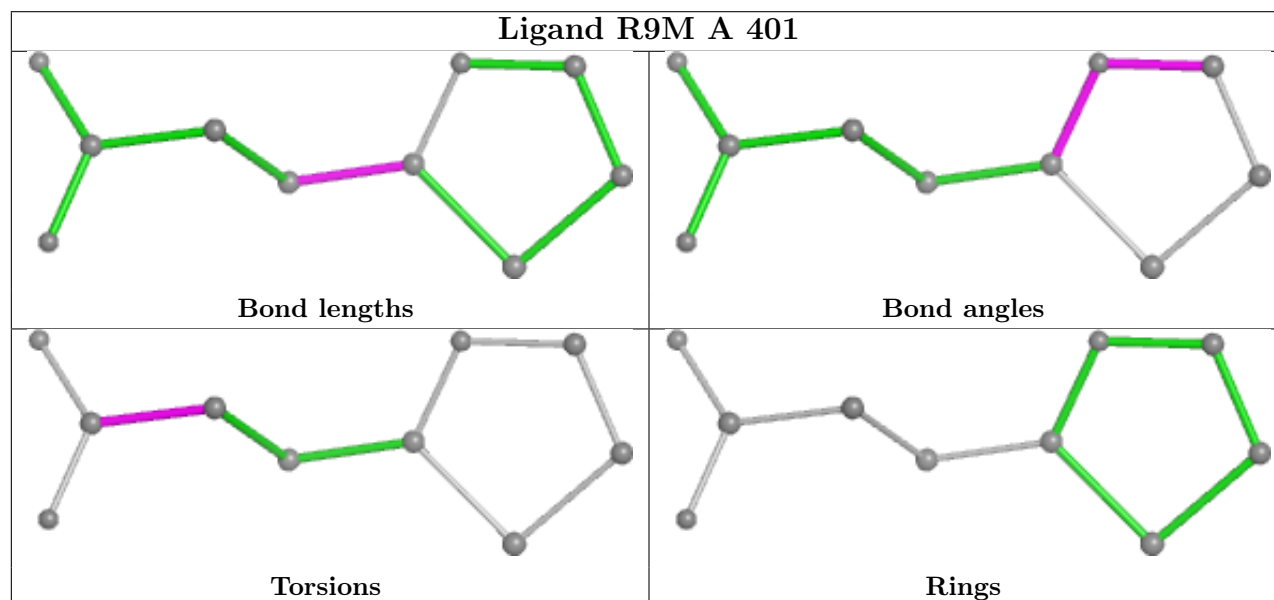
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	409	EDO	1	0
3	A	403	EDO	1	0
3	B	411	EDO	1	0
3	D	405	EDO	1	0
4	C	404	DMS	3	0
3	C	403	EDO	1	0
4	C	405	DMS	1	0
3	C	411	EDO	1	0
4	D	406	DMS	1	0
3	A	404	EDO	2	0
3	D	404	EDO	4	0
3	B	404	EDO	1	0
3	A	409	EDO	1	0
4	B	406	DMS	2	0
4	A	405	DMS	3	0
3	E	406	EDO	1	0

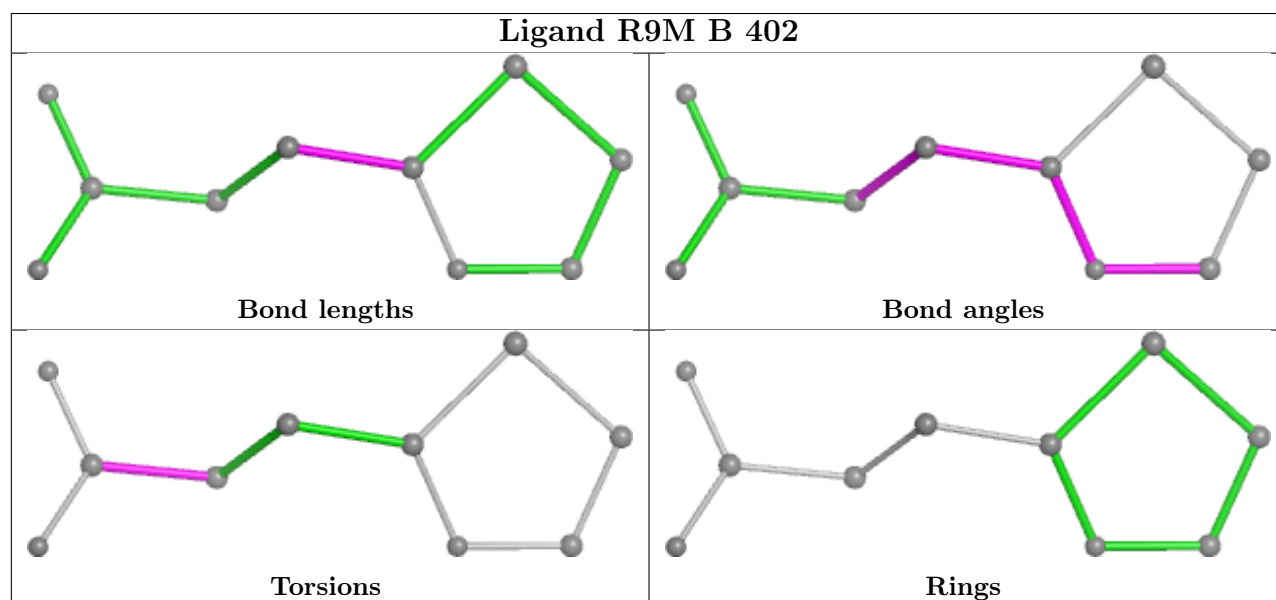
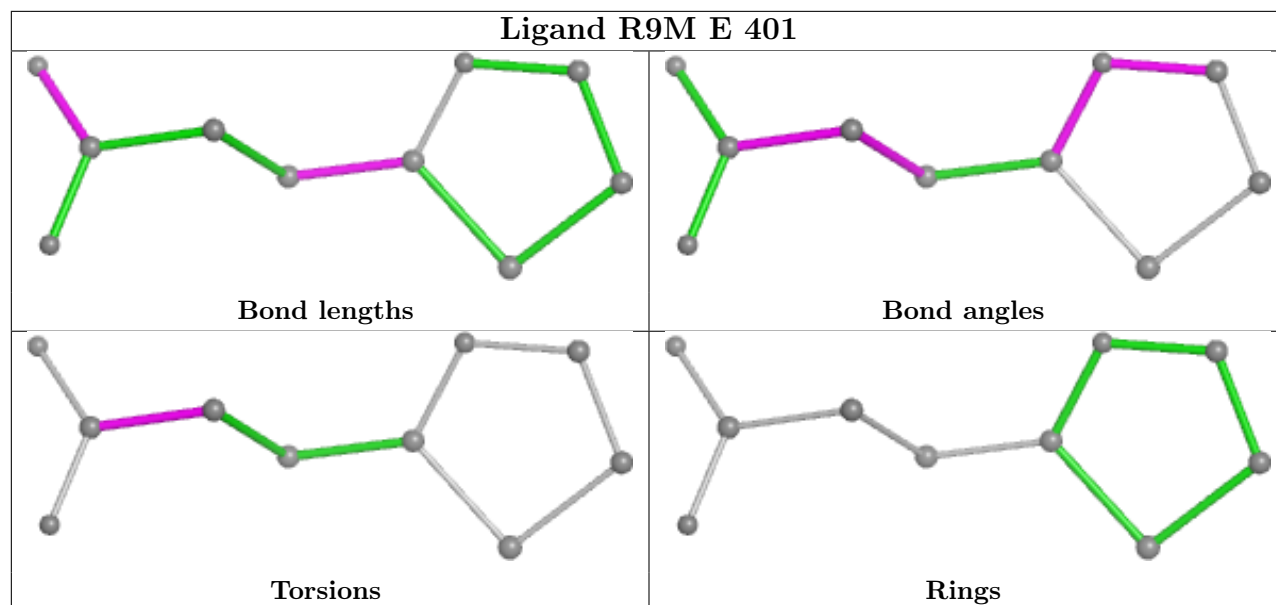
Continued on next page...

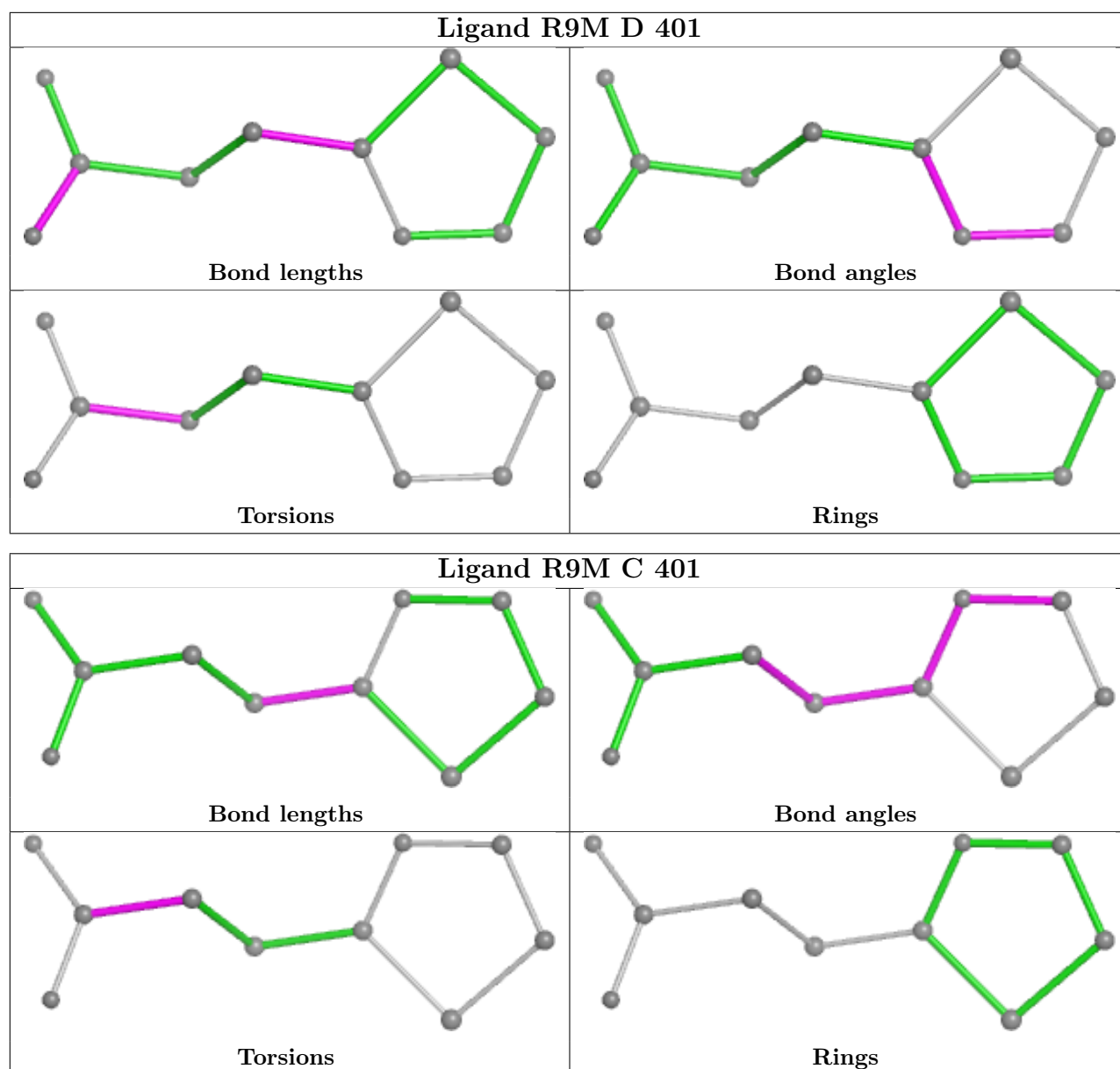
Continued from previous page...

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	405	EDO	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 in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







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	273/342 (79%)	0.30	19 (6%)	24 26	16, 38, 74, 119	4 (1%)
1	B	278/342 (81%)	0.22	18 (6%)	26 28	16, 37, 67, 111	5 (1%)
1	C	273/342 (79%)	0.26	21 (7%)	21 22	15, 36, 79, 174	3 (1%)
1	D	275/342 (80%)	0.18	16 (5%)	30 32	16, 36, 62, 92	4 (1%)
1	E	273/342 (79%)	0.38	22 (8%)	19 20	16, 39, 77, 171	4 (1%)
All	All	1372/1710 (80%)	0.27	96 (6%)	24 26	15, 37, 73, 174	20 (1%)

All (96) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	42	VAL	8.6
1	C	39	LEU	7.8
1	E	37	MET	6.9
1	C	40	ALA	6.3
1	D	143	VAL	6.3
1	B	-3	TYR	6.0
1	E	143	VAL	5.8
1	E	39	LEU	5.3
1	B	143	VAL	5.2
1	E	41	THR	5.0
1	C	41	THR	4.9
1	C	42	VAL	4.7
1	E	40	ALA	4.6
1	B	-4	LEU	4.6
1	D	273	ASP	4.3
1	E	273	ASP	4.2
1	A	273	ASP	4.1
1	A	42	VAL	4.1
1	A	272	ALA	4.0
1	C	37	MET	3.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	272	ALA	3.9
1	E	43	SER	3.9
1	B	0	SER	3.8
1	D	270	SER	3.8
1	B	271	MET	3.8
1	A	271	MET	3.6
1	E	270	SER	3.5
1	B	40	ALA	3.4
1	E	272	ALA	3.4
1	D	271	MET	3.4
1	A	40	ALA	3.4
1	A	39	LEU	3.3
1	E	38	ASP	3.3
1	B	-2	PHE	3.2
1	D	-1	GLN	3.2
1	C	38	ASP	3.2
1	C	273	ASP	3.2
1	A	52	LEU	3.1
1	B	273	ASP	3.1
1	E	26	ALA	3.1
1	E	46	ARG	3.0
1	C	46	ARG	2.9
1	E	35	PRO	2.9
1	B	270	SER	2.9
1	E	44	ALA	2.9
1	A	269	GLN	2.8
1	C	272	ALA	2.8
1	D	77	PHE	2.8
1	A	270	SER	2.8
1	A	35	PRO	2.7
1	C	44	ALA	2.7
1	D	70	VAL	2.7
1	C	144	GLU	2.7
1	A	143	VAL	2.7
1	D	269	GLN	2.7
1	C	45	LEU	2.7
1	E	45	LEU	2.7
1	A	62	HIS	2.7
1	C	43	SER	2.6
1	B	39	LEU	2.6
1	A	8	ALA	2.6
1	B	35	PRO	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	62	HIS	2.6
1	D	39	LEU	2.6
1	E	36	ASP	2.5
1	E	8	ALA	2.5
1	A	37	MET	2.5
1	B	37	MET	2.5
1	C	8	ALA	2.5
1	C	271	MET	2.5
1	E	271	MET	2.5
1	D	7	GLY	2.5
1	C	52	LEU	2.4
1	B	269	GLN	2.4
1	B	77	PHE	2.3
1	D	72	PRO	2.3
1	E	47	LYS	2.3
1	B	272	ALA	2.3
1	A	70	VAL	2.2
1	A	77	PHE	2.2
1	B	-1	GLN	2.2
1	E	269	GLN	2.2
1	D	11	LEU	2.2
1	A	110	ALA	2.2
1	B	8	ALA	2.2
1	A	44	ALA	2.1
1	A	49	GLY	2.1
1	D	71	LYS	2.1
1	C	270	SER	2.1
1	C	26	ALA	2.1
1	D	8	ALA	2.1
1	B	70	VAL	2.1
1	C	143	VAL	2.1
1	E	62	HIS	2.0
1	C	27	ALA	2.0
1	D	40	ALA	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.

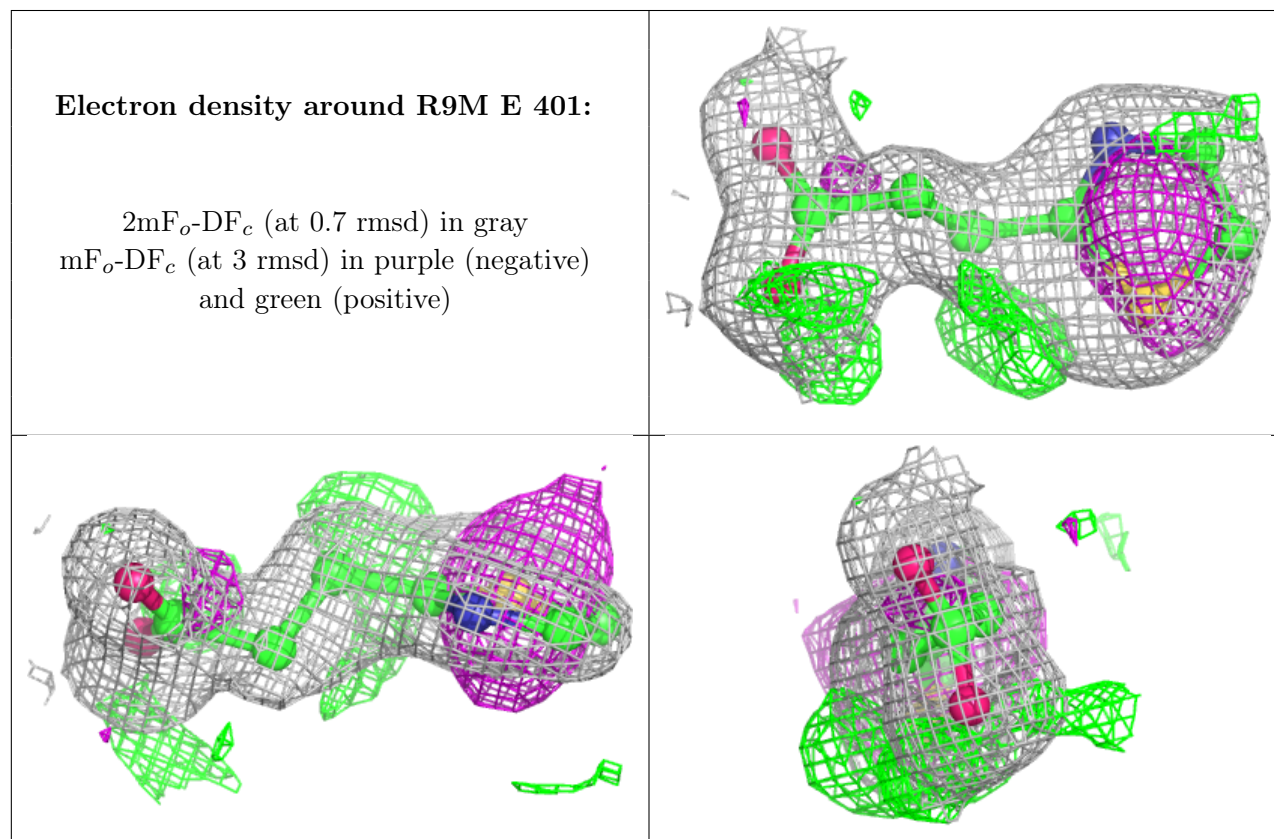
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	EDO	C	408	4/4	0.65	0.24	51,53,66,70	0
4	DMS	C	405	4/4	0.65	0.26	58,59,71,76	0
3	EDO	B	405	4/4	0.71	0.22	51,56,62,65	0
4	DMS	C	404	4/4	0.72	0.22	45,54,60,90	0
3	EDO	B	408	4/4	0.74	0.17	86,92,94,98	0
3	EDO	A	404	4/4	0.76	0.17	48,51,58,61	0
4	DMS	A	405	4/4	0.76	0.18	51,52,60,89	0
3	EDO	A	407	4/4	0.77	0.16	55,60,62,67	0
3	EDO	C	409	4/4	0.77	0.16	48,49,57,59	0
3	EDO	C	410	4/4	0.77	0.15	53,56,59,64	0
3	EDO	C	412	4/4	0.79	0.15	56,63,67,69	0
3	EDO	C	413	4/4	0.80	0.18	47,49,61,63	0
3	EDO	D	408	4/4	0.80	0.19	43,52,62,65	0
3	EDO	E	406	4/4	0.80	0.17	45,56,62,65	0
3	EDO	D	404	4/4	0.81	0.18	37,44,44,74	0
4	DMS	D	406	4/4	0.81	0.20	59,59,70,76	0
3	EDO	E	405	4/4	0.82	0.15	50,53,64,76	0
3	EDO	B	409	4/4	0.83	0.15	52,57,63,67	0
3	EDO	D	407	4/4	0.83	0.16	55,56,58,62	0
3	EDO	A	406	4/4	0.83	0.15	39,40,58,61	0
3	EDO	C	407	4/4	0.84	0.16	46,54,64,65	0
4	DMS	B	406	4/4	0.85	0.15	56,57,59,93	0
3	EDO	B	407	4/4	0.85	0.14	49,58,60,71	0
3	EDO	A	410	4/4	0.85	0.13	45,48,62,68	0
2	R9M	E	401	10/10	0.85	0.16	32,46,50,50	0
2	R9M	B	402	10/10	0.86	0.16	34,43,48,51	0
3	EDO	B	410	4/4	0.86	0.15	43,59,67,71	0
3	EDO	C	411	4/4	0.86	0.12	46,59,60,67	0
3	EDO	A	409	4/4	0.87	0.12	43,49,53,73	0
2	R9M	C	401	10/10	0.87	0.15	31,42,46,50	0
3	EDO	B	411	4/4	0.88	0.12	51,51,55,63	0
3	EDO	A	408	4/4	0.88	0.13	50,54,56,57	0

Continued on next page...

Continued from previous page...

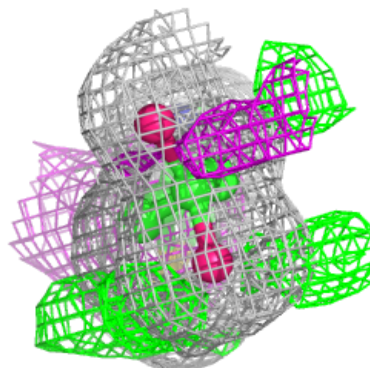
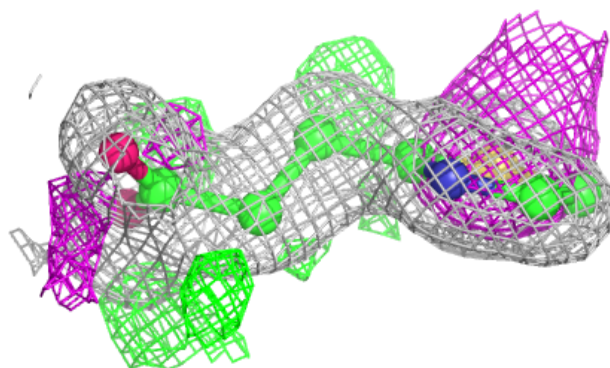
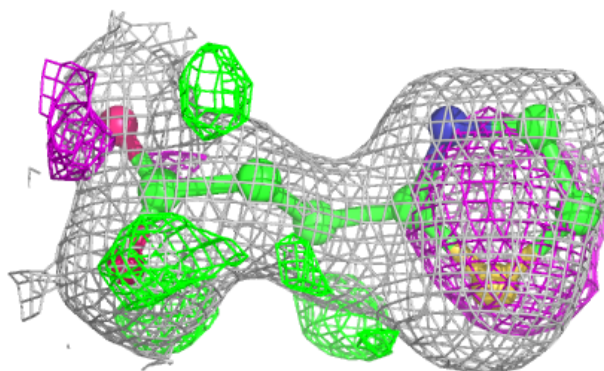
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	EDO	B	401	4/4	0.89	0.12	50,55,57,58	0
2	R9M	D	401	10/10	0.89	0.14	30,42,48,55	0
3	EDO	C	402	4/4	0.90	0.13	53,53,56,58	0
3	EDO	E	403	4/4	0.90	0.14	44,47,50,58	0
3	EDO	B	403	4/4	0.90	0.13	49,49,52,52	0
3	EDO	E	404	4/4	0.91	0.12	34,46,53,64	0
2	R9M	A	401	10/10	0.91	0.12	32,41,45,48	0
3	EDO	A	402	4/4	0.91	0.12	50,52,54,57	0
3	EDO	B	404	4/4	0.91	0.13	44,47,51,51	0
3	EDO	D	405	4/4	0.92	0.11	41,41,47,53	0
3	EDO	D	402	4/4	0.92	0.10	44,47,49,49	0
3	EDO	D	403	4/4	0.92	0.12	39,42,45,47	0
3	EDO	C	406	4/4	0.92	0.11	38,45,53,58	0
3	EDO	C	403	4/4	0.94	0.10	39,43,43,45	0
3	EDO	A	403	4/4	0.94	0.11	39,41,43,48	0
3	EDO	E	402	4/4	0.96	0.09	53,54,56,57	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.

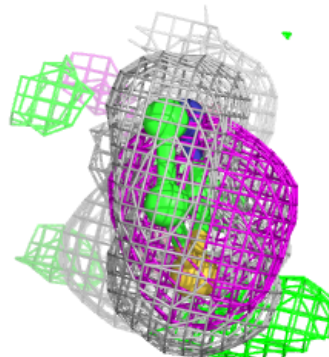
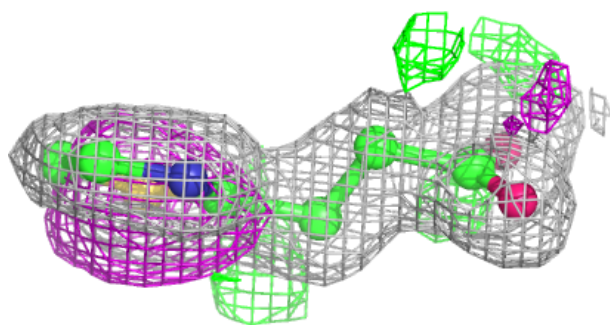
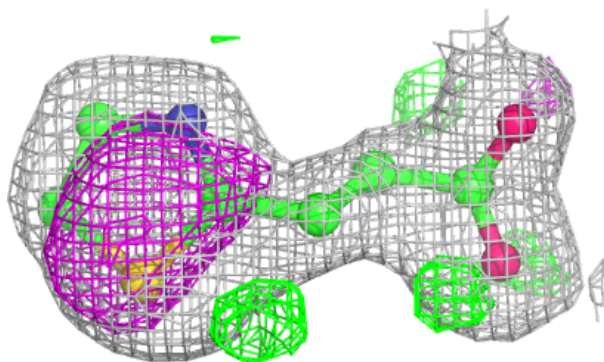


Electron density around R9M B 402:

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

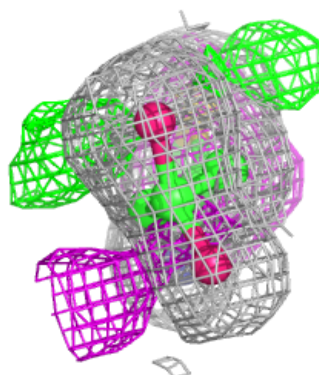
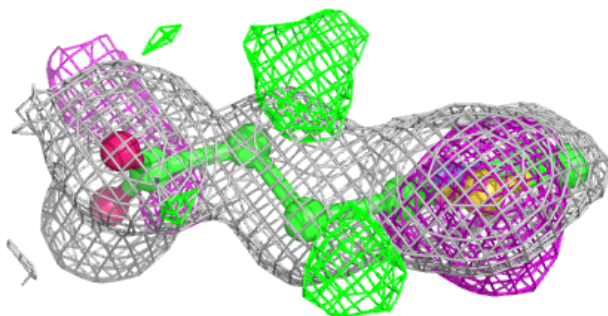
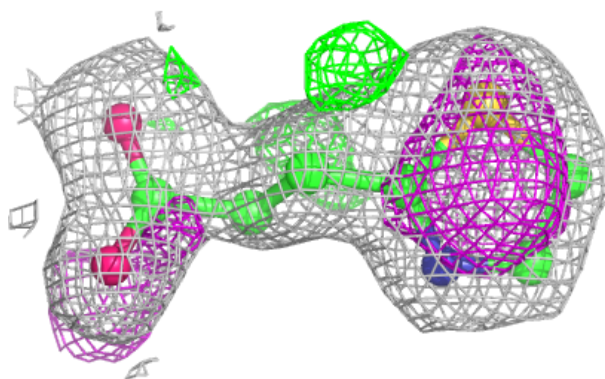
**Electron density around R9M C 401:**

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

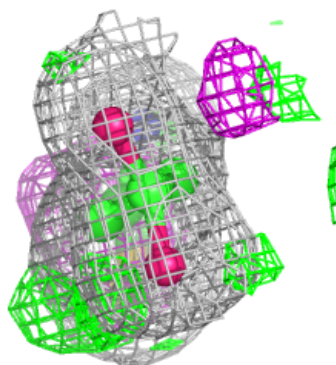
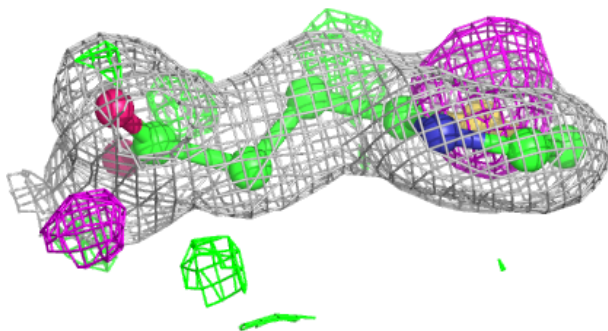
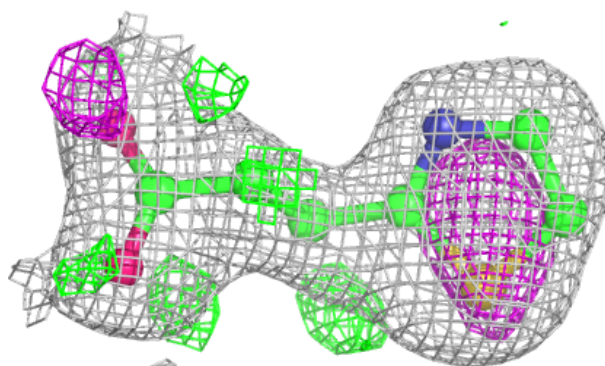


Electron density around R9M D 401:

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

**Electron density around R9M A 401:**

$2mF_o-DF_c$ (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.