



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 23, 2021 – 08:44 AM EDT

PDB ID : 1FUR  
Title : FUMARASE MUTANT H188N WITH BOUND SUBSTRATE L-MALATE  
AT PUTATIVE ACTIVATOR SITE  
Authors : Weaver, T.M.; Lees, M.; Banaszak, L.J.  
Deposited on : 1997-01-09  
Resolution : 1.95 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.23.2  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.23.2

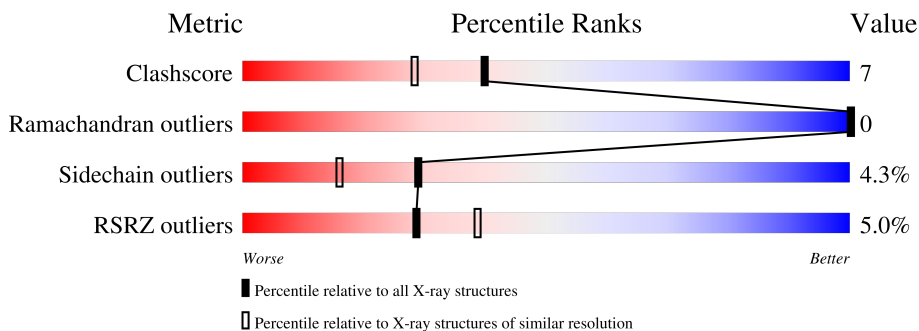
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.95 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))
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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  $\geq 3$ , 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	467	
1	B	467	

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
2	MLT	A	468	X	-	X	-
2	MLT	B	468	X	-	-	-

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 9695 atoms, of which 2368 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.

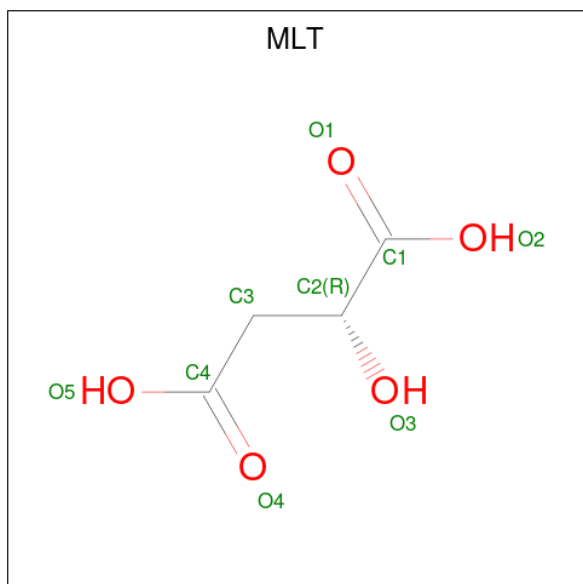
- Molecule 1 is a protein called FUMARASE C.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	456	4235	2156	781	616	659	23	0	0	0
1	B	455	4236	2155	783	616	658	24	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	188	ASN	HIS	engineered mutation	UNP P05042
B	188	ASN	HIS	engineered mutation	UNP P05042

- Molecule 2 is D-MALATE (three-letter code: MLT) (formula: C<sub>4</sub>H<sub>6</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	9	4	5	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	B	1	9	4	5	0	0

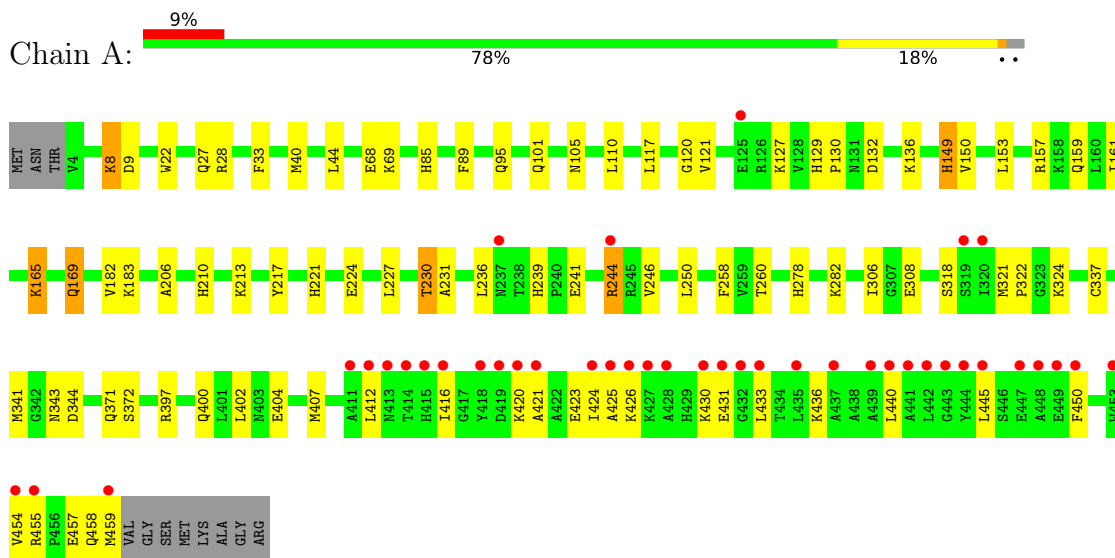
- Molecule 3 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	H	O		
3	A	215	645	430	215	0	0
3	B	187	561	374	187	0	0

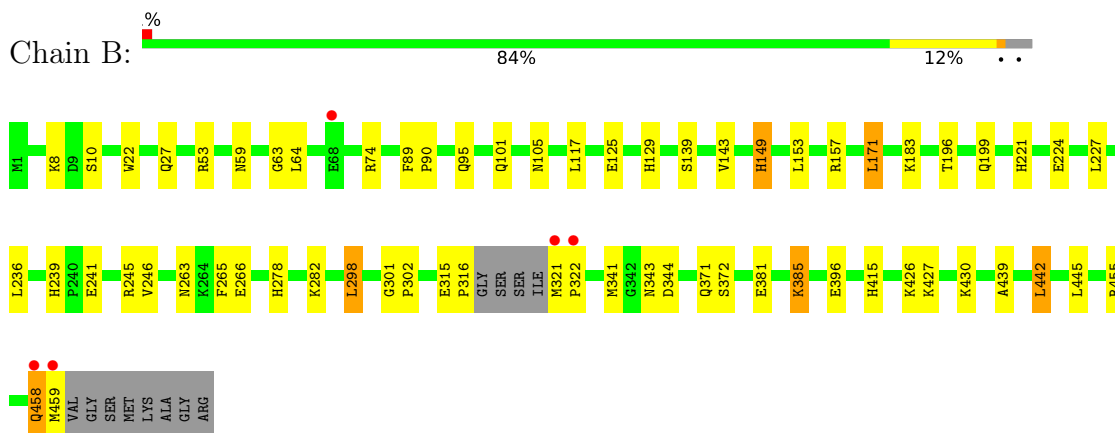
### 3 Residue-property plots

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: FUMARASE C



- Molecule 1: FUMARASE C



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	104.03Å 219.32Å 86.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 1.95 14.50 – 1.82	Depositor EDS
% Data completeness (in resolution range)	87.6 (8.00-1.95) 80.3 (14.50-1.82)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.25 (at 1.82Å)	Xtrriage
Refinement program	X-PLOR 3.1	Depositor
R, $R_{free}$	0.170 , 0.209 0.180 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	15.3	Xtrriage
Anisotropy	0.328	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 58.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	9695	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	16.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.51% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: MLT

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	0.33	0/3511	0.58	0/4756
1	B	0.34	0/3509	0.58	0/4752
All	All	0.33	0/7020	0.58	0/9508

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

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	3454	781	3472	61	1
1	B	3453	783	3472	43	1
2	A	9	0	3	4	0
2	B	9	0	3	2	0
3	A	215	430	0	7	1
3	B	187	374	0	5	1
All	All	7327	2368	6950	101	2

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 (101) 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:B:89:PHE:HA	1:B:105:ASN:HD21	1.52	0.74
1:A:343:ASN:HD21	1:A:371:GLN:HE21	1.34	0.73
1:B:263:ASN:HD22	1:B:266:GLU:H	1.38	0.71
1:B:439:ALA:HB1	1:B:445:LEU:HD21	1.70	0.71
1:A:407:MET:SD	1:B:321:MET:HG2	2.30	0.71
1:A:431:GLU:HB3	1:A:433:LEU:HD13	1.73	0.69
1:A:230:THR:HG23	1:A:231:ALA:H	1.58	0.68
1:A:230:THR:CG2	1:A:231:ALA:H	2.07	0.67
1:A:423:GLU:HA	1:A:426:LYS:HD2	1.76	0.67
1:A:85:HIS:HD2	3:A:535:HOH:O	1.80	0.65
1:A:455:ARG:HH11	1:A:458:GLN:HE22	1.44	0.64
1:B:426:LYS:O	1:B:430:LYS:HD3	1.97	0.64
1:B:343:ASN:HD21	1:B:371:GLN:HE21	1.45	0.63
1:B:125:GLU:HG2	3:B:575:HOH:O	1.97	0.62
1:B:415:HIS:HD2	3:B:557:HOH:O	1.82	0.62
1:B:415:HIS:HE1	3:B:556:HOH:O	1.82	0.61
1:B:196:THR:H	1:B:199:GLN:HE21	1.50	0.60
1:A:445:LEU:HD13	1:A:450:PHE:HB2	1.83	0.60
1:A:230:THR:HG21	1:A:236:LEU:O	2.02	0.60
1:A:244:ARG:NH1	1:A:260:THR:H	2.00	0.59
1:A:95:GLN:HB2	1:A:101:GLN:OE1	2.04	0.58
1:A:230:THR:HG22	3:A:486:HOH:O	2.02	0.57
1:B:427:LYS:HE2	1:B:442:LEU:HD13	1.85	0.57
1:A:244:ARG:HH12	1:A:260:THR:H	1.53	0.57
1:A:230:THR:CG2	1:A:231:ALA:N	2.68	0.56
1:B:139:SER:O	1:B:143:VAL:HG23	2.05	0.56
1:B:22:TRP:HE1	1:B:27:GLN:NE2	2.03	0.56
1:A:89:PHE:HA	1:A:105:ASN:HD21	1.71	0.56
1:B:439:ALA:HB1	1:B:445:LEU:CD2	2.36	0.56
1:A:149:HIS:HD2	1:A:224:GLU:O	1.90	0.54
1:A:343:ASN:HD21	1:A:371:GLN:NE2	2.05	0.54
1:B:196:THR:H	1:B:199:GLN:NE2	2.06	0.54
1:B:263:ASN:HD21	1:B:265:PHE:HB2	1.72	0.54
1:B:95:GLN:HB2	1:B:101:GLN:NE2	2.23	0.54
1:B:227:LEU:HD23	1:B:246:VAL:HG21	1.90	0.54
1:B:278:HIS:HE1	1:B:372:SER:OG	1.91	0.54
1:A:9:ASP:HB2	3:A:604:HOH:O	2.07	0.53
1:A:129:HIS:HD2	2:A:468:MLT:O4	1.92	0.53
1:B:59:ASN:HD22	1:B:64:LEU:HD12	1.72	0.53
1:A:455:ARG:NH1	1:A:458:GLN:HE22	2.08	0.52
1:A:22:TRP:HE1	1:A:27:GLN:NE2	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:412:LEU:O	1:A:416:ILE:HG22	2.09	0.52
1:A:33:PHE:HE2	1:A:101:GLN:HE22	1.52	0.51
1:B:455:ARG:O	1:B:458:GLN:HG3	2.11	0.50
1:B:22:TRP:HE1	1:B:27:GLN:HE21	1.58	0.50
1:A:182:VAL:HG21	1:A:457:GLU:HA	1.92	0.50
1:A:400:GLN:O	1:A:404:GLU:HB2	2.11	0.50
1:A:22:TRP:HE1	1:A:27:GLN:HE21	1.60	0.49
1:A:40:MET:SD	1:A:44:LEU:HD23	2.51	0.49
1:A:306:ILE:HG22	1:A:308:GLU:HG3	1.93	0.49
1:A:318:SER:HB2	1:A:321:MET:HB2	1.93	0.49
1:A:121:VAL:HG12	3:A:582:HOH:O	2.12	0.49
1:B:149:HIS:HD2	1:B:224:GLU:O	1.94	0.49
1:B:278:HIS:HD2	1:B:344:ASP:OD1	1.96	0.49
1:B:241:GLU:O	1:B:245:ARG:HG3	2.13	0.49
1:A:421:ALA:O	1:A:424:ILE:HG12	2.13	0.48
1:A:169:GLN:NE2	3:A:570:HOH:O	2.46	0.48
1:A:129:HIS:CD2	2:A:468:MLT:H32	2.49	0.48
1:B:239:HIS:HD2	1:B:241:GLU:H	1.61	0.48
1:A:161:ILE:HG22	1:A:165:LYS:HE2	1.96	0.47
1:A:8:LYS:HB2	1:A:8:LYS:NZ	2.29	0.47
1:A:454:VAL:HG13	1:A:454:VAL:O	2.15	0.47
1:B:63:GLY:HA2	3:B:567:HOH:O	2.14	0.47
1:A:278:HIS:HD2	1:A:344:ASP:OD1	1.97	0.46
1:A:239:HIS:CD2	1:A:241:GLU:H	2.34	0.46
1:B:427:LYS:HE2	1:B:442:LEU:CD1	2.45	0.46
1:B:53:ARG:HG3	1:B:74:ARG:HG3	1.98	0.45
1:A:149:HIS:HE1	3:A:500:HOH:O	1.99	0.45
1:A:450:PHE:O	1:A:454:VAL:HG12	2.17	0.45
1:B:149:HIS:HE1	3:B:488:HOH:O	2.00	0.45
1:A:425:ALA:HB1	1:B:322:PRO:HG2	1.99	0.45
1:A:206:ALA:O	1:A:210:HIS:HD2	2.00	0.45
1:A:132:ASP:O	1:A:136:LYS:HE3	2.17	0.44
1:A:210:HIS:CE1	3:A:588:HOH:O	2.71	0.44
1:A:227:LEU:HD13	1:A:246:VAL:HG21	1.99	0.44
1:A:420:LYS:HA	1:A:420:LYS:HD3	1.80	0.44
1:A:436:LYS:O	1:A:440:LEU:HG	2.18	0.43
1:B:239:HIS:CD2	1:B:241:GLU:H	2.36	0.43
1:A:68:GLU:HG3	1:A:69:LYS:N	2.32	0.43
1:A:120:GLY:CA	1:A:127:LYS:HE2	2.49	0.43
1:A:213:LYS:HG2	1:A:217:TYR:CE2	2.54	0.43
1:A:337:CYS:O	1:A:341:MET:HG2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:263:ASN:ND2	1:B:266:GLU:H	2.11	0.42
1:A:321:MET:CE	1:A:324:LYS:HD3	2.50	0.42
1:A:130:PRO:HG2	2:A:468:MLT:O1	2.18	0.42
1:A:244:ARG:HD3	1:A:244:ARG:HA	1.72	0.42
1:B:315:GLU:HB3	1:B:316:PRO:HD2	2.01	0.42
1:A:322:PRO:HG3	1:B:426:LYS:HA	2.02	0.42
1:A:120:GLY:HA2	1:A:127:LYS:HE2	2.01	0.42
1:B:171:LEU:HD12	1:B:171:LEU:HA	1.92	0.41
1:A:278:HIS:HE1	1:A:372:SER:OG	2.03	0.41
1:A:150:VAL:HA	1:A:258:PHE:CZ	2.55	0.41
1:A:153:LEU:HG	1:A:157:ARG:HD3	2.02	0.41
1:B:301:GLY:HA3	1:B:302:PRO:HA	1.91	0.41
1:B:381:GLU:HG2	1:B:385:LYS:HE2	2.02	0.41
1:A:129:HIS:CD2	2:A:468:MLT:O4	2.73	0.40
1:B:298:LEU:HD12	1:B:298:LEU:HA	1.92	0.40
1:B:8:LYS:HE3	1:B:8:LYS:HB2	1.76	0.40
1:B:129:HIS:CG	2:B:468:MLT:H32	2.56	0.40
1:B:129:HIS:HB3	2:B:468:MLT:O2	2.21	0.40
1:B:153:LEU:HG	1:B:157:ARG:HD3	2.04	0.40

All (2) 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)
1:A:183:LYS:HZ2	3:B:477:HOH:H1[3_756]	1.21	0.39
1:B:183:LYS:HZ2	3:A:473:HOH:H2[3_756]	1.31	0.29

## 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	454/467 (97%)	441 (97%)	13 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	451/467 (97%)	439 (97%)	12 (3%)	0	100	100
All	All	905/934 (97%)	880 (97%)	25 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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	368/376 (98%)	351 (95%)	17 (5%)	27	14
1	B	368/376 (98%)	353 (96%)	15 (4%)	30	18
All	All	736/752 (98%)	704 (96%)	32 (4%)	29	16

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	LYS
1	A	28	ARG
1	A	110	LEU
1	A	117	LEU
1	A	149	HIS
1	A	159	GLN
1	A	165	LYS
1	A	169	GLN
1	A	221	HIS
1	A	230	THR
1	A	244	ARG
1	A	250	LEU
1	A	282	LYS
1	A	397	ARG
1	A	402	LEU
1	A	430	LYS
1	A	459	MET
1	B	10	SER

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Mol	Chain	Res	Type
1	B	90	PRO
1	B	117	LEU
1	B	149	HIS
1	B	171	LEU
1	B	221	HIS
1	B	236	LEU
1	B	282	LYS
1	B	298	LEU
1	B	341	MET
1	B	385	LYS
1	B	396	GLU
1	B	442	LEU
1	B	458	GLN
1	B	459	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (39) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	A	85	HIS
1	A	105	ASN
1	A	129	HIS
1	A	149	HIS
1	A	163	GLN
1	A	169	GLN
1	A	190	GLN
1	A	210	HIS
1	A	239	HIS
1	A	278	HIS
1	A	329	GLN
1	A	339	GLN
1	A	368	ASN
1	A	371	GLN
1	A	384	ASN
1	A	399	ASN
1	A	413	ASN
1	A	458	GLN
1	B	25	GLN
1	B	27	GLN
1	B	59	ASN
1	B	101	GLN
1	B	105	ASN

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Mol	Chain	Res	Type
1	B	149	HIS
1	B	163	GLN
1	B	199	GLN
1	B	239	HIS
1	B	263	ASN
1	B	278	HIS
1	B	314	ASN
1	B	329	GLN
1	B	339	GLN
1	B	368	ASN
1	B	371	GLN
1	B	399	ASN
1	B	415	HIS
1	B	429	HIS
1	B	458	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 monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

2 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
2	MLT	A	468	-	2,8,8	0.91	0	3,10,10	7.34	3 (100%)
2	MLT	B	468	-	2,8,8	0.72	0	3,10,10	7.39	3 (100%)

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	MLT	A	468	-	1/1/3/3	1/2/8/8	-
2	MLT	B	468	-	1/1/3/3	1/2/8/8	-

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	468	MLT	O3-C2-C3	11.90	134.65	108.50
2	A	468	MLT	O3-C2-C3	11.69	134.20	108.50
2	A	468	MLT	C3-C2-C1	4.21	116.45	111.10
2	B	468	MLT	C3-C2-C1	3.86	116.01	111.10
2	B	468	MLT	O3-C2-C1	-2.70	103.53	111.66
2	A	468	MLT	O3-C2-C1	-2.70	103.55	111.66

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	468	MLT	C2
2	B	468	MLT	C2

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	468	MLT	C1-C2-C3-C4
2	A	468	MLT	C1-C2-C3-C4

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	468	MLT	4	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	468	MLT	2	0

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	456/467 (97%)	0.15	41 (8%) <span style="border: 1px solid red; padding: 2px;">9</span> <span style="border: 1px solid red; padding: 2px;">15</span>	4, 16, 78, 92	0
1	B	455/467 (97%)	-0.20	5 (1%) <span style="border: 1px solid blue; padding: 2px;">80</span> <span style="border: 1px solid blue; padding: 2px;">85</span>	4, 15, 38, 83	0
All	All	911/934 (97%)	-0.03	46 (5%) <span style="border: 1px solid red; padding: 2px;">28</span> <span style="border: 1px solid red; padding: 2px;">39</span>	4, 15, 65, 92	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	442	LEU	6.6
1	A	424	ILE	6.3
1	A	453	TRP	6.3
1	A	418	TYR	6.1
1	A	412	LEU	5.7
1	B	459	MET	5.5
1	A	414	THR	5.3
1	A	415	HIS	5.0
1	A	416	ILE	4.9
1	A	441	ALA	4.7
1	B	322	PRO	4.4
1	B	321	MET	4.4
1	B	458	GLN	4.1
1	A	425	ALA	4.0
1	A	419	ASP	4.0
1	A	420	LYS	3.9
1	A	450	PHE	3.7
1	A	433	LEU	3.7
1	A	439	ALA	3.7
1	A	426	LYS	3.7
1	A	448	ALA	3.6
1	A	444	TYR	3.5
1	A	455	ARG	3.4
1	A	445	LEU	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	459	MET	3.0
1	A	319	SER	3.0
1	A	430	LYS	2.8
1	A	431	GLU	2.8
1	A	447	GLU	2.7
1	A	413	ASN	2.7
1	A	435	LEU	2.7
1	A	449	GLU	2.5
1	A	427	LYS	2.5
1	A	125	GLU	2.5
1	A	437	ALA	2.4
1	A	244	ARG	2.4
1	A	428	ALA	2.3
1	A	432	GLY	2.3
1	A	421	ALA	2.3
1	B	68	GLU	2.3
1	A	443	GLY	2.2
1	A	320	ILE	2.1
1	A	237	ASN	2.1
1	A	454	VAL	2.1
1	A	411	ALA	2.1
1	A	440	LEU	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [\(i\)](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates [\(i\)](#)

There are no monosaccharides 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	MLT	B	468	9/9	0.79	0.22	38,42,43,49	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	MLT	A	468	9/9	0.88	0.16	34,38,41,42	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.