



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:10 PM UTC

PDB ID : 7HT3 / pdb\_00007ht3  
Title : PanDDA analysis group deposition – Crystal Structure of FatA in complex with Z31478129  
Authors : Kot, E.; Ni, X.; Tomlinson, C.W.E.; Fearon, D.; Aschenbrenner, J.C.; Fairhead, M.; Koekemoer, L.; Marx, M.L.; Wright, N.D.; Mulholland, N.P.; Montgomery, M.G.; von Delft, F.  
Deposited on : 2024-12-23  
Resolution : 1.89 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references](#) ①) 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

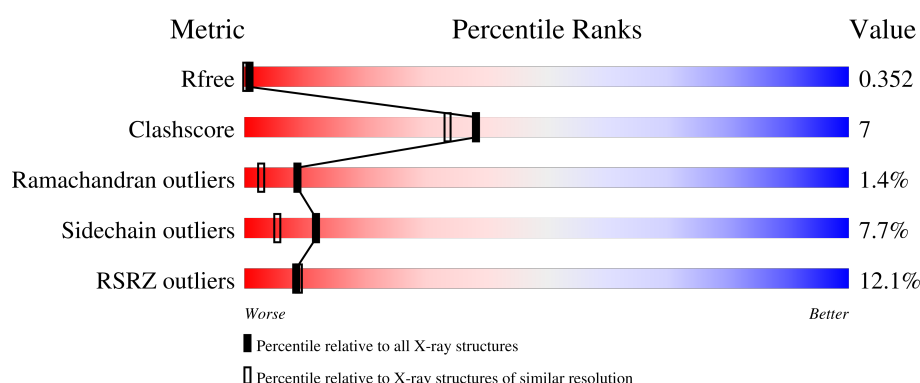
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.89 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	295	<div> <div>6%</div> <div>68%</div> <div>19%</div> <div>•</div> <div>10%</div> </div>
1	B	295	<div> <div>16%</div> <div>68%</div> <div>19%</div> <div>•</div> <div>11%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4612 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.

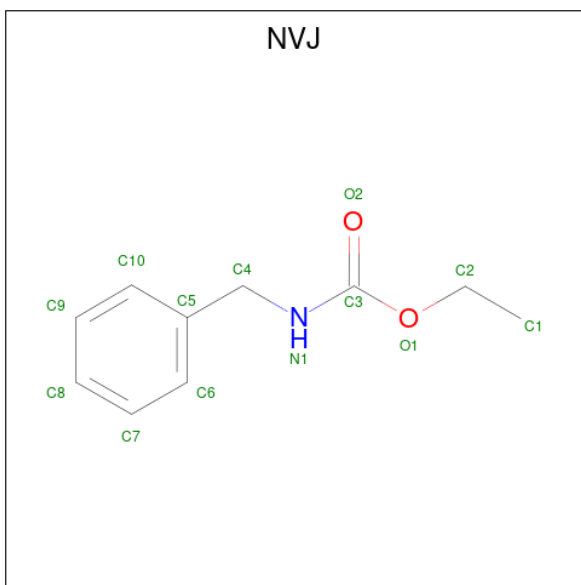
- Molecule 1 is a protein called Oleoyl-acyl carrier protein thioesterase 1, chloroplastic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	0	4	0
			2192	1363	388	427	14			
1	B	264	Total	C	N	O	S	0	3	0
			2190	1363	390	427	10			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	74	MET	-	initiating methionine	UNP Q42561
A	363	HIS	-	expression tag	UNP Q42561
A	364	HIS	-	expression tag	UNP Q42561
A	365	HIS	-	expression tag	UNP Q42561
A	366	HIS	-	expression tag	UNP Q42561
A	367	HIS	-	expression tag	UNP Q42561
A	368	HIS	-	expression tag	UNP Q42561
B	74	MET	-	initiating methionine	UNP Q42561
B	363	HIS	-	expression tag	UNP Q42561
B	364	HIS	-	expression tag	UNP Q42561
B	365	HIS	-	expression tag	UNP Q42561
B	366	HIS	-	expression tag	UNP Q42561
B	367	HIS	-	expression tag	UNP Q42561
B	368	HIS	-	expression tag	UNP Q42561

- Molecule 2 is ethyl benzylcarbamate (CCD ID: NVJ) (formula:  $C_{10}H_{13}NO_2$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			13	10	1	2		
2	B	1	Total	C	N	O	0	0
			13	10	1	2		

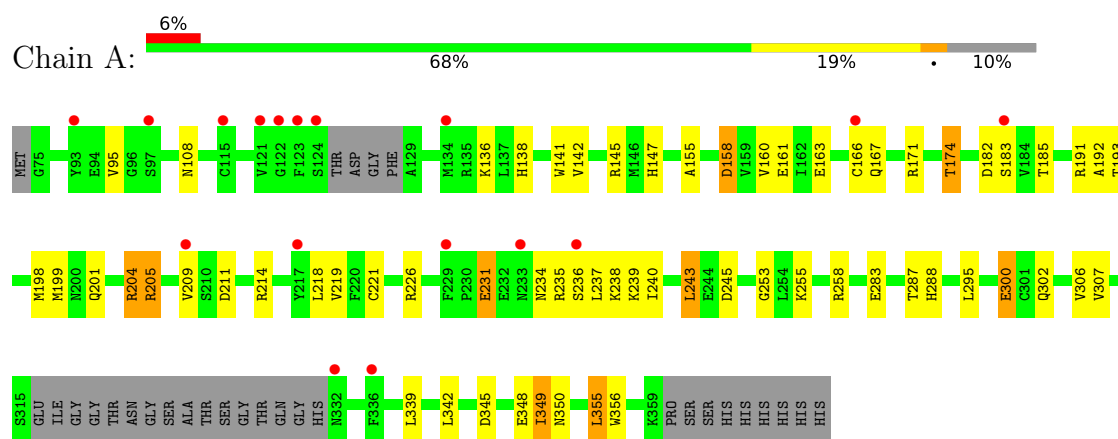
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	106	Total	O	0	0
			106	106		
3	B	98	Total	O	0	0
			98	98		

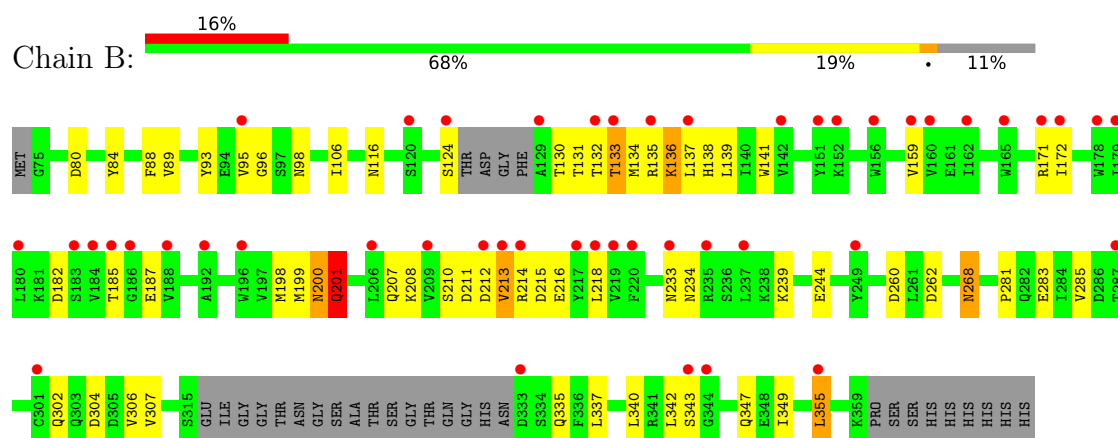
### 3 Residue-property plots [i](#)

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: Oleoyl-acyl carrier protein thioesterase 1, chloroplastic



- Molecule 1: Oleoyl-acyl carrier protein thioesterase 1, chloroplastic



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	98.70Å 98.61Å 128.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	78.24 – 1.89 78.24 – 1.89	Depositor EDS
% Data completeness (in resolution range)	94.5 (78.24-1.89) 94.6 (78.24-1.89)	Depositor EDS
$R_{merge}$	0.29	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	29.27 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.8.0267, REFMAC5	Depositor
R, $R_{free}$	0.312 , 0.377 0.317 , 0.352	Depositor DCC
$R_{free}$ test set	2428 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	8.3	Xtriage
Anisotropy	1.011	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 39.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.460 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.81	EDS
Total number of atoms	4612	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.76% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NVJ

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	1.25	5/2230 (0.2%)	1.54	10/3017 (0.3%)
1	B	1.18	6/2228 (0.3%)	1.47	9/3014 (0.3%)
All	All	1.22	11/4458 (0.2%)	1.50	19/6031 (0.3%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	142	VAL	C-O	7.56	1.31	1.24
1	B	96	GLY	C-O	7.25	1.30	1.23
1	A	95	VAL	C-O	7.07	1.31	1.24
1	A	307	VAL	C-O	6.98	1.31	1.24
1	B	307	VAL	C-O	6.42	1.30	1.24
1	A	243	LEU	C-O	6.31	1.31	1.23
1	B	340	LEU	C-O	6.01	1.31	1.24
1	B	116	ASN	C-O	5.64	1.30	1.24
1	B	96	GLY	N-CA	5.22	1.49	1.44
1	A	138	HIS	C-O	5.10	1.30	1.23
1	B	84	TYR	C-O	5.00	1.29	1.23

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	95	VAL	CA-C-N	8.01	130.20	122.27
1	A	95	VAL	C-N-CA	8.01	130.20	122.27
1	B	268	ASN	CB-CA-C	7.19	121.95	109.65
1	B	281	PRO	CA-C-N	6.47	128.95	120.28
1	B	281	PRO	C-N-CA	6.47	128.95	120.28
1	A	158	ASP	CA-CB-CG	6.31	118.91	112.60
1	A	201	GLN	CB-CA-C	-5.77	101.20	110.79
1	A	342	LEU	CA-C-O	-5.65	114.48	120.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	88	PHE	CA-C-O	-5.59	114.11	120.32
1	B	80	ASP	CA-CB-CG	5.50	118.10	112.60
1	A	95	VAL	O-C-N	5.39	128.69	123.03
1	A	193	THR	CB-CA-C	5.30	120.46	109.38
1	A	161	GLU	CB-CG-CD	5.28	121.58	112.60
1	A	245	ASP	CB-CA-C	5.26	117.74	109.27
1	B	349	ILE	CA-C-O	-5.13	115.48	120.30
1	A	300	GLU	CB-CA-C	5.10	118.23	109.51
1	B	337	LEU	N-CA-C	-5.02	101.97	109.95
1	B	95	VAL	CA-C-N	5.01	126.89	122.43
1	B	95	VAL	C-N-CA	5.01	126.89	122.43

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	2192	0	2152	32	0
1	B	2190	0	2149	30	0
2	A	13	0	0	0	0
2	B	13	0	0	0	0
3	A	106	0	0	4	0
3	B	98	0	0	2	0
All	All	4612	0	4301	60	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 (60) 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:205:ARG:HD3	3:A:510:HOH:O	1.38	1.22
1:A:141:TRP:CZ3	1:A:174:THR:OG1	2.12	1.03
1:B:138:HIS:O	1:B:200:ASN:O	2.01	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:ARG:NE	3:A:501:HOH:O	2.08	0.74
1:A:205:ARG:HG2	1:A:205:ARG:HH11	1.51	0.73
1:A:211:ASP:OD1	1:A:214:ARG:NH2	2.21	0.73
1:B:137:LEU:HB2	3:B:557:HOH:O	1.90	0.69
1:A:167:GLN:O	1:A:174:THR:HG22	1.99	0.61
1:B:214:ARG:O	1:B:216:GLU:N	2.28	0.60
1:A:199:MET:HE1	1:A:204:ARG:CZ	2.32	0.60
1:B:200:ASN:O	1:B:201:GLN:HB3	2.06	0.55
1:A:355:LEU:HD12	1:A:356:TRP:N	2.22	0.54
1:A:237:LEU:HD11	3:A:512:HOH:O	2.08	0.54
1:A:355:LEU:HD12	1:A:355:LEU:C	2.33	0.54
1:B:210:SER:O	1:B:214:ARG:HB2	2.07	0.53
1:A:258:ARG:CZ	1:B:285:VAL:HG21	2.39	0.52
1:B:207:GLN:NE2	3:B:505:HOH:O	2.43	0.51
1:A:163:GLU:CD	1:A:226:ARG:HH22	2.18	0.51
1:A:198:MET:HE2	1:A:209:VAL:HA	1.92	0.51
1:B:306:VAL:HG12	1:B:343:SER:OG	2.11	0.50
1:A:147:HIS:ND1	1:A:236:SER:HA	2.26	0.50
1:A:199:MET:HE1	1:A:204:ARG:NH2	2.26	0.50
1:A:283:GLU:O	1:A:287:THR:HG23	2.12	0.49
1:B:171:ARG:O	1:B:208:LYS:HB2	2.13	0.49
1:B:182:ASP:HB3	1:B:185:THR:OG1	2.12	0.49
1:B:260:ASP:O	1:B:268:ASN:OD1	2.31	0.49
1:B:342:LEU:HD12	1:B:347:GLN:HB3	1.95	0.48
1:B:132:THR:O	1:B:216:GLU:OE1	2.30	0.48
1:B:134:MET:HE1	1:B:141:TRP:CD2	2.47	0.48
1:A:287:THR:HG1	1:A:288:HIS:CE1	2.31	0.48
1:B:335:GLN:HG2	1:B:355:LEU:HD13	1.96	0.47
1:B:214:ARG:C	1:B:216:GLU:H	2.21	0.47
1:A:287:THR:OG1	1:A:288:HIS:ND1	2.48	0.47
1:A:166[C]:CYS:HG	1:A:221:CYS:HG	1.61	0.47
1:A:155:ALA:O	1:A:158:ASP:HB2	2.15	0.46
1:A:166[A]:CYS:HG	1:A:221:CYS:HG	1.61	0.46
1:A:253:GLY:HA2	1:A:306:VAL:CG1	2.46	0.45
1:B:133:THR:HA	1:B:136:LYS:HB2	1.99	0.45
1:B:134:MET:HE1	1:B:141:TRP:CE3	2.52	0.45
1:B:210:SER:HB3	1:B:213:VAL:HG23	1.98	0.44
1:B:200:ASN:O	1:B:201:GLN:CB	2.66	0.44
1:B:304:ASP:N	1:B:304:ASP:OD1	2.48	0.43
1:A:349:ILE:HG13	1:A:350:ASN:OD1	2.19	0.43
1:A:145:ARG:NH2	1:A:240:ILE:HD12	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:182:ASP:HB3	1:A:185:THR:OG1	2.18	0.43
1:A:108:ASN:HB3	1:B:93:TYR:CE2	2.54	0.43
1:B:134:MET:CE	1:B:198:MET:HE1	2.47	0.43
1:B:139:LEU:HD23	1:B:200:ASN:HA	2.00	0.43
1:B:212:ASP:O	1:B:216:GLU:CD	2.62	0.43
1:B:89:VAL:HG22	1:B:159:VAL:HG22	2.00	0.43
1:A:147:HIS:O	1:A:192:ALA:HA	2.19	0.42
1:B:141:TRP:CD1	1:B:198:MET:CE	3.03	0.42
1:B:210:SER:OG	1:B:211:ASP:N	2.53	0.42
3:A:585:HOH:O	1:B:262:ASP:HB2	2.19	0.41
1:A:198:MET:CE	1:A:209:VAL:HA	2.51	0.41
1:B:214:ARG:C	1:B:216:GLU:N	2.78	0.41
1:A:339:LEU:HD11	1:A:348:GLU:HG3	2.02	0.41
1:A:218:LEU:HD13	1:A:218:LEU:C	2.46	0.41
1:A:231:GLU:HB2	1:A:234:ASN:HB3	2.02	0.40
1:A:295:LEU:HD11	1:A:350:ASN:HB3	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	267/295 (90%)	254 (95%)	11 (4%)	2 (1%)	18	10
1	B	264/295 (90%)	250 (95%)	9 (3%)	5 (2%)	6	1
All	All	531/590 (90%)	504 (95%)	20 (4%)	7 (1%)	9	3

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	231	GLU
1	A	345	ASP

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Mol	Chain	Res	Type
1	B	201	GLN
1	B	233	ASN
1	B	200	ASN
1	B	215	ASP
1	B	172	ILE

### 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	248/263 (94%)	230 (93%)	18 (7%)	13	6
1	B	245/263 (93%)	226 (92%)	19 (8%)	11	5
All	All	493/526 (94%)	456 (92%)	37 (8%)	12	6

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	136	LYS
1	A	160	VAL
1	A	171	ARG
1	A	174	THR
1	A	183	SER
1	A	191	ARG
1	A	204	ARG
1	A	205	ARG
1	A	219	VAL
1	A	235	ARG
1	A	238	LYS
1	A	239	LYS
1	A	243	LEU
1	A	255	LYS
1	A	300	GLU
1	A	302	GLN
1	A	349	ILE
1	A	355	LEU

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Mol	Chain	Res	Type
1	B	98	ASN
1	B	106	ILE
1	B	124	SER
1	B	130	THR
1	B	131	THR
1	B	133	THR
1	B	135	ARG
1	B	136	LYS
1	B	187	GLU
1	B	199	MET
1	B	201	GLN
1	B	213	VAL
1	B	218	LEU
1	B	234	ASN
1	B	239	LYS
1	B	244	GLU
1	B	283	GLU
1	B	302	GLN
1	B	355	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	ASN
1	A	119	GLN
1	A	223	GLN
1	A	291	GLN
1	A	302	GLN
1	B	119	GLN
1	B	207	GLN
1	B	264	ASN
1	B	268	ASN
1	B	303	GLN

### 5.3.3 RNA ⓘ

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](#)

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	NVJ	B	401	-	13,13,13	0.21	0	15,15,15	0.29	0
2	NVJ	A	401	-	13,13,13	0.27	0	15,15,15	0.36	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
2	NVJ	B	401	-	-	2/8/8/8	0/1/1/1
2	NVJ	A	401	-	-	4/8/8/8	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	401	NVJ	N1-C3-O1-C2
2	A	401	NVJ	N1-C3-O1-C2

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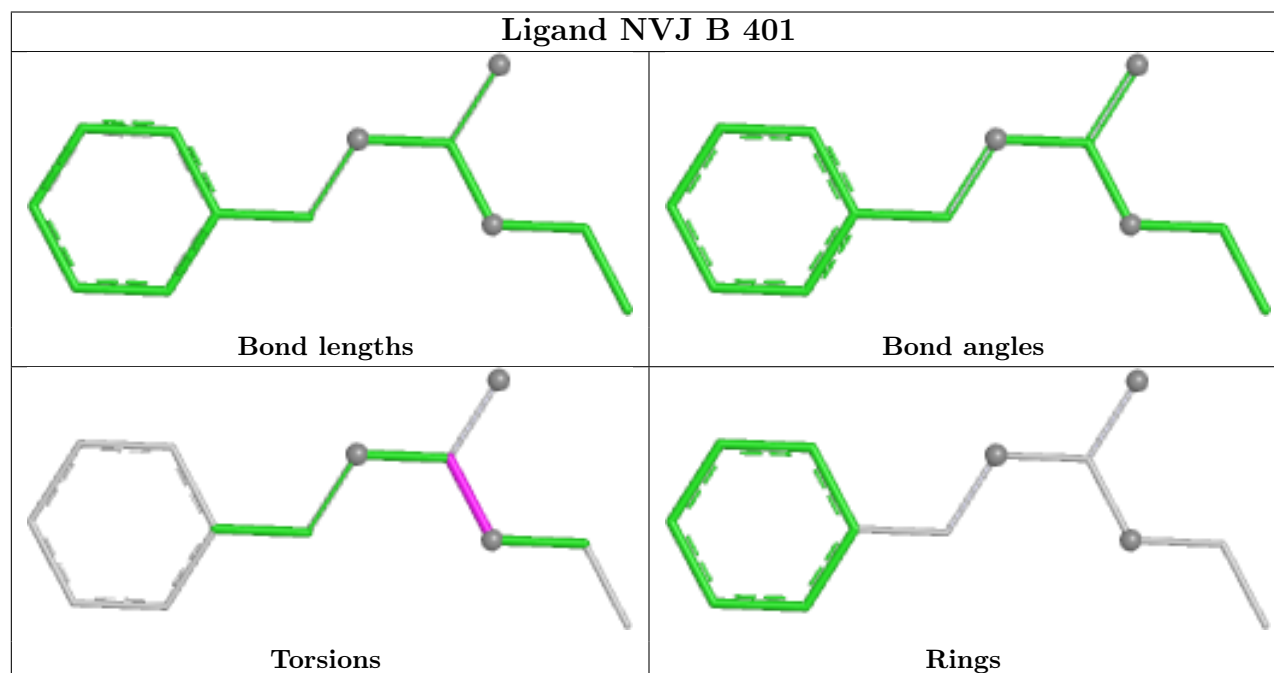
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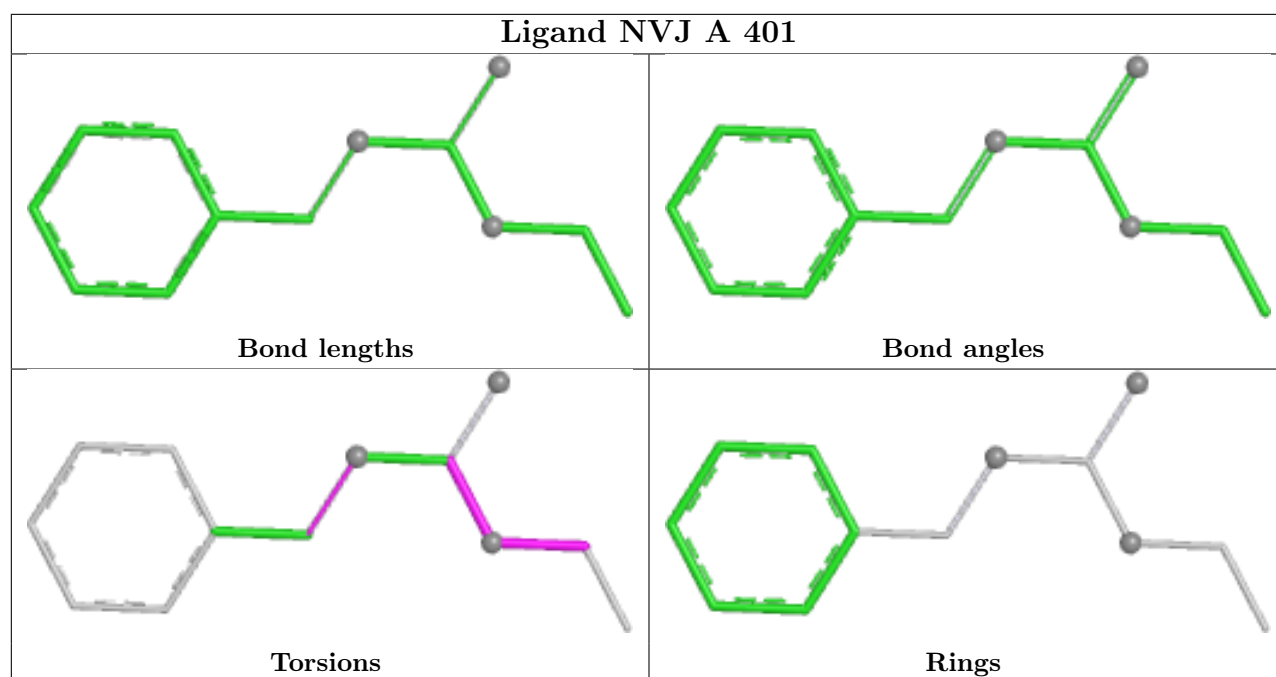
Mol	Chain	Res	Type	Atoms
2	A	401	NVJ	O2-C3-O1-C2
2	B	401	NVJ	O2-C3-O1-C2
2	A	401	NVJ	C1-C2-O1-C3
2	A	401	NVJ	C5-C4-N1-C3

There are no ring outliers.

No monomer is involved in short contacts.

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, 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	265/295 (89%)	0.58	17 (6%) 25 27	6, 37, 62, 96	13 (4%)
1	B	264/295 (89%)	1.08	47 (17%) 4 3	5, 32, 58, 100	85 (32%)
All	All	529/590 (89%)	0.83	64 (12%) 8 9	5, 35, 61, 100	98 (18%)

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	213	VAL	4.5
1	A	217	TYR	3.9
1	A	97[A]	SER	3.7
1	A	166[A]	CYS	3.6
1	B	184	VAL	3.6
1	A	123	PHE	3.5
1	B	179	ILE	3.0
1	B	137	LEU	3.0
1	B	219	VAL	3.0
1	A	229	PHE	3.0
1	B	344	GLY	3.0
1	B	172	ILE	2.9
1	B	188	VAL	2.8
1	B	206	LEU	2.8
1	B	217	TYR	2.8
1	A	121	VAL	2.8
1	A	336	PHE	2.8
1	B	156	TRP	2.8
1	B	209	VAL	2.7
1	B	192	ALA	2.7
1	B	218	LEU	2.7
1	B	186	GLY	2.7
1	B	233	ASN	2.7
1	B	214	ARG	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	142	VAL	2.6
1	B	129	ALA	2.6
1	B	235	ARG	2.6
1	B	133	THR	2.6
1	B	151	TYR	2.5
1	B	180	LEU	2.5
1	A	233	ASN	2.5
1	B	185	THR	2.5
1	B	237	LEU	2.4
1	B	120	SER	2.4
1	B	212	ASP	2.4
1	B	165	TRP	2.4
1	B	196	TRP	2.4
1	B	301	CYS	2.4
1	B	178	TRP	2.4
1	A	122	GLY	2.4
1	B	249	TYR	2.4
1	B	287	THR	2.3
1	B	95	VAL	2.3
1	B	343	SER	2.3
1	B	152	LYS	2.3
1	B	183	SER	2.3
1	B	160	VAL	2.3
1	B	135	ARG	2.3
1	A	115	CYS	2.3
1	A	134[A]	MET	2.2
1	B	355	LEU	2.2
1	B	220	PHE	2.2
1	B	159	VAL	2.2
1	A	124	SER	2.2
1	A	93	TYR	2.1
1	B	162	ILE	2.1
1	B	333	ASP	2.1
1	A	183	SER	2.1
1	A	236	SER	2.1
1	B	124	SER	2.1
1	A	209	VAL	2.1
1	B	171	ARG	2.1
1	A	332	ASN	2.1
1	B	132	THR	2.1

## 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](#)

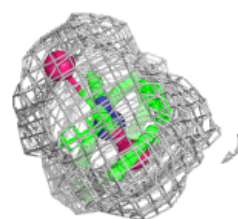
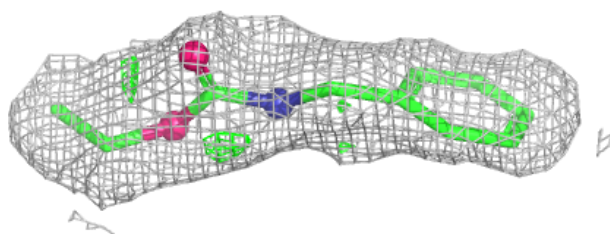
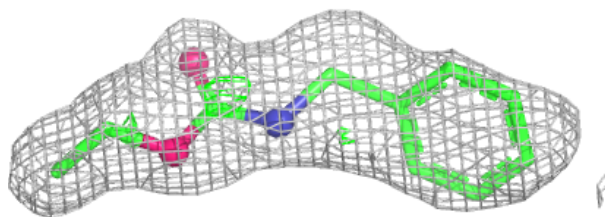
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	NVJ	A	401	13/13	0.95	0.11	20,21,22,25	13
2	NVJ	B	401	13/13	0.96	0.08	20,22,27,27	13

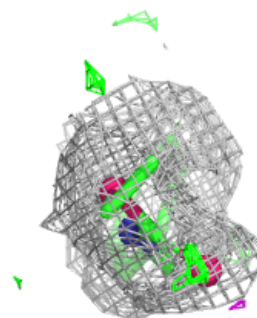
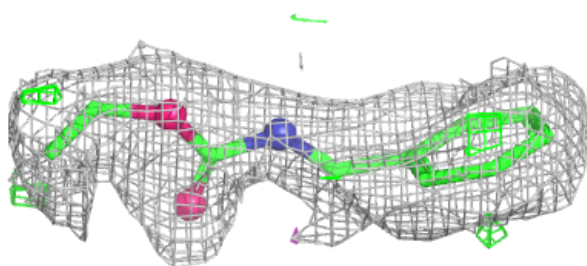
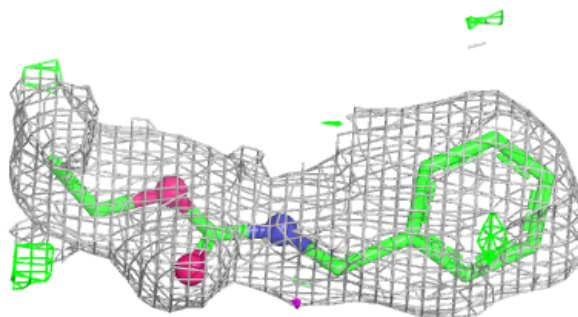
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around NVJ 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)

**Electron density around NVJ B 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.