



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:02 PM UTC

PDB ID : 7NBD / pdb_00007nbd
Title : Crystal structure of human serine racemase in complex with DSiP fragment Z235449082, XChem fragment screen.
Authors : Koulouris, C.R.; Roe, S.M.
Deposited on : 2021-01-26
Resolution : 1.86 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

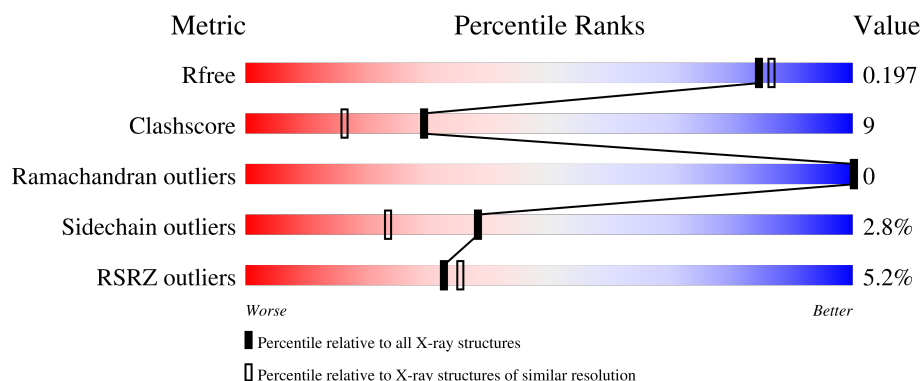
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.86 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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	AAA	346	2% 79% 13% • 7%
1	BBB	346	9% 80% 9% • 9%
1	CCC	346	3% 83% 10% 7%
1	DDD	346	4% 80% 11% • 8%

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	GOL	AAA	403	-	-	X	-
3	GOL	AAA	404	-	-	X	-
3	GOL	AAA	409	-	-	X	-
3	GOL	BBB	402	-	-	X	-

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 11477 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 Serine racemase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	323	Total	C	N	O	S	0	35	0
			2688	1703	452	521	12			
1	BBB	315	Total	C	N	O	S	0	13	1
			2412	1535	400	466	11			
1	CCC	323	Total	C	N	O	S	5	22	1
			2573	1636	427	498	12			
1	DDD	318	Total	C	N	O	S	0	20	0
			2527	1603	434	477	13			

There are 32 discrepancies between the modelled and reference sequences:

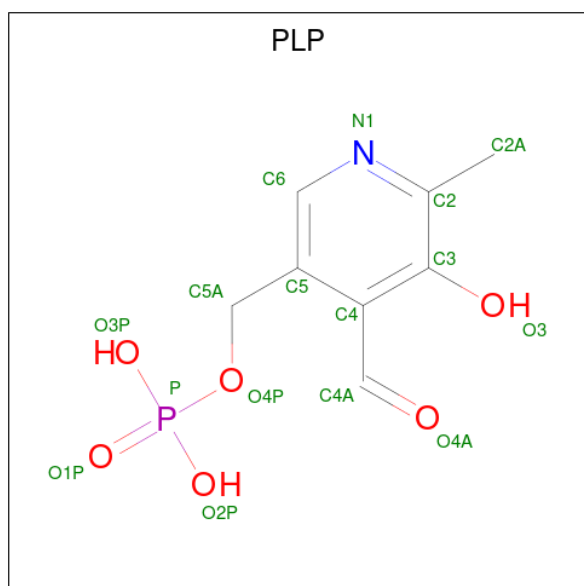
Chain	Residue	Modelled	Actual	Comment	Reference
AAA	2	ASP	CYS	engineered mutation	UNP Q9GZT4
AAA	6	ASP	CYS	engineered mutation	UNP Q9GZT4
AAA	341	HIS	-	expression tag	UNP Q9GZT4
AAA	342	HIS	-	expression tag	UNP Q9GZT4
AAA	343	HIS	-	expression tag	UNP Q9GZT4
AAA	344	HIS	-	expression tag	UNP Q9GZT4
AAA	345	HIS	-	expression tag	UNP Q9GZT4
AAA	346	HIS	-	expression tag	UNP Q9GZT4
BBB	2	ASP	CYS	engineered mutation	UNP Q9GZT4
BBB	6	ASP	CYS	engineered mutation	UNP Q9GZT4
BBB	341	HIS	-	expression tag	UNP Q9GZT4
BBB	342	HIS	-	expression tag	UNP Q9GZT4
BBB	343	HIS	-	expression tag	UNP Q9GZT4
BBB	344	HIS	-	expression tag	UNP Q9GZT4
BBB	345	HIS	-	expression tag	UNP Q9GZT4
BBB	346	HIS	-	expression tag	UNP Q9GZT4
CCC	2	ASP	CYS	engineered mutation	UNP Q9GZT4
CCC	6	ASP	CYS	engineered mutation	UNP Q9GZT4
CCC	341	HIS	-	expression tag	UNP Q9GZT4
CCC	342	HIS	-	expression tag	UNP Q9GZT4
CCC	343	HIS	-	expression tag	UNP Q9GZT4

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
CCC	344	HIS	-	expression tag	UNP Q9GZT4
CCC	345	HIS	-	expression tag	UNP Q9GZT4
CCC	346	HIS	-	expression tag	UNP Q9GZT4
DDD	2	ASP	CYS	engineered mutation	UNP Q9GZT4
DDD	6	ASP	CYS	engineered mutation	UNP Q9GZT4
DDD	362	HIS	-	expression tag	UNP Q9GZT4
DDD	363	HIS	-	expression tag	UNP Q9GZT4
DDD	364	HIS	-	expression tag	UNP Q9GZT4
DDD	365	HIS	-	expression tag	UNP Q9GZT4
DDD	366	HIS	-	expression tag	UNP Q9GZT4
DDD	367	HIS	-	expression tag	UNP Q9GZT4

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: $C_8H_{10}NO_6P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	AAA	1	Total	C	N	O	P	0	1
			30	16	2	10	2		
2	BBB	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	CCC	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	DDD	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	AAA	1	Total	C	O	0	0
			6	3	3		
3	AAA	1	Total	C	O	0	0
			6	3	3		
3	AAA	1	Total	C	O	0	0
			6	3	3		
3	AAA	1	Total	C	O	0	0
			6	3	3		
3	AAA	1	Total	C	O	0	0
			6	3	3		
3	BBB	1	Total	C	O	0	0
			6	3	3		
3	BBB	1	Total	C	O	0	0
			6	3	3		
3	CCC	1	Total	C	O	0	0
			6	3	3		

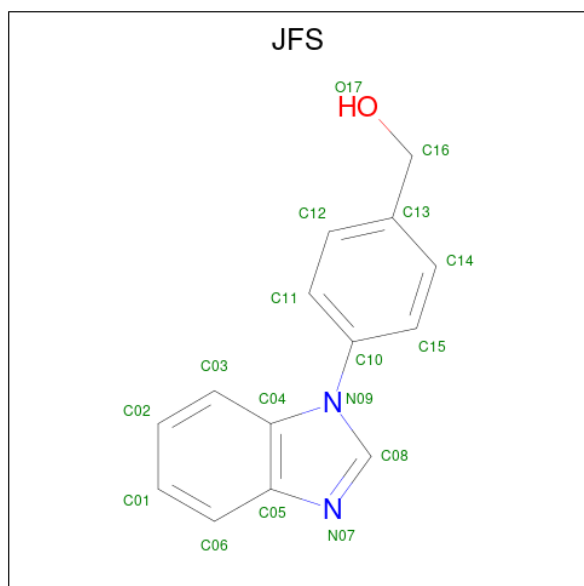
- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	AAA	1	Total	Ca	0	0
			1	1		
4	BBB	2	Total	Ca	0	0
			2	2		
4	CCC	2	Total	Ca	0	1
			3	3		
4	DDD	2	Total	Ca	0	1
			3	3		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

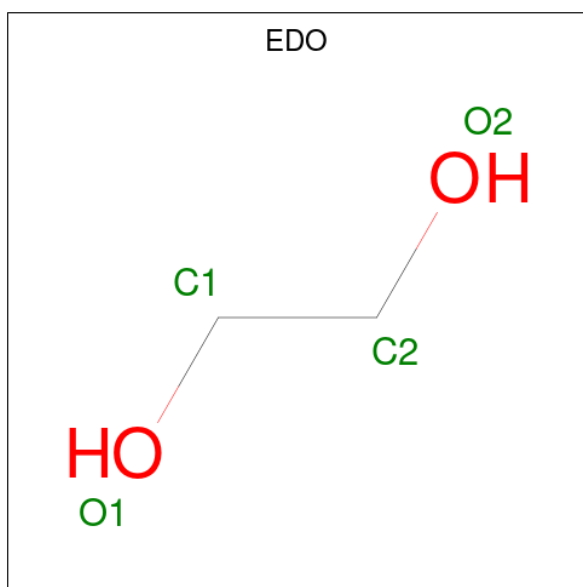
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	AAA	1	Total	Mg	0	0
			1	1		
5	DDD	1	Total	Mg	0	0
			1	1		

- Molecule 6 is [4-(1H-benzimidazol-1-yl)phenyl]methanol (CCD ID: JFS) (formula: C₁₄H₁₂N₂O) (labeled as "Ligand of Interest" by depositor).



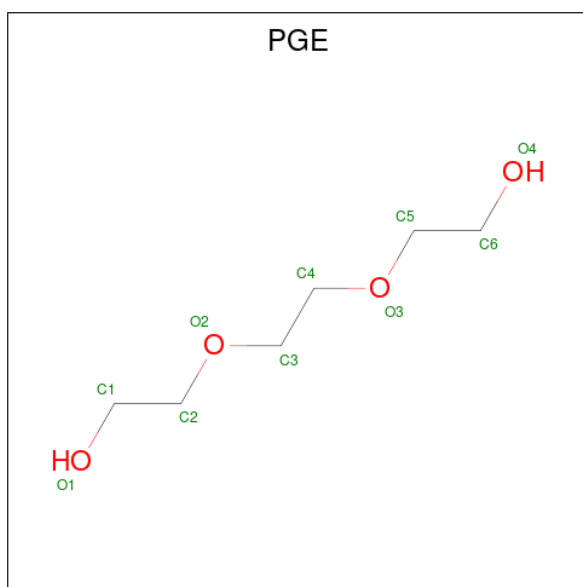
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	AAA	1	Total	C	N	O	0	1
			34	28	4	2		
6	BBB	1	Total	C	N	O	0	0
			17	14	2	1		

- Molecule 7 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	BBB	1	Total	C	O	0	0
			4	2	2		
7	CCC	1	Total	C	O	0	0
			4	2	2		
7	DDD	1	Total	C	O	0	0
			4	2	2		

- Molecule 8 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	CCC	1	Total	C	O	0	0
			10	6	4		

- Molecule 9 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	DDD	1	Total	Na	0	0
			1	1		

- Molecule 10 is water.

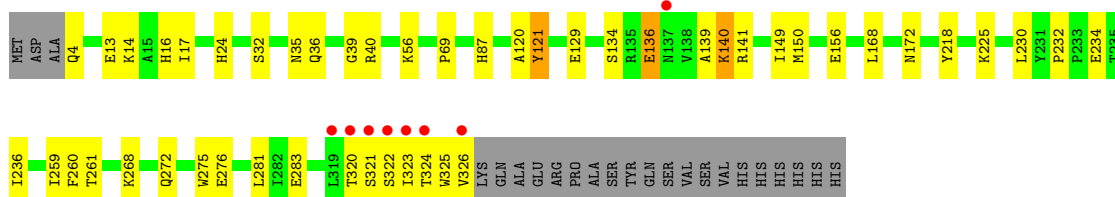
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	AAA	323	Total	O	0	4
			327	327		
10	BBB	182	Total	O	0	2
			184	184		
10	CCC	262	Total	O	0	6
			268	268		
10	DDD	286	Total	O	0	4
			290	290		

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.

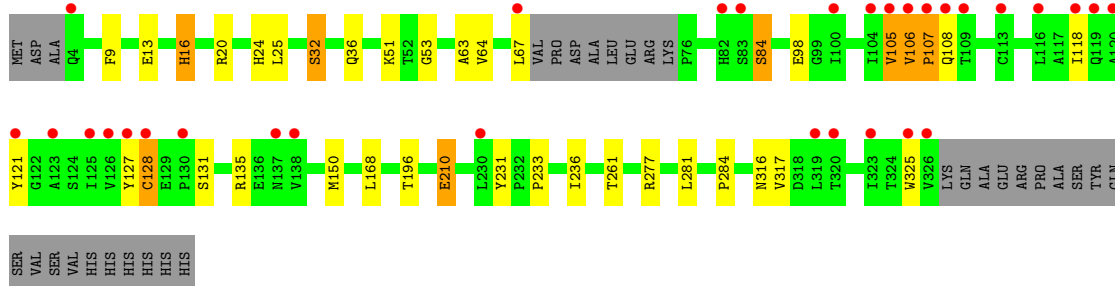
- Molecule 1: Serine racemase

Chain AAA: 



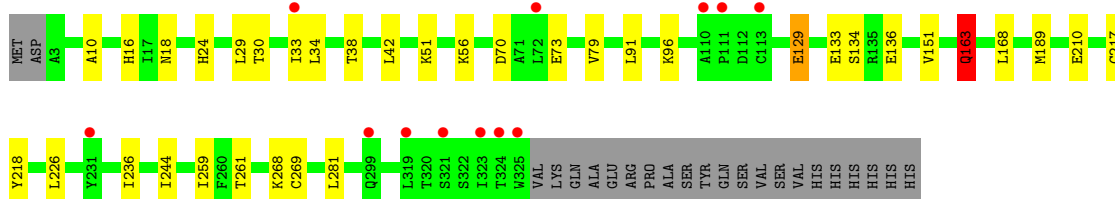
- Molecule 1: Serine racemase

Chain BBB: 



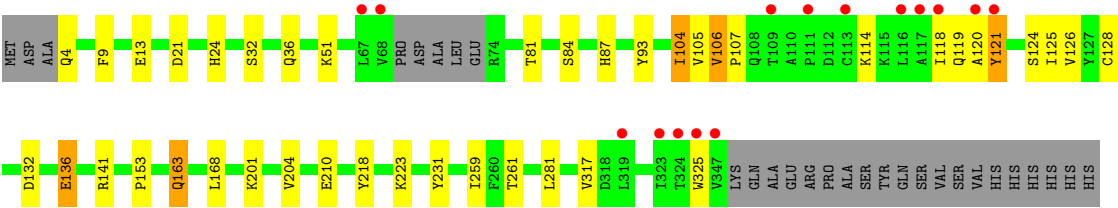
- Molecule 1: Serine racemase

Chain CCC: 



- Molecule 1: Serine racemase

Chain DDD: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	48.01Å 154.85Å 85.45Å 90.00° 98.04° 90.00°	Depositor
Resolution (Å)	84.60 – 1.86 84.60 – 1.86	Depositor EDS
% Data completeness (in resolution range)	93.9 (84.60-1.86) 93.9 (84.60-1.86)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.37 (at 1.87Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.176 , 0.207 (Not available) , 0.197	Depositor DCC
R_{free} test set	4830 reflections (4.71%)	wwPDB-VP
Wilson B-factor (Å ²)	28.5	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 50.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	11477	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.26% 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: CA, JFS, PGE, NA, GOL, PLP, EDO, MG

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	AAA	0.75	0/2735	1.02	3/3728 (0.1%)
1	BBB	0.65	0/2454	1.05	9/3352 (0.3%)
1	CCC	0.69	1/2619 (0.0%)	1.05	3/3576 (0.1%)
1	DDD	0.74	1/2571 (0.0%)	1.09	3/3505 (0.1%)
All	All	0.71	2/10379 (0.0%)	1.05	18/14161 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	CCC	16	HIS	CE1-NE2	5.85	1.38	1.32
1	DDD	325	TRP	C-N	5.38	1.40	1.33

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AAA	283	GLU	CB-CA-C	7.96	122.08	109.27
1	CCC	163	GLN	CB-CG-CD	7.83	125.91	112.60
1	DDD	121	TYR	CB-CA-C	7.70	124.51	109.72
1	BBB	128	CYS	N-CA-CB	-7.21	99.98	110.87
1	AAA	121	TYR	CB-CA-C	6.45	121.62	110.01
1	DDD	163	GLN	CB-CG-CD	6.45	123.56	112.60
1	BBB	121	TYR	CB-CA-C	6.32	121.85	109.72
1	BBB	107	PRO	N-CA-C	-6.16	101.54	111.03
1	CCC	129[A]	GLU	CB-CA-C	5.69	117.06	108.86
1	CCC	129[B]	GLU	CB-CA-C	5.69	117.06	108.86
1	BBB	196	THR	CA-CB-OG1	-5.64	101.14	109.60
1	BBB	210	GLU	N-CA-CB	-5.61	101.63	111.17
1	BBB	105	VAL	CA-C-N	-5.60	117.27	123.43
1	BBB	105	VAL	C-N-CA	-5.60	117.27	123.43
1	BBB	277	ARG	CG-CD-NE	-5.50	99.91	112.00

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BBB	127	TYR	CB-CA-C	5.18	118.17	109.89
1	DDD	106	VAL	CB-CA-C	5.16	115.14	110.13
1	AAA	276	GLU	CB-CG-CD	5.07	121.22	112.60

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	AAA	2688	0	2666	73	1
1	BBB	2412	0	2371	28	0
1	CCC	2573	0	2556	38	0
1	DDD	2527	0	2519	36	1
2	AAA	30	0	12	0	0
2	BBB	15	0	6	0	0
2	CCC	15	0	6	1	0
2	DDD	15	0	6	0	0
3	AAA	30	0	40	21	0
3	BBB	12	0	16	5	0
3	CCC	6	0	8	2	0
4	AAA	1	0	0	0	0
4	BBB	2	0	0	0	0
4	CCC	3	0	0	0	0
4	DDD	3	0	0	0	0
5	AAA	1	0	0	0	0
5	DDD	1	0	0	0	0
6	AAA	34	0	0	0	0
6	BBB	17	0	0	0	0
7	BBB	4	0	6	0	0
7	CCC	4	0	5	2	0
7	DDD	4	0	6	0	0
8	CCC	10	0	14	5	0
9	DDD	1	0	0	0	0
10	AAA	327	0	0	21	1
10	BBB	184	0	0	3	1

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	CCC	268	0	0	9	1
10	DDD	290	0	0	5	1
All	All	11477	0	10237	179	5

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 9.

All (179) 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)
3:AAA:403:GOL:H12	10:AAA:521:HOH:O	1.17	1.32
1:AAA:140:LYS:CE	1:AAA:150[B]:MET:HE1	1.59	1.29
1:AAA:322[B]:SER:O	1:AAA:326[B]:VAL:HG12	1.39	1.18
1:AAA:261:THR:H	3:AAA:403:GOL:H32	1.13	1.08
1:AAA:140:LYS:HE3	1:AAA:150[B]:MET:HE1	1.20	1.06
1:AAA:14[B]:LYS:HE3	10:AAA:609:HOH:O	1.55	1.04
1:DDD:81[B]:THR:CG2	1:DDD:104[B]:ILE:CD1	2.37	1.02
1:DDD:81[B]:THR:CG2	1:DDD:104[B]:ILE:HD13	1.97	0.95
1:CCC:51[B]:LYS:NZ	10:CCC:501[B]:HOH:O	1.98	0.95
1:AAA:140:LYS:HE3	1:AAA:150[B]:MET:CE	1.98	0.94
1:DDD:223:LYS:HG2	10:DDD:506:HOH:O	1.68	0.94
1:AAA:136:GLU:HA	10:AAA:506:HOH:O	1.66	0.93
1:CCC:30:THR:OG1	8:CCC:402:PGE:H12	1.66	0.92
1:DDD:81[B]:THR:HG22	1:DDD:104[B]:ILE:HD13	1.51	0.92
1:AAA:140:LYS:HE2	1:AAA:150[B]:MET:HE1	1.53	0.88
1:AAA:172:ASN:ND2	10:AAA:501:HOH:O	2.06	0.87
1:CCC:51[A]:LYS:NZ	10:CCC:501[A]:HOH:O	2.05	0.87
1:CCC:79[A]:VAL:HG11	1:CCC:91:LEU:HG	1.56	0.87
1:AAA:149:ILE:HA	3:AAA:404:GOL:H32	1.57	0.86
1:DDD:81[B]:THR:HG23	1:DDD:104[B]:ILE:CD1	2.04	0.86
1:AAA:272:GLN:HE21	1:AAA:325[B]:TRP:HE1	1.24	0.86
1:AAA:35[B]:ASN:O	1:AAA:39[B]:GLY:N	2.10	0.85
8:CCC:402:PGE:H32	10:CCC:718:HOH:O	1.76	0.84
1:CCC:30:THR:O	8:CCC:402:PGE:H42	1.78	0.84
1:CCC:24:HIS:NE2	1:CCC:51[B]:LYS:HE3	1.93	0.83
1:AAA:322[B]:SER:C	1:AAA:326[B]:VAL:HG12	2.02	0.83
1:DDD:21[B]:ASP:OD1	10:DDD:762[B]:HOH:O	1.96	0.83
1:AAA:16[B]:HIS:CD2	3:AAA:409:GOL:HO2	1.96	0.83
1:AAA:322[B]:SER:O	1:AAA:326[B]:VAL:CG1	2.26	0.82
1:CCC:24:HIS:CE1	1:CCC:51[B]:LYS:HE3	2.15	0.81
1:AAA:16[B]:HIS:CD2	3:AAA:409:GOL:O2	2.34	0.81

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:DDD:81[B]:THR:HG23	1:DDD:104[B]:ILE:HD12	1.63	0.80
1:AAA:261:THR:H	3:AAA:403:GOL:C3	1.95	0.80
1:DDD:104[A]:ILE:HG21	1:DDD:118:ILE:HG21	1.62	0.80
1:AAA:261:THR:N	3:AAA:403:GOL:H32	1.96	0.79
1:DDD:51:LYS:NZ	10:DDD:502:HOH:O	2.03	0.78
1:AAA:168:LEU:HD13	3:AAA:409:GOL:H31	1.65	0.76
1:CCC:18[A]:ASN:OD1	3:CCC:403:GOL:H11	1.85	0.76
3:AAA:402:GOL:H31	10:AAA:602:HOH:O	1.87	0.74
3:AAA:403:GOL:H31	10:AAA:718:HOH:O	1.88	0.74
1:AAA:322[A]:SER:O	1:AAA:324[A]:THR:N	2.20	0.74
1:CCC:129[A]:GLU:HG2	1:CCC:134:SER:CB	2.20	0.72
1:AAA:139:ALA:HB1	1:AAA:150[A]:MET:HE1	1.71	0.72
1:DDD:132:ASP:O	1:DDD:136:GLU:HG2	1.90	0.72
1:CCC:18[A]:ASN:OD1	3:CCC:403:GOL:C1	2.37	0.71
1:AAA:4[B]:GLN:CD	1:AAA:4[B]:GLN:N	2.49	0.70
1:AAA:39[B]:GLY:O	1:AAA:40[B]:ARG:HG3	1.92	0.70
1:BBB:67[B]:LEU:CD1	1:BBB:67[B]:LEU:N	2.55	0.70
1:AAA:325[A]:TRP:O	1:AAA:325[A]:TRP:HE3	1.74	0.69
1:DDD:32:SER:O	1:DDD:36[A]:GLN:HG3	1.92	0.69
1:CCC:217[B]:CYS:HB2	1:CCC:244[B]:ILE:HD11	1.75	0.68
1:BBB:281[B]:LEU:HD13	1:DDD:281[B]:LEU:HD21	1.77	0.66
1:BBB:107:PRO:HA	1:BBB:128:CYS:HB3	1.76	0.66
1:AAA:325[A]:TRP:O	1:AAA:325[A]:TRP:CE3	2.48	0.66
1:BBB:24:HIS:CE1	1:BBB:51:LYS:HE3	2.30	0.66
1:AAA:87:HIS:ND1	10:AAA:507:HOH:O	2.29	0.65
1:BBB:84:SER:HB3	1:BBB:118:ILE:HD11	1.78	0.65
1:AAA:69:PRO:O	10:AAA:503:HOH:O	2.14	0.65
1:AAA:234[B]:GLU:OE1	10:AAA:502:HOH:O	2.14	0.65
3:AAA:404:GOL:H31	10:AAA:639:HOH:O	1.96	0.65
1:DDD:223:LYS:HE3	10:DDD:506:HOH:O	1.96	0.64
1:CCC:236[B]:ILE:CD1	1:CCC:268:LYS:HG3	2.28	0.64
1:CCC:129[A]:GLU:HG2	1:CCC:134:SER:HB2	1.78	0.63
1:AAA:275:TRP:CH2	1:AAA:325[B]:TRP:CH2	2.87	0.62
1:AAA:149:ILE:CA	3:AAA:404:GOL:H32	2.29	0.62
1:AAA:232:PRO:HA	3:AAA:402:GOL:H32	1.82	0.62
1:CCC:236[B]:ILE:HD13	1:CCC:268:LYS:HG3	1.82	0.61
3:AAA:404:GOL:O3	10:AAA:505:HOH:O	2.16	0.61
1:BBB:67[B]:LEU:N	1:BBB:67[B]:LEU:HD12	2.15	0.61
1:AAA:140:LYS:CE	1:AAA:150[B]:MET:CE	2.55	0.60
1:DDD:105:VAL:HG22	1:DDD:126[B]:VAL:CG2	2.31	0.60
1:AAA:156[B]:GLU:OE1	10:AAA:504:HOH:O	2.16	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BBB:231[A]:TYR:H	3:BBB:402:GOL:HO1	1.48	0.60
10:AAA:501:HOH:O	1:DDD:13:GLU:OE2	2.17	0.59
1:BBB:231[B]:TYR:N	3:BBB:402:GOL:O1	2.27	0.59
1:CCC:29:LEU:HB3	8:CCC:402:PGE:H5	1.85	0.59
1:AAA:322[A]:SER:O	1:AAA:323[A]:ILE:C	2.45	0.58
1:BBB:105:VAL:HG12	1:BBB:128:CYS:HB2	1.86	0.58
1:AAA:32:SER:O	1:AAA:36[A]:GLN:HG3	2.05	0.57
1:AAA:320[B]:THR:O	1:AAA:323[B]:ILE:CB	2.52	0.57
1:AAA:14[B]:LYS:HG3	10:AAA:631:HOH:O	2.05	0.57
1:BBB:231[B]:TYR:H	3:BBB:402:GOL:HO1	1.52	0.56
1:AAA:56[B]:LYS:HD2	10:AAA:507:HOH:O	2.04	0.56
1:AAA:322[A]:SER:C	1:AAA:324[A]:THR:N	2.62	0.56
3:BBB:402:GOL:H11	10:BBB:621:HOH:O	2.04	0.56
1:DDD:104[A]:ILE:HD13	1:DDD:106:VAL:HG12	1.87	0.56
1:DDD:105:VAL:HG22	1:DDD:126[B]:VAL:HG21	1.87	0.56
1:BBB:281[B]:LEU:HG	1:BBB:317:VAL:HG11	1.88	0.55
1:BBB:32:SER:O	1:BBB:36:GLN:HG3	2.06	0.55
1:CCC:29:LEU:HD13	8:CCC:402:PGE:H5	1.88	0.55
1:AAA:139:ALA:HB1	1:AAA:150[A]:MET:CE	2.36	0.55
1:BBB:236[A]:ILE:HD12	1:BBB:284:PRO:HB3	1.88	0.55
1:AAA:13[B]:GLU:OE1	3:AAA:409:GOL:O3	2.24	0.54
1:CCC:24:HIS:NE2	1:CCC:51[B]:LYS:CE	2.67	0.54
1:DDD:201:LYS:HG2	1:DDD:204:VAL:HG23	1.88	0.54
1:DDD:259[A]:ILE:HG22	10:DDD:623:HOH:O	2.08	0.53
1:CCC:269[B]:CYS:SG	10:CCC:752:HOH:O	2.59	0.53
1:DDD:104[B]:ILE:HG22	1:DDD:106:VAL:HG12	1.89	0.53
1:DDD:105:VAL:HA	1:DDD:126[B]:VAL:HG23	1.91	0.53
1:CCC:70:ASP:OD2	10:CCC:502:HOH:O	2.19	0.53
1:AAA:120:ALA:HB3	10:AAA:580:HOH:O	2.09	0.52
1:AAA:140:LYS:CD	1:AAA:150[B]:MET:HE1	2.33	0.52
1:DDD:118:ILE:C	1:DDD:120:ALA:H	2.17	0.51
1:AAA:13[A]:GLU:OE2	1:DDD:13:GLU:HG2	2.11	0.51
1:AAA:136:GLU:O	10:AAA:506:HOH:O	2.19	0.51
1:BBB:233:PRO:O	10:BBB:501:HOH:O	2.19	0.51
1:AAA:139:ALA:HB3	10:AAA:506:HOH:O	2.10	0.51
10:BBB:584:HOH:O	1:DDD:24:HIS:HE1	1.94	0.51
1:BBB:281[A]:LEU:HD21	1:DDD:281[A]:LEU:HD13	1.92	0.50
1:AAA:272:GLN:NE2	1:AAA:325[B]:TRP:HE1	2.02	0.50
1:BBB:53:GLY:HA3	1:BBB:316:ASN:HD21	1.76	0.50
1:AAA:321[B]:SER:C	1:AAA:323[B]:ILE:N	2.67	0.50
1:BBB:107:PRO:CA	1:BBB:128:CYS:HB3	2.42	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:DDD:218:TYR:N	1:DDD:259[A]:ILE:HD13	2.27	0.50
1:DDD:259[A]:ILE:O	1:DDD:259[A]:ILE:HG23	2.12	0.49
1:AAA:218:TYR:CE1	3:AAA:403:GOL:H11	2.47	0.49
1:AAA:272:GLN:HE21	1:AAA:325[B]:TRP:NE1	2.03	0.49
1:AAA:281[B]:LEU:HD21	1:CCC:281[B]:LEU:HD13	1.95	0.49
1:AAA:39[B]:GLY:C	1:AAA:40[B]:ARG:HG3	2.37	0.48
1:AAA:39[B]:GLY:O	1:AAA:40[B]:ARG:CG	2.62	0.48
1:DDD:107:PRO:HA	1:DDD:128[B]:CYS:O	2.14	0.48
1:AAA:236:ILE:HD11	1:AAA:325[A]:TRP:CD1	2.49	0.47
1:CCC:226:LEU:HD11	1:CCC:244[A]:ILE:HG13	1.96	0.47
1:DDD:9:PHE:CE1	1:DDD:168[B]:LEU:HD21	2.49	0.47
1:AAA:268:LYS:HE2	1:AAA:325[B]:TRP:HB2	1.96	0.47
1:CCC:79[A]:VAL:CG1	1:CCC:91:LEU:HG	2.38	0.47
1:DDD:118:ILE:O	1:DDD:119:GLN:HB2	2.14	0.47
1:BBB:16:HIS:O	1:BBB:20:ARG:HB3	2.16	0.46
1:BBB:236[B]:ILE:HD11	1:BBB:325:TRP:CE2	2.50	0.46
1:CCC:38:THR:HG21	1:CCC:42:LEU:HD11	1.96	0.46
1:DDD:106:VAL:HG13	1:DDD:125:ILE:HG23	1.97	0.46
1:CCC:236[B]:ILE:HD11	1:CCC:268:LYS:HG3	1.97	0.46
1:BBB:13:GLU:HG3	1:CCC:10:ALA:HA	1.98	0.46
1:CCC:218:TYR:HA	1:CCC:259[B]:ILE:HG21	1.96	0.46
1:AAA:321[B]:SER:C	1:AAA:323[B]:ILE:H	2.22	0.46
1:AAA:150[B]:MET:HE2	1:AAA:150[B]:MET:HB2	1.60	0.46
1:CCC:33[B]:ILE:HG23	10:CCC:575:HOH:O	2.16	0.46
1:BBB:63:ALA:O	1:BBB:67[A]:LEU:HB2	2.15	0.45
1:BBB:64:VAL:HG12	1:BBB:98:GLU:HG3	1.98	0.45
1:BBB:128:CYS:SG	1:BBB:135:ARG:HA	2.57	0.45
1:BBB:210:GLU:O	1:BBB:261:THR:HA	2.16	0.45
1:AAA:275:TRP:CH2	1:AAA:325[B]:TRP:HH2	2.32	0.45
1:CCC:226:LEU:CD1	1:CCC:244[A]:ILE:HG13	2.47	0.45
1:DDD:51:LYS:HD2	1:DDD:93:TYR:OH	2.16	0.45
1:DDD:87:HIS:CG	1:DDD:153:PRO:HB3	2.51	0.44
1:BBB:106:VAL:O	1:BBB:107:PRO:C	2.60	0.44
1:CCC:236[B]:ILE:HD13	1:CCC:268:LYS:CG	2.47	0.44
1:CCC:56:LYS:NZ	2:CCC:401:PLP:O3	2.50	0.44
1:BBB:67[B]:LEU:H	1:BBB:67[B]:LEU:HD13	1.82	0.44
1:CCC:210[A]:GLU:HB2	10:CCC:518[A]:HOH:O	2.17	0.44
1:AAA:322[A]:SER:O	1:AAA:325[A]:TRP:N	2.51	0.44
1:CCC:151:VAL:O	7:CCC:406:EDO:H21	2.17	0.44
1:DDD:84:SER:HA	1:DDD:121:TYR:OH	2.19	0.43
1:CCC:217[B]:CYS:HA	1:CCC:244[B]:ILE:HD12	2.01	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CCC:217[B]:CYS:CB	1:CCC:244[B]:ILE:HD11	2.47	0.43
1:AAA:129[B]:GLU:HB2	1:AAA:134:SER:CB	2.49	0.43
1:AAA:322[A]:SER:C	1:AAA:324[A]:THR:H	2.26	0.43
1:DDD:126[B]:VAL:HG23	1:DDD:126[B]:VAL:O	2.18	0.43
1:AAA:150[A]:MET:H	3:AAA:404:GOL:C3	2.31	0.43
1:AAA:14[B]:LYS:HD2	10:AAA:798:HOH:O	2.18	0.43
1:AAA:168:LEU:HD22	3:AAA:409:GOL:H31	2.01	0.43
1:AAA:259[B]:ILE:O	1:AAA:259[B]:ILE:HG23	2.18	0.42
1:CCC:236[B]:ILE:HG21	1:CCC:268:LYS:HE3	2.01	0.42
1:AAA:150[B]:MET:H	3:AAA:404:GOL:C3	2.33	0.42
1:AAA:322[B]:SER:O	1:AAA:326[B]:VAL:N	2.49	0.42
1:AAA:230:LEU:HA	1:AAA:230:LEU:HD23	1.80	0.41
1:AAA:13[A]:GLU:O	1:AAA:17[A]:ILE:HG12	2.20	0.41
3:AAA:404:GOL:C3	10:AAA:639:HOH:O	2.64	0.41
1:BBB:9:PHE:HE1	1:BBB:168:LEU:HD21	1.84	0.41
1:CCC:217[B]:CYS:HB2	1:CCC:244[B]:ILE:CD1	2.47	0.41
1:BBB:231[B]:TYR:N	3:BBB:402:GOL:HO1	2.13	0.41
1:AAA:24:HIS:CE1	10:CCC:603:HOH:O	2.73	0.41
1:DDD:210:GLU:O	1:DDD:261:THR:HA	2.21	0.41
1:AAA:260:PHE:HA	3:AAA:403:GOL:O3	2.21	0.41
7:CCC:406:EDO:H12	10:CCC:727:HOH:O	2.20	0.41
1:CCC:163:GLN:HB3	1:CCC:189:MET:HA	2.03	0.40
1:AAA:14[B]:LYS:HE2	1:AAA:14[B]:LYS:HB2	1.78	0.40
1:CCC:210[A]:GLU:O	1:CCC:261:THR:HA	2.22	0.40
1:AAA:136:GLU:C	10:AAA:506:HOH:O	2.64	0.40

All (5) 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)
10:DDD:530:HOH:O	10:DDD:597:HOH:O[1_655]	1.82	0.38
10:AAA:693:HOH:O	10:AAA:764:HOH:O[1_455]	2.07	0.13
10:CCC:521:HOH:O	10:CCC:707:HOH:O[1_455]	2.07	0.13
1:AAA:129[B]:GLU:OE2	1:DDD:231:TYR:OH[1_554]	2.09	0.11
10:BBB:572:HOH:O	10:BBB:601:HOH:O[1_655]	2.16	0.04

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	AAA	355/346 (103%)	339 (96%)	16 (4%)	0	100	100
1	BBB	323/346 (93%)	314 (97%)	9 (3%)	0	100	100
1	CCC	343/346 (99%)	334 (97%)	9 (3%)	0	100	100
1	DDD	334/346 (96%)	326 (98%)	8 (2%)	0	100	100
All	All	1355/1384 (98%)	1313 (97%)	42 (3%)	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	AAA	293/291 (101%)	288 (98%)	5 (2%)	53	42
1	BBB	257/291 (88%)	249 (97%)	8 (3%)	35	20
1	CCC	279/291 (96%)	270 (97%)	9 (3%)	34	19
1	DDD	272/291 (94%)	262 (96%)	10 (4%)	30	15
All	All	1101/1164 (95%)	1069 (97%)	32 (3%)	38	22

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

Mol	Chain	Res	Type
1	AAA	121	TYR
1	AAA	136	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	AAA	140	LYS
1	AAA	141	ARG
1	AAA	225	LYS
1	BBB	16	HIS
1	BBB	25	LEU
1	BBB	32	SER
1	BBB	84	SER
1	BBB	106	VAL
1	BBB	108	GLN
1	BBB	131	SER
1	BBB	150	MET
1	CCC	34[A]	LEU
1	CCC	34[B]	LEU
1	CCC	73	GLU
1	CCC	96	LYS
1	CCC	133	GLU
1	CCC	136[A]	GLU
1	CCC	136[B]	GLU
1	CCC	163	GLN
1	CCC	168	LEU
1	DDD	4	GLN
1	DDD	104[A]	ILE
1	DDD	104[B]	ILE
1	DDD	114	LYS
1	DDD	124	SER
1	DDD	136	GLU
1	DDD	141[A]	ARG
1	DDD	141[B]	ARG
1	DDD	163	GLN
1	DDD	317	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 32 ligands modelled in this entry, 12 are monoatomic - leaving 20 for Mogul analysis.

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	GOL	AAA	409	-	5,5,5	0.14	0	5,5,5	0.30	0
2	PLP	AAA	401[B]	1	15,15,16	0.87	1 (6%)	21,22,23	0.81	0
3	GOL	AAA	408	-	5,5,5	0.18	0	5,5,5	0.41	0
7	EDO	CCC	406	-	3,3,3	0.80	0	2,2,2	0.28	0
8	PGE	CCC	402	-	9,9,9	0.17	0	8,8,8	0.25	0
6	JFS	AAA	407[B]	-	19,19,19	0.34	0	26,26,26	0.55	0
6	JFS	BBB	407	-	19,19,19	0.39	0	26,26,26	0.63	0
3	GOL	BBB	403	-	5,5,5	0.10	0	5,5,5	0.31	0
3	GOL	BBB	402	-	5,5,5	0.15	0	5,5,5	0.13	0
3	GOL	AAA	402	-	5,5,5	0.15	0	5,5,5	0.26	0
2	PLP	AAA	401[A]	1	15,15,16	0.87	1 (6%)	21,22,23	1.31	2 (9%)
3	GOL	AAA	403	-	5,5,5	0.21	0	5,5,5	0.64	0
3	GOL	AAA	404	-	5,5,5	0.17	0	5,5,5	0.53	0
2	PLP	DDD	401	1	15,15,16	0.79	1 (6%)	21,22,23	1.07	1 (4%)
2	PLP	BBB	401	1	15,15,16	0.79	1 (6%)	21,22,23	0.76	1 (4%)
6	JFS	AAA	407[A]	-	19,19,19	0.32	0	26,26,26	0.50	0
7	EDO	DDD	406	-	3,3,3	0.09	0	2,2,2	0.25	0
7	EDO	BBB	406	-	3,3,3	0.06	0	2,2,2	0.15	0
3	GOL	CCC	403	-	5,5,5	0.09	0	5,5,5	0.31	0
2	PLP	CCC	401	1	15,15,16	0.93	0	21,22,23	1.13	2 (9%)

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	GOL	AAA	409	-	-	0/4/4/4	-
2	PLP	AAA	401[B]	1	-	3/6/6/8	0/1/1/1
3	GOL	AAA	408	-	-	2/4/4/4	-
7	EDO	CCC	406	-	-	1/1/1/1	-
8	PGE	CCC	402	-	-	6/7/7/7	-
6	JFS	AAA	407[B]	-	-	0/6/6/6	0/3/3/3
6	JFS	BBB	407	-	-	0/6/6/6	0/3/3/3
3	GOL	BBB	403	-	-	2/4/4/4	-
3	GOL	BBB	402	-	-	2/4/4/4	-
3	GOL	AAA	402	-	-	2/4/4/4	-
2	PLP	AAA	401[A]	1	-	0/6/6/8	0/1/1/1
3	GOL	AAA	403	-	-	1/4/4/4	-
3	GOL	AAA	404	-	-	2/4/4/4	-
2	PLP	DDD	401	1	-	0/6/6/8	0/1/1/1
2	PLP	BBB	401	1	-	0/6/6/8	0/1/1/1
6	JFS	AAA	407[A]	-	-	0/6/6/6	0/3/3/3
7	EDO	DDD	406	-	-	1/1/1/1	-
7	EDO	BBB	406	-	-	0/1/1/1	-
3	GOL	CCC	403	-	-	0/4/4/4	-
2	PLP	CCC	401	1	-	0/6/6/8	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	AAA	401[B]	PLP	C4A-C4	-2.72	1.46	1.51
2	AAA	401[A]	PLP	C4A-C4	-2.64	1.46	1.51
2	DDD	401	PLP	C4A-C4	-2.60	1.46	1.51
2	BBB	401	PLP	C4A-C4	-2.32	1.47	1.51

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	AAA	401[A]	PLP	O4P-C5A-C5	4.55	117.89	109.36
2	DDD	401	PLP	O4P-C5A-C5	3.40	115.73	109.36
2	CCC	401	PLP	C4A-C4-C5	3.07	124.11	120.94
2	AAA	401[A]	PLP	C4A-C4-C5	2.32	123.33	120.94
2	CCC	401	PLP	O4P-C5A-C5	2.20	113.48	109.36
2	BBB	401	PLP	O4P-C5A-C5	2.02	113.14	109.36

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	AAA	401[B]	PLP	C4-C5-C5A-O4P
2	AAA	401[B]	PLP	C6-C5-C5A-O4P
3	AAA	404	GOL	C1-C2-C3-O3
3	AAA	404	GOL	O2-C2-C3-O3
3	AAA	408	GOL	O1-C1-C2-O2
3	AAA	408	GOL	O1-C1-C2-C3
8	CCC	402	PGE	O2-C3-C4-O3
3	AAA	402	GOL	C1-C2-C3-O3
3	BBB	403	GOL	C1-C2-C3-O3
8	CCC	402	PGE	O1-C1-C2-O2
3	BBB	403	GOL	O2-C2-C3-O3
8	CCC	402	PGE	C3-C4-O3-C5
2	AAA	401[B]	PLP	C5A-O4P-P-O1P
7	CCC	406	EDO	O1-C1-C2-O2
3	BBB	402	GOL	O1-C1-C2-C3
3	BBB	402	GOL	C1-C2-C3-O3
7	DDD	406	EDO	O1-C1-C2-O2
3	AAA	402	GOL	O2-C2-C3-O3
3	AAA	403	GOL	O2-C2-C3-O3
8	CCC	402	PGE	C6-C5-O3-C4
8	CCC	402	PGE	C4-C3-O2-C2
8	CCC	402	PGE	C1-C2-O2-C3

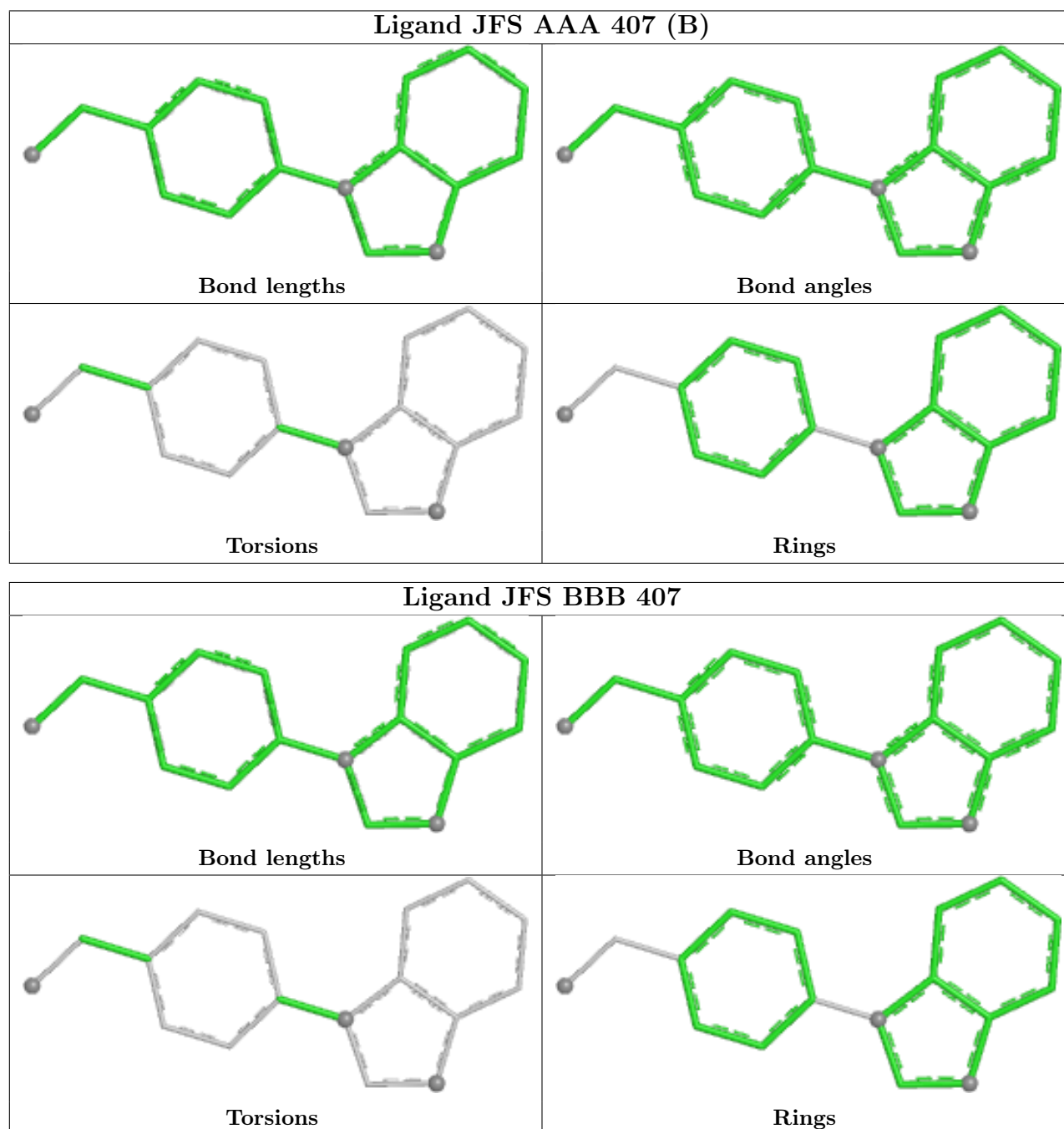
There are no ring outliers.

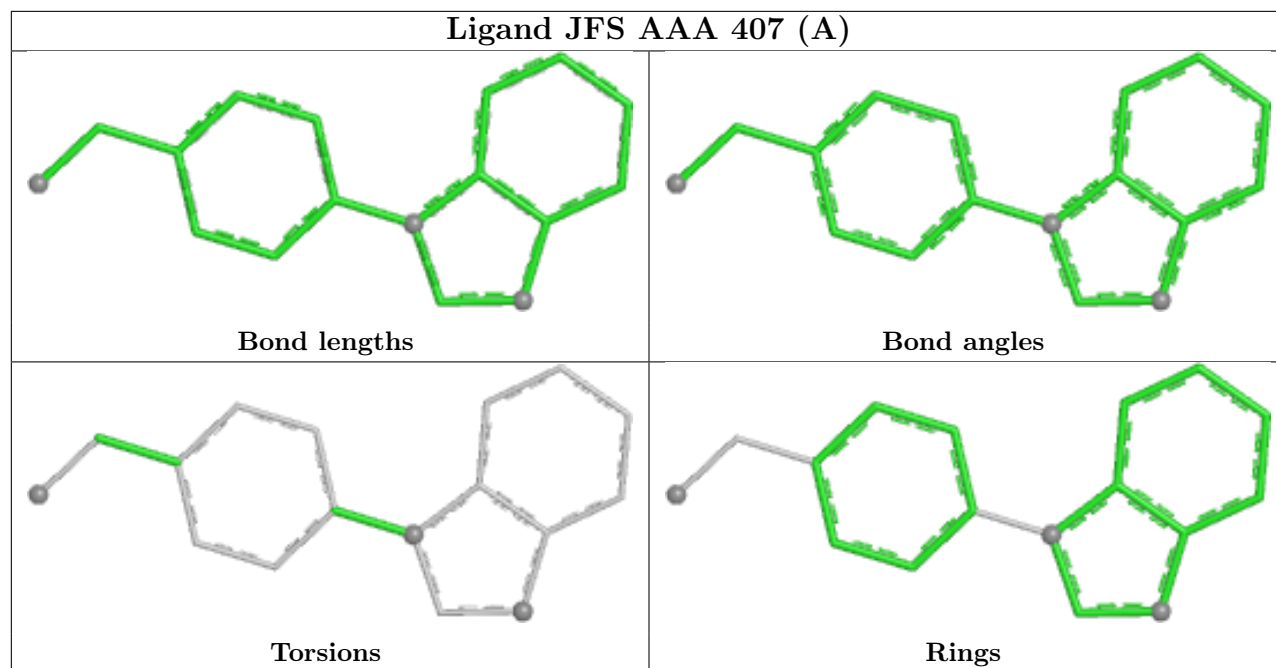
9 monomers are involved in 36 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	AAA	409	GOL	5	0
7	CCC	406	EDO	2	0
8	CCC	402	PGE	5	0
3	BBB	402	GOL	5	0
3	AAA	402	GOL	2	0
3	AAA	403	GOL	7	0
3	AAA	404	GOL	7	0
3	CCC	403	GOL	2	0
2	CCC	401	PLP	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	AAA	323/346 (93%)	-0.21	8 (2%) 58 62	8, 24, 55, 84	35 (10%)
1	BBB	315/346 (91%)	0.39	31 (9%) 13 13	13, 35, 97, 123	13 (4%)
1	CCC	323/346 (93%)	0.10	12 (3%) 45 49	12, 32, 68, 89	22 (6%)
1	DDD	318/346 (91%)	-0.04	15 (4%) 36 39	12, 27, 77, 108	20 (6%)
All	All	1279/1384 (92%)	0.06	66 (5%) 33 35	8, 30, 79, 123	90 (7%)

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	AAA	323[A]	ILE	7.2
1	BBB	105	VAL	6.1
1	BBB	106	VAL	4.8
1	BBB	125	ILE	4.1
1	BBB	67[A]	LEU	4.1
1	AAA	326[A]	VAL	3.9
1	BBB	326	VAL	3.8
1	CCC	325	TRP	3.5
1	BBB	107	PRO	3.5
1	BBB	323	ILE	3.4
1	BBB	120	ALA	3.3
1	AAA	320[A]	THR	3.3
1	BBB	109	THR	3.3
1	BBB	130	PRO	3.3
1	BBB	82	HIS	3.3
1	BBB	126	VAL	3.2
1	DDD	67	LEU	3.2
1	CCC	33[A]	ILE	3.2
1	CCC	111	PRO	3.1
1	BBB	230	LEU	3.0
1	DDD	323	ILE	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	BBB	319	LEU	2.9
1	CCC	113	CYS	2.9
1	BBB	104	ILE	2.9
1	DDD	319	LEU	2.8
1	CCC	110	ALA	2.8
1	DDD	347	VAL	2.8
1	DDD	117	ALA	2.7
1	BBB	123	ALA	2.7
1	CCC	231[A]	TYR	2.7
1	BBB	127	TYR	2.6
1	DDD	118	ILE	2.6
1	DDD	68	VAL	2.6
1	AAA	322[A]	SER	2.6
1	AAA	319[A]	LEU	2.6
1	BBB	100	ILE	2.6
1	BBB	320	THR	2.6
1	BBB	108	GLN	2.5
1	DDD	121	TYR	2.5
1	BBB	325	TRP	2.5
1	BBB	116	LEU	2.5
1	DDD	109	THR	2.5
1	AAA	321[A]	SER	2.4
1	CCC	321	SER	2.4
1	AAA	137	ASN	2.4
1	BBB	137	ASN	2.4
1	BBB	138	VAL	2.4
1	CCC	319	LEU	2.3
1	BBB	119	GLN	2.3
1	DDD	116	LEU	2.3
1	BBB	121	TYR	2.3
1	CCC	324	THR	2.3
1	DDD	111	PRO	2.3
1	BBB	128	CYS	2.3
1	DDD	113	CYS	2.2
1	AAA	324[A]	THR	2.2
1	BBB	4	GLN	2.1
1	DDD	325	TRP	2.1
1	CCC	299	GLN	2.1
1	BBB	118	ILE	2.1
1	CCC	323	ILE	2.1
1	BBB	83	SER	2.1
1	BBB	113	CYS	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	DDD	120	ALA	2.1
1	CCC	72	LEU	2.0
1	DDD	324	THR	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.

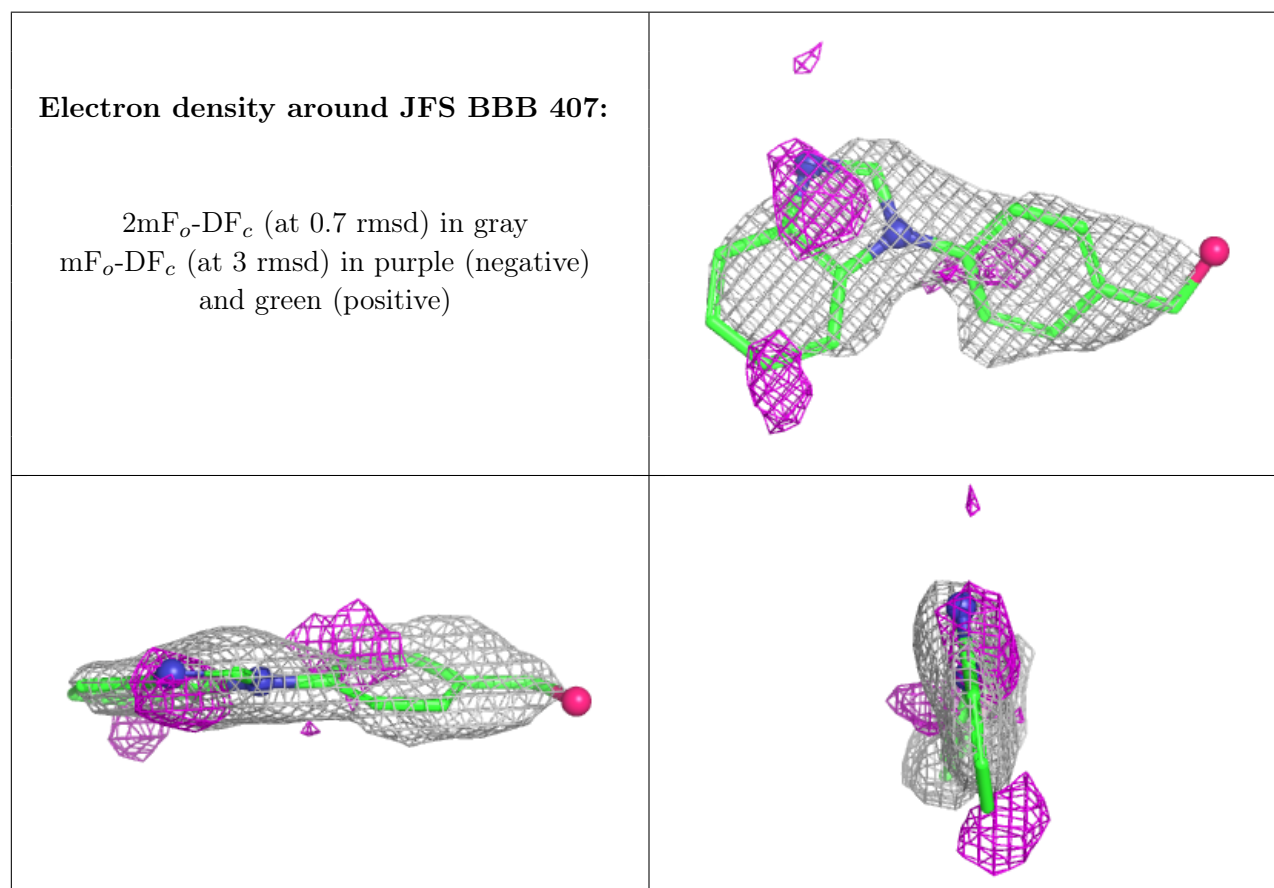
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	JFS	BBB	407	17/17	0.76	0.17	53,58,63,71	0
3	GOL	BBB	403	6/6	0.77	0.18	66,78,84,100	0
7	EDO	BBB	406	4/4	0.79	0.18	57,59,63,66	0
6	JFS	AAA	407[B]	17/17	0.81	0.17	47,50,57,62	17
3	GOL	AAA	404	6/6	0.81	0.18	38,48,55,62	0
6	JFS	AAA	407[A]	17/17	0.81	0.17	54,56,62,65	17
3	GOL	AAA	408	6/6	0.82	0.17	46,67,74,75	0
8	PGE	CCC	402	10/10	0.82	0.14	48,55,61,64	0
7	EDO	CCC	406	4/4	0.83	0.18	34,39,47,52	0
3	GOL	BBB	402	6/6	0.83	0.19	60,66,74,75	0
3	GOL	CCC	403	6/6	0.84	0.18	58,67,75,80	0
3	GOL	AAA	409	6/6	0.86	0.11	51,54,55,55	0
4	CA	DDD	404[A]	1/1	0.86	0.18	64,64,64,64	1
7	EDO	DDD	406	4/4	0.86	0.15	56,58,59,62	0
4	CA	DDD	404[B]	1/1	0.86	0.18	44,44,44,44	1
3	GOL	AAA	403	6/6	0.89	0.14	31,39,43,51	0
3	GOL	AAA	402	6/6	0.89	0.12	52,56,64,66	0
9	NA	DDD	405	1/1	0.89	0.10	54,54,54,54	0
4	CA	CCC	405	1/1	0.90	0.14	70,70,70,70	0
5	MG	DDD	403	1/1	0.96	0.05	45,45,45,45	0

Continued on next page...

Continued from previous page...

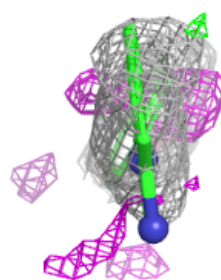
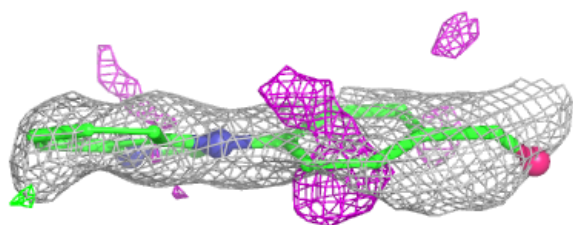
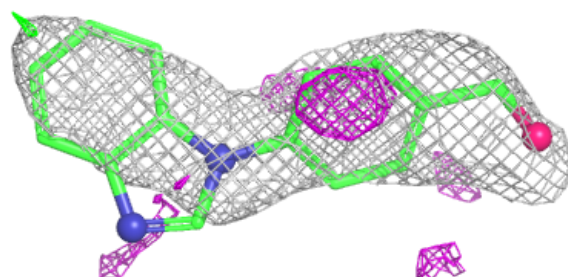
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PLP	AAA	401[B]	15/16	0.98	0.05	19,19,21,21	15
2	PLP	BBB	401	15/16	0.98	0.04	21,23,27,28	0
5	MG	AAA	406	1/1	0.98	0.03	31,31,31,31	0
2	PLP	CCC	401	15/16	0.98	0.05	21,23,24,26	0
4	CA	BBB	404	1/1	0.98	0.04	33,33,33,33	0
2	PLP	AAA	401[A]	15/16	0.98	0.05	16,17,19,20	15
4	CA	BBB	405	1/1	0.99	0.08	48,48,48,48	0
2	PLP	DDD	401	15/16	0.99	0.04	19,20,22,23	0
4	CA	DDD	402	1/1	0.99	0.03	25,25,25,25	0
4	CA	CCC	404[B]	1/1	1.00	0.02	29,29,29,29	1
4	CA	AAA	405	1/1	1.00	0.01	22,22,22,22	0
4	CA	CCC	404[A]	1/1	1.00	0.02	22,22,22,22	1

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.

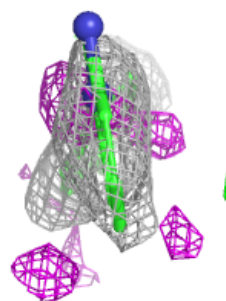
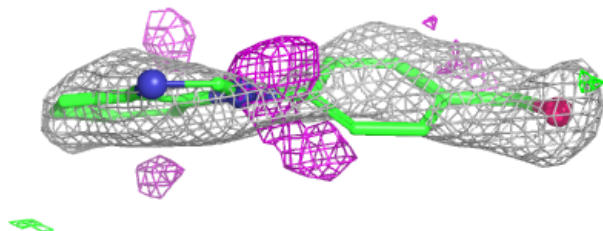
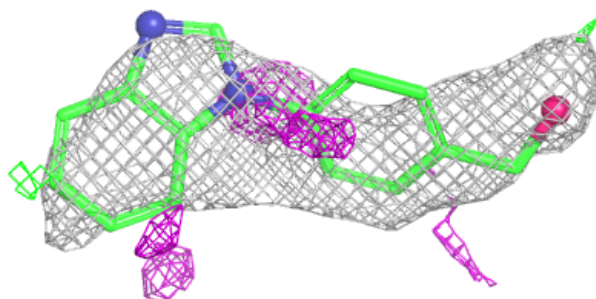


Electron density around JFS AAA 407 (B):

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

**Electron density around JFS AAA 407 (A):**

$2mF_o-DF_c$ (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.