



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 10:36 PM UTC

PDB ID : 1XET / pdb_00001xet
Title : Crystal structure of stilbene synthase from *Pinus sylvestris*, complexed with methylmalonyl CoA
Authors : Ng, S.H.; Chirgadze, D.; Spiteller, D.; Li, T.L.; Spencer, J.B.; Blundell, T.L.
Deposited on : 2004-09-12
Resolution : 2.00 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

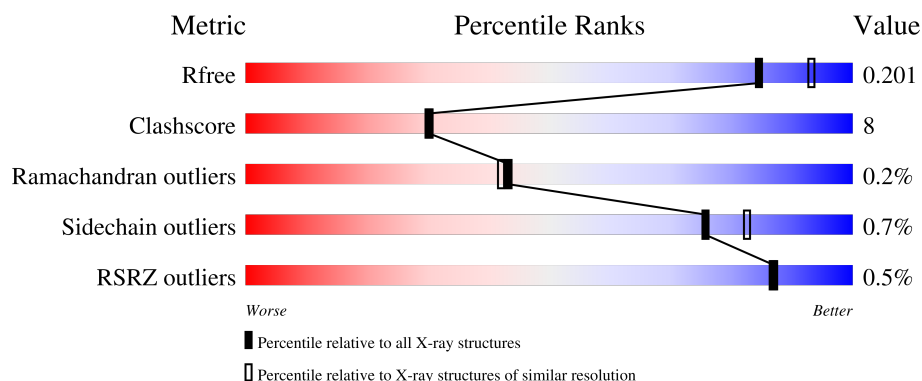
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	413	 79% 14% • 6%
1	B	413	 74% 19% 6%
1	C	413	 77% 15% • 8%
1	D	413	 73% 19% • 7%

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	3IO	D	4000	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12508 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 Dihydropinosylvin synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	387	Total	C	N	O	S	0	0	0
			2923	1857	503	546	17			
1	B	387	Total	C	N	O	S	0	0	0
			2921	1854	503	547	17			
1	C	382	Total	C	N	O	S	0	0	0
			2857	1815	494	531	17			
1	D	384	Total	C	N	O	S	4	0	0
			2860	1819	495	529	17			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	cloning artifact	UNP Q02323
A	-18	GLY	-	cloning artifact	UNP Q02323
A	-17	SER	-	cloning artifact	UNP Q02323
A	-16	SER	-	cloning artifact	UNP Q02323
A	-15	HIS	-	cloning artifact	UNP Q02323
A	-14	HIS	-	cloning artifact	UNP Q02323
A	-13	HIS	-	cloning artifact	UNP Q02323
A	-12	HIS	-	cloning artifact	UNP Q02323
A	-11	HIS	-	cloning artifact	UNP Q02323
A	-10	HIS	-	cloning artifact	UNP Q02323
A	-9	SER	-	cloning artifact	UNP Q02323
A	-8	SER	-	cloning artifact	UNP Q02323
A	-7	GLY	-	cloning artifact	UNP Q02323
A	-6	LEU	-	cloning artifact	UNP Q02323
A	-5	VAL	-	cloning artifact	UNP Q02323
A	-4	PRO	-	cloning artifact	UNP Q02323
A	-3	ARG	-	cloning artifact	UNP Q02323
A	-2	GLY	-	cloning artifact	UNP Q02323
A	-1	SER	-	cloning artifact	UNP Q02323
A	0	HIS	-	cloning artifact	UNP Q02323
B	-19	MET	-	cloning artifact	UNP Q02323

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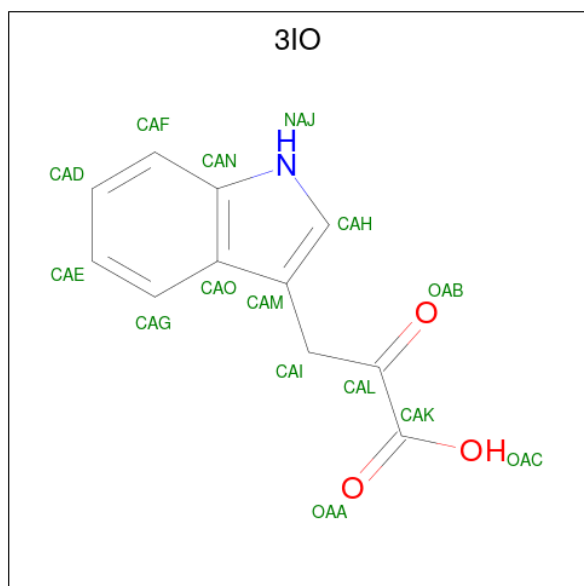
Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	cloning artifact	UNP Q02323
B	-17	SER	-	cloning artifact	UNP Q02323
B	-16	SER	-	cloning artifact	UNP Q02323
B	-15	HIS	-	cloning artifact	UNP Q02323
B	-14	HIS	-	cloning artifact	UNP Q02323
B	-13	HIS	-	cloning artifact	UNP Q02323
B	-12	HIS	-	cloning artifact	UNP Q02323
B	-11	HIS	-	cloning artifact	UNP Q02323
B	-10	HIS	-	cloning artifact	UNP Q02323
B	-9	SER	-	cloning artifact	UNP Q02323
B	-8	SER	-	cloning artifact	UNP Q02323
B	-7	GLY	-	cloning artifact	UNP Q02323
B	-6	LEU	-	cloning artifact	UNP Q02323
B	-5	VAL	-	cloning artifact	UNP Q02323
B	-4	PRO	-	cloning artifact	UNP Q02323
B	-3	ARG	-	cloning artifact	UNP Q02323
B	-2	GLY	-	cloning artifact	UNP Q02323
B	-1	SER	-	cloning artifact	UNP Q02323
B	0	HIS	-	cloning artifact	UNP Q02323
C	-19	MET	-	cloning artifact	UNP Q02323
C	-18	GLY	-	cloning artifact	UNP Q02323
C	-17	SER	-	cloning artifact	UNP Q02323
C	-16	SER	-	cloning artifact	UNP Q02323
C	-15	HIS	-	cloning artifact	UNP Q02323
C	-14	HIS	-	cloning artifact	UNP Q02323
C	-13	HIS	-	cloning artifact	UNP Q02323
C	-12	HIS	-	cloning artifact	UNP Q02323
C	-11	HIS	-	cloning artifact	UNP Q02323
C	-10	HIS	-	cloning artifact	UNP Q02323
C	-9	SER	-	cloning artifact	UNP Q02323
C	-8	SER	-	cloning artifact	UNP Q02323
C	-7	GLY	-	cloning artifact	UNP Q02323
C	-6	LEU	-	cloning artifact	UNP Q02323
C	-5	VAL	-	cloning artifact	UNP Q02323
C	-4	PRO	-	cloning artifact	UNP Q02323
C	-3	ARG	-	cloning artifact	UNP Q02323
C	-2	GLY	-	cloning artifact	UNP Q02323
C	-1	SER	-	cloning artifact	UNP Q02323
C	0	HIS	-	cloning artifact	UNP Q02323
D	-19	MET	-	cloning artifact	UNP Q02323
D	-18	GLY	-	cloning artifact	UNP Q02323
D	-17	SER	-	cloning artifact	UNP Q02323

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	cloning artifact	UNP Q02323
D	-15	HIS	-	cloning artifact	UNP Q02323
D	-14	HIS	-	cloning artifact	UNP Q02323
D	-13	HIS	-	cloning artifact	UNP Q02323
D	-12	HIS	-	cloning artifact	UNP Q02323
D	-11	HIS	-	cloning artifact	UNP Q02323
D	-10	HIS	-	cloning artifact	UNP Q02323
D	-9	SER	-	cloning artifact	UNP Q02323
D	-8	SER	-	cloning artifact	UNP Q02323
D	-7	GLY	-	cloning artifact	UNP Q02323
D	-6	LEU	-	cloning artifact	UNP Q02323
D	-5	VAL	-	cloning artifact	UNP Q02323
D	-4	PRO	-	cloning artifact	UNP Q02323
D	-3	ARG	-	cloning artifact	UNP Q02323
D	-2	GLY	-	cloning artifact	UNP Q02323
D	-1	SER	-	cloning artifact	UNP Q02323
D	0	HIS	-	cloning artifact	UNP Q02323

- Molecule 2 is 3-(1H-INDOL-3-YL)-2-OXOPROPANOIC ACID (CCD ID: 3IO) (formula: C₁₁H₉NO₃).



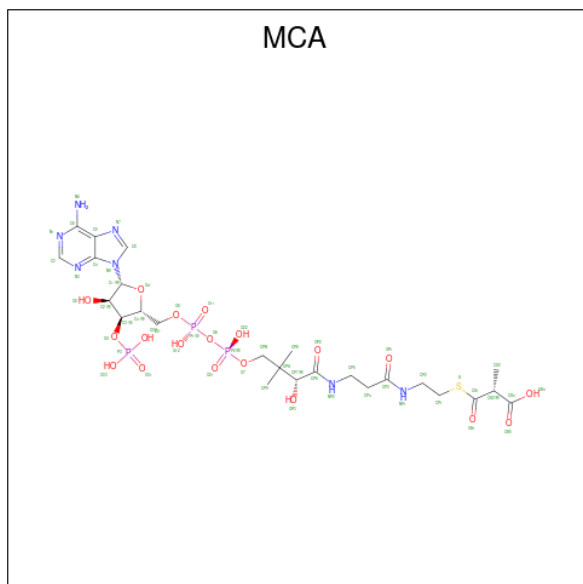
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			15	11	1	3		
2	C	1	Total	C	N	O	0	0
			15	11	1	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	D	1	Total	C	N	O	0	0
			15	11	1	3		

- Molecule 3 is METHYLMALONYL-COENZYME A (CCD ID: MCA) (formula: $C_{25}H_{40}N_7O_{19}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	D	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		

- Molecule 4 is water.

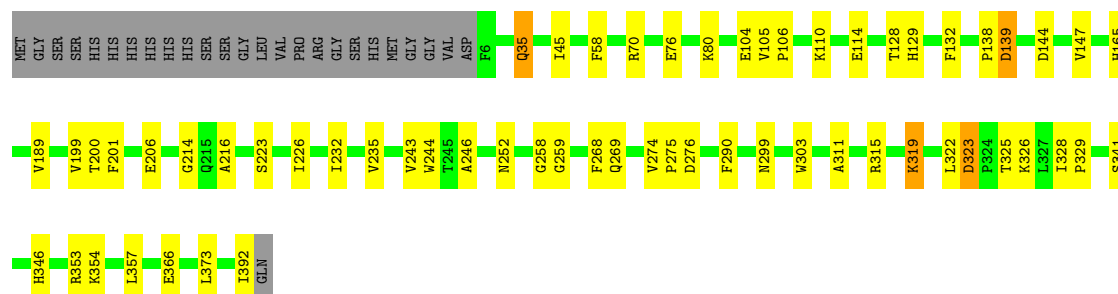
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	204	Total	O	0	0
			204	204		
4	B	214	Total	O	0	0
			214	214		
4	C	241	Total	O	0	0
			241	241		
4	D	195	Total	O	0	0
			195	195		

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.

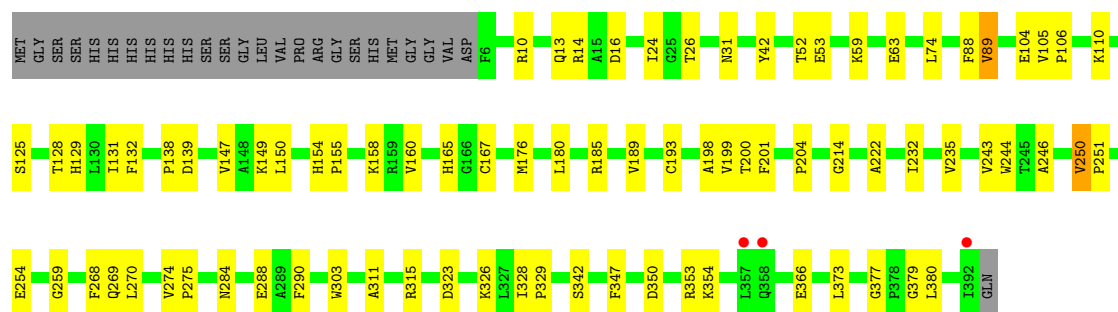
• Molecule 1: Dihydropinosylvin synthase

Chain A: 



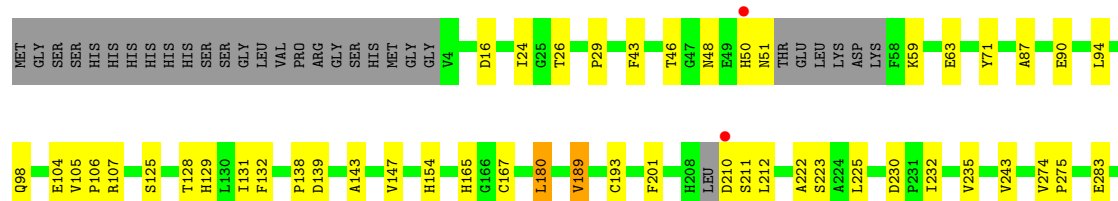
• Molecule 1: Dihydropinosylvin synthase

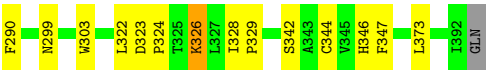
Chain B: 



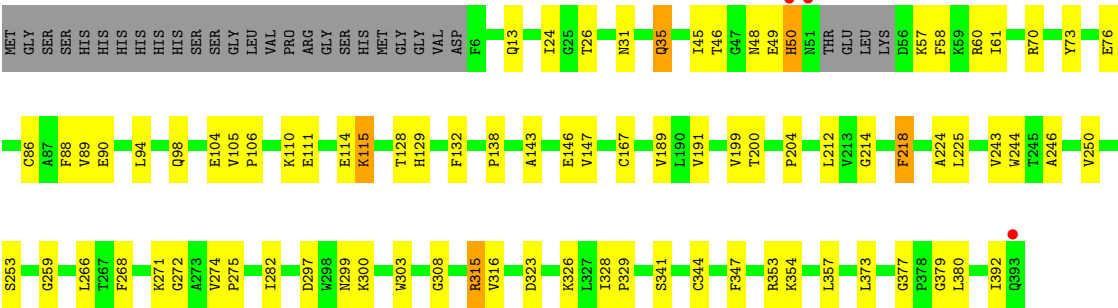
• Molecule 1: Dihydropinosylvin synthase

Chain C: 





● Molecule 1: Dihydropinosylvin synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.47Å 111.19Å 132.01Å 90.00° 93.13° 90.00°	Depositor
Resolution (Å)	22.58 – 2.00 22.58 – 2.00	Depositor EDS
% Data completeness (in resolution range)	91.9 (22.58-2.00) 91.2 (22.58-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.16 (at 1.99Å)	Xtriage
Refinement program	CNS 1.1, REFMAC 5.1.24	Depositor
R, R_{free}	0.195 , 0.223 0.197 , 0.201	Depositor DCC
R_{free} test set	4809 reflections (4.54%)	wwPDB-VP
Wilson B-factor (Å ²)	25.9	Xtriage
Anisotropy	0.028	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12508	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.55% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 3IO, MCA

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	0.39	0/2983	0.91	12/4050 (0.3%)
1	B	0.38	0/2980	0.95	14/4045 (0.3%)
1	C	0.41	0/2914	0.90	10/3957 (0.3%)
1	D	0.39	0/2919	0.93	14/3967 (0.4%)
All	All	0.39	0/11796	0.92	50/16019 (0.3%)

There are no bond length outliers.

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	243	VAL	N-CA-C	9.96	120.48	110.82
1	A	243	VAL	N-CA-C	8.52	118.50	110.74
1	D	243	VAL	N-CA-C	8.08	118.66	110.82
1	C	243	VAL	N-CA-C	8.00	118.58	110.82
1	B	104	GLU	N-CA-C	7.80	119.86	111.36
1	D	104	GLU	N-CA-C	7.78	119.75	111.28
1	C	104	GLU	N-CA-C	7.12	119.12	111.36
1	A	104	GLU	N-CA-C	7.10	119.10	111.36
1	B	16	ASP	N-CA-C	6.81	119.70	111.40
1	C	290	PHE	N-CA-C	6.35	121.19	113.38
1	B	165	HIS	N-CA-C	-6.32	104.28	112.68
1	D	303	TRP	N-CA-C	6.29	119.94	109.95
1	C	303	TRP	N-CA-C	6.27	119.75	109.72
1	A	138	PRO	N-CA-C	6.21	121.70	113.57
1	B	138	PRO	N-CA-C	6.16	121.64	113.57
1	B	303	TRP	N-CA-C	6.10	119.47	109.72
1	B	250	VAL	CA-C-N	6.03	126.05	119.90
1	B	250	VAL	C-N-CA	6.03	126.05	119.90
1	B	326	LYS	N-CA-C	5.99	118.30	111.11
1	A	303	TRP	N-CA-C	5.92	119.19	109.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	138	PRO	N-CA-C	5.90	121.31	113.40
1	A	189	VAL	N-CA-C	5.89	116.59	107.99
1	D	138	PRO	N-CA-C	5.75	121.11	113.57
1	C	324	PRO	N-CA-CB	5.72	109.26	103.25
1	C	342	SER	N-CA-C	5.59	118.14	111.71
1	A	226	ILE	N-CA-C	-5.57	99.63	107.75
1	A	290	PHE	N-CA-C	5.50	120.15	113.38
1	C	189	VAL	N-CA-C	5.46	115.97	107.99
1	A	139	ASP	N-CA-C	5.36	115.47	108.07
1	C	165	HIS	N-CA-C	-5.32	104.37	112.04
1	B	189	VAL	N-CA-C	5.29	115.72	107.99
1	B	342	SER	N-CA-C	5.29	117.73	111.33
1	A	165	HIS	N-CA-C	-5.28	104.44	112.04
1	D	218	PHE	N-CA-C	5.27	117.87	109.96
1	C	326	LYS	N-CA-C	5.22	116.65	111.07
1	B	377	GLY	CA-C-N	5.19	125.49	119.93
1	B	377	GLY	C-N-CA	5.19	125.49	119.93
1	A	341	SER	CB-CA-C	-5.19	110.57	116.54
1	D	189	VAL	N-CA-C	5.16	115.52	107.99
1	D	344	CYS	N-CA-C	5.13	117.27	111.02
1	D	73	TYR	N-CA-C	-5.12	105.78	111.36
1	D	323	ASP	CA-C-N	5.09	126.20	119.84
1	D	323	ASP	C-N-CA	5.09	126.20	119.84
1	D	341	SER	CB-CA-C	-5.08	110.73	116.63
1	D	31	ASN	N-CA-C	5.06	116.72	108.63
1	B	290	PHE	N-CA-C	5.05	120.29	113.72
1	D	377	GLY	CA-C-N	5.05	125.33	119.93
1	D	377	GLY	C-N-CA	5.05	125.33	119.93
1	A	323	ASP	CA-C-N	5.01	126.10	119.84
1	A	323	ASP	C-N-CA	5.01	126.10	119.84

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	2923	0	2897	41	0
1	B	2921	0	2892	45	0
1	C	2857	0	2803	41	0
1	D	2860	0	2807	66	0
2	B	15	0	8	2	0
2	C	15	0	8	3	0
2	D	15	0	8	8	0
3	D	48	0	31	5	0
4	A	204	0	0	6	0
4	B	214	0	0	3	0
4	C	241	0	0	7	0
4	D	195	0	0	2	0
All	All	12508	0	11454	193	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 8.

All (193) 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:D:297:ASP:HB3	1:D:300:LYS:HE2	1.49	0.95
1:C:167:CYS:SG	2:C:3000:3IO:CAL	2.58	0.92
1:B:10:ARG:HE	1:B:13:GLN:HE21	1.22	0.87
1:B:167:CYS:SG	2:B:2000:3IO:CAL	2.67	0.83
4:A:592:HOH:O	1:B:149:LYS:HE2	1.80	0.81
1:D:35:GLN:HA	1:D:35:GLN:HE21	1.48	0.79
1:A:35:GLN:HE21	1:A:35:GLN:HA	1.51	0.74
1:D:167:CYS:SG	2:D:4000:3IO:CAL	2.76	0.74
1:D:297:ASP:HB3	1:D:300:LYS:CE	2.19	0.72
1:D:46:THR:HG21	1:D:204:PRO:HD2	1.70	0.72
1:C:210:ASP:O	1:C:212:LEU:N	2.22	0.71
1:D:357:LEU:HD23	1:D:392:ILE:HD13	1.72	0.70
1:A:357:LEU:HD23	1:A:392:ILE:HD13	1.74	0.70
1:B:105:VAL:HB	1:B:106:PRO:HD3	1.74	0.70
1:D:266:LEU:HD21	2:D:4000:3IO:HAD	1.74	0.68
1:D:49:GLU:O	1:D:50:HIS:HB2	1.93	0.67
1:D:132:PHE:CE2	1:D:147:VAL:HG21	2.29	0.67
1:D:250:VAL:HG23	1:D:253:SER:OG	1.94	0.67
1:D:46:THR:HG22	1:D:48:ASN:ND2	2.10	0.66
1:A:128:THR:OG1	1:A:129:HIS:HD2	1.78	0.66
1:A:366:GLU:HG3	4:A:544:HOH:O	1.95	0.66
1:B:366:GLU:HG2	4:B:2168:HOH:O	1.93	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:GLN:HE22	1:A:70:ARG:HH21	1.43	0.66
1:B:10:ARG:HE	1:B:13:GLN:NE2	1.91	0.66
1:C:167:CYS:HG	2:C:3000:3IO:CAL	2.09	0.64
1:B:128:THR:OG1	1:B:129:HIS:HD2	1.81	0.63
1:D:105:VAL:HB	1:D:106:PRO:HD3	1.80	0.62
2:D:4000:3IO:CAK	3:D:4001:MCA:S	2.88	0.62
1:B:131:ILE:HD11	1:B:180:LEU:HD22	1.81	0.61
1:D:46:THR:HG22	1:D:48:ASN:HD22	1.65	0.60
1:D:250:VAL:CG2	1:D:380:LEU:HB3	2.31	0.60
1:D:214:GLY:HA3	1:D:268:PHE:CZ	2.37	0.60
1:C:180:LEU:HD11	4:C:3179:HOH:O	2.01	0.60
1:C:105:VAL:HB	1:C:106:PRO:HD3	1.84	0.60
1:D:167:CYS:SG	2:D:4000:3IO:CAK	2.90	0.59
1:A:299:ASN:HD21	1:A:323:ASP:H	1.49	0.59
1:C:328:ILE:HB	1:C:329:PRO:HD3	1.85	0.59
1:B:10:ARG:NE	1:B:13:GLN:HE21	1.97	0.59
1:D:46:THR:HG23	1:D:86:CYS:O	2.03	0.58
1:C:299:ASN:HD21	1:C:322:LEU:HA	1.68	0.58
1:D:49:GLU:O	1:D:50:HIS:CB	2.50	0.58
1:D:61:ILE:HD12	1:D:212:LEU:HD12	1.85	0.58
1:D:35:GLN:HE22	1:D:70:ARG:HH11	1.52	0.57
1:A:105:VAL:HB	1:A:106:PRO:HD3	1.87	0.56
1:D:218:PHE:CE1	2:D:4000:3IO:CAH	2.89	0.56
1:A:110:LYS:HE2	1:A:114:GLU:OE2	2.05	0.56
1:C:299:ASN:HD21	1:C:323:ASP:H	1.53	0.56
1:B:42:TYR:CZ	1:B:204:PRO:HD3	2.42	0.55
1:D:299:ASN:HA	1:D:326:LYS:HE3	1.87	0.55
1:D:282:ILE:HD13	1:D:316:VAL:HG13	1.89	0.55
1:A:328:ILE:HB	1:A:329:PRO:HD3	1.87	0.55
1:D:46:THR:HG22	1:D:46:THR:O	2.05	0.55
4:A:535:HOH:O	1:B:158:LYS:HE2	2.06	0.55
1:D:297:ASP:CB	1:D:300:LYS:HE2	2.31	0.55
1:D:24:ILE:HG12	1:D:225:LEU:HD22	1.87	0.55
1:D:274:VAL:HB	1:D:275:PRO:HD3	1.88	0.55
1:A:110:LYS:O	1:A:114:GLU:HG3	2.07	0.54
1:C:132:PHE:CE2	1:C:147:VAL:HG21	2.42	0.54
1:D:128:THR:OG1	1:D:129:HIS:HD2	1.90	0.54
1:A:259:GLY:O	1:B:139:ASP:HB3	2.08	0.53
1:A:199:VAL:HG13	1:A:200:THR:HG23	1.90	0.53
1:D:143:ALA:O	1:D:147:VAL:HG23	2.08	0.53
1:B:110:LYS:HD2	1:B:150:LEU:HB3	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:167:CYS:SG	2:C:3000:3IO:CAK	2.97	0.53
1:A:274:VAL:HB	1:A:275:PRO:HD3	1.91	0.53
3:D:4001:MCA:OP2	3:D:4001:MCA:CP4	2.54	0.52
1:C:43:PHE:HA	1:C:46:THR:HG22	1.91	0.52
1:A:319:LYS:HB2	1:A:319:LYS:NZ	2.24	0.52
1:D:167:CYS:SG	2:D:4000:3IO:CAH	2.98	0.52
1:B:125:SER:HB2	1:B:154:HIS:NE2	2.25	0.52
1:C:299:ASN:HA	1:C:326:LYS:HE3	1.90	0.52
1:B:269:GLN:O	1:B:270:LEU:HD12	2.10	0.51
3:D:4001:MCA:OP2	3:D:4001:MCA:H91	2.11	0.51
1:C:201:PHE:C	1:C:201:PHE:CD1	2.89	0.51
1:D:373:LEU:C	1:D:373:LEU:HD23	2.36	0.50
1:D:94:LEU:O	1:D:98:GLN:HG2	2.12	0.50
1:C:143:ALA:O	1:C:147:VAL:HG23	2.11	0.50
1:C:223:SER:OG	1:C:346:HIS:HD2	1.94	0.50
1:C:26:THR:HB	1:C:347:PHE:CZ	2.47	0.50
1:C:180:LEU:HD21	4:C:3179:HOH:O	2.11	0.50
1:D:218:PHE:CE1	2:D:4000:3IO:HAH	2.47	0.50
1:B:59:LYS:O	1:B:63:GLU:HG3	2.12	0.49
1:A:139:ASP:HB3	1:B:259:GLY:O	2.12	0.49
1:C:46:THR:HG23	1:C:48:ASN:HD22	1.77	0.49
1:A:214:GLY:HA3	1:A:268:PHE:CZ	2.47	0.49
1:C:373:LEU:C	1:C:373:LEU:HD23	2.37	0.49
1:B:24:ILE:HG21	1:B:350:ASP:HB2	1.95	0.49
1:B:328:ILE:HB	1:B:329:PRO:HD3	1.95	0.49
1:C:232:ILE:HB	1:C:235:VAL:HB	1.94	0.49
1:C:299:ASN:ND2	1:C:323:ASP:H	2.10	0.49
1:D:49:GLU:O	1:D:50:HIS:ND1	2.45	0.49
1:D:250:VAL:HG22	1:D:380:LEU:HB3	1.93	0.48
1:B:353:ARG:HG3	1:B:354:LYS:N	2.29	0.48
1:C:129:HIS:HE1	4:C:3160:HOH:O	1.95	0.48
1:A:373:LEU:HD23	1:A:373:LEU:C	2.39	0.48
1:C:125:SER:HB2	1:C:154:HIS:NE2	2.28	0.48
1:D:88:PHE:CD2	1:D:89:VAL:HG22	2.49	0.48
1:B:74:LEU:HD11	1:B:198:ALA:HA	1.96	0.48
1:B:88:PHE:CD2	1:B:89:VAL:HG22	2.48	0.48
1:B:132:PHE:CE2	1:B:147:VAL:HG21	2.48	0.48
1:B:155:PRO:HD2	4:B:2161:HOH:O	2.14	0.47
1:D:90:GLU:CD	1:D:90:GLU:H	2.21	0.47
1:A:276:ASP:OD1	1:A:315:ARG:NH1	2.42	0.47
1:A:315:ARG:HD2	4:A:529:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:308:GLY:C	3:D:4001:MCA:HN1	2.22	0.47
1:B:14:ARG:O	1:B:185:ARG:HB2	2.13	0.47
1:D:191:VAL:O	1:D:224:ALA:HA	2.14	0.47
1:D:297:ASP:HB3	1:D:300:LYS:CD	2.43	0.47
1:A:201:PHE:C	1:A:201:PHE:CD1	2.92	0.47
1:B:373:LEU:C	1:B:373:LEU:HD23	2.40	0.47
1:A:223:SER:OG	1:A:346:HIS:HD2	1.98	0.47
1:A:319:LYS:HB2	1:A:319:LYS:HZ3	1.80	0.46
1:D:244:TRP:CH2	1:D:246:ALA:HB2	2.50	0.46
1:B:269:GLN:C	1:B:270:LEU:HD12	2.40	0.46
1:C:87:ALA:HB1	1:C:90:GLU:OE1	2.15	0.46
1:A:214:GLY:HA3	1:A:268:PHE:CE2	2.50	0.46
1:A:353:ARG:HG3	1:A:354:LYS:N	2.30	0.46
1:B:167:CYS:SG	2:B:2000:3IO:CAK	3.04	0.46
1:D:45:ILE:HD13	1:D:76:GLU:HG3	1.97	0.46
1:A:45:ILE:HD13	1:A:76:GLU:HG3	1.97	0.46
1:C:274:VAL:HB	1:C:275:PRO:HD3	1.96	0.46
1:D:353:ARG:HG3	1:D:354:LYS:N	2.31	0.46
1:B:284:ASN:O	1:B:288:GLU:HG3	2.16	0.45
1:D:46:THR:CG2	1:D:48:ASN:HD22	2.26	0.45
1:B:274:VAL:HB	1:B:275:PRO:HD3	1.98	0.45
1:D:13:GLN:HG2	4:D:4140:HOH:O	2.16	0.45
2:D:4000:3IO:CAL	3:D:4001:MCA:S	3.05	0.45
1:C:46:THR:CG2	1:C:48:ASN:HD22	2.30	0.45
1:A:311:ALA:O	1:A:315:ARG:HG2	2.17	0.45
1:D:271:LYS:O	1:D:274:VAL:HG23	2.17	0.45
1:D:110:LYS:O	1:D:114:GLU:HG2	2.17	0.45
1:B:199:VAL:HG13	1:B:200:THR:HG23	1.97	0.44
1:D:272:GLY:O	1:D:275:PRO:HD2	2.17	0.44
1:B:193:CYS:O	1:B:222:ALA:HA	2.16	0.44
1:C:139:ASP:HB3	1:D:259:GLY:O	2.17	0.44
1:C:128:THR:OG1	1:C:129:HIS:HD2	2.01	0.44
1:D:24:ILE:HG12	1:D:225:LEU:CD2	2.48	0.44
1:A:325:THR:HA	1:A:328:ILE:HG12	1.99	0.44
1:B:160:VAL:HG11	1:B:176:MET:HE1	1.99	0.44
1:D:328:ILE:HB	1:D:329:PRO:HD3	1.99	0.44
1:C:193:CYS:O	1:C:222:ALA:HA	2.17	0.43
1:D:24:ILE:HG23	1:D:225:LEU:HD23	2.00	0.43
1:D:111:GLU:O	1:D:115:LYS:HD2	2.18	0.43
1:A:58:PHE:CE1	1:A:216:ALA:HB2	2.54	0.43
1:B:128:THR:OG1	1:B:129:HIS:CD2	2.68	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:214:GLY:HA3	1:B:268:PHE:CE1	2.53	0.43
1:A:80:LYS:HE3	1:A:80:LYS:HB2	1.85	0.43
1:A:132:PHE:CE2	1:A:147:VAL:HG21	2.54	0.43
1:D:212:LEU:C	1:D:212:LEU:HD13	2.44	0.43
1:A:35:GLN:HE22	1:A:70:ARG:NH2	2.12	0.42
1:A:232:ILE:HB	1:A:235:VAL:HB	2.02	0.42
1:D:58:PHE:HD2	1:D:212:LEU:HD11	1.84	0.42
1:B:201:PHE:CD1	1:B:201:PHE:C	2.97	0.42
1:D:214:GLY:HA3	1:D:268:PHE:CE1	2.55	0.42
1:D:379:GLY:N	1:D:380:LEU:HA	2.35	0.42
1:A:132:PHE:HB3	1:A:144:ASP:HB3	2.01	0.42
1:A:244:TRP:CH2	1:A:246:ALA:HB2	2.54	0.42
1:D:61:ILE:HD12	1:D:212:LEU:CD1	2.49	0.42
1:D:90:GLU:CD	1:D:90:GLU:N	2.77	0.42
1:D:315:ARG:HH11	1:D:315:ARG:HG3	1.84	0.42
1:B:379:GLY:N	1:B:380:LEU:HA	2.35	0.42
1:D:58:PHE:HD2	1:D:212:LEU:CD1	2.33	0.42
1:C:16:ASP:HB2	1:C:230:ASP:OD2	2.20	0.42
1:D:26:THR:HB	1:D:347:PHE:CZ	2.55	0.42
1:A:299:ASN:HD21	1:A:322:LEU:HA	1.85	0.41
1:B:244:TRP:CH2	1:B:246:ALA:HB2	2.54	0.41
1:D:199:VAL:HG13	1:D:200:THR:HG23	2.03	0.41
1:B:31:ASN:HB3	4:B:2129:HOH:O	2.21	0.41
4:C:3139:HOH:O	1:D:129:HIS:HE1	2.02	0.41
1:A:299:ASN:ND2	1:A:323:ASP:H	2.17	0.41
1:A:258:GLY:C	1:A:269:GLN:NE2	2.79	0.41
1:B:250:VAL:HA	1:B:251:PRO:HD3	1.92	0.41
1:B:323:ASP:C	1:B:323:ASP:OD2	2.63	0.41
1:C:131:ILE:HD11	1:C:180:LEU:HD22	2.01	0.41
1:A:206:GLU:HG2	4:A:546:HOH:O	2.20	0.41
1:A:252:ASN:OD1	1:D:50:HIS:HB3	2.21	0.41
1:C:50:HIS:HB3	1:C:51:ASN:H	1.30	0.41
1:C:283:GLU:HB3	4:C:3163:HOH:O	2.21	0.41
1:B:311:ALA:O	1:B:315:ARG:HG2	2.20	0.41
1:B:52:THR:HG23	1:B:53:GLU:N	2.35	0.41
1:C:24:ILE:HG12	1:C:225:LEU:HD22	2.02	0.41
1:B:232:ILE:HB	1:B:235:VAL:HB	2.03	0.41
1:C:344:CYS:HB2	4:C:3001:HOH:O	2.20	0.41
1:D:146:GLU:HB2	4:D:4108:HOH:O	2.21	0.41
1:C:29:PRO:HD2	1:C:71:TYR:HB3	2.03	0.40
1:C:299:ASN:HD21	1:C:323:ASP:N	2.19	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:LYS:NZ	4:A:581:HOH:O	2.53	0.40
1:B:26:THR:HB	1:B:347:PHE:CZ	2.56	0.40
1:C:59:LYS:O	1:C:63:GLU:HG3	2.22	0.40
1:C:94:LEU:O	1:C:98:GLN:HG2	2.21	0.40
1:C:107:ARG:HD2	4:C:3217:HOH:O	2.21	0.40
1:C:129:HIS:HB2	1:C:189:VAL:HG22	2.03	0.40
1:A:323:ASP:O	1:A:326:LYS:HG2	2.21	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	385/413 (93%)	370 (96%)	15 (4%)	0	100	100
1	B	385/413 (93%)	372 (97%)	12 (3%)	1 (0%)	36	35
1	C	376/413 (91%)	362 (96%)	13 (4%)	1 (0%)	36	35
1	D	380/413 (92%)	366 (96%)	13 (3%)	1 (0%)	36	35
All	All	1526/1652 (92%)	1470 (96%)	53 (4%)	3 (0%)	43	42

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	211	SER
1	D	50	HIS
1	B	89	VAL

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	310/340 (91%)	308 (99%)	2 (1%)	78	85
1	B	309/340 (91%)	308 (100%)	1 (0%)	86	91
1	C	298/340 (88%)	297 (100%)	1 (0%)	86	91
1	D	297/340 (87%)	292 (98%)	5 (2%)	53	60
All	All	1214/1360 (89%)	1205 (99%)	9 (1%)	76	82

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	GLN
1	A	319	LYS
1	B	254	GLU
1	C	180	LEU
1	D	35	GLN
1	D	57	LYS
1	D	60	ARG
1	D	115	LYS
1	D	315	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	GLN
1	A	69	GLN
1	A	118	GLN
1	A	129	HIS
1	A	269	GLN
1	A	281	ASN
1	A	299	ASN
1	A	346	HIS
1	A	358	GLN
1	B	13	GLN
1	B	51	ASN
1	B	69	GLN
1	B	129	HIS
1	B	208	HIS

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Mol	Chain	Res	Type
1	B	215	GLN
1	B	284	ASN
1	B	346	HIS
1	B	359	ASN
1	C	31	ASN
1	C	48	ASN
1	C	69	GLN
1	C	129	HIS
1	C	208	HIS
1	C	281	ASN
1	C	299	ASN
1	C	346	HIS
1	D	31	ASN
1	D	35	GLN
1	D	48	ASN
1	D	129	HIS
1	D	154	HIS
1	D	234	GLN
1	D	284	ASN
1	D	346	HIS
1	D	351	GLN
1	D	358	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](#)

4 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	3IO	C	3000	-	16,16,16	1.58	5 (31%)	21,22,22	1.10	2 (9%)
2	3IO	D	4000	-	16,16,16	1.58	5 (31%)	21,22,22	1.02	2 (9%)
2	3IO	B	2000	-	16,16,16	1.58	5 (31%)	21,22,22	1.16	2 (9%)
3	MCA	D	4001	-	47,50,57	2.04	9 (19%)	69,75,85	1.88	14 (20%)

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	3IO	C	3000	-	-	1/8/8/8	0/2/2/2
2	3IO	D	4000	-	-	1/8/8/8	0/2/2/2
2	3IO	B	2000	-	-	1/8/8/8	0/2/2/2
3	MCA	D	4001	-	-	6/48/64/75	0/3/3/3

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	4001	MCA	OP1-CP3	6.64	1.36	1.23
3	D	4001	MCA	CP5-NP2	-6.04	1.32	1.46
3	D	4001	MCA	CP4-CP3	-5.95	1.39	1.51
3	D	4001	MCA	CP5-CP4	-3.66	1.39	1.51
3	D	4001	MCA	P3-O31	3.52	1.61	1.50
3	D	4001	MCA	C5-N7	-3.23	1.33	1.39
3	D	4001	MCA	CP1-S	-3.11	1.67	1.80
2	B	2000	3IO	CAN-NAJ	-2.79	1.33	1.37
2	D	4000	3IO	CAN-NAJ	-2.78	1.33	1.37
2	C	3000	3IO	CAO-CAM	-2.74	1.39	1.44
2	C	3000	3IO	CAN-NAJ	-2.72	1.33	1.37
2	B	2000	3IO	CAO-CAM	-2.71	1.39	1.44
2	D	4000	3IO	CAO-CAM	-2.68	1.39	1.44
2	C	3000	3IO	CAI-CAM	2.66	1.53	1.50
2	B	2000	3IO	CAI-CAM	2.61	1.53	1.50
2	D	4000	3IO	CAI-CAM	2.54	1.53	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	4001	MCA	C8-N9	-2.32	1.33	1.37
2	C	3000	3IO	CAH-CAM	-2.23	1.33	1.36
2	B	2000	3IO	CAH-CAM	-2.23	1.33	1.36
2	D	4000	3IO	CAH-CAM	-2.20	1.33	1.36
2	D	4000	3IO	CAH-NAJ	-2.18	1.33	1.37
2	C	3000	3IO	CAH-NAJ	-2.15	1.33	1.37
2	B	2000	3IO	CAH-NAJ	-2.09	1.33	1.37
3	D	4001	MCA	C4-N9	-2.01	1.33	1.37

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	4001	MCA	CP2-NP1-CP3	-6.98	109.82	122.82
3	D	4001	MCA	C5-C4-N3	-5.22	119.53	126.72
3	D	4001	MCA	CP4-CP5-NP2	4.75	122.10	112.00
3	D	4001	MCA	N1-C2-N3	-4.49	121.79	128.58
3	D	4001	MCA	CP5-CP4-CP3	4.36	119.66	112.39
3	D	4001	MCA	N3-C4-N9	3.68	133.42	127.17
3	D	4001	MCA	C2-N3-C4	3.52	120.42	111.83
2	B	2000	3IO	OAA-CAK-CAL	-3.29	117.72	121.81
2	C	3000	3IO	OAA-CAK-CAL	-2.95	118.14	121.81
3	D	4001	MCA	N9-C8-N7	-2.92	109.79	113.94
2	D	4000	3IO	OAA-CAK-CAL	-2.75	118.39	121.81
3	D	4001	MCA	O4'-C1'-C2'	-2.72	100.78	106.62
3	D	4001	MCA	C5-N7-C8	2.66	107.63	103.45
3	D	4001	MCA	C4-C5-N7	-2.57	107.65	110.58
2	B	2000	3IO	CAF-CAN-CAO	-2.50	119.77	122.19
2	C	3000	3IO	CAF-CAN-CAO	-2.36	119.91	122.19
3	D	4001	MCA	C4'-O4'-C1'	-2.25	104.49	109.47
3	D	4001	MCA	CP4-CP3-NP1	2.15	120.26	116.34
2	D	4000	3IO	CAF-CAN-CAO	-2.13	120.13	122.19
3	D	4001	MCA	C4-N9-C8	2.04	107.88	105.74

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	2000	3IO	OAC-CAK-CAL-CAI
2	C	3000	3IO	OAC-CAK-CAL-CAI
2	D	4000	3IO	CAM-CAI-CAL-OAB
3	D	4001	MCA	CPB-O7-P2-O6
3	D	4001	MCA	CPB-O7-P2-O21

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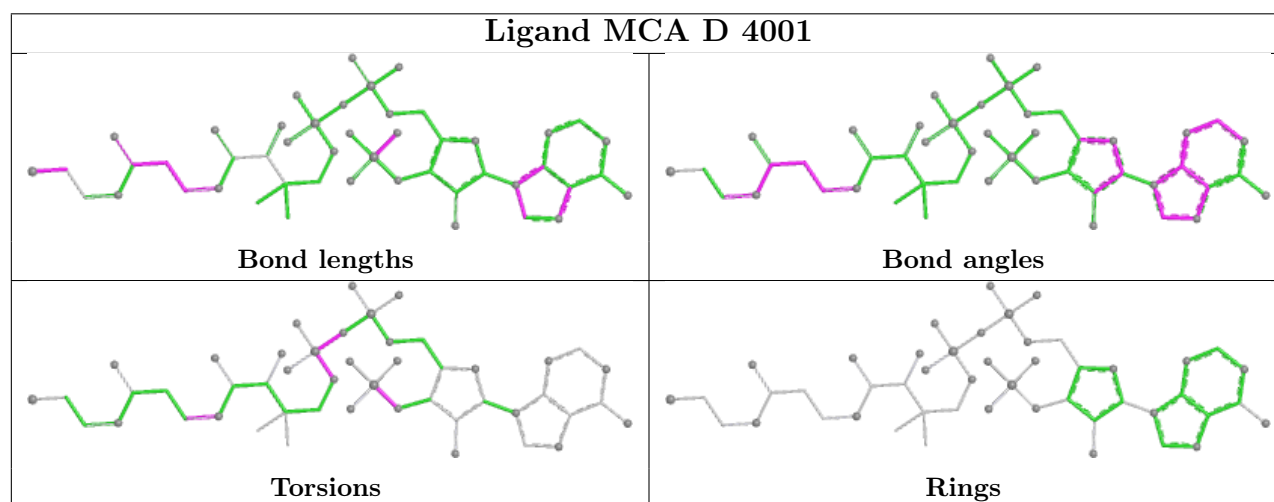
Mol	Chain	Res	Type	Atoms
3	D	4001	MCA	CPB-O7-P2-O22
3	D	4001	MCA	CP4-CP5-NP2-CP6
3	D	4001	MCA	C3'-O3'-P3-O33
3	D	4001	MCA	P1-O6-P2-O22

There are no ring outliers.

4 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	3000	3IO	3	0
2	D	4000	3IO	8	0
2	B	2000	3IO	2	0
3	D	4001	MCA	5	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	387/413 (93%)	-0.28	0 100 100	15, 24, 36, 47	4 (1%)
1	B	387/413 (93%)	-0.19	3 (0%) 82 82	15, 25, 40, 58	3 (0%)
1	C	382/413 (92%)	-0.20	2 (0%) 87 87	16, 25, 40, 60	5 (1%)
1	D	384/413 (92%)	-0.12	3 (0%) 82 82	16, 27, 40, 57	3 (0%)
All	All	1540/1652 (93%)	-0.20	8 (0%) 87 87	15, 25, 40, 60	15 (0%)

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	210	ASP	4.2
1	D	50	HIS	3.3
1	D	51	ASN	3.2
1	C	50	HIS	2.6
1	B	392	ILE	2.4
1	B	357	LEU	2.2
1	B	358	GLN	2.1
1	D	393	GLN	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](#)

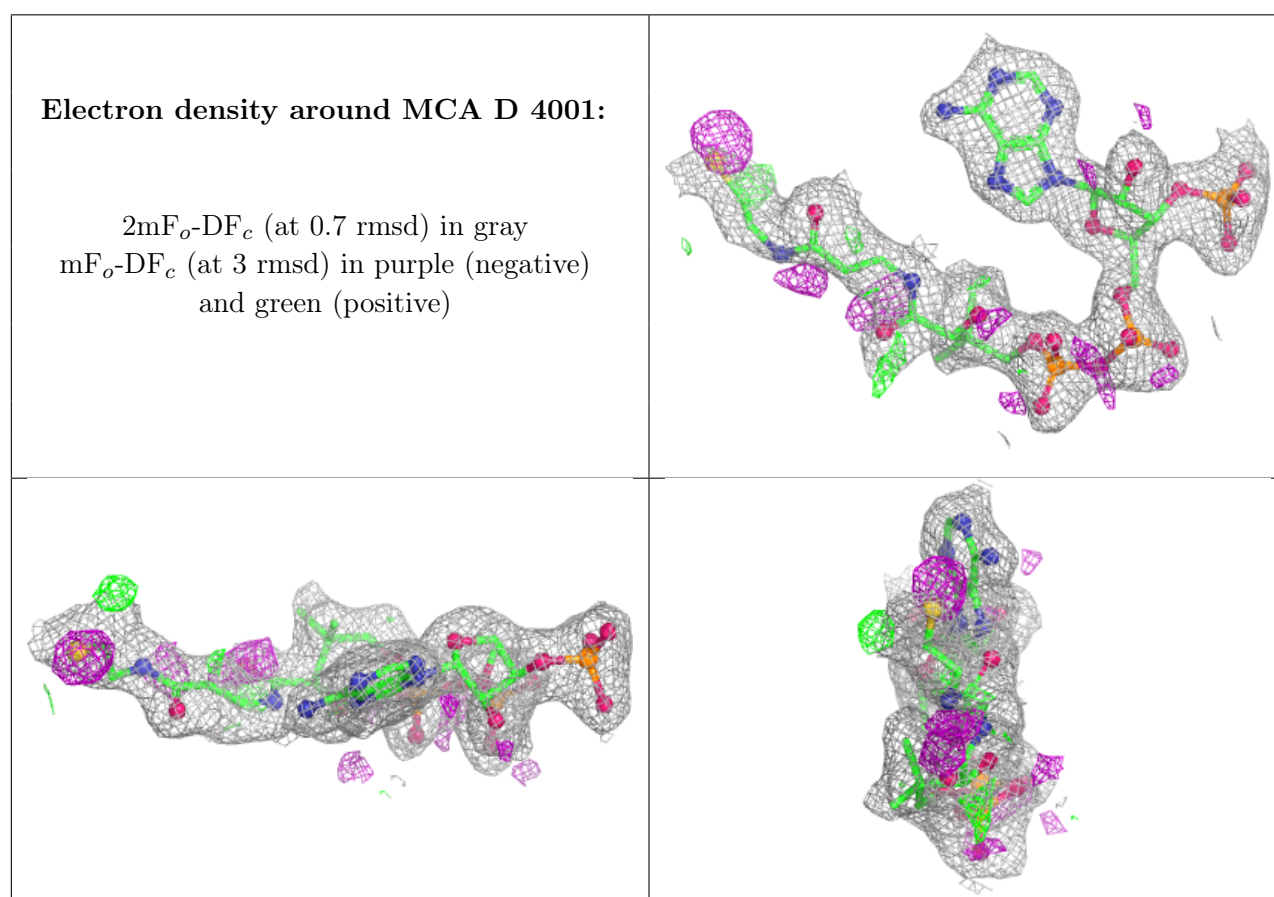
There are no oligosaccharides in this entry.

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($\text{\AA}^2$)	Q<0.9
2	3IO	D	4000	15/15	0.70	0.24	18,23,24,25	15
2	3IO	C	3000	15/15	0.84	0.12	27,35,36,37	0
2	3IO	B	2000	15/15	0.84	0.10	25,31,33,34	0
3	MCA	D	4001	48/55	0.91	0.10	25,31,41,57	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.