



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 04:58 PM UTC

PDB ID : 4QJE / pdb_00004qje
Title : 1.85 Angstrom resolution crystal structure of apo betaine aldehyde dehydrogenase (betB) G234S mutant from Staphylococcus aureus (IDP00699) with BME-free sulfinic acid form of Cys289
Authors : Halavaty, A.S.; Minasov, G.; Chen, C.; Joo, J.C.; Yakunin, A.F.; Anderson, W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2014-06-03
Resolution : 1.85 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](#) 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

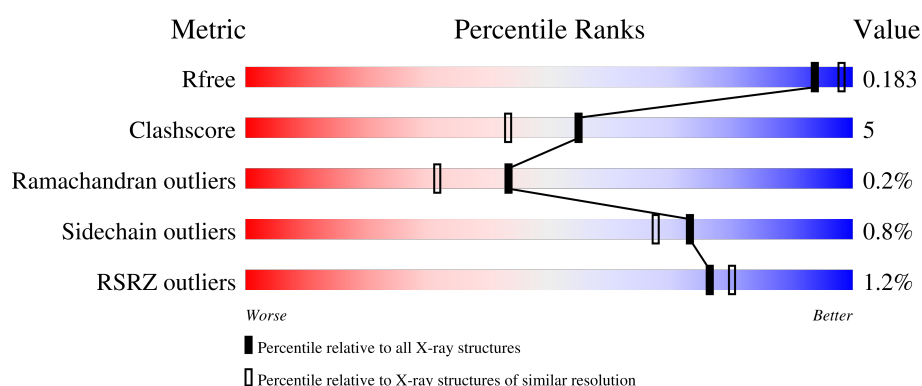
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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	A	517	88% 7% 5%
1	C	517	2% 87% 8% . .
1	D	517	2% 86% 9% . .
2	B	517	% 88% 7% . 5%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 18985 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 Betaine aldehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	493	Total	C	N	O	S	0	34	0
			4125	2592	695	823	15			
1	C	496	Total	C	N	O	S	0	25	0
			4055	2554	684	801	16			
1	D	496	Total	C	N	O	S	0	29	0
			4089	2568	696	809	16			

There are 66 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	expression tag	UNP Q5HCU0
A	-19	GLY	-	expression tag	UNP Q5HCU0
A	-18	SER	-	expression tag	UNP Q5HCU0
A	-17	SER	-	expression tag	UNP Q5HCU0
A	-16	HIS	-	expression tag	UNP Q5HCU0
A	-15	HIS	-	expression tag	UNP Q5HCU0
A	-14	HIS	-	expression tag	UNP Q5HCU0
A	-13	HIS	-	expression tag	UNP Q5HCU0
A	-12	HIS	-	expression tag	UNP Q5HCU0
A	-11	HIS	-	expression tag	UNP Q5HCU0
A	-10	SER	-	expression tag	UNP Q5HCU0
A	-9	SER	-	expression tag	UNP Q5HCU0
A	-8	GLY	-	expression tag	UNP Q5HCU0
A	-7	ARG	-	expression tag	UNP Q5HCU0
A	-6	GLU	-	expression tag	UNP Q5HCU0
A	-5	ASN	-	expression tag	UNP Q5HCU0
A	-4	LEU	-	expression tag	UNP Q5HCU0
A	-3	TYR	-	expression tag	UNP Q5HCU0
A	-2	PHE	-	expression tag	UNP Q5HCU0
A	-1	GLN	-	expression tag	UNP Q5HCU0
A	0	GLY	-	expression tag	UNP Q5HCU0
A	234	SER	GLY	engineered mutation	UNP Q5HCU0
C	-20	MET	-	expression tag	UNP Q5HCU0

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-19	GLY	-	expression tag	UNP Q5HCU0
C	-18	SER	-	expression tag	UNP Q5HCU0
C	-17	SER	-	expression tag	UNP Q5HCU0
C	-16	HIS	-	expression tag	UNP Q5HCU0
C	-15	HIS	-	expression tag	UNP Q5HCU0
C	-14	HIS	-	expression tag	UNP Q5HCU0
C	-13	HIS	-	expression tag	UNP Q5HCU0
C	-12	HIS	-	expression tag	UNP Q5HCU0
C	-11	HIS	-	expression tag	UNP Q5HCU0
C	-10	SER	-	expression tag	UNP Q5HCU0
C	-9	SER	-	expression tag	UNP Q5HCU0
C	-8	GLY	-	expression tag	UNP Q5HCU0
C	-7	ARG	-	expression tag	UNP Q5HCU0
C	-6	GLU	-	expression tag	UNP Q5HCU0
C	-5	ASN	-	expression tag	UNP Q5HCU0
C	-4	LEU	-	expression tag	UNP Q5HCU0
C	-3	TYR	-	expression tag	UNP Q5HCU0
C	-2	PHE	-	expression tag	UNP Q5HCU0
C	-1	GLN	-	expression tag	UNP Q5HCU0
C	0	GLY	-	expression tag	UNP Q5HCU0
C	234	SER	GLY	engineered mutation	UNP Q5HCU0
D	-20	MET	-	expression tag	UNP Q5HCU0
D	-19	GLY	-	expression tag	UNP Q5HCU0
D	-18	SER	-	expression tag	UNP Q5HCU0
D	-17	SER	-	expression tag	UNP Q5HCU0
D	-16	HIS	-	expression tag	UNP Q5HCU0
D	-15	HIS	-	expression tag	UNP Q5HCU0
D	-14	HIS	-	expression tag	UNP Q5HCU0
D	-13	HIS	-	expression tag	UNP Q5HCU0
D	-12	HIS	-	expression tag	UNP Q5HCU0
D	-11	HIS	-	expression tag	UNP Q5HCU0
D	-10	SER	-	expression tag	UNP Q5HCU0
D	-9	SER	-	expression tag	UNP Q5HCU0
D	-8	GLY	-	expression tag	UNP Q5HCU0
D	-7	ARG	-	expression tag	UNP Q5HCU0
D	-6	GLU	-	expression tag	UNP Q5HCU0
D	-5	ASN	-	expression tag	UNP Q5HCU0
D	-4	LEU	-	expression tag	UNP Q5HCU0
D	-3	TYR	-	expression tag	UNP Q5HCU0
D	-2	PHE	-	expression tag	UNP Q5HCU0
D	-1	GLN	-	expression tag	UNP Q5HCU0
D	0	GLY	-	expression tag	UNP Q5HCU0

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Chain	Residue	Modelled	Actual	Comment	Reference
D	234	SER	GLY	engineered mutation	UNP Q5HCU0

- Molecule 2 is a protein called Betaine aldehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	493	Total	C	N	O	S	0	30	0
			4079	2571	688	804	16			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	MET	-	expression tag	UNP Q5HCU0
B	-19	GLY	-	expression tag	UNP Q5HCU0
B	-18	SER	-	expression tag	UNP Q5HCU0
B	-17	SER	-	expression tag	UNP Q5HCU0
B	-16	HIS	-	expression tag	UNP Q5HCU0
B	-15	HIS	-	expression tag	UNP Q5HCU0
B	-14	HIS	-	expression tag	UNP Q5HCU0
B	-13	HIS	-	expression tag	UNP Q5HCU0
B	-12	HIS	-	expression tag	UNP Q5HCU0
B	-11	HIS	-	expression tag	UNP Q5HCU0
B	-10	SER	-	expression tag	UNP Q5HCU0
B	-9	SER	-	expression tag	UNP Q5HCU0
B	-8	GLY	-	expression tag	UNP Q5HCU0
B	-7	ARG	-	expression tag	UNP Q5HCU0
B	-6	GLU	-	expression tag	UNP Q5HCU0
B	-5	ASN	-	expression tag	UNP Q5HCU0
B	-4	LEU	-	expression tag	UNP Q5HCU0
B	-3	TYR	-	expression tag	UNP Q5HCU0
B	-2	PHE	-	expression tag	UNP Q5HCU0
B	-1	GLN	-	expression tag	UNP Q5HCU0
B	0	GLY	-	expression tag	UNP Q5HCU0
B	234	SER	GLY	engineered mutation	UNP Q5HCU0

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

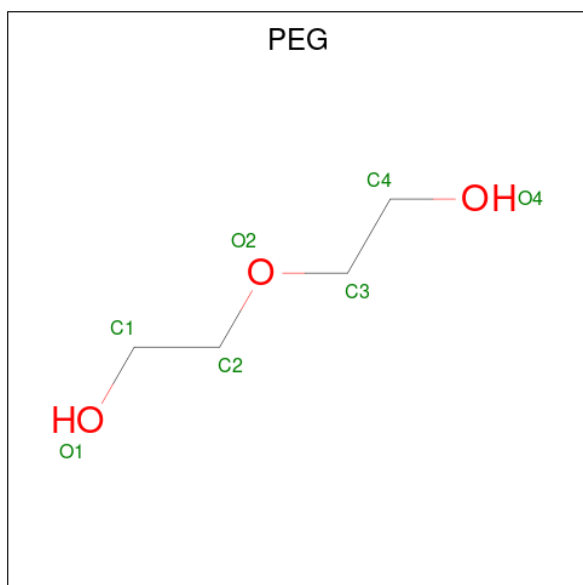
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Na	0	0
			2	2		
3	B	2	Total	Na	0	0
			2	2		

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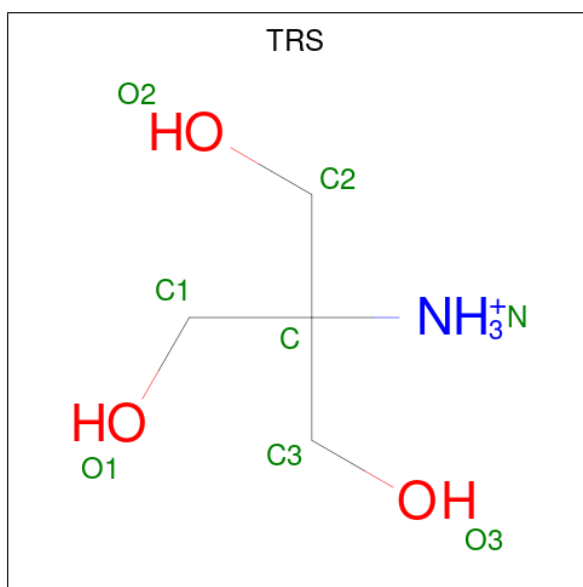
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	2	Total	Na	0	0
			2	2		
3	D	2	Total	Na	0	0
			2	2		

- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



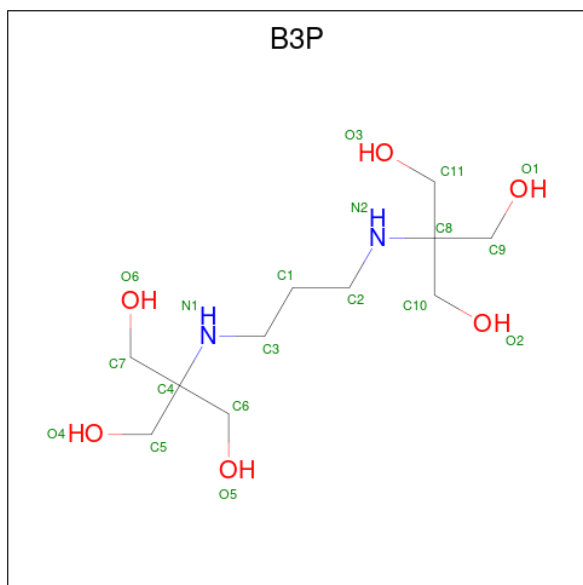
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		
4	A	1	Total	C	O	0	0
			7	4	3		
4	B	1	Total	C	O	0	0
			7	4	3		
4	C	1	Total	C	O	0	0
			7	4	3		
4	C	1	Total	C	O	0	0
			7	4	3		
4	C	1	Total	C	O	0	0
			7	4	3		
4	D	1	Total	C	O	0	0
			7	4	3		
4	D	1	Total	C	O	0	0
			7	4	3		

- Molecule 5 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			8	4	1	3		
5	C	1	Total	C	N	O	0	0
			8	4	1	3		
5	C	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 6 is 2-[3-(2-HYDROXY-1,1-DIHYDROXYMETHYL-ETHYLAMINO)-PROPYLAMINO]-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: B3P) (formula: C₁₁H₂₆N₂O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	C	1	Total	C	N	O	0	1
			38	22	4	12		

- Molecule 7 is water.

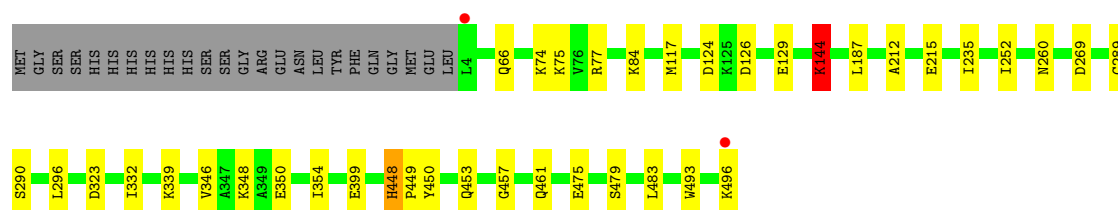
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	625	Total	O	0	29
			640	640		
7	B	592	Total	O	0	16
			598	598		
7	C	623	Total	O	0	23
			632	632		
7	D	632	Total	O	0	20
			641	641		

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.

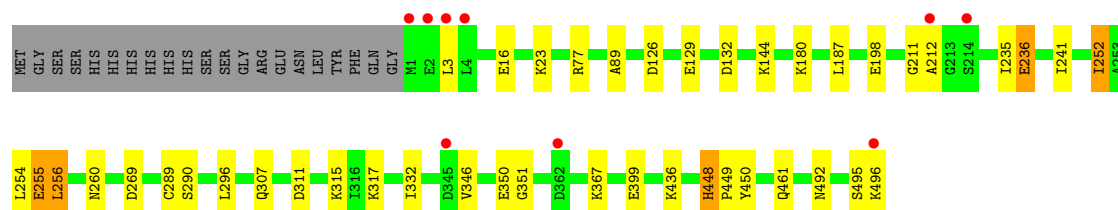
- Molecule 1: Betaine aldehyde dehydrogenase

Chain A: 



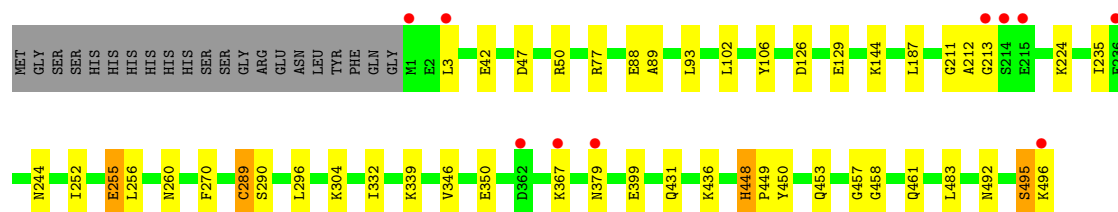
- Molecule 1: Betaine aldehyde dehydrogenase

Chain C: 



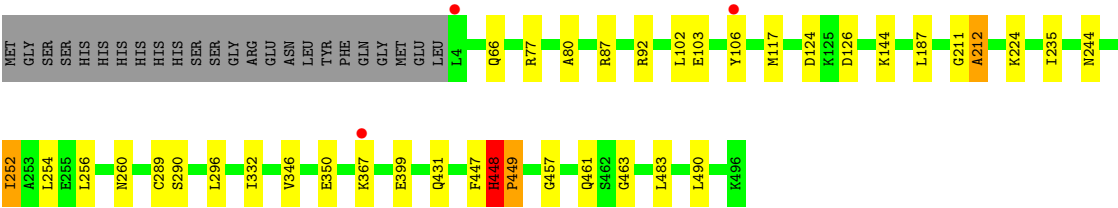
- Molecule 1: Betaine aldehyde dehydrogenase

Chain D: 



- Molecule 2: Betaine aldehyde dehydrogenase

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	218.66Å 102.93Å 118.07Å 90.00° 101.42° 90.00°	Depositor
Resolution (Å)	29.83 – 1.85 29.83 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.83-1.85) 99.7 (29.83-1.85)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 1.85Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.150 , 0.175 0.161 , 0.183	Depositor DCC
R_{free} test set	10934 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	16.9	Xtriage
Anisotropy	0.613	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	18985	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% 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: PEG, TRS, CSD, NA, B3P

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	0.65	0/4184	0.89	10/5656 (0.2%)
1	C	0.70	0/4112	0.89	9/5561 (0.2%)
1	D	0.70	0/4147	0.88	10/5605 (0.2%)
2	B	0.67	0/4153	0.88	11/5614 (0.2%)
All	All	0.68	0/16596	0.88	40/22436 (0.2%)

There are no bond length outliers.

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	457	GLY	CA-C-N	-7.28	117.16	122.18
1	D	457	GLY	C-N-CA	-7.28	117.16	122.18
2	B	457	GLY	CA-C-N	-7.08	117.29	122.18
2	B	457	GLY	C-N-CA	-7.08	117.29	122.18
1	A	457	GLY	CA-C-N	-7.08	117.30	122.18
1	A	457	GLY	C-N-CA	-7.08	117.30	122.18
1	D	332	ILE	N-CA-C	6.40	119.09	111.09
2	B	212	ALA	N-CA-C	-6.38	102.64	110.61
2	B	449	PRO	N-CA-C	6.37	120.84	111.03
1	C	212	ALA	N-CA-C	-6.21	102.85	110.61
1	D	212	ALA	N-CA-C	-6.16	102.91	110.61
1	A	187	LEU	N-CA-C	5.88	117.36	111.07
1	C	256[A]	LEU	N-CA-C	-5.78	100.83	109.96
1	C	256[B]	LEU	N-CA-C	-5.78	100.83	109.96
1	C	187	LEU	N-CA-C	5.74	117.21	111.07
1	D	495	SER	N-CA-C	5.68	117.62	110.24
1	C	269	ASP	N-CA-C	-5.65	101.01	109.15
2	B	256[A]	LEU	N-CA-C	-5.61	101.10	109.96
2	B	256[B]	LEU	N-CA-C	-5.61	101.10	109.96
1	D	187	LEU	N-CA-C	5.56	117.02	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	458	GLY	N-CA-C	5.52	120.84	111.19
1	C	449	PRO	N-CA-C	5.47	119.46	111.03
1	D	448	HIS	CA-C-N	-5.47	114.16	119.80
1	D	448	HIS	C-N-CA	-5.47	114.16	119.80
1	C	332	ILE	N-CA-C	5.45	118.74	111.44
2	B	448	HIS	CA-C-N	-5.30	114.11	119.83
2	B	448	HIS	C-N-CA	-5.30	114.11	119.83
1	A	144	LYS	CB-CG-CD	5.28	123.45	111.30
1	A	269	ASP	N-CA-C	-5.21	101.65	109.15
1	C	448	HIS	CA-C-N	-5.19	114.22	119.83
1	C	448	HIS	C-N-CA	-5.19	114.22	119.83
1	A	449	PRO	N-CA-C	5.18	119.17	111.09
2	B	187	LEU	N-CA-C	5.17	116.61	111.07
1	A	448	HIS	CA-C-N	-5.15	114.49	119.80
1	A	448	HIS	C-N-CA	-5.15	114.49	119.80
2	B	332	ILE	N-CA-C	5.14	118.32	111.44
1	A	332	ILE	N-CA-C	5.13	118.32	111.44
1	A	212	ALA	N-CA-C	5.07	117.69	110.24
2	B	463	GLY	N-CA-C	5.06	117.86	110.63
1	D	449	PRO	N-CA-C	5.03	118.94	111.09

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	4125	0	4008	32	0
1	C	4055	0	3985	38	0
1	D	4089	0	4004	49	0
2	B	4079	0	3992	37	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
4	A	14	0	20	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	7	0	10	1	0
4	C	21	0	30	1	0
4	D	14	0	20	0	0
5	A	8	0	12	0	0
5	C	16	0	24	3	0
6	C	38	0	52	4	0
7	A	640	0	0	14	0
7	B	598	0	0	16	0
7	C	632	0	0	27	0
7	D	641	0	0	24	0
All	All	18985	0	16157	158	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 5.

All (158) 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:450[B]:TYR:HE2	7:D:1123:HOH:O	1.42	1.02
1:C:450[B]:TYR:HE2	7:C:764:HOH:O	1.43	0.98
2:B:126[B]:ASP:HB3	7:B:1180[B]:HOH:O	1.64	0.96
1:A:450[B]:TYR:HE2	7:A:806:HOH:O	1.50	0.95
1:D:126[A]:ASP:HB3	7:D:1193[A]:HOH:O	1.67	0.94
1:C:126[B]:ASP:HB3	7:C:1174[B]:HOH:O	1.72	0.88
2:B:92[A]:ARG:HD2	7:B:839:HOH:O	1.74	0.88
2:B:289[A]:CYS:SG	7:B:1108:HOH:O	2.36	0.84
1:C:450[A]:TYR:HE2	7:C:913[A]:HOH:O	1.59	0.84
2:B:126[B]:ASP:CB	7:B:1180[B]:HOH:O	2.26	0.81
2:B:117[B]:MET:HA	2:B:117[B]:MET:CE	2.10	0.81
1:A:144:LYS:HE3	1:A:479:SER:OG	1.80	0.80
1:D:450[B]:TYR:CE2	7:D:1123:HOH:O	2.24	0.80
2:B:126[B]:ASP:CG	7:B:1180[B]:HOH:O	2.25	0.80
2:B:77[B]:ARG:HG2	2:B:117[B]:MET:HE1	1.64	0.79
1:A:450[B]:TYR:CE2	7:A:806:HOH:O	2.30	0.79
6:C:508[A]:B3P:H31	6:C:508[A]:B3P:O6	1.83	0.77
1:A:323:ASP:HB3	7:A:1051:HOH:O	1.83	0.77
5:C:507:TRS:H21	7:C:1029:HOH:O	1.84	0.76
1:A:126[A]:ASP:HB3	7:A:1225[A]:HOH:O	1.86	0.74
2:B:77[B]:ARG:HG2	2:B:117[B]:MET:CE	2.17	0.73
1:A:339:LYS:HD2	7:A:846[B]:HOH:O	1.89	0.72
2:B:117[B]:MET:HA	2:B:117[B]:MET:HE3	1.71	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:450[B]:TYR:CE2	7:C:764:HOH:O	2.28	0.71
1:A:129:GLU:HG3	7:A:1011:HOH:O	1.92	0.70
1:C:450[A]:TYR:CE2	7:C:913[A]:HOH:O	2.40	0.69
1:D:3:LEU:CD2	1:D:93:LEU:HD22	2.24	0.68
1:C:144:LYS:CE	7:C:1119:HOH:O	2.41	0.67
1:D:483[B]:LEU:C	1:D:483[B]:LEU:HD23	2.19	0.67
1:D:339:LYS:HD2	7:D:1134:HOH:O	1.94	0.66
2:B:77[B]:ARG:CG	2:B:117[B]:MET:HE1	2.25	0.66
1:C:317[A]:LYS:NZ	1:D:495:SER:OG	2.22	0.64
1:D:77[B]:ARG:NH2	7:D:1155:HOH:O	2.29	0.64
1:C:144:LYS:HD2	7:C:798:HOH:O	1.97	0.63
1:A:453:GLN:HG2	7:B:809:HOH:O	1.99	0.63
1:D:224:LYS:NZ	7:D:1064:HOH:O	2.31	0.63
1:C:315:LYS:NZ	1:D:496:LYS:O	2.25	0.62
1:D:431[A]:GLN:NE2	7:D:874:HOH:O	2.32	0.62
1:C:144:LYS:HE2	7:C:1119:HOH:O	1.98	0.62
1:A:450[A]:TYR:HD2	7:A:977:HOH:O	1.83	0.61
1:C:77[B]:ARG:NH1	7:C:762:HOH:O	2.34	0.61
1:A:215[A]:GLU:OE2	1:A:215[A]:GLU:HA	2.00	0.61
1:D:3:LEU:HG	7:D:920:HOH:O	1.98	0.61
2:B:346:VAL:O	2:B:350:GLU:HG2	2.02	0.60
1:A:346:VAL:O	1:A:350[A]:GLU:HG2	2.02	0.60
1:C:317[A]:LYS:HE2	7:C:907:HOH:O	2.00	0.60
1:A:346:VAL:O	1:A:350[B]:GLU:HG2	2.02	0.59
1:D:346:VAL:O	1:D:350:GLU:HG2	2.03	0.59
1:A:348:LYS:HE2	1:A:354[B]:ILE:HD12	1.85	0.58
2:B:77[A]:ARG:NH1	7:B:1107:HOH:O	2.36	0.58
1:D:436:LYS:HD2	7:D:1139:HOH:O	2.02	0.58
1:A:77[A]:ARG:NH1	7:A:791:HOH:O	2.36	0.58
1:D:88:GLU:OE1	1:D:106[A]:TYR:OH	2.21	0.58
1:C:346:VAL:O	1:C:350:GLU:HG2	2.04	0.57
1:A:215[A]:GLU:OE2	1:A:215[A]:GLU:CA	2.53	0.57
1:D:42:GLU:HG3	7:D:1159:HOH:O	2.04	0.56
1:D:379[A]:ASN:HB2	7:D:1217[A]:HOH:O	2.04	0.56
1:C:255[A]:GLU:OE1	1:C:256[A]:LEU:O	2.24	0.56
1:D:47:ASP:OD1	1:D:50[B]:ARG:NH1	2.38	0.56
6:C:508[A]:B3P:O6	6:C:508[A]:B3P:C3	2.52	0.56
1:A:144:LYS:CE	1:A:479:SER:OG	2.54	0.55
1:D:255[B]:GLU:CD	7:D:1079:HOH:O	2.49	0.55
1:D:255[A]:GLU:OE1	1:D:256[A]:LEU:O	2.25	0.55
1:A:75[B]:LYS:NZ	7:A:1038:HOH:O	2.29	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:1154:HOH:O	2:B:431:GLN:HG2	2.07	0.54
2:B:66:GLN:HB2	4:B:503:PEG:H12	1.89	0.54
1:D:379[B]:ASN:O	1:D:379[B]:ASN:CG	2.50	0.53
1:D:88:GLU:CD	1:D:106[A]:TYR:OH	2.52	0.53
1:C:241[B]:ILE:HD13	7:C:920:HOH:O	2.08	0.53
1:D:3:LEU:HD21	1:D:93:LEU:HD22	1.89	0.53
1:C:198:GLU:OE2	6:C:508[B]:B3P:H91	2.09	0.53
1:C:16:GLU:HG3	7:C:932:HOH:O	2.07	0.53
1:D:88:GLU:CD	1:D:106[A]:TYR:HH	2.18	0.52
1:D:492:ASN:OD1	1:D:496:LYS:HE3	2.10	0.52
1:D:102:LEU:HG	1:D:106[A]:TYR:CE2	2.44	0.52
1:C:351:GLY:C	7:C:882:HOH:O	2.52	0.52
1:D:244[B]:ASN:ND2	7:D:906:HOH:O	2.40	0.51
1:C:436:LYS:HD3	7:D:1161:HOH:O	2.11	0.51
1:A:493:TRP:CZ3	2:B:103[B]:GLU:HG2	2.46	0.51
2:B:102[B]:LEU:O	2:B:106[B]:TYR:CD2	2.64	0.51
2:B:490[B]:LEU:HD23	7:B:695:HOH:O	2.11	0.50
2:B:483[B]:LEU:HD23	2:B:483[B]:LEU:C	2.36	0.50
1:A:124:ASP:HB2	7:C:787:HOH:O	2.11	0.50
1:D:213:GLY:CA	7:D:1159:HOH:O	2.59	0.50
1:C:492:ASN:OD1	1:C:496:LYS:HE3	2.12	0.50
1:C:3:LEU:HD13	1:C:89:ALA:HB1	1.94	0.50
5:C:506:TRS:H12	7:C:1140:HOH:O	2.11	0.49
1:D:88:GLU:OE2	1:D:106[A]:TYR:OH	2.29	0.49
1:C:132:ASP:O	7:C:787:HOH:O	2.20	0.49
1:D:144:LYS:HD2	7:D:789:HOH:O	2.11	0.49
1:A:483[B]:LEU:C	1:A:483[B]:LEU:HD23	2.37	0.49
1:C:180:LYS:HE2	7:C:1165:HOH:O	2.13	0.49
1:C:126[B]:ASP:CB	7:C:1174[B]:HOH:O	2.47	0.48
1:C:3:LEU:HB3	7:C:1139:HOH:O	2.13	0.48
1:C:495:SER:N	7:C:1146:HOH:O	2.46	0.48
1:D:379[A]:ASN:CB	7:D:1217[A]:HOH:O	2.59	0.48
2:B:212:ALA:HA	7:B:770:HOH:O	2.14	0.48
1:A:475:GLU:HG3	7:A:1153:HOH:O	2.14	0.47
1:D:450[B]:TYR:CD2	1:D:450[B]:TYR:C	2.87	0.47
1:D:304[A]:LYS:HD3	7:D:1225[A]:HOH:O	2.13	0.47
2:B:117[B]:MET:HA	2:B:117[B]:MET:HE2	1.90	0.47
1:C:23:LYS:O	7:C:1054:HOH:O	2.21	0.46
7:C:787:HOH:O	1:D:453:GLN:HG2	2.14	0.46
1:D:270:PHE:CE1	1:D:304[A]:LYS:HD2	2.51	0.46
2:B:252[B]:ILE:HD11	2:B:254:LEU:HD21	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:270:PHE:CZ	1:D:304[A]:LYS:HD2	2.51	0.46
1:A:117[B]:MET:CE	1:D:77[B]:ARG:NH1	2.79	0.45
1:A:496:LYS:HD2	7:B:1099:HOH:O	2.16	0.45
1:A:74:LYS:HE2	7:A:808:HOH:O	2.16	0.45
1:C:252[B]:ILE:HD11	1:C:254:LEU:HD21	1.99	0.45
1:A:339:LYS:HD2	7:A:846[A]:HOH:O	2.17	0.45
2:B:244[B]:ASN:ND2	7:B:1033:HOH:O	2.49	0.45
1:D:483[B]:LEU:C	1:D:483[B]:LEU:CD2	2.89	0.45
2:B:80:ALA:HB1	2:B:117[B]:MET:HG2	1.99	0.45
2:B:296:LEU:HD23	2:B:399:GLU:HB2	1.98	0.44
1:D:296:LEU:HD23	1:D:399:GLU:HB2	1.99	0.44
1:A:66:GLN:HB2	4:A:503:PEG:H22	1.98	0.44
2:B:124:ASP:HB2	7:D:823:HOH:O	2.17	0.44
1:D:3:LEU:CD1	7:D:920:HOH:O	2.65	0.44
1:D:260:ASN:ND2	1:D:290:SER:HA	2.32	0.44
1:D:50[B]:ARG:NH2	1:D:50[B]:ARG:HB2	2.33	0.44
2:B:235:ILE:HD11	2:B:461:GLN:OE1	2.18	0.43
1:C:211:GLY:HA3	7:C:911:HOH:O	2.18	0.43
1:A:296:LEU:HD23	1:A:399:GLU:HB2	1.99	0.43
1:A:84:LYS:HE2	7:A:1135:HOH:O	2.17	0.43
1:C:492:ASN:OD1	1:C:496:LYS:CE	2.67	0.43
1:C:235:ILE:HD11	1:C:461:GLN:OE1	2.19	0.43
1:D:3:LEU:HD13	1:D:89:ALA:HB1	2.00	0.43
1:C:296:LEU:HD23	1:C:399:GLU:HB2	2.00	0.43
2:B:447:PHE:C	2:B:449:PRO:HD3	2.44	0.43
6:C:508[A]:B3P:H22	6:C:508[A]:B3P:H112	1.76	0.43
1:D:289[A]:CSD:OD2	7:D:1140:HOH:O	2.20	0.42
2:B:87:ARG:NH2	7:B:714:HOH:O	2.51	0.42
2:B:260:ASN:ND2	2:B:290:SER:HA	2.34	0.42
5:C:507:TRS:C2	7:C:1029:HOH:O	2.55	0.42
2:B:102[B]:LEU:HB3	2:B:106[B]:TYR:CE2	2.55	0.42
2:B:447:PHE:O	2:B:448:HIS:HB2	2.20	0.42
1:C:236:GLU:H	1:C:236:GLU:CD	2.26	0.42
1:C:252[B]:ILE:HG13	1:C:254:LEU:HG	2.02	0.42
1:A:354[B]:ILE:HD13	1:A:354[B]:ILE:HG21	1.76	0.42
2:B:117[B]:MET:HE2	2:B:117[B]:MET:CA	2.50	0.42
1:C:346:VAL:HG21	4:C:504:PEG:H42	2.01	0.42
2:B:224:LYS:NZ	7:B:927:HOH:O	2.53	0.41
2:B:144:LYS:HD2	7:B:880:HOH:O	2.21	0.41
1:C:311[B]:ASP:OD1	7:C:855[B]:HOH:O	2.21	0.41
2:B:103[B]:GLU:HA	2:B:106[B]:TYR:HD2	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:307[B]:GLN:HG3	7:C:752:HOH:O	2.20	0.41
1:D:492:ASN:OD1	1:D:496:LYS:CE	2.68	0.41
1:D:211:GLY:HA3	7:D:777:HOH:O	2.20	0.41
1:A:235:ILE:HD11	1:A:461:GLN:OE1	2.20	0.41
2:B:211:GLY:HA3	7:B:665:HOH:O	2.21	0.41
1:D:235:ILE:HD11	1:D:461:GLN:OE1	2.20	0.41
1:D:304[A]:LYS:NZ	7:D:691:HOH:O	2.32	0.41
1:A:450[B]:TYR:CD2	1:A:450[B]:TYR:C	2.94	0.40
2:B:461:GLN:NE2	7:B:1136:HOH:O	2.34	0.40
1:C:260:ASN:ND2	1:C:290:SER:HA	2.36	0.40
1:D:3:LEU:CG	7:D:920:HOH:O	2.65	0.40
1:A:260:ASN:ND2	1:A:290:SER:HA	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	524/517 (101%)	514 (98%)	9 (2%)	1 (0%)	43	31
1	C	518/517 (100%)	508 (98%)	9 (2%)	1 (0%)	43	31
1	D	521/517 (101%)	510 (98%)	10 (2%)	1 (0%)	43	31
2	B	522/517 (101%)	511 (98%)	10 (2%)	1 (0%)	43	31
All	All	2085/2068 (101%)	2043 (98%)	38 (2%)	4 (0%)	43	31

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	448	HIS
2	B	448	HIS
1	C	448	HIS

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Mol	Chain	Res	Type
1	D	448	HIS

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	440/429 (103%)	437 (99%)	3 (1%)	76	70
1	C	434/429 (101%)	426 (98%)	8 (2%)	51	40
1	D	437/429 (102%)	432 (99%)	5 (1%)	65	57
2	B	438/430 (102%)	435 (99%)	3 (1%)	76	70
All	All	1749/1717 (102%)	1730 (99%)	19 (1%)	73	57

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	144	LYS
1	A	252[A]	ILE
1	A	252[B]	ILE
2	B	252[A]	ILE
2	B	252[B]	ILE
2	B	367	LYS
1	C	129[A]	GLU
1	C	129[B]	GLU
1	C	236	GLU
1	C	252[A]	ILE
1	C	252[B]	ILE
1	C	255[A]	GLU
1	C	255[B]	GLU
1	C	367	LYS
1	D	129	GLU
1	D	252	ILE
1	D	255[A]	GLU
1	D	255[B]	GLU
1	D	367	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	GLN
1	A	260	ASN
1	A	279	ASN
2	B	34	GLN
2	B	260	ASN
2	B	279	ASN
2	B	431	GLN
1	C	34	GLN
1	C	260	ASN
1	C	279	ASN
1	D	34	GLN
1	D	66	GLN
1	D	260	ASN
1	D	279	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

6 non-standard protein/DNA/RNA residues 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
1	CSD	C	289[A]	1	4,7,8	1.69	1 (25%)	1,8,10	3.37	1 (100%)
1	CSD	A	289[A]	1	4,7,8	2.15	1 (25%)	1,8,10	0.76	0
1	CSD	D	289[A]	1	4,7,8	1.65	1 (25%)	1,8,10	3.42	1 (100%)
1	CSD	C	289[B]	1	4,7,8	2.57	1 (25%)	1,8,10	2.40	1 (100%)
1	CSD	A	289[B]	1	4,7,8	1.16	0	1,8,10	4.03	1 (100%)
1	CSD	D	289[B]	1	4,7,8	2.20	1 (25%)	1,8,10	3.32	1 (100%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	CSD	C	289[A]	1	-	1/2/6/8	-
1	CSD	A	289[A]	1	-	1/2/6/8	-
1	CSD	D	289[A]	1	-	1/2/6/8	-
1	CSD	C	289[B]	1	-	1/2/6/8	-
1	CSD	A	289[B]	1	-	1/2/6/8	-
1	CSD	D	289[B]	1	-	1/2/6/8	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	289[B]	CSD	OD1-SG	-4.59	1.43	1.47
1	D	289[B]	CSD	OD1-SG	-3.76	1.44	1.47
1	A	289[A]	CSD	OD1-SG	-3.71	1.44	1.47
1	C	289[A]	CSD	OD1-SG	-2.60	1.45	1.47
1	D	289[A]	CSD	OD1-SG	-2.25	1.45	1.47

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	289[B]	CSD	OD1-SG-CB	4.03	113.03	105.60
1	D	289[A]	CSD	OD1-SG-CB	3.42	111.90	105.60
1	C	289[A]	CSD	OD1-SG-CB	3.37	111.80	105.60
1	D	289[B]	CSD	OD1-SG-CB	3.32	111.71	105.60
1	C	289[B]	CSD	OD1-SG-CB	2.40	110.02	105.60

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	289[A]	CSD	CA-CB-SG-OD1
1	A	289[B]	CSD	CA-CB-SG-OD1
1	C	289[A]	CSD	CA-CB-SG-OD1
1	C	289[B]	CSD	CA-CB-SG-OD1
1	D	289[A]	CSD	CA-CB-SG-OD1
1	D	289[B]	CSD	CA-CB-SG-OD1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	289[A]	CSD	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 21 ligands modelled in this entry, 8 are monoatomic - leaving 13 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	$\# Z > 2$
4	PEG	C	503	-	6,6,6	0.42	0	5,5,5	0.29	0
4	PEG	D	504	-	6,6,6	0.46	0	5,5,5	0.29	0
6	B3P	C	508[A]	-	18,18,18	0.71	0	23,23,23	1.66	4 (17%)
4	PEG	C	504	-	6,6,6	0.43	0	5,5,5	0.68	0
5	TRS	A	505	-	7,7,7	0.61	0	9,9,9	0.82	0
5	TRS	C	507	-	7,7,7	0.37	0	9,9,9	0.73	0
5	TRS	C	506	-	7,7,7	0.42	0	9,9,9	0.91	1 (11%)
4	PEG	B	503	-	6,6,6	0.37	0	5,5,5	0.25	0
4	PEG	A	503	-	6,6,6	0.46	0	5,5,5	0.44	0
4	PEG	A	504	-	6,6,6	0.42	0	5,5,5	0.27	0
6	B3P	C	508[B]	-	18,18,18	0.72	0	23,23,23	1.19	2 (8%)
4	PEG	D	503	-	6,6,6	0.41	0	5,5,5	0.56	0
4	PEG	C	505	-	6,6,6	0.42	0	5,5,5	0.30	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
4	PEG	C	503	-	-	1/4/4/4	-
4	PEG	D	504	-	-	2/4/4/4	-
6	B3P	C	508[A]	-	-	12/28/28/28	-
4	PEG	C	504	-	-	3/4/4/4	-
5	TRS	A	505	-	-	5/9/9/9	-
5	TRS	C	507	-	-	3/9/9/9	-
5	TRS	C	506	-	-	3/9/9/9	-
4	PEG	B	503	-	-	3/4/4/4	-
4	PEG	A	503	-	-	2/4/4/4	-
4	PEG	A	504	-	-	1/4/4/4	-
6	B3P	C	508[B]	-	-	12/28/28/28	-
4	PEG	D	503	-	-	3/4/4/4	-
4	PEG	C	505	-	-	3/4/4/4	-

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	508[A]	B3P	C3-N1-C4	5.87	124.76	116.17
6	C	508[B]	B3P	C3-N1-C4	3.78	121.70	116.17
6	C	508[B]	B3P	C2-N2-C8	3.06	120.64	116.17
6	C	508[A]	B3P	C11-C8-C10	-2.60	104.32	110.02
6	C	508[A]	B3P	C7-C4-N1	2.31	115.92	109.02
6	C	508[A]	B3P	C2-N2-C8	2.19	119.38	116.17
5	C	506	TRS	C3-C-C2	-2.10	105.07	110.66

There are no chirality outliers.

All (53) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	C	508[A]	B3P	C5-C4-N1-C3
6	C	508[A]	B3P	C6-C4-N1-C3
6	C	508[A]	B3P	C7-C4-N1-C3
6	C	508[A]	B3P	N1-C4-C6-O5
6	C	508[A]	B3P	C5-C4-C6-O5
6	C	508[A]	B3P	N1-C4-C7-O6
6	C	508[A]	B3P	C5-C4-C7-O6
6	C	508[A]	B3P	C6-C4-C7-O6
6	C	508[B]	B3P	N1-C4-C7-O6

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Mol	Chain	Res	Type	Atoms
6	C	508[B]	B3P	C5-C4-C7-O6
6	C	508[B]	B3P	C6-C4-C7-O6
6	C	508[B]	B3P	O3-C11-C8-N2
6	C	508[B]	B3P	O3-C11-C8-C9
6	C	508[B]	B3P	O3-C11-C8-C10
4	A	503	PEG	O2-C3-C4-O4
6	C	508[A]	B3P	C7-C4-C6-O5
4	A	504	PEG	O2-C3-C4-O4
4	C	503	PEG	O1-C1-C2-O2
4	C	505	PEG	O2-C3-C4-O4
4	D	503	PEG	O1-C1-C2-O2
4	D	504	PEG	O2-C3-C4-O4
5	A	505	TRS	C1-C-C2-O2
5	C	507	TRS	C1-C-C2-O2
6	C	508[B]	B3P	C6-C4-C5-O4
6	C	508[B]	B3P	C7-C4-C5-O4
4	A	503	PEG	C1-C2-O2-C3
4	D	503	PEG	O2-C3-C4-O4
4	B	503	PEG	O2-C3-C4-O4
6	C	508[B]	B3P	C10-C8-N2-C2
4	C	504	PEG	C1-C2-O2-C3
6	C	508[B]	B3P	C1-C3-N1-C4
4	D	503	PEG	C4-C3-O2-C2
4	C	504	PEG	O1-C1-C2-O2
4	B	503	PEG	O1-C1-C2-O2
5	A	505	TRS	N-C-C2-O2
5	C	507	TRS	N-C-C2-O2
6	C	508[A]	B3P	C2-C1-C3-N1
4	C	505	PEG	O1-C1-C2-O2
4	C	505	PEG	C1-C2-O2-C3
6	C	508[A]	B3P	C11-C8-N2-C2
6	C	508[B]	B3P	C9-C8-N2-C2
5	A	505	TRS	C3-C-C2-O2
5	C	506	TRS	C3-C-C1-O1
5	C	507	TRS	C3-C-C2-O2
4	C	504	PEG	C4-C3-O2-C2
6	C	508[B]	B3P	N1-C4-C5-O4
4	D	504	PEG	C1-C2-O2-C3
4	B	503	PEG	C4-C3-O2-C2
5	A	505	TRS	C2-C-C1-O1
5	C	506	TRS	C1-C-C3-O3
6	C	508[A]	B3P	O3-C11-C8-C10

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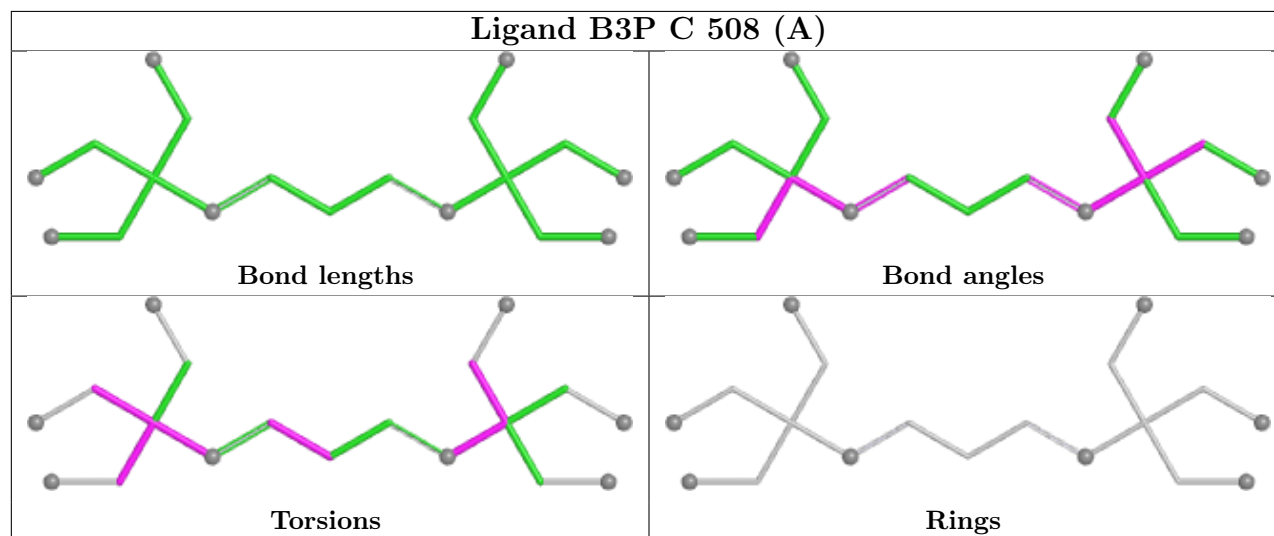
Mol	Chain	Res	Type	Atoms
5	A	505	TRS	N-C-C1-O1
5	C	506	TRS	N-C-C3-O3

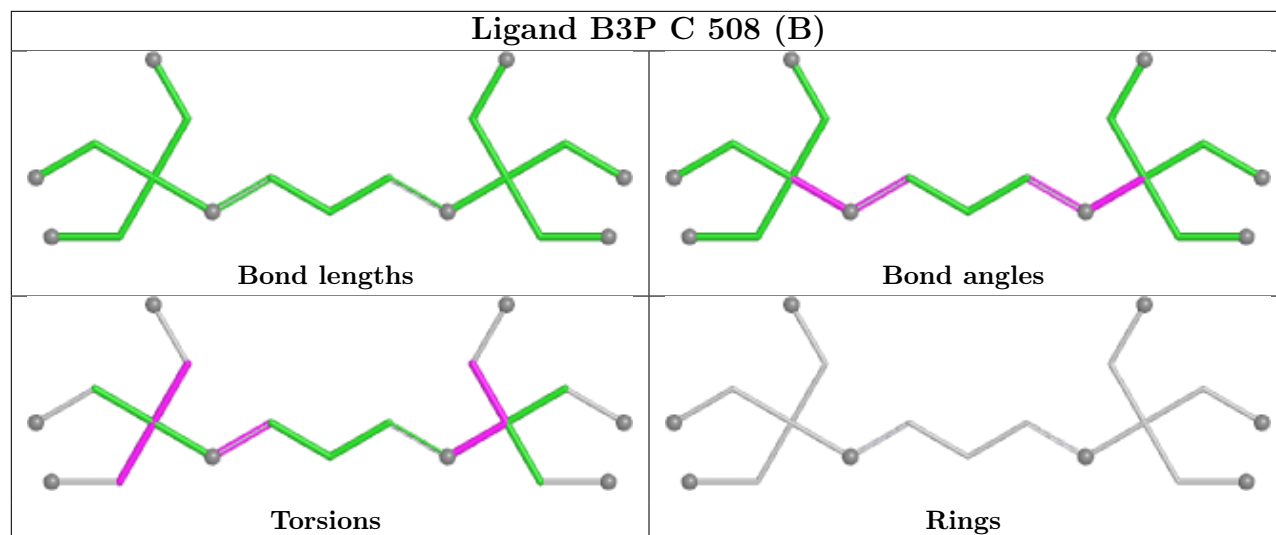
There are no ring outliers.

7 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	508[A]	B3P	3	0
4	C	504	PEG	1	0
5	C	507	TRS	2	0
5	C	506	TRS	1	0
4	B	503	PEG	1	0
4	A	503	PEG	1	0
6	C	508[B]	B3P	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	492/517 (95%)	-0.42	2 (0%) 88 91	5, 18, 31, 69	33 (6%)
1	C	495/517 (95%)	-0.35	9 (1%) 67 71	4, 17, 35, 93	24 (4%)
1	D	495/517 (95%)	-0.40	10 (2%) 65 68	6, 16, 32, 85	28 (5%)
2	B	493/517 (95%)	-0.30	3 (0%) 85 88	6, 20, 36, 67	30 (6%)
All	All	1975/2068 (95%)	-0.37	24 (1%) 76 80	4, 17, 34, 93	115 (5%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	3	LEU	5.2
1	C	3	LEU	4.9
1	D	1	MET	3.9
1	C	1	MET	3.6
2	B	4	LEU	3.1
1	D	379[A]	ASN	2.8
1	C	4	LEU	2.8
1	D	367	LYS	2.8
2	B	106[A]	TYR	2.7
1	D	362	ASP	2.6
1	A	4	LEU	2.5
1	D	496	LYS	2.5
1	C	496	LYS	2.4
1	C	345	ASP	2.3
1	C	2	GLU	2.3
1	C	212	ALA	2.3
1	D	214	SER	2.3
1	C	362	ASP	2.2
1	C	214	SER	2.2
2	B	367	LYS	2.2
1	A	496	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	215	GLU	2.0
1	D	236	GLU	2.0
1	D	213	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	CSD	A	289[A]	8/9	0.84	0.13	18,20,25,26	8
1	CSD	A	289[B]	8/9	0.84	0.13	19,22,32,33	8
1	CSD	D	289[A]	8/9	0.84	0.16	21,23,29,31	8
1	CSD	D	289[B]	8/9	0.84	0.16	21,24,30,30	8
1	CSD	C	289[A]	8/9	0.88	0.13	22,24,31,32	8
1	CSD	C	289[B]	8/9	0.88	0.13	22,24,30,31	8

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
4	PEG	D	504	7/7	0.75	0.17	65,68,72,72	0
5	TRS	C	507	8/8	0.79	0.14	38,39,43,45	0
4	PEG	C	503	7/7	0.81	0.14	53,56,59,60	0
4	PEG	C	505	7/7	0.81	0.16	50,56,59,62	0
4	PEG	B	503	7/7	0.82	0.16	46,52,58,62	0
4	PEG	D	503	7/7	0.82	0.17	47,49,63,78	0
5	TRS	A	505	8/8	0.83	0.13	36,42,43,49	0
4	PEG	C	504	7/7	0.85	0.13	52,54,56,61	0
5	TRS	C	506	8/8	0.87	0.10	32,34,36,41	0

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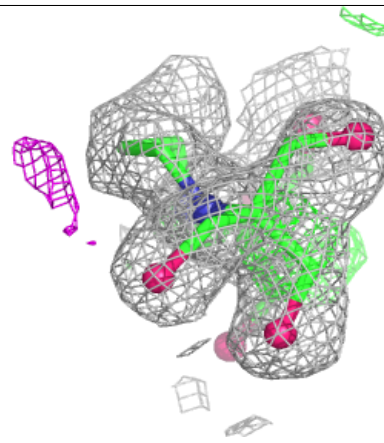
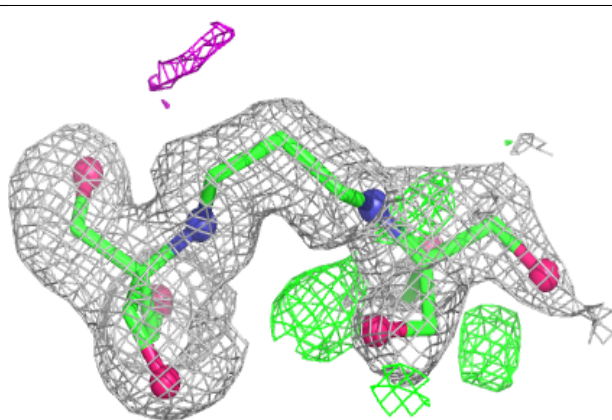
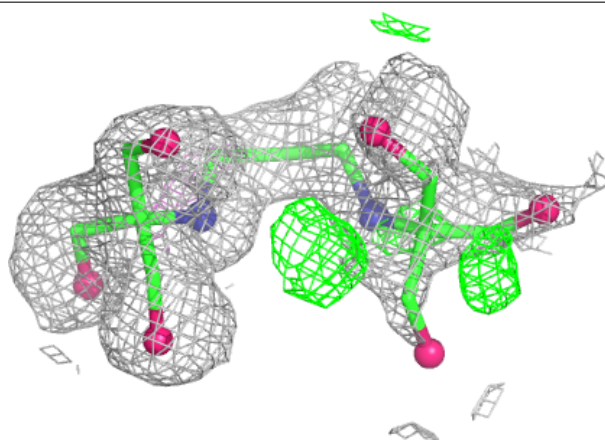
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	PEG	A	504	7/7	0.88	0.13	52,54,57,58	0
6	B3P	C	508[A]	19/19	0.88	0.11	14,18,25,26	19
6	B3P	C	508[B]	19/19	0.88	0.11	23,27,40,42	19
4	PEG	A	503	7/7	0.89	0.13	40,41,55,56	0
3	NA	C	501	1/1	0.96	0.06	19,19,19,19	0
3	NA	A	501	1/1	0.97	0.06	21,21,21,21	0
3	NA	D	501	1/1	0.98	0.04	19,19,19,19	0
3	NA	B	501	1/1	0.98	0.07	21,21,21,21	0
3	NA	A	502	1/1	0.99	0.03	15,15,15,15	0
3	NA	D	502	1/1	0.99	0.03	13,13,13,13	0
3	NA	C	502	1/1	0.99	0.06	12,12,12,12	0
3	NA	B	502	1/1	1.00	0.03	15,15,15,15	0

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

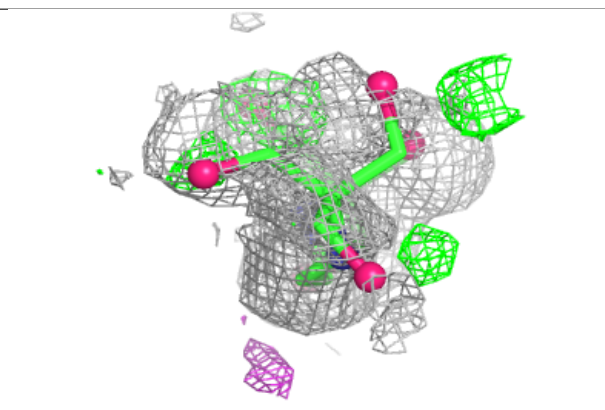
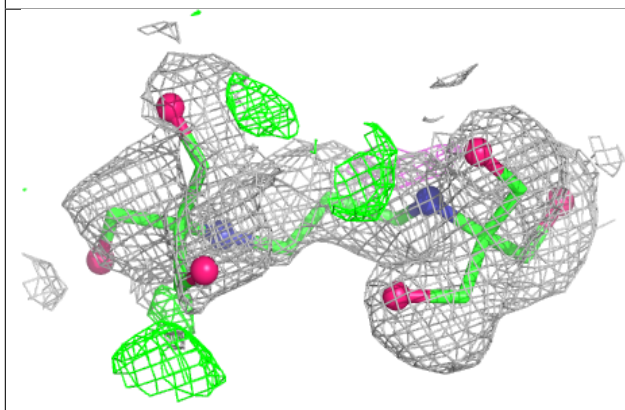
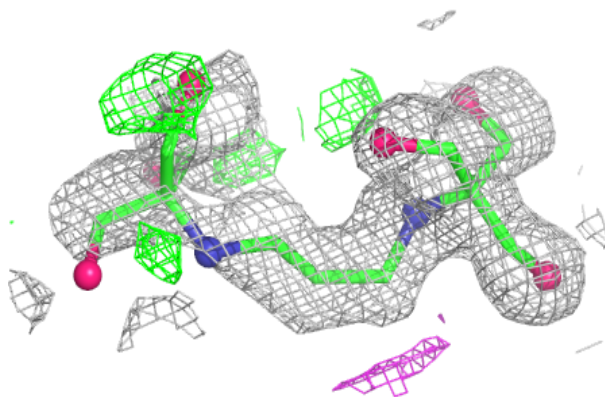
Electron density around B3P C 508 (A):

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 B3P C 508 (B):

$2mF_o-DF_c$ (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.