



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:26 PM GMT

PDB ID : 1AT1  
Title : CRYSTAL STRUCTURES OF PHOSPHONOACETAMIDE LIGATED T AND PHOSPHONOACETAMIDE AND MALONATE LIGATED R STATES OF ASPARTATE CARBAMOYLTRANSFERASE AT 2.8-ANGSTROMS RESOLUTION AND NEUTRAL P\*H  
Authors : Gouaux, J.E.; Lipscomb, W.N.  
Deposited on : 1989-09-22  
Resolution : 2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

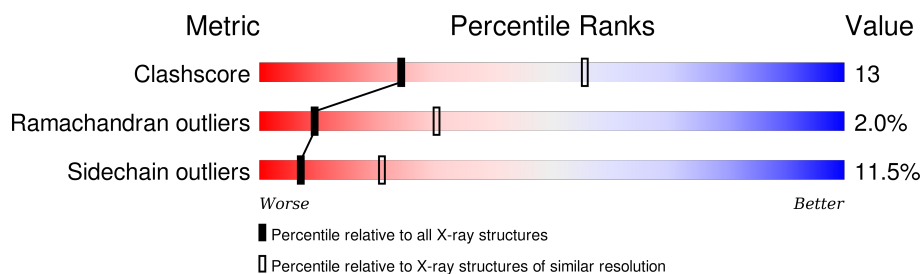
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	310	
1	C	310	
2	B	153	
2	D	153	

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7138 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 ASPARTATE CARBAMOYLTRANSFERASE, CATALYTIC CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	310	Total	C	N	O	S	0	0	0
			2415	1527	423	456	9			
1	C	310	Total	C	N	O	S	0	0	0
			2415	1527	423	456	9			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	60	GLN	GLU	CONFLICT	UNP P0A786
A	147	GLN	GLU	CONFLICT	UNP P0A786
A	149	GLU	GLN	CONFLICT	UNP P0A786
A	196	GLU	GLN	CONFLICT	UNP P0A786
C	60	GLN	GLU	CONFLICT	UNP P0A786
C	147	GLN	GLU	CONFLICT	UNP P0A786
C	149	GLU	GLN	CONFLICT	UNP P0A786
C	196	GLU	GLN	CONFLICT	UNP P0A786

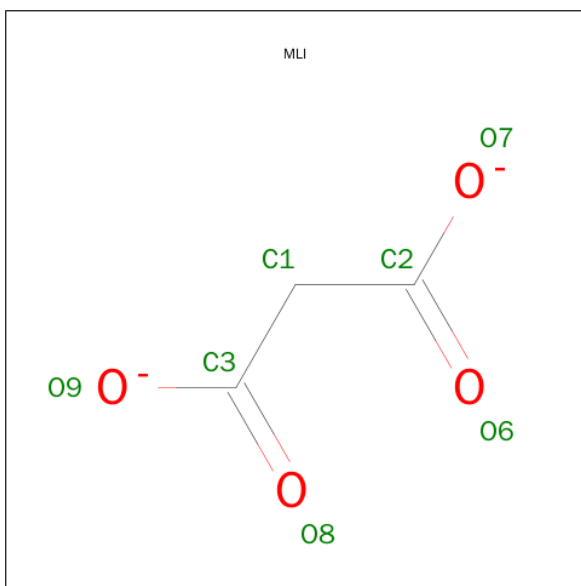
- Molecule 2 is a protein called ASPARTATE CARBAMOYLTRANSFERASE REGULATORY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	146	Total	C	N	O	S	0	0	0
			1138	714	201	218	5			
2	D	146	Total	C	N	O	S	0	0	0
			1138	714	201	218	5			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	8	GLY	GLN	CONFLICT	UNP P0A7F3
D	8	GLY	GLN	CONFLICT	UNP P0A7F3

- Molecule 3 is MALONATE ION (three-letter code: MLI) (formula:  $C_3H_2O_4$ ).

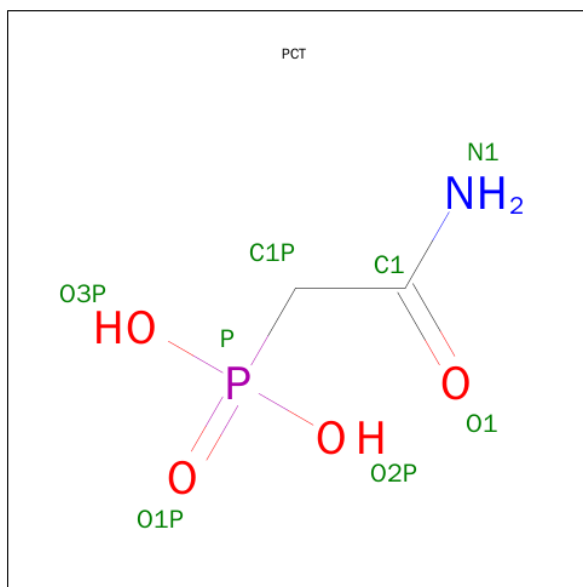


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	3	4		
3	C	1	Total	C	O	0	0
			7	3	4		

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula:  $Zn$ ).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Zn	0	0
			1	1		
4	D	1	Total	Zn	0	0
			1	1		

- Molecule 5 is PHOSPHONOACETAMIDE (three-letter code: PCT) (formula:  $C_2H_6NO_4P$ ).



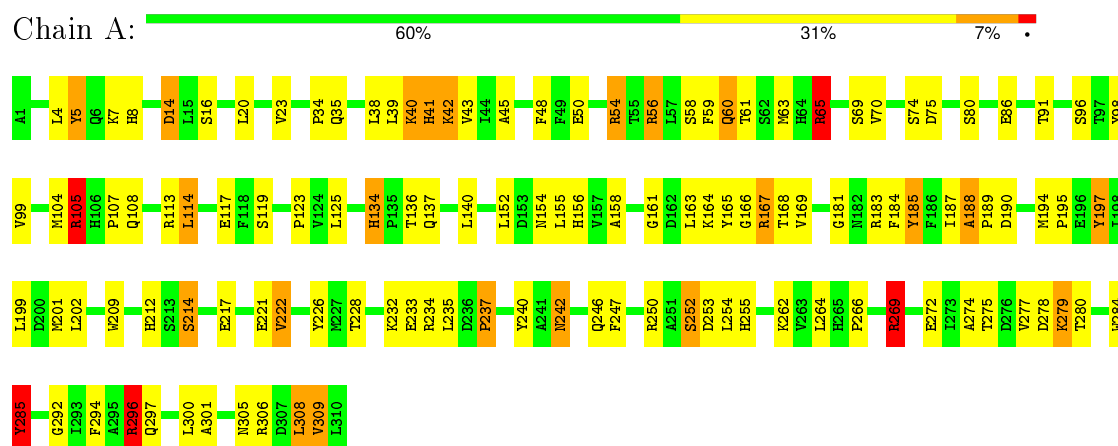
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			8	2	1	4	1		
5	C	1	Total	C	N	O	P	0	0
			8	2	1	4	1		

### 3 Residue-property plots

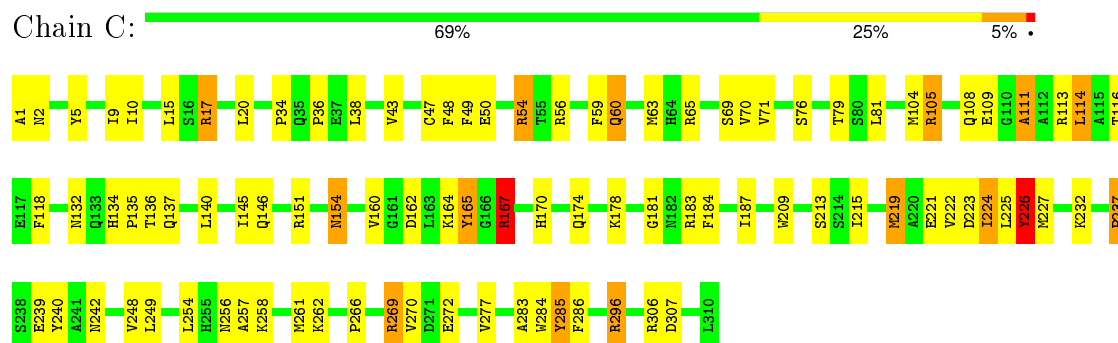
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

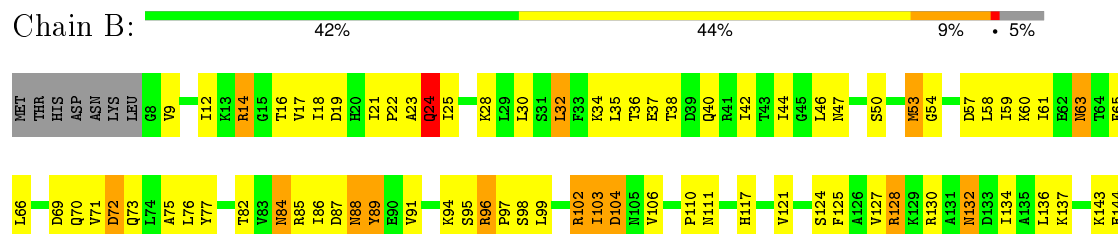
- Molecule 1: ASPARTATE CARBAMOYLTRANSFERASE, CATALYTIC CHAIN



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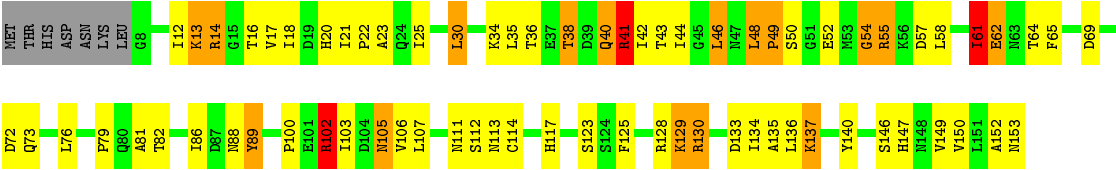


- Molecule 2: ASPARTATE CARBAMOYLTRANSFERASE REGULATORY CHAIN





● Molecule 2: ASPARTATE CARBAMOYLTRANSFERASE REGULATORY CHAIN



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	122.30 Å   122.30 Å   156.40 Å 90.00°   90.00°   120.00°	Depositor
Resolution (Å)	6.00 – 2.80	Depositor
% Data completeness (in resolution range)	(Not available) (6.00-2.80)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.164 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7138	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, MLI, PCT

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	0.89	0/2461	1.69	44/3339 (1.3%)
1	C	0.92	0/2461	1.72	39/3339 (1.2%)
2	B	0.82	1/1155 (0.1%)	1.58	8/1561 (0.5%)
2	D	0.80	0/1155	1.62	11/1561 (0.7%)
All	All	0.88	1/7232 (0.0%)	1.67	102/9800 (1.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	C	0	6
2	B	0	2
2	D	0	2
All	All	0	13

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	106	VAL	CA-CB	5.39	1.66	1.54

All (102) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	65	ARG	NE-CZ-NH2	13.52	127.06	120.30
1	A	56	ARG	NE-CZ-NH1	-13.15	113.73	120.30
2	B	130	ARG	NE-CZ-NH2	11.54	126.07	120.30
1	C	54	ARG	NE-CZ-NH2	10.77	125.69	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	102	ARG	NE-CZ-NH2	10.61	125.61	120.30
2	D	14	ARG	NE-CZ-NH2	10.54	125.57	120.30
1	C	56	ARG	NE-CZ-NH2	10.41	125.51	120.30
1	C	105	ARG	NE-CZ-NH2	10.27	125.43	120.30
1	A	234	ARG	NE-CZ-NH1	-9.95	115.33	120.30
1	A	98	TYR	CB-CG-CD2	-9.92	115.05	121.00
1	C	269	ARG	NE-CZ-NH2	9.72	125.16	120.30
1	A	234	ARG	NE-CZ-NH2	9.31	124.96	120.30
1	C	54	ARG	NE-CZ-NH1	-9.19	115.70	120.30
1	C	17	ARG	NE-CZ-NH2	9.02	124.81	120.30
1	C	209	TRP	CD1-CG-CD2	8.74	113.29	106.30
1	A	54	ARG	NE-CZ-NH2	8.65	124.63	120.30
2	B	96	ARG	NE-CZ-NH1	-8.63	115.99	120.30
1	C	284	TRP	CD1-CG-CD2	8.61	113.18	106.30
1	C	113	ARG	NE-CZ-NH1	-8.47	116.06	120.30
1	C	114	LEU	CA-CB-CG	8.29	134.38	115.30
1	A	65	ARG	NE-CZ-NH2	8.14	124.37	120.30
1	C	284	TRP	CE2-CD2-CG	-8.12	100.81	107.30
1	C	209	TRP	CE2-CD2-CG	-8.09	100.83	107.30
1	C	151	ARG	NE-CZ-NH2	8.07	124.33	120.30
2	B	89	TYR	CB-CG-CD2	-7.83	116.30	121.00
1	C	113	ARG	NE-CZ-NH2	7.80	124.20	120.30
1	A	209	TRP	CD1-CG-CD2	7.57	112.36	106.30
1	A	284	TRP	CD1-CG-CD2	7.47	112.27	106.30
1	A	105	ARG	NE-CZ-NH2	7.40	124.00	120.30
1	C	60	GLN	CA-CB-CG	-7.30	97.33	113.40
1	A	309	VAL	CA-C-N	-7.27	101.20	117.20
1	A	56	ARG	NE-CZ-NH2	7.17	123.89	120.30
1	C	167	ARG	CB-CG-CD	7.10	130.06	111.60
2	D	102	ARG	NE-CZ-NH1	-7.08	116.76	120.30
1	C	269	ARG	CA-CB-CG	7.05	128.91	113.40
1	C	105	ARG	NE-CZ-NH1	-7.04	116.78	120.30
1	C	65	ARG	NE-CZ-NH1	-6.97	116.82	120.30
2	D	152	ALA	CA-C-N	-6.91	102.00	117.20
1	A	91	THR	CA-CB-CG2	6.81	121.94	112.40
1	C	226	TYR	CB-CG-CD2	-6.80	116.92	121.00
1	A	284	TRP	CE2-CD2-CG	-6.70	101.94	107.30
1	A	309	VAL	O-C-N	6.64	133.32	122.70
1	A	114	LEU	CA-CB-CG	6.56	130.39	115.30
1	C	43	VAL	CG1-CB-CG2	-6.55	100.41	110.90
1	C	167	ARG	NE-CZ-NH1	-6.50	117.05	120.30
1	A	269	ARG	NE-CZ-NH2	6.46	123.53	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	209	TRP	CB-CG-CD1	-6.44	118.63	127.00
2	B	152	ALA	N-CA-C	6.41	128.31	111.00
1	A	209	TRP	CE2-CD2-CG	-6.41	102.17	107.30
1	A	185	TYR	CB-CG-CD1	-6.31	117.21	121.00
1	A	201	MET	CA-CB-CG	6.26	123.94	113.30
1	A	105	ARG	CA-CB-CG	-6.11	99.95	113.40
1	A	285	TYR	CB-CG-CD2	-6.11	117.34	121.00
2	B	102	ARG	NE-CZ-NH2	6.05	123.33	120.30
1	A	104	MET	CG-SD-CE	-6.04	90.53	100.20
2	D	61	ILE	CA-CB-CG2	-6.00	98.91	110.90
2	B	128	ARG	NE-CZ-NH2	5.97	123.28	120.30
1	A	91	THR	CA-CB-OG1	-5.96	96.50	109.00
1	A	98	TYR	CA-CB-CG	5.91	124.63	113.40
1	A	5	TYR	CB-CG-CD1	-5.88	117.47	121.00
1	A	104	MET	N-CA-CB	-5.81	100.14	110.60
1	C	219	MET	CG-SD-CE	-5.76	90.98	100.20
1	A	167	ARG	NE-CZ-NH1	-5.75	117.42	120.30
1	C	269	ARG	CG-CD-NE	-5.75	99.73	111.80
1	C	104	MET	CA-CB-CG	5.70	122.98	113.30
1	A	14	ASP	CA-C-N	5.68	129.71	117.20
2	D	105	ASN	N-CA-C	5.68	126.34	111.00
2	D	130	ARG	NE-CZ-NH2	5.67	123.13	120.30
1	C	269	ARG	NE-CZ-NH1	-5.65	117.47	120.30
1	A	23	VAL	CG1-CB-CG2	-5.64	101.87	110.90
1	A	86	GLU	CA-CB-CG	5.63	125.78	113.40
1	A	232	LYS	N-CA-CB	-5.62	100.49	110.60
1	C	165	TYR	CB-CG-CD1	-5.53	117.68	121.00
1	A	296	ARG	NE-CZ-NH2	5.52	123.06	120.30
2	B	37	GLU	CA-CB-CG	-5.52	101.25	113.40
1	C	209	TRP	CG-CD1-NE1	-5.51	104.58	110.10
1	A	14	ASP	O-C-N	-5.45	113.98	122.70
1	A	169	VAL	CA-CB-CG1	-5.41	102.78	110.90
1	A	308	LEU	CA-CB-CG	5.39	127.70	115.30
1	C	284	TRP	CB-CG-CD1	-5.34	120.05	127.00
1	A	54	ARG	NE-CZ-NH1	-5.34	117.63	120.30
2	D	46	LEU	CA-C-N	-5.33	105.48	117.20
1	C	56	ARG	NH1-CZ-NH2	-5.30	113.57	119.40
1	C	296	ARG	NE-CZ-NH1	-5.28	117.66	120.30
1	A	269	ARG	CB-CG-CD	-5.27	97.89	111.60
2	D	41	ARG	NE-CZ-NH2	5.27	122.94	120.30
1	A	242	ASN	CA-C-N	-5.27	105.61	117.20
2	B	14	ARG	NE-CZ-NH2	5.23	122.92	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	111	ALA	CA-C-N	-5.20	105.75	117.20
1	C	47	CYS	CA-CB-SG	-5.15	104.74	114.00
1	C	222	VAL	CA-C-N	-5.14	105.89	117.20
1	A	65	ARG	NE-CZ-NH1	-5.12	117.74	120.30
1	C	48	PHE	CA-C-N	-5.11	105.95	117.20
2	D	114	CYS	CA-CB-SG	-5.11	104.81	114.00
1	A	169	VAL	CA-CB-CG2	5.09	118.53	110.90
1	A	222	VAL	N-CA-C	5.08	124.71	111.00
2	D	14	ARG	NE-CZ-NH1	-5.06	117.77	120.30
1	A	58	SER	CA-C-N	5.04	128.29	117.20
1	A	136	THR	CA-CB-CG2	-5.04	105.34	112.40
1	C	162	ASP	CB-CG-OD1	5.03	122.83	118.30
1	A	60	GLN	O-C-N	-5.03	114.66	122.70
1	C	225	LEU	CA-CB-CG	5.01	126.81	115.30

There are no chirality outliers.

All (13) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	188	ALA	Peptide
1	A	197	TYR	Sidechain
1	A	285	TYR	Sidechain
2	B	77	TYR	Sidechain
2	B	89	TYR	Sidechain
1	C	118	PHE	Sidechain
1	C	226	TYR	Sidechain
1	C	240	TYR	Sidechain
1	C	285	TYR	Sidechain
1	C	286	PHE	Sidechain
1	C	5	TYR	Sidechain
2	D	140	TYR	Sidechain
2	D	89	TYR	Sidechain

## 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	2415	0	2422	57	0
1	C	2415	0	2422	41	0
2	B	1138	0	1154	50	0
2	D	1138	0	1154	46	0
3	A	7	0	2	0	0
3	C	7	0	2	0	0
4	B	1	0	0	0	0
4	D	1	0	0	0	0
5	A	8	0	4	1	0
5	C	8	0	4	1	0
All	All	7138	0	7164	186	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 13.

All (186) 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)
2:B:14:ARG:HG3	2:B:87:ASP:HA	1.53	0.91
2:D:146:SER:HB3	2:D:149:VAL:HG23	1.55	0.88
2:B:22:PRO:HB2	2:B:25:ILE:HD12	1.63	0.81
2:D:102:ARG:HH21	2:D:102:ARG:HB3	1.45	0.81
1:C:145:ILE:HG12	1:C:224:ILE:HD13	1.61	0.80
1:A:292:GLY:O	1:A:296:ARG:HB2	1.84	0.77
2:D:49:PRO:HA	2:D:55:ARG:HA	1.67	0.75
1:C:154:ASN:HA	1:C:181:GLY:O	1.89	0.72
2:D:13:LYS:HG3	2:D:88:ASN:HA	1.73	0.71
2:D:130:ARG:HD2	2:D:135:ALA:HB2	1.74	0.69
2:B:84:ASN:HD21	2:B:91:VAL:HG13	1.59	0.67
2:D:106:VAL:HG23	2:D:107:LEU:HD12	1.76	0.66
1:A:43:VAL:HG22	1:A:69:SER:HB2	1.78	0.66
1:A:154:ASN:HA	1:A:181:GLY:O	1.96	0.66
1:C:10:ILE:HD11	1:C:116:THR:HG21	1.79	0.65
2:B:42:ILE:HB	2:D:46:LEU:HB2	1.79	0.65
2:B:30:LEU:HD11	2:B:59:ILE:HG21	1.77	0.65
1:A:4:LEU:HD23	1:A:7:LYS:HD3	1.78	0.64
1:C:54:ARG:HE	5:C:311:PCT:H1P2	1.61	0.64
2:B:66:LEU:HD12	2:B:71:VAL:HG22	1.81	0.63
2:B:16:THR:OG1	2:B:65:PHE:HA	1.98	0.63
2:B:75:ALA:HB2	2:B:97:PRO:HB2	1.79	0.63
2:B:53:MET:SD	2:B:54:GLY:N	2.72	0.63
2:B:30:LEU:HD12	2:B:35:LEU:HD13	1.79	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:72:ASP:HB3	2:D:100:PRO:HG3	1.80	0.62
2:B:134:ILE:O	2:B:147:HIS:HB3	1.99	0.61
1:C:248:VAL:HG22	1:C:272:GLU:HA	1.82	0.61
1:C:9:ILE:HD12	1:C:15:LEU:HD11	1.84	0.60
1:C:76:SER:HA	1:C:79:THR:HG23	1.83	0.60
2:D:111:ASN:O	2:D:117:HIS:HE1	1.84	0.60
2:B:44:ILE:HB	2:D:44:ILE:HD13	1.83	0.60
1:A:63:MET:SD	1:A:70:VAL:HG22	2.43	0.59
1:A:254:LEU:HD12	1:A:280:THR:HG21	1.84	0.59
1:C:137:GLN:O	1:C:140:LEU:HG	2.03	0.58
2:B:46:LEU:HB2	2:D:42:ILE:HB	1.85	0.58
1:A:137:GLN:HG2	1:A:168:THR:HG22	1.85	0.57
2:D:38:THR:HG23	2:D:40:GLN:H	1.71	0.56
2:B:44:ILE:HG23	2:B:59:ILE:HG13	1.85	0.56
1:A:114:LEU:HD22	2:B:121:VAL:HG11	1.88	0.56
2:B:104:ASP:HA	2:B:124:SER:HA	1.87	0.55
1:A:35:GLN:HB3	1:A:38:LEU:HD13	1.88	0.55
1:C:50:GLU:HB3	1:C:105:ARG:HG2	1.88	0.55
1:A:235:LEU:HB2	1:A:240:TYR:HE2	1.72	0.55
2:D:14:ARG:HG3	2:D:65:PHE:CE1	2.41	0.55
1:C:50:GLU:HB3	1:C:105:ARG:CG	2.37	0.55
1:C:154:ASN:HD22	1:C:181:GLY:HA3	1.73	0.54
1:C:63:MET:SD	1:C:70:VAL:HG22	2.48	0.53
1:C:160:VAL:HB	1:C:227:MET:SD	2.49	0.53
1:A:4:LEU:HA	1:A:7:LYS:HD2	1.90	0.52
2:B:136:LEU:HD12	2:B:150:VAL:HG21	1.92	0.52
1:A:185:TYR:HD2	1:A:212:HIS:NE2	2.07	0.52
1:C:49:PHE:HD2	1:C:76:SER:HB3	1.75	0.52
2:D:125:PHE:HA	2:D:137:LYS:O	2.10	0.52
2:D:107:LEU:HD23	2:D:150:VAL:HG11	1.92	0.52
2:D:14:ARG:HA	2:D:86:ILE:O	2.10	0.51
2:B:103:ILE:HD11	2:B:127:VAL:HG22	1.92	0.51
2:D:14:ARG:HG3	2:D:65:PHE:HE1	1.75	0.51
2:B:50:SER:HB3	2:B:53:MET:O	2.10	0.51
1:C:254:LEU:HD22	1:C:261:MET:HE1	1.92	0.51
2:B:72:ASP:HA	2:B:97:PRO:HB3	1.92	0.51
1:A:185:TYR:HB3	1:A:212:HIS:CD2	2.46	0.51
1:C:1:ALA:HA	1:C:306:ARG:HG2	1.91	0.51
1:C:134:HIS:HB2	1:C:167:ARG:HD3	1.92	0.50
2:B:38:THR:HG21	2:B:42:ILE:HD11	1.92	0.50
1:A:301:ALA:O	1:A:305:ASN:HB2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:56:ARG:HE	1:A:60:GLN:NE2	2.08	0.50
2:B:23:ALA:O	2:B:24:GLN:HB2	2.11	0.50
2:B:91:VAL:HG11	2:B:94:LYS:NZ	2.27	0.49
1:A:214:SER:HB3	1:A:217:GLU:HG3	1.93	0.49
2:D:102:ARG:HA	2:D:125:PHE:O	2.12	0.49
1:A:161:GLY:HA3	1:A:228:THR:OG1	2.12	0.49
2:D:12:ILE:HG23	2:D:41:ARG:NH1	2.28	0.49
2:B:18:ILE:HD12	2:B:18:ILE:N	2.27	0.49
1:A:163:LEU:HG	1:A:188:ALA:HB2	1.94	0.49
2:B:21:ILE:HD13	2:B:58:LEU:HA	1.93	0.49
1:A:189:PRO:HD3	1:A:247:PHE:HE1	1.78	0.49
1:A:137:GLN:O	1:A:140:LEU:HG	2.13	0.49
1:C:257:ALA:HB1	1:C:261:MET:HE3	1.95	0.49
2:D:16:THR:OG1	2:D:65:PHE:HA	2.13	0.49
1:C:249:LEU:HD21	1:C:254:LEU:HD21	1.94	0.49
1:C:136:THR:HB	1:C:296:ARG:NH2	2.28	0.48
1:C:239:GLU:O	1:C:242:ASN:HB2	2.13	0.48
1:A:278:ASP:HB2	1:A:279:LYS:NZ	2.27	0.48
2:D:12:ILE:HD11	2:D:17:VAL:HG23	1.95	0.48
1:C:170:HIS:O	1:C:174:GLN:HG3	2.12	0.48
1:A:183:ARG:NH2	1:A:184:PHE:O	2.46	0.48
2:D:13:LYS:CG	2:D:88:ASN:HA	2.42	0.48
1:C:219:MET:HE3	1:C:254:LEU:HD23	1.96	0.48
2:B:40:GLN:NE2	2:B:63:ASN:HB2	2.28	0.48
2:B:18:ILE:HD12	2:B:18:ILE:H	1.78	0.48
2:D:22:PRO:O	2:D:25:ILE:HB	2.13	0.47
2:D:134:ILE:HB	2:D:147:HIS:HB3	1.96	0.47
2:D:130:ARG:CD	2:D:135:ALA:HB2	2.44	0.47
1:C:10:ILE:HD11	1:C:116:THR:CG2	2.45	0.47
2:B:14:ARG:HA	2:B:86:ILE:HG22	1.95	0.47
1:C:60:GLN:HG2	1:C:70:VAL:HG21	1.96	0.47
1:A:269:ARG:HA	1:A:272:GLU:OE2	2.15	0.47
1:A:134:HIS:CD2	1:A:137:GLN:HB2	2.50	0.46
1:C:10:ILE:O	1:C:135:PRO:HG3	2.16	0.46
1:C:160:VAL:HG22	1:C:187:ILE:HB	1.96	0.46
1:A:254:LEU:HD11	1:A:277:VAL:HG13	1.97	0.46
1:A:39:LEU:O	1:A:42:LYS:HB2	2.16	0.46
1:A:194:MET:SD	1:A:195:PRO:HD2	2.55	0.46
1:A:275:THR:O	1:A:279:LYS:NZ	2.46	0.46
1:A:158:ALA:HB2	1:A:222:VAL:HG11	1.98	0.46
1:C:20:LEU:HD11	1:C:178:LYS:HD2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:254:LEU:HD22	1:C:261:MET:CE	2.45	0.46
1:A:38:LEU:HD11	1:A:309:VAL:HG21	1.98	0.46
2:B:70:GLN:HA	2:B:73:GLN:OE1	2.16	0.46
1:A:187:ILE:HG12	1:A:212:HIS:HB2	1.97	0.46
2:D:61:ILE:CG2	2:D:64:THR:HB	2.46	0.46
1:C:164:LYS:HD3	1:C:165:TYR:CE2	2.51	0.46
2:B:136:LEU:HD12	2:B:150:VAL:HG11	1.98	0.46
1:A:250:ARG:HG2	1:A:274:ALA:CB	2.45	0.46
2:B:23:ALA:HA	2:B:57:ASP:CG	2.35	0.45
1:A:294:PHE:HA	1:A:297:GLN:HG2	1.97	0.45
1:A:61:THR:O	1:A:65:ARG:HG3	2.16	0.45
2:B:69:ASP:O	2:B:73:GLN:HG3	2.17	0.45
2:B:110:PRO:HG2	2:B:145:PHE:CG	2.51	0.45
2:D:23:ALA:HB3	2:D:55:ARG:NH2	2.31	0.45
1:A:269:ARG:NH1	1:A:278:ASP:OD2	2.49	0.45
1:C:226:TYR:OH	1:C:266:PRO:HG3	2.16	0.45
2:B:36:THR:HG23	2:D:46:LEU:HD13	1.97	0.45
2:D:100:PRO:HG2	2:D:103:ILE:HD13	1.98	0.45
1:A:156:HIS:HA	1:A:183:ARG:HB3	1.99	0.45
1:A:199:LEU:O	1:A:202:LEU:HB2	2.17	0.45
2:B:84:ASN:ND2	2:B:91:VAL:HG13	2.29	0.45
1:C:223:ASP:O	1:C:261:MET:HA	2.16	0.45
2:B:102:ARG:HA	2:B:125:PHE:O	2.17	0.45
1:C:183:ARG:NH2	1:C:184:PHE:O	2.50	0.45
2:B:16:THR:HG21	2:B:66:LEU:HG	1.99	0.45
1:A:164:LYS:HD3	1:A:165:TYR:CE2	2.51	0.45
1:C:49:PHE:CD2	1:C:76:SER:HB3	2.52	0.44
2:D:30:LEU:HD13	2:D:35:LEU:HB2	1.99	0.44
1:A:5:TYR:CZ	1:A:306:ARG:HG2	2.52	0.44
2:D:49:PRO:HA	2:D:54:GLY:O	2.17	0.44
1:A:96:SER:OG	1:A:119:SER:HA	2.17	0.44
2:B:137:LYS:HG3	2:B:144:GLU:HG2	1.99	0.44
2:B:99:LEU:HD23	2:B:99:LEU:HA	1.84	0.44
2:B:17:VAL:HG13	2:B:60:LYS:HG2	2.00	0.44
2:B:84:ASN:ND2	2:B:91:VAL:HG22	2.33	0.43
2:D:48:LEU:O	2:D:55:ARG:HA	2.19	0.43
2:D:12:ILE:HG22	2:D:62:GLU:HG3	1.99	0.43
1:A:40:LYS:O	1:A:41:HIS:HB2	2.18	0.43
1:A:4:LEU:HD23	1:A:7:LYS:CD	2.46	0.43
2:B:66:LEU:HB2	2:B:71:VAL:HG23	2.00	0.43
2:D:38:THR:HG21	2:D:42:ILE:HD11	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:108:GLN:HA	2:D:113:ASN:ND2	2.32	0.43
1:A:16:SER:O	1:A:20:LEU:HG	2.19	0.43
2:D:69:ASP:O	2:D:73:GLN:HG3	2.19	0.43
2:D:20:HIS:HB2	2:D:81:ALA:HA	2.01	0.43
1:C:34:PRO:O	1:C:36:PRO:HD3	2.19	0.43
1:A:48:PHE:CD2	1:A:105:ARG:HB2	2.54	0.43
2:D:13:LYS:HB2	2:D:89:TYR:OH	2.19	0.42
1:A:262:LYS:HA	1:A:262:LYS:HD3	1.90	0.42
2:D:129:LYS:N	2:D:129:LYS:HD2	2.35	0.42
1:C:277:VAL:O	1:C:283:ALA:HB2	2.19	0.42
1:A:8:HIS:ND1	1:A:123:PRO:HA	2.35	0.42
1:A:152:LEU:HA	1:A:155:LEU:HD11	2.01	0.42
2:D:18:ILE:O	2:D:58:LEU:HD12	2.20	0.42
2:B:127:VAL:HA	2:B:136:LEU:HD23	2.00	0.41
1:A:233:GLU:H	1:A:233:GLU:CD	2.23	0.41
1:A:252:SER:O	1:A:255:HIS:HB3	2.21	0.41
2:B:136:LEU:O	2:B:144:GLU:HA	2.19	0.41
2:D:61:ILE:HG22	2:D:64:THR:HB	2.02	0.41
2:B:110:PRO:HG2	2:B:145:PHE:CD2	2.54	0.41
1:A:226:TYR:CZ	1:A:266:PRO:HD3	2.55	0.41
2:D:125:PHE:HB3	2:D:136:LEU:HD22	2.02	0.41
2:B:28:LYS:O	2:B:32:LEU:HB2	2.20	0.41
1:A:50:GLU:HG3	1:A:107:PRO:HD3	2.02	0.41
1:C:109:GLU:OE2	2:D:113:ASN:HB3	2.20	0.41
2:B:111:ASN:O	2:B:117:HIS:CE1	2.74	0.41
1:A:197:TYR:HE2	2:B:143:LYS:HD3	1.85	0.41
1:A:45:ALA:HB2	1:A:99:VAL:HG21	2.02	0.41
1:C:160:VAL:HG11	1:C:215:ILE:HD11	2.02	0.41
1:A:54:ARG:HE	5:A:311:PCT:H1P2	1.86	0.41
2:D:21:ILE:HB	2:D:57:ASP:O	2.20	0.41
1:A:296:ARG:O	1:A:300:LEU:HG	2.21	0.41
2:B:30:LEU:HA	2:B:35:LEU:HD12	2.02	0.41
2:B:96:ARG:HA	2:B:97:PRO:HD3	1.89	0.41
1:A:237:PRO:HA	1:A:240:TYR:CD1	2.57	0.40
1:C:256:ASN:OD1	1:C:256:ASN:N	2.54	0.40
1:C:111:ALA:O	1:C:114:LEU:HB3	2.21	0.40
1:A:189:PRO:HD3	1:A:247:PHE:CE1	2.56	0.40
2:D:50:SER:N	2:D:54:GLY:O	2.54	0.40
2:D:16:THR:HG1	2:D:65:PHE:HA	1.87	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	308/310 (99%)	283 (92%)	23 (8%)	2 (1%)	30	65
1	C	308/310 (99%)	285 (92%)	19 (6%)	4 (1%)	15	44
2	B	144/153 (94%)	120 (83%)	18 (12%)	6 (4%)	3	11
2	D	144/153 (94%)	113 (78%)	25 (17%)	6 (4%)	3	11
All	All	904/926 (98%)	801 (89%)	85 (9%)	18 (2%)	9	30

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	34	LYS
2	D	105	ASN
2	B	24	GLN
2	B	132	ASN
2	D	54	GLY
2	B	9	VAL
2	B	88	ASN
2	B	47	ASN
1	C	154	ASN
2	D	48	LEU
2	D	133	ASP
1	A	41	HIS
2	D	13	LYS
1	C	132	ASN
2	D	49	PRO
1	C	237	PRO
1	A	166	GLY
1	C	270	VAL

### 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	261/261 (100%)	231 (88%)	30 (12%)	7	21
1	C	261/261 (100%)	242 (93%)	19 (7%)	17	44
2	B	129/136 (95%)	109 (84%)	20 (16%)	3	10
2	D	129/136 (95%)	108 (84%)	21 (16%)	3	8
All	All	780/794 (98%)	690 (88%)	90 (12%)	7	21

All (90) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ASP
1	A	34	PRO
1	A	40	LYS
1	A	42	LYS
1	A	59	PHE
1	A	65	ARG
1	A	74	SER
1	A	75	ASP
1	A	80	SER
1	A	105	ARG
1	A	108	GLN
1	A	113	ARG
1	A	117	GLU
1	A	125	LEU
1	A	134	HIS
1	A	167	ARG
1	A	190	ASP
1	A	214	SER
1	A	221	GLU
1	A	237	PRO
1	A	242	ASN
1	A	246	GLN
1	A	252	SER
1	A	253	ASP

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Mol	Chain	Res	Type
1	A	264	LEU
1	A	269	ARG
1	A	279	LYS
1	A	285	TYR
1	A	296	ARG
1	A	308	LEU
2	B	12	ILE
2	B	19	ASP
2	B	24	GLN
2	B	32	LEU
2	B	53	MET
2	B	61	ILE
2	B	63	ASN
2	B	72	ASP
2	B	76	LEU
2	B	82	THR
2	B	84	ASN
2	B	85	ARG
2	B	88	ASN
2	B	95	SER
2	B	98	SER
2	B	103	ILE
2	B	104	ASP
2	B	128	ARG
2	B	132	ASN
2	B	153	ASN
1	C	2	ASN
1	C	17	ARG
1	C	38	LEU
1	C	59	PHE
1	C	69	SER
1	C	71	VAL
1	C	81	LEU
1	C	146	GLN
1	C	167	ARG
1	C	213	SER
1	C	221	GLU
1	C	224	ILE
1	C	232	LYS
1	C	237	PRO
1	C	258	LYS
1	C	262	LYS

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Mol	Chain	Res	Type
1	C	269	ARG
1	C	285	TYR
1	C	307	ASP
2	D	30	LEU
2	D	34	LYS
2	D	36	THR
2	D	38	THR
2	D	40	GLN
2	D	41	ARG
2	D	43	THR
2	D	52	GLU
2	D	55	ARG
2	D	61	ILE
2	D	62	GLU
2	D	76	LEU
2	D	79	PRO
2	D	82	THR
2	D	102	ARG
2	D	112	SER
2	D	123	SER
2	D	128	ARG
2	D	129	LYS
2	D	137	LYS
2	D	153	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	GLN
1	A	154	ASN
1	A	231	GLN
2	B	63	ASN
2	B	117	HIS
2	B	153	ASN
1	C	13	ASN
1	C	60	GLN
1	C	154	ASN
1	C	305	ASN
2	D	117	HIS

### 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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	$\# Z  > 2$
5	PCT	A	311	-	7,7,7	3.01	3 (42%)	9,10,10	2.03	5 (55%)
3	MLI	A	312	-	0,6,6	0.00	-	0,7,7	0.00	-
5	PCT	C	311	-	7,7,7	3.13	4 (57%)	9,10,10	1.79	2 (22%)
3	MLI	C	312	-	0,6,6	0.00	-	0,7,7	0.00	-

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
5	PCT	A	311	-	-	0/4/5/5	0/0/0/0
3	MLI	A	312	-	-	0/0/4/4	0/0/0/0
5	PCT	C	311	-	-	0/4/5/5	0/0/0/0
3	MLI	C	312	-	-	0/0/4/4	0/0/0/0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	311	PCT	P-O2P	2.15	1.60	1.54
5	A	311	PCT	C1P-C1	2.52	1.54	1.51
5	C	311	PCT	P-O3P	2.95	1.62	1.54
5	A	311	PCT	P-O1P	4.49	1.60	1.50
5	C	311	PCT	P-O1P	4.99	1.61	1.50
5	C	311	PCT	P-C1P	5.11	1.87	1.79
5	A	311	PCT	P-C1P	5.59	1.87	1.79

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	311	PCT	O1P-P-C1P	-3.00	103.54	110.92
5	C	311	PCT	O1P-P-C1P	-2.56	104.62	110.92
5	A	311	PCT	O3P-P-C1P	2.11	111.66	107.01
5	A	311	PCT	O1-C1-C1P	2.18	122.43	119.90
5	A	311	PCT	C1P-C1-N1	2.46	118.23	115.25
5	A	311	PCT	O3P-P-O2P	2.81	116.37	108.13
5	C	311	PCT	O1-C1-C1P	3.63	124.11	119.90

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	311	PCT	1	0
5	C	311	PCT	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

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

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.