



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:41 AM GMT

PDB ID : 2I4N
Title : Rhodopseudomonas palustris prolyl-tRNA synthetase in complex with CysAMS
Authors : Crepin, T.; Yaremchuk, A.; Tukalo, M.; Cusack, S.
Deposited on : 2006-08-22
Resolution : 2.85 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

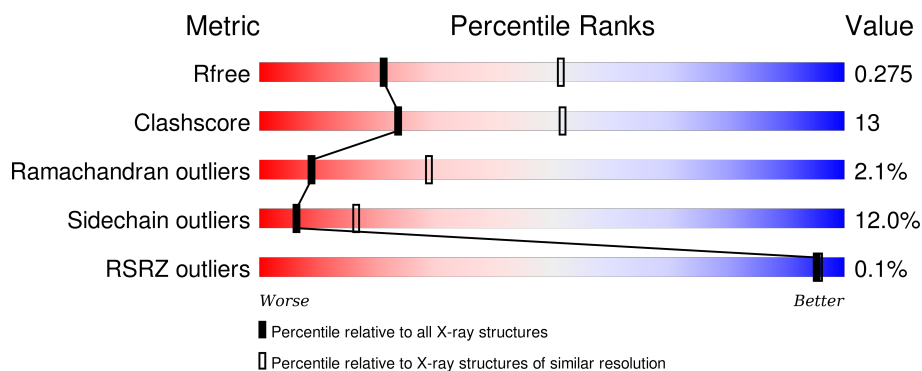
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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}	91344	2228 (2.90-2.82)
Clashscore	102246	2499 (2.90-2.82)
Ramachandran outliers	100387	2439 (2.90-2.82)
Sidechain outliers	100360	2442 (2.90-2.82)
RSRZ outliers	91569	2236 (2.90-2.82)

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 ≥ 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.

Mol	Chain	Length	Quality of chain
1	A	458	
1	B	458	
1	C	458	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10672 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 Proline-tRNA ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	441	Total	C	N	O	S	0	0	0
			3495	2211	620	650	14			
1	B	442	Total	C	N	O	S	0	0	0
			3506	2217	624	651	14			
1	C	442	Total	C	N	O	S	0	0	0
			3506	2217	624	651	14			

There are 60 discrepancies between the modelled and reference sequences:

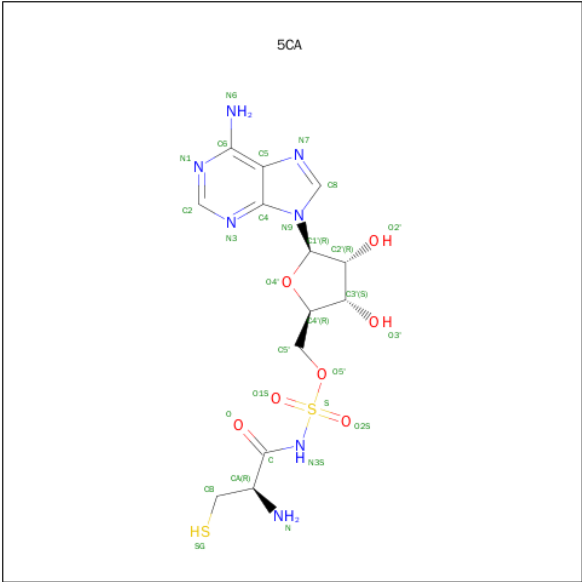
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP Q6N5P6
A	-18	GLY	-	EXPRESSION TAG	UNP Q6N5P6
A	-17	SER	-	EXPRESSION TAG	UNP Q6N5P6
A	-16	SER	-	EXPRESSION TAG	UNP Q6N5P6
A	-15	HIS	-	EXPRESSION TAG	UNP Q6N5P6
A	-14	HIS	-	EXPRESSION TAG	UNP Q6N5P6
A	-13	HIS	-	EXPRESSION TAG	UNP Q6N5P6
A	-12	HIS	-	EXPRESSION TAG	UNP Q6N5P6
A	-11	HIS	-	EXPRESSION TAG	UNP Q6N5P6
A	-10	HIS	-	EXPRESSION TAG	UNP Q6N5P6
A	-9	SER	-	EXPRESSION TAG	UNP Q6N5P6
A	-8	SER	-	EXPRESSION TAG	UNP Q6N5P6
A	-7	GLY	-	EXPRESSION TAG	UNP Q6N5P6
A	-6	LEU	-	EXPRESSION TAG	UNP Q6N5P6
A	-5	VAL	-	EXPRESSION TAG	UNP Q6N5P6
A	-4	PRO	-	EXPRESSION TAG	UNP Q6N5P6
A	-3	ARG	-	EXPRESSION TAG	UNP Q6N5P6
A	-2	GLY	-	EXPRESSION TAG	UNP Q6N5P6
A	-1	SER	-	EXPRESSION TAG	UNP Q6N5P6
A	0	HIS	-	EXPRESSION TAG	UNP Q6N5P6
B	-19	MET	-	EXPRESSION TAG	UNP Q6N5P6
B	-18	GLY	-	EXPRESSION TAG	UNP Q6N5P6
B	-17	SER	-	EXPRESSION TAG	UNP Q6N5P6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-16	SER	-	EXPRESSION TAG	UNP Q6N5P6
B	-15	HIS	-	EXPRESSION TAG	UNP Q6N5P6
B	-14	HIS	-	EXPRESSION TAG	UNP Q6N5P6
B	-13	HIS	-	EXPRESSION TAG	UNP Q6N5P6
B	-12	HIS	-	EXPRESSION TAG	UNP Q6N5P6
B	-11	HIS	-	EXPRESSION TAG	UNP Q6N5P6
B	-10	HIS	-	EXPRESSION TAG	UNP Q6N5P6
B	-9	SER	-	EXPRESSION TAG	UNP Q6N5P6
B	-8	SER	-	EXPRESSION TAG	UNP Q6N5P6
B	-7	GLY	-	EXPRESSION TAG	UNP Q6N5P6
B	-6	LEU	-	EXPRESSION TAG	UNP Q6N5P6
B	-5	VAL	-	EXPRESSION TAG	UNP Q6N5P6
B	-4	PRO	-	EXPRESSION TAG	UNP Q6N5P6
B	-3	ARG	-	EXPRESSION TAG	UNP Q6N5P6
B	-2	GLY	-	EXPRESSION TAG	UNP Q6N5P6
B	-1	SER	-	EXPRESSION TAG	UNP Q6N5P6
B	0	HIS	-	EXPRESSION TAG	UNP Q6N5P6
C	-19	MET	-	EXPRESSION TAG	UNP Q6N5P6
C	-18	GLY	-	EXPRESSION TAG	UNP Q6N5P6
C	-17	SER	-	EXPRESSION TAG	UNP Q6N5P6
C	-16	SER	-	EXPRESSION TAG	UNP Q6N5P6
C	-15	HIS	-	EXPRESSION TAG	UNP Q6N5P6
C	-14	HIS	-	EXPRESSION TAG	UNP Q6N5P6
C	-13	HIS	-	EXPRESSION TAG	UNP Q6N5P6
C	-12	HIS	-	EXPRESSION TAG	UNP Q6N5P6
C	-11	HIS	-	EXPRESSION TAG	UNP Q6N5P6
C	-10	HIS	-	EXPRESSION TAG	UNP Q6N5P6
C	-9	SER	-	EXPRESSION TAG	UNP Q6N5P6
C	-8	SER	-	EXPRESSION TAG	UNP Q6N5P6
C	-7	GLY	-	EXPRESSION TAG	UNP Q6N5P6
C	-6	LEU	-	EXPRESSION TAG	UNP Q6N5P6
C	-5	VAL	-	EXPRESSION TAG	UNP Q6N5P6
C	-4	PRO	-	EXPRESSION TAG	UNP Q6N5P6
C	-3	ARG	-	EXPRESSION TAG	UNP Q6N5P6
C	-2	GLY	-	EXPRESSION TAG	UNP Q6N5P6
C	-1	SER	-	EXPRESSION TAG	UNP Q6N5P6
C	0	HIS	-	EXPRESSION TAG	UNP Q6N5P6

- Molecule 2 is 5'-O-(N-(L-CYSTEINYL)-SULFAMOYL)ADENOSINE (three-letter code: 5CA) (formula: C₁₃H₁₉N₇O₇S₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			29	13	7	7	2		
2	B	1	Total	C	N	O	S	0	0
			29	13	7	7	2		
2	C	1	Total	C	N	O	S	0	0
			29	13	7	7	2		

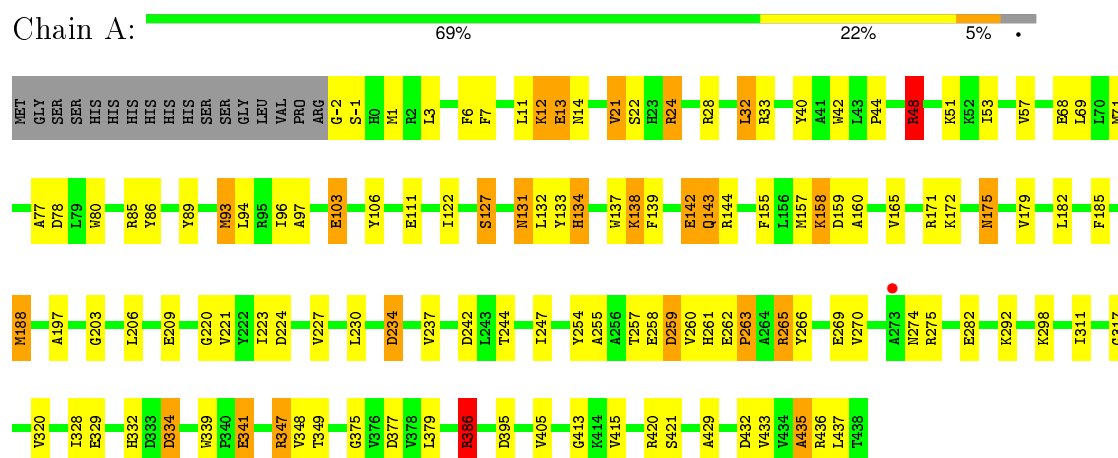
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	28	Total	O	0	0
			28	28		
3	B	30	Total	O	0	0
			30	30		
3	C	20	Total	O	0	0
			20	20		

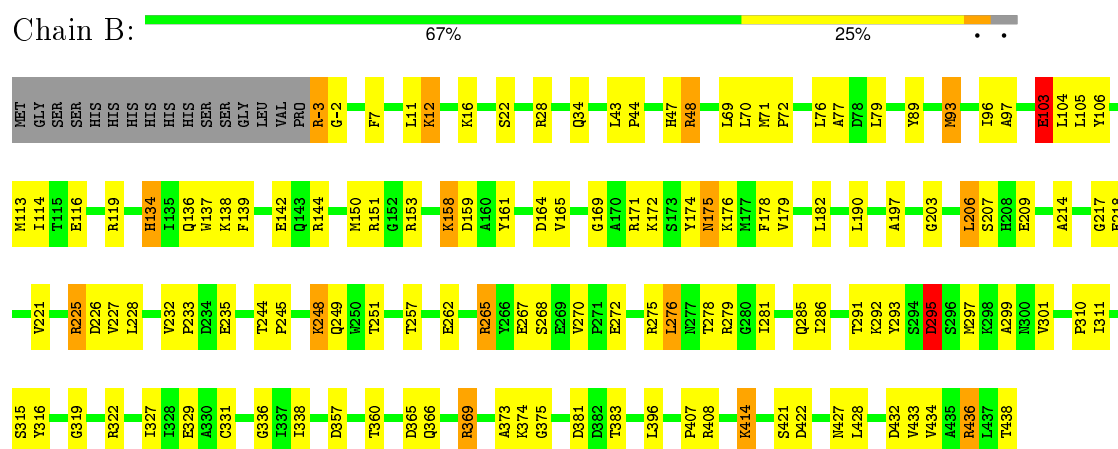
3 Residue-property plots

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.

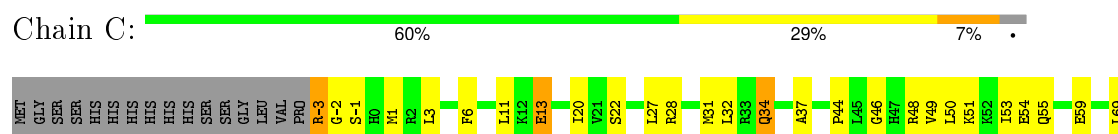
• Molecule 1: Proline-tRNA ligase

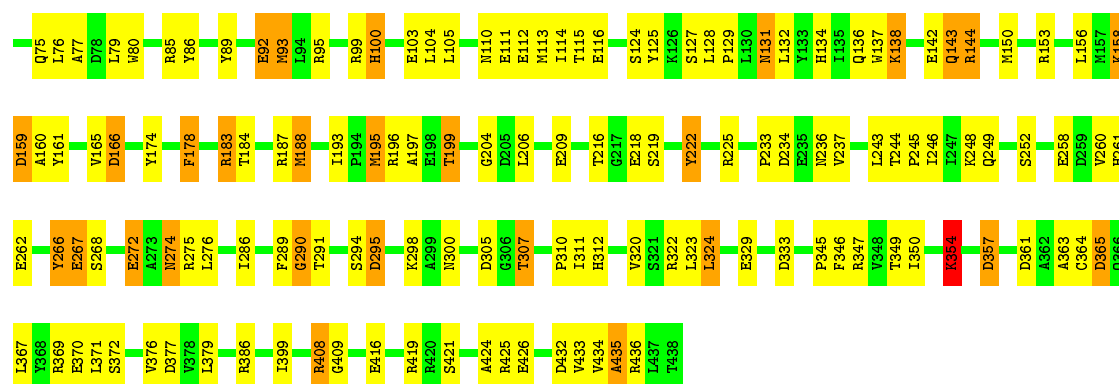


• Molecule 1: Proline-tRNA ligase



• Molecule 1: Proline-tRNA ligase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	217.90Å 107.47Å 110.66Å 90.00° 120.43° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.85 29.84 – 2.70	Depositor EDS
% Data completeness (in resolution range)	94.6 ((Not available)-2.85) 94.6 (29.84-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.01	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.20 (at 2.68Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.211 , 0.277 0.216 , 0.275	Depositor DCC
R_{free} test set	2903 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	63.7	Xtriage
Anisotropy	0.144	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 33.6	EDS
Estimated twinning fraction	0.047 for -h+k-l,-l,-k 0.034 for -h-k-l,l,k 0.064 for -h-2*l,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	0 of 57190 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10672	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 18.69% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 5CA

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	0.63	0/3570	0.76	2/4828 (0.0%)
1	B	0.63	0/3581	0.75	1/4842 (0.0%)
1	C	0.64	0/3581	0.74	0/4842
All	All	0.64	0/10732	0.75	3/14512 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	48	ARG	NE-CZ-NH1	5.22	122.91	120.30
1	B	79	LEU	CA-CB-CG	5.12	127.08	115.30
1	A	386	ARG	NE-CZ-NH1	5.09	122.84	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3495	0	3466	94	0
1	B	3506	0	3479	89	0
1	C	3506	0	3479	99	0
2	A	29	0	19	1	0
2	B	29	0	19	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	29	0	19	1	0
3	A	28	0	0	3	0
3	B	30	0	0	0	0
3	C	20	0	0	1	0
All	All	10672	0	10481	278	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 (278) 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:C:-3:ARG:HH11	1:C:-3:ARG:HG3	1.10	1.13
1:B:436:ARG:HH11	1:B:436:ARG:HG2	0.95	1.09
1:B:48:ARG:HH11	1:B:48:ARG:HG2	1.18	1.00
1:A:-2:GLY:HA3	1:A:375:GLY:HA2	1.48	0.96
1:B:175:ASN:HD22	1:B:175:ASN:N	1.63	0.96
1:A:12:LYS:HD3	1:A:12:LYS:H	1.29	0.95
1:A:12:LYS:N	1:A:12:LYS:HD3	1.84	0.92
1:C:188:MET:CE	1:C:323:LEU:HD13	2.00	0.92
1:B:436:ARG:NH1	1:B:436:ARG:HG2	1.74	0.92
1:B:436:ARG:HH11	1:B:436:ARG:CG	1.84	0.91
1:C:-3:ARG:NH1	1:C:-3:ARG:HG3	1.86	0.88
1:C:199:THR:HG21	1:C:204:GLY:O	1.72	0.88
1:B:175:ASN:HD22	1:B:175:ASN:H	1.18	0.86
1:A:259:ASP:O	1:A:261:HIS:N	2.08	0.86
1:C:272:GLU:HA	1:C:275:ARG:HD2	1.59	0.85
1:A:386:ARG:HH11	1:A:386:ARG:H	1.22	0.85
1:B:72:PRO:HD2	1:B:113:MET:HE1	1.59	0.84
1:C:-3:ARG:HH11	1:C:-3:ARG:CG	1.89	0.83
1:B:48:ARG:HH11	1:B:48:ARG:CG	1.91	0.83
1:C:188:MET:HE3	1:C:323:LEU:HD13	1.62	0.80
1:B:171:ARG:O	1:B:175:ASN:ND2	2.13	0.80
1:B:22:SER:HB3	1:B:329:GLU:HG2	1.63	0.79
1:B:175:ASN:H	1:B:175:ASN:ND2	1.80	0.79
1:A:48:ARG:CG	1:A:48:ARG:HH11	1.96	0.78
1:C:76:LEU:HD23	1:C:104:LEU:HD23	1.63	0.78
1:B:293:TYR:O	1:B:297:MET:HB2	1.85	0.77
1:B:267:GLU:HA	1:B:275:ARG:HH12	1.49	0.76
1:C:158:LYS:HD2	1:C:158:LYS:C	2.06	0.76
1:A:435:ALA:O	1:A:437:LEU:N	2.19	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:LEU:HD11	1:A:40:TYR:HB3	1.69	0.73
1:A:48:ARG:HG2	1:A:48:ARG:HH11	1.53	0.73
1:A:14:ASN:HD22	1:A:24:ARG:HG3	1.53	0.73
1:C:138:LYS:NZ	1:C:159:ASP:OD1	2.21	0.73
1:C:305:ASP:OD1	1:C:307:THR:HG23	1.88	0.72
1:A:265:ARG:NH2	1:A:269:GLU:OE1	2.21	0.71
1:B:225:ARG:HH11	1:B:225:ARG:HG3	1.53	0.71
1:A:259:ASP:C	1:A:261:HIS:H	1.93	0.71
1:A:22:SER:HB3	1:A:329:GLU:HG2	1.72	0.70
1:A:-2:GLY:CA	1:A:375:GLY:HA2	2.22	0.70
1:B:72:PRO:HD2	1:B:113:MET:CE	2.23	0.69
1:B:71:MET:HB3	1:B:113:MET:CE	2.25	0.66
1:B:175:ASN:N	1:B:175:ASN:ND2	2.35	0.66
1:A:220:GLY:O	1:A:258:GLU:HB2	1.96	0.66
1:A:386:ARG:NH1	1:A:386:ARG:H	1.93	0.66
1:C:188:MET:HE1	1:C:323:LEU:HD13	1.78	0.65
1:A:11:LEU:HB2	1:A:28:ARG:HA	1.79	0.64
1:A:234:ASP:O	1:A:237:VAL:HG23	1.96	0.64
1:A:175:ASN:O	1:A:179:VAL:HG23	1.98	0.64
1:A:68:GLU:HB2	1:A:133:TYR:CZ	2.34	0.63
1:C:131:ASN:HD22	1:C:131:ASN:C	2.01	0.63
1:C:51:LYS:HD2	1:C:54:GLU:OE2	1.99	0.63
1:B:76:LEU:HD23	1:B:104:LEU:HD22	1.81	0.62
1:B:77:ALA:HB3	1:B:103:GLU:HB2	1.81	0.62
1:A:138:LYS:HD3	1:A:157:MET:HB3	1.81	0.62
1:C:408:ARG:HG3	1:C:409:GLY:H	1.64	0.62
1:C:158:LYS:HD2	1:C:158:LYS:O	2.00	0.62
1:B:71:MET:HB3	1:B:113:MET:HE3	1.82	0.62
1:C:432:ASP:O	1:C:435:ALA:HB2	1.99	0.61
1:A:224:ASP:HB2	1:A:274:ASN:HB3	1.81	0.61
1:B:272:GLU:HA	1:B:275:ARG:HE	1.64	0.61
1:C:294:SER:HB3	1:C:311:ILE:HG13	1.82	0.61
1:B:267:GLU:HA	1:B:275:ARG:NH1	2.15	0.61
1:B:207:SER:HA	1:B:286:ILE:O	2.00	0.61
1:A:89:TYR:HB2	1:A:93:MET:HG3	1.83	0.61
1:A:-1:SER:O	1:A:377:ASP:HA	2.00	0.61
1:B:138:LYS:HE3	1:B:159:ASP:OD1	2.01	0.61
1:C:354:LYS:HE2	1:C:386:ARG:HH12	1.66	0.60
1:A:257:THR:O	1:A:261:HIS:HB2	2.01	0.59
1:A:21:VAL:HA	1:A:24:ARG:HD2	1.83	0.59
1:C:11:LEU:HB2	1:C:28:ARG:HA	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:ILE:HG22	1:A:255:ALA:HA	1.84	0.59
1:B:48:ARG:NH1	1:B:48:ARG:HG2	1.97	0.59
1:B:365:ASP:O	1:B:369:ARG:HD3	2.03	0.59
1:C:77:ALA:HB3	1:C:103:GLU:HB3	1.84	0.59
1:A:106:TYR:CD2	1:A:139:PHE:HB2	2.38	0.58
1:B:267:GLU:CA	1:B:275:ARG:HH12	2.15	0.58
1:C:363:ALA:O	1:C:367:LEU:HB2	2.02	0.58
1:A:224:ASP:HB3	1:A:270:VAL:HG11	1.85	0.57
1:C:99:ARG:HG3	1:C:100:HIS:ND1	2.20	0.57
1:A:158:LYS:O	1:A:317:GLY:HA2	2.04	0.57
1:C:76:LEU:HD23	1:C:104:LEU:CD2	2.34	0.57
1:C:44:PRO:O	1:C:48:ARG:HG3	2.05	0.57
1:C:361:ASP:O	1:C:365:ASP:HB2	2.06	0.56
1:C:196:ARG:HB3	3:C:451:HOH:O	2.05	0.56
1:C:349:THR:HA	1:C:379:LEU:O	2.05	0.56
1:A:395:ASP:HA	1:A:420:ARG:HE	1.71	0.56
1:A:197:ALA:HB2	1:A:209:GLU:HB2	1.87	0.56
1:C:433:VAL:C	1:C:435:ALA:H	2.08	0.56
1:A:265:ARG:O	1:A:265:ARG:HD3	2.06	0.55
1:C:346:PHE:O	1:C:399:ILE:HG23	2.06	0.55
1:B:48:ARG:NH1	1:B:48:ARG:CG	2.58	0.55
1:A:1:MET:O	1:A:379:LEU:HD12	2.06	0.55
1:A:421:SER:HB3	3:A:453:HOH:O	2.07	0.55
1:C:225:ARG:HG3	1:C:274:ASN:O	2.07	0.55
1:C:-1:SER:O	1:C:377:ASP:HA	2.06	0.54
1:B:217:GLY:O	1:B:279:ARG:NH1	2.35	0.54
1:B:226:ASP:C	1:B:228:LEU:H	2.11	0.54
1:A:137:TRP:CZ2	1:B:70:LEU:HD21	2.41	0.54
1:B:244:THR:HB	1:B:245:PRO:HD3	1.88	0.54
1:A:263:PRO:C	1:A:265:ARG:H	2.11	0.54
1:B:144:ARG:NH1	1:B:144:ARG:HB2	2.22	0.54
1:B:262:GLU:HB3	1:B:265:ARG:HB2	1.89	0.54
1:C:49:VAL:HG22	1:C:345:PRO:HD3	1.90	0.54
1:A:185:PHE:HA	1:A:188:MET:HG3	1.90	0.54
1:A:311:ILE:HD12	1:A:311:ILE:C	2.28	0.53
1:C:294:SER:CB	1:C:311:ILE:HG13	2.39	0.53
1:C:234:ASP:O	1:C:237:VAL:HG23	2.09	0.53
1:C:1:MET:HG2	1:C:6:PHE:CD1	2.43	0.53
1:C:93:MET:SD	1:C:105:LEU:HD22	2.48	0.53
1:A:7:PHE:HB3	1:A:44:PRO:HG2	1.91	0.53
1:A:12:LYS:N	1:A:12:LYS:CD	2.66	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:ASN:C	1:A:131:ASN:HD22	2.12	0.52
1:B:139:PHE:HA	1:B:153:ARG:O	2.08	0.52
1:A:53:ILE:HG22	1:A:320:VAL:HG13	1.91	0.52
1:B:11:LEU:HB2	1:B:28:ARG:HA	1.91	0.52
1:B:214:ALA:HB2	1:B:281:ILE:HD12	1.91	0.52
1:A:429:ALA:O	1:A:432:ASP:HB2	2.08	0.52
1:C:75:GLN:HG2	1:C:113:MET:CE	2.40	0.52
1:C:89:TYR:O	1:C:92:GLU:HB2	2.09	0.51
1:C:291:THR:O	1:C:295:ASP:HB2	2.10	0.51
1:B:179:VAL:HG13	1:B:233:PRO:HG2	1.93	0.51
1:A:421:SER:CB	3:A:453:HOH:O	2.59	0.51
1:C:246:ILE:HA	1:C:249:GLN:HB3	1.92	0.50
1:C:93:MET:HE3	1:C:93:MET:HA	1.93	0.50
1:A:203:GLY:HA3	1:A:292:LYS:HE3	1.92	0.50
1:A:334:ASP:OD2	1:A:334:ASP:N	2.43	0.50
1:C:244:THR:O	1:C:248:LYS:HG3	2.12	0.50
1:A:32:LEU:HD12	1:A:33:ARG:N	2.26	0.50
2:B:439:5CA:O	2:B:439:5CA:O1S	2.28	0.50
1:C:110:ASN:O	1:C:114:ILE:HG22	2.12	0.50
1:C:22:SER:HB3	1:C:329:GLU:HG2	1.94	0.50
1:C:142:GLU:H	1:C:153:ARG:HG2	1.77	0.50
1:A:-1:SER:HB3	1:A:377:ASP:OD1	2.12	0.49
1:C:31:MET:HE3	1:C:46:GLY:HA2	1.94	0.49
1:A:386:ARG:N	1:A:386:ARG:HH11	1.99	0.49
1:B:144:ARG:HH11	1:B:144:ARG:HB2	1.77	0.49
1:C:183:ARG:HD2	1:C:233:PRO:O	2.12	0.49
1:C:34:GLN:OE1	1:C:37:ALA:HA	2.11	0.49
1:C:233:PRO:HB3	1:C:237:VAL:HG21	1.93	0.49
1:C:89:TYR:HB2	1:C:93:MET:HG2	1.95	0.49
1:C:50:LEU:HD13	1:C:324:LEU:CD1	2.43	0.48
1:B:291:THR:O	1:B:295:ASP:HB2	2.14	0.48
1:A:13:GLU:CD	1:A:13:GLU:H	2.16	0.48
1:C:-2:GLY:HA3	1:C:372:SER:O	2.14	0.48
1:B:114:ILE:HG21	1:B:161:TYR:CG	2.48	0.48
1:B:209:GLU:HB3	1:B:285:GLN:HG3	1.96	0.48
1:B:299:ALA:O	1:B:310:PRO:HA	2.13	0.48
1:A:48:ARG:NH1	1:A:48:ARG:HG2	2.23	0.48
1:C:128:LEU:HB3	1:C:129:PRO:HA	1.94	0.48
1:B:373:ALA:C	1:B:375:GLY:H	2.16	0.48
1:B:336:GLY:HA3	1:B:396:LEU:HA	1.95	0.48
1:B:174:TYR:CZ	1:B:206:LEU:HD22	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:131:ASN:HD21	1:C:160:ALA:HB1	1.79	0.48
1:C:416:GLU:OE2	1:C:425:ARG:HD3	2.14	0.48
1:C:174:TYR:CE2	1:C:286:ILE:HD12	2.48	0.47
1:B:244:THR:HG22	1:B:248:LYS:HD3	1.95	0.47
1:B:76:LEU:HD23	1:B:104:LEU:CD2	2.43	0.47
1:A:122:ILE:HD12	1:A:127:SER:OG	2.13	0.47
1:B:217:GLY:HA3	1:B:279:ARG:HB3	1.95	0.47
1:A:311:ILE:HD12	1:A:311:ILE:O	2.15	0.47
1:C:93:MET:HE1	1:C:105:LEU:HB2	1.96	0.47
1:B:174:TYR:OH	1:B:206:LEU:HD22	2.15	0.47
1:B:221:VAL:HG22	1:B:278:THR:O	2.15	0.47
1:A:242:ASP:OD1	1:A:244:THR:HG22	2.15	0.47
1:A:131:ASN:HD21	1:A:160:ALA:HB1	1.79	0.47
1:C:159:ASP:OD1	1:C:159:ASP:N	2.47	0.47
1:C:13:GLU:H	1:C:13:GLU:CD	2.18	0.47
1:A:96:ILE:HG22	1:A:97:ALA:N	2.30	0.47
1:B:96:ILE:HG22	1:B:97:ALA:N	2.29	0.47
1:C:136:GLN:HG3	1:C:137:TRP:N	2.30	0.47
1:C:50:LEU:CD1	1:C:320:VAL:HG12	2.46	0.46
1:C:20:ILE:HD12	1:C:329:GLU:OE2	2.15	0.46
1:A:341:GLU:OE1	1:A:347:ARG:NH1	2.48	0.46
1:B:182:LEU:HB3	1:B:232:VAL:HG13	1.97	0.46
1:B:275:ARG:O	1:B:276:LEU:C	2.54	0.46
1:C:323:LEU:O	1:C:324:LEU:C	2.53	0.46
1:C:289:PHE:O	1:C:290:GLY:C	2.54	0.46
1:B:93:MET:HE1	1:B:105:LEU:HD13	1.97	0.46
1:A:254:TYR:CE2	1:A:266:TYR:HD2	2.34	0.46
1:B:311:ILE:C	1:B:311:ILE:HD12	2.36	0.46
1:B:432:ASP:O	1:B:433:VAL:C	2.54	0.46
1:C:137:TRP:CE2	1:C:156:LEU:HD13	2.51	0.45
1:A:103:GLU:HG3	3:A:450:HOH:O	2.16	0.45
1:C:244:THR:N	1:C:245:PRO:HD2	2.31	0.45
1:A:413:GLY:O	1:A:429:ALA:HA	2.16	0.45
1:A:48:ARG:HG3	1:A:48:ARG:HH11	1.80	0.45
1:C:85:ARG:O	1:C:86:TYR:C	2.53	0.45
1:C:346:PHE:HA	1:C:377:ASP:HB3	1.99	0.45
1:B:357:ASP:HB3	1:B:360:THR:OG1	2.17	0.45
1:A:22:SER:HA	1:A:328:ILE:HG21	1.98	0.45
1:C:77:ALA:O	1:C:80:TRP:N	2.35	0.45
1:B:190:LEU:HD21	1:B:327:ILE:HA	1.99	0.45
1:A:111:GLU:OE2	2:A:439:5CA:N	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:TRP:HZ2	1:B:70:LEU:HD21	1.82	0.45
1:C:93:MET:HE3	1:C:93:MET:CA	2.47	0.45
1:C:183:ARG:O	1:C:184:THR:C	2.56	0.44
1:B:197:ALA:HB2	1:B:209:GLU:HG2	1.98	0.44
1:A:175:ASN:OD1	1:A:247:ILE:HG13	2.17	0.44
1:A:94:LEU:HD12	1:A:94:LEU:HA	1.73	0.44
1:A:80:TRP:HB3	1:A:86:TYR:HA	1.99	0.44
1:B:151:ARG:HD3	1:B:218:GLU:OE2	2.17	0.44
1:A:77:ALA:HB2	1:A:103:GLU:HB2	1.99	0.44
1:B:267:GLU:HG2	1:B:275:ARG:HH12	1.82	0.44
1:B:136:GLN:HG3	1:B:137:TRP:N	2.33	0.44
1:A:227:VAL:O	1:A:230:LEU:HD12	2.18	0.44
1:B:150:MET:O	1:B:322:ARG:NH1	2.51	0.44
1:B:169:GLY:O	1:B:172:LYS:HB2	2.17	0.44
1:A:68:GLU:HG2	1:B:47:HIS:CD2	2.54	0.43
1:C:110:ASN:HB2	1:C:161:TYR:OH	2.18	0.43
1:C:354:LYS:O	1:C:357:ASP:HB3	2.19	0.43
1:C:188:MET:HE1	1:C:323:LEU:CD1	2.46	0.43
1:C:114:ILE:HD11	1:C:132:LEU:HD12	2.01	0.43
1:B:7:PHE:HB3	1:B:44:PRO:HG2	2.00	0.43
1:B:381:ASP:OD1	1:B:383:THR:OG1	2.30	0.43
1:A:42:TRP:CD1	1:A:42:TRP:N	2.86	0.43
1:B:315:SER:C	1:B:316:TYR:CG	2.91	0.43
1:B:69:LEU:O	1:B:134:HIS:HB3	2.19	0.43
1:B:225:ARG:NH1	1:B:225:ARG:HG3	2.28	0.43
1:A:71:MET:HB2	1:A:134:HIS:CE1	2.53	0.43
1:A:348:VAL:HG22	1:A:349:THR:N	2.34	0.43
1:B:209:GLU:OE2	1:B:257:THR:HG21	2.18	0.42
1:C:55:GLN:NE2	1:C:59:GLU:OE2	2.45	0.42
1:C:193:ILE:HB	1:C:195:MET:HE3	2.00	0.42
1:C:419:ARG:HB3	1:C:424:ALA:HB3	2.00	0.42
1:A:262:GLU:HA	1:A:263:PRO:HD3	1.82	0.42
1:B:225:ARG:HH11	1:B:225:ARG:CG	2.28	0.42
1:A:254:TYR:HB3	1:A:270:VAL:HG21	2.02	0.42
1:C:131:ASN:C	1:C:131:ASN:ND2	2.72	0.42
1:C:244:THR:N	1:C:245:PRO:CD	2.82	0.42
1:C:27:LEU:HD23	1:C:32:LEU:HB3	2.01	0.42
1:A:405:VAL:HG22	1:A:415:VAL:HG11	2.01	0.42
1:A:142:GLU:O	1:A:143:GLN:C	2.57	0.42
1:C:112:GLU:O	1:C:115:THR:HB	2.19	0.42
1:B:203:GLY:HA3	1:B:292:LYS:HD2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:339:TRP:CE3	1:A:339:TRP:HA	2.55	0.42
1:B:-3:ARG:N	1:B:-3:ARG:HE	2.18	0.42
1:A:171:ARG:O	1:A:175:ASN:ND2	2.53	0.41
1:C:347:ARG:H	1:C:377:ASP:HB2	1.85	0.41
1:A:137:TRP:HA	1:A:155:PHE:O	2.20	0.41
1:A:137:TRP:HZ2	1:B:70:LEU:CD2	2.33	0.41
1:B:331:CYS:HB3	1:B:338:ILE:O	2.20	0.41
1:A:57:VAL:HG11	1:A:158:LYS:HG3	2.02	0.41
1:C:222:TYR:CD2	1:C:222:TYR:N	2.88	0.41
1:B:436:ARG:NH1	1:B:436:ARG:CG	2.54	0.41
1:C:350:ILE:HD11	1:C:371:LEU:HD12	2.03	0.41
1:B:175:ASN:O	1:B:179:VAL:HG23	2.20	0.41
1:C:197:ALA:HB2	1:C:209:GLU:HB2	2.03	0.41
1:C:178:PHE:C	1:C:178:PHE:CD2	2.93	0.41
1:B:434:VAL:O	1:B:438:THR:HG23	2.21	0.41
1:A:22:SER:HA	1:A:328:ILE:CG2	2.51	0.41
1:C:261:HIS:CD2	1:C:261:HIS:C	2.94	0.41
1:A:234:ASP:OD1	1:A:234:ASP:C	2.60	0.41
1:C:131:ASN:HD22	1:C:132:LEU:N	2.19	0.41
1:A:-1:SER:CB	1:A:377:ASP:OD1	2.69	0.41
1:C:93:MET:HE2	1:C:93:MET:HB3	1.91	0.41
1:A:131:ASN:HD22	1:A:132:LEU:N	2.19	0.41
1:A:85:ARG:NH2	1:A:111:GLU:HB2	2.35	0.41
1:B:43:LEU:HB3	1:B:44:PRO:CD	2.51	0.41
1:B:158:LYS:HD2	1:B:158:LYS:C	2.41	0.41
1:B:175:ASN:O	1:B:176:LYS:C	2.58	0.41
1:A:57:VAL:CG1	1:A:158:LYS:HG3	2.50	0.40
1:B:106:TYR:CD2	1:B:139:PHE:HB2	2.55	0.40
1:B:89:TYR:HB2	1:B:93:MET:HG3	2.01	0.40
1:B:311:ILE:HD12	1:B:311:ILE:O	2.21	0.40
1:C:93:MET:CE	1:C:105:LEU:HD22	2.52	0.40
1:C:53:ILE:HG23	1:C:188:MET:HE3	2.02	0.40
1:A:13:GLU:CD	1:A:13:GLU:N	2.75	0.40
1:C:266:TYR:CE1	1:C:275:ARG:HG2	2.57	0.40
1:A:48:ARG:NH1	1:A:48:ARG:CG	2.66	0.40
1:A:32:LEU:HA	1:A:32:LEU:HD13	1.80	0.40
1:C:364:CYS:O	1:C:367:LEU:N	2.54	0.40
1:A:137:TRP:CZ2	1:B:70:LEU:CD2	3.04	0.40
1:C:166:ASP:O	1:C:312:HIS:CG	2.75	0.40
1:A:182:LEU:HD23	1:A:182:LEU:HA	1.91	0.40
1:C:333:ASP:OD1	1:C:333:ASP:C	2.60	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:414:LYS:HB3	1:B:414:LYS:HE2	1.89	0.40
1:C:111:GLU:OE2	2:C:439:5CA:N	2.55	0.40
1:C:143:GLN:HB3	1:C:144:ARG:HG2	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	439/458 (96%)	393 (90%)	41 (9%)	5 (1%)	17	47
1	B	440/458 (96%)	392 (89%)	36 (8%)	12 (3%)	6	22
1	C	440/458 (96%)	374 (85%)	55 (12%)	11 (2%)	7	24
All	All	1319/1374 (96%)	1159 (88%)	132 (10%)	28 (2%)	9	29

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	260	VAL
1	A	436	ARG
1	B	12	LYS
1	B	103	GLU
1	C	408	ARG
1	A	175	ASN
1	A	435	ALA
1	B	-2	GLY
1	B	227	VAL
1	B	276	LEU
1	B	421	SER
1	C	267	GLU
1	C	354	LYS
1	C	434	VAL

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Mol	Chain	Res	Type
1	C	435	ALA
1	B	142	GLU
1	B	235	GLU
1	C	266	TYR
1	C	252	SER
1	A	143	GLN
1	B	295	ASP
1	B	374	LYS
1	C	166	ASP
1	C	258	GLU
1	C	290	GLY
1	B	407	PRO
1	C	310	PRO
1	B	319	GLY

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	361/376 (96%)	322 (89%)	39 (11%)	8	21
1	B	362/376 (96%)	329 (91%)	33 (9%)	12	31
1	C	362/376 (96%)	304 (84%)	58 (16%)	3	7
All	All	1085/1128 (96%)	955 (88%)	130 (12%)	6	16

All (130) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	6	PHE
1	A	12	LYS
1	A	13	GLU
1	A	21	VAL
1	A	24	ARG
1	A	32	LEU
1	A	48	ARG

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Mol	Chain	Res	Type
1	A	51	LYS
1	A	69	LEU
1	A	78	ASP
1	A	93	MET
1	A	103	GLU
1	A	127	SER
1	A	131	ASN
1	A	134	HIS
1	A	138	LYS
1	A	142	GLU
1	A	144	ARG
1	A	158	LYS
1	A	159	ASP
1	A	165	VAL
1	A	172	LYS
1	A	188	MET
1	A	206	LEU
1	A	221	VAL
1	A	234	ASP
1	A	259	ASP
1	A	263	PRO
1	A	265	ARG
1	A	275	ARG
1	A	282	GLU
1	A	298	LYS
1	A	332	HIS
1	A	334	ASP
1	A	341	GLU
1	A	347	ARG
1	A	386	ARG
1	A	433	VAL
1	B	-3	ARG
1	B	12	LYS
1	B	16	LYS
1	B	34	GLN
1	B	48	ARG
1	B	93	MET
1	B	103	GLU
1	B	116	GLU
1	B	119	ARG
1	B	134	HIS
1	B	158	LYS

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Mol	Chain	Res	Type
1	B	164	ASP
1	B	165	VAL
1	B	175	ASN
1	B	178	PHE
1	B	206	LEU
1	B	225	ARG
1	B	248	LYS
1	B	249	GLN
1	B	251	THR
1	B	265	ARG
1	B	268	SER
1	B	270	VAL
1	B	295	ASP
1	B	301	VAL
1	B	366	GLN
1	B	369	ARG
1	B	408	ARG
1	B	414	LYS
1	B	422	ASP
1	B	427	ASN
1	B	428	LEU
1	B	436	ARG
1	C	-3	ARG
1	C	3	LEU
1	C	13	GLU
1	C	34	GLN
1	C	69	LEU
1	C	79	LEU
1	C	92	GLU
1	C	93	MET
1	C	95	ARG
1	C	100	HIS
1	C	116	GLU
1	C	124	SER
1	C	125	TYR
1	C	127	SER
1	C	131	ASN
1	C	134	HIS
1	C	138	LYS
1	C	143	GLN
1	C	144	ARG
1	C	150	MET

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Mol	Chain	Res	Type
1	C	158	LYS
1	C	159	ASP
1	C	165	VAL
1	C	178	PHE
1	C	183	ARG
1	C	187	ARG
1	C	188	MET
1	C	195	MET
1	C	199	THR
1	C	206	LEU
1	C	216	THR
1	C	218	GLU
1	C	219	SER
1	C	222	TYR
1	C	236	ASN
1	C	243	LEU
1	C	260	VAL
1	C	262	GLU
1	C	267	GLU
1	C	268	SER
1	C	272	GLU
1	C	274	ASN
1	C	276	LEU
1	C	295	ASP
1	C	298	LYS
1	C	300	ASN
1	C	307	THR
1	C	322	ARG
1	C	324	LEU
1	C	354	LYS
1	C	357	ASP
1	C	365	ASP
1	C	369	ARG
1	C	370	GLU
1	C	376	VAL
1	C	421	SER
1	C	426	GLU
1	C	436	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	ASN
1	A	131	ASN
1	A	285	GLN
1	A	335	ASN
1	B	47	HIS
1	B	61	GLN
1	B	175	ASN
1	B	236	ASN
1	B	285	GLN
1	B	300	ASN
1	C	0	HIS
1	C	23	HIS
1	C	131	ASN
1	C	143	GLN
1	C	229	ASN
1	C	236	ASN
1	C	249	GLN
1	C	274	ASN
1	C	285	GLN
1	C	332	HIS
1	C	355	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

3 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	5CA	A	439	-	24,31,31	2.68	4 (16%)	29,46,46	2.59	7 (24%)
2	5CA	B	439	-	24,31,31	2.16	2 (8%)	29,46,46	3.03	10 (34%)
2	5CA	C	439	-	24,31,31	2.43	4 (16%)	29,46,46	2.48	7 (24%)

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	5CA	A	439	-	-	0/15/37/37	0/3/3/3
2	5CA	B	439	-	-	0/15/37/37	0/3/3/3
2	5CA	C	439	-	-	0/15/37/37	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	439	5CA	CB-CA	-2.78	1.49	1.52
2	A	439	5CA	CB-SG	-2.09	1.76	1.81
2	C	439	5CA	C2-N3	2.03	1.35	1.32
2	C	439	5CA	O4'-C1'	2.24	1.44	1.41
2	B	439	5CA	O1S-S	6.22	1.47	1.42
2	C	439	5CA	O1S-S	7.38	1.48	1.42
2	B	439	5CA	O2S-S	7.74	1.49	1.42
2	C	439	5CA	O2S-S	7.98	1.49	1.42
2	A	439	5CA	O1S-S	8.45	1.49	1.42
2	A	439	5CA	O2S-S	8.74	1.49	1.42

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	439	5CA	N3-C2-N1	-9.36	121.72	128.89
2	A	439	5CA	N3-C2-N1	-9.12	121.91	128.89
2	C	439	5CA	O2S-S-O1S	-8.27	110.86	120.77
2	B	439	5CA	O2S-S-O1S	-7.96	111.24	120.77
2	C	439	5CA	N3-C2-N1	-7.14	123.43	128.89
2	A	439	5CA	C2'-C1'-N9	-5.53	105.84	114.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	439	5CA	CA-CB-SG	-5.15	102.69	114.48
2	A	439	5CA	O2S-S-O1S	-5.00	114.78	120.77
2	B	439	5CA	O1S-S-N3S	-4.84	102.03	108.50
2	B	439	5CA	CA-CB-SG	-4.34	104.53	114.48
2	B	439	5CA	C-N3S-S	-4.01	118.43	124.05
2	B	439	5CA	C2'-C1'-N9	-3.89	108.35	114.29
2	C	439	5CA	C-N3S-S	-3.26	119.49	124.05
2	B	439	5CA	C1'-N9-C4	-2.92	122.53	126.94
2	C	439	5CA	CA-CB-SG	-2.60	108.52	114.48
2	A	439	5CA	CB-CA-C	-2.55	103.98	109.70
2	B	439	5CA	C4-C5-N7	-2.34	107.33	109.48
2	C	439	5CA	C4'-O4'-C1'	-2.31	107.19	109.72
2	A	439	5CA	C4-C5-N7	-2.01	107.63	109.48
2	C	439	5CA	O-C-N3S	2.18	125.44	121.62
2	A	439	5CA	O1S-S-N3S	2.46	111.78	108.50
2	B	439	5CA	C5'-O5'-S	2.69	123.17	118.02
2	C	439	5CA	O1S-S-N3S	2.73	112.14	108.50
2	B	439	5CA	O2S-S-N3S	3.14	112.69	108.50

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	439	5CA	1	0
2	B	439	5CA	1	0
2	C	439	5CA	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	441/458 (96%)	-0.26	1 (0%) 95 95	44, 55, 77, 84	0
1	B	442/458 (96%)	-0.33	0 100 100	45, 58, 69, 81	0
1	C	442/458 (96%)	-0.59	0 100 100	33, 63, 86, 101	0
All	All	1325/1374 (96%)	-0.39	1 (0%) 95 95	33, 58, 81, 101	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	273	ALA	2.5

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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	5CA	C	439	29/29	0.98	0.12	-0.72	37,43,46,48	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	5CA	A	439	29/29	0.97	0.14	-1.91	46,48,50,51	0
2	5CA	B	439	29/29	0.97	0.13	-1.93	44,50,53,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.