Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 09:59 PM GMT

PDB ID	:	1RLC
Title	:	CRYSTAL STRUCTURE OF THE UNACTIVATED RIBULOSE 1, 5-BISPHOSPHATE CARBOXYLASE(SLASH)OXYGENASE COMPLEXED WITH A TRANSITION STATE ANALOG, 2-CARBOXY-D-ARABINITOL 1,5-BISPHOSPHATE
Authors	:	Zhang, K.Y.J.; Cascio, D.; Eisenberg, D.
Deposited on	:	1993-08-04
Resolution	:	2.70 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 (i) symbol.

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

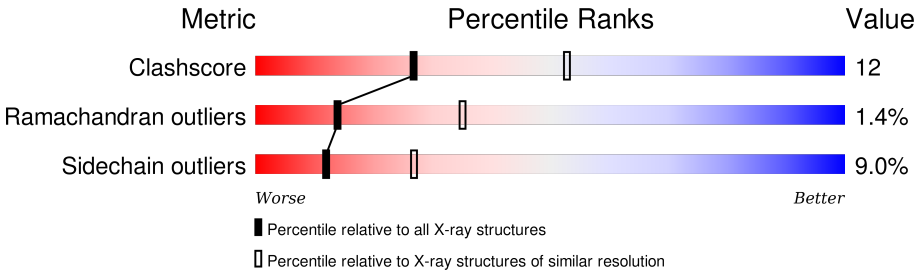
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)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

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)

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 <=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	L	477	56% 29% 7% 8%
2	S	123	59% 32% 8% •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4505 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.

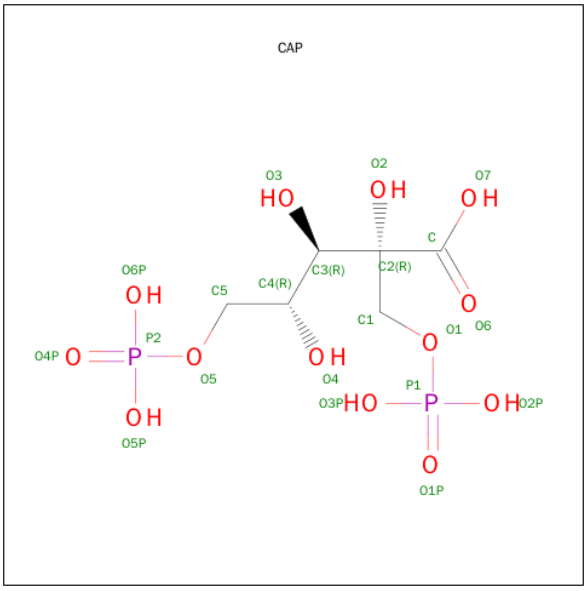
- Molecule 1 is a protein called RIBULOSE 1,5 BISPHOSPHATE CARBOXYLASE/OXYGENASE (LARGE CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	L	441	3455	2194	612	633	16	0	0	0

- Molecule 2 is a protein called RIBULOSE 1,5 BISPHOSPHATE CARBOXYLASE/OXYGENASE (SMALL CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	S	123	1029	672	163	188	6	0	0	0

- Molecule 3 is 2-CARBOXYARABINITOL-1,5-DIPHOSPHATE (three-letter code: CAP) (formula: $C_6H_{14}O_{13}P_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	P		
3	L	1	21	6	13	2	0	0

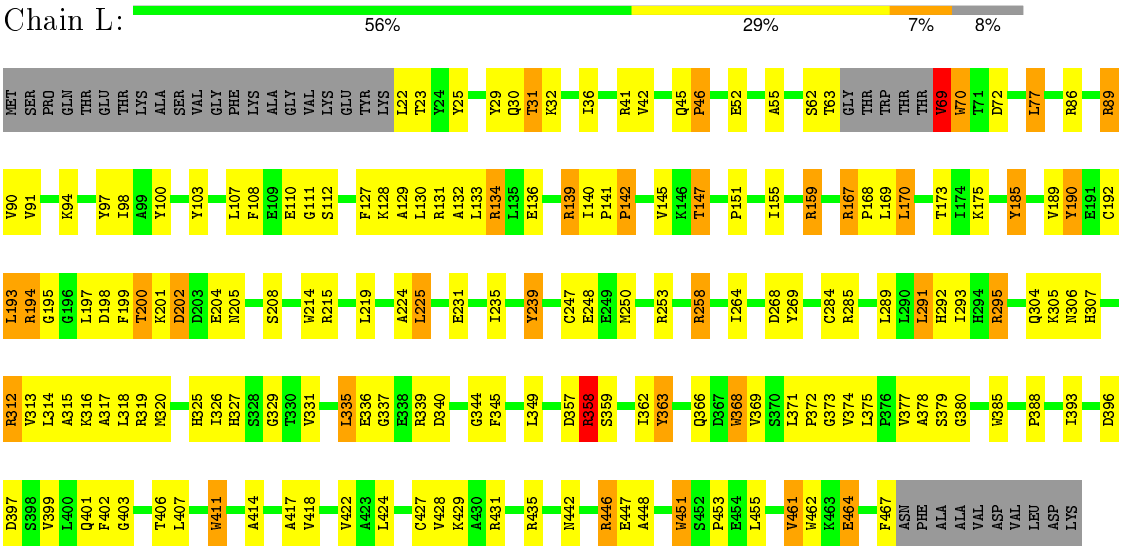
3 Residue-property plots (i)

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.

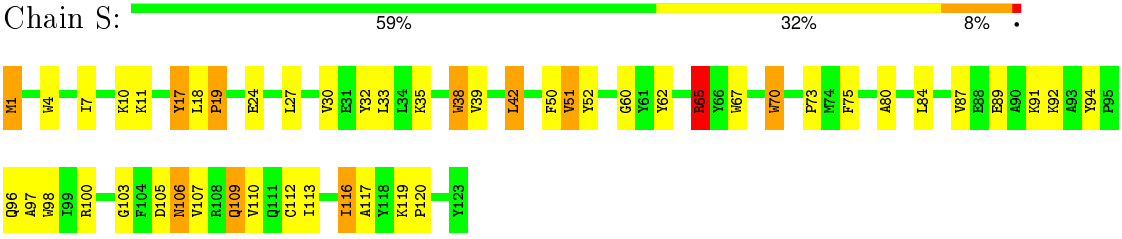
• Molecule 1: RIBULOSE 1,5 BISPHOSPHATE CARBOXYLASE/OXYGENASE (LARGE CHAIN)

Chain L:



• Molecule 2: RIBULOSE 1,5 BISPHOSPHATE CARBOXYLASE/OXYGENASE (SMALL CHAIN)

Chain S:



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α, β, γ	149.00Å 149.00Å 136.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.70	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.70)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.196 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4505	wwPDB-VP
Average B, all atoms (Å^2)	34.0	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: CAP

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	L	1.02	2/3538 (0.1%)	1.90	86/4796 (1.8%)
2	S	0.96	0/1062	1.82	26/1442 (1.8%)
All	All	1.01	2/4600 (0.0%)	1.88	112/6238 (1.8%)

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	L	0	1
2	S	0	1
All	All	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	248	GLU	CD-OE1	-6.04	1.19	1.25
1	L	208	SER	CA-CB	-5.43	1.44	1.52

All (112) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	L	215	ARG	NE-CZ-NH2	-22.87	108.86	120.30
1	L	215	ARG	NE-CZ-NH1	16.47	128.53	120.30
1	L	285	ARG	NE-CZ-NH1	12.63	126.61	120.30
2	S	70	TRP	CD1-CG-CD2	9.25	113.70	106.30
1	L	159	ARG	NE-CZ-NH1	9.21	124.90	120.30
2	S	67	TRP	CG-CD2-CE3	9.12	142.10	133.90
1	L	411	TRP	CD1-CG-CD2	9.01	113.50	106.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	S	98	TRP	CD1-CG-CD2	8.90	113.42	106.30
1	L	159	ARG	NE-CZ-NH2	-8.63	115.98	120.30
1	L	112	SER	CA-CB-OG	8.52	134.20	111.20
1	L	214	TRP	CD1-CG-CD2	8.51	113.11	106.30
1	L	70	TRP	N-CA-C	8.45	133.80	111.00
1	L	462	TRP	CD1-CG-CD2	8.26	112.91	106.30
2	S	38	TRP	CD1-CG-CD2	8.25	112.90	106.30
1	L	385	TRP	CD1-CG-CD2	8.16	112.83	106.30
1	L	320	MET	CG-SD-CE	8.00	113.00	100.20
1	L	435	ARG	NE-CZ-NH2	7.85	124.22	120.30
2	S	65	ARG	CB-CG-CD	-7.84	91.20	111.60
2	S	70	TRP	CE2-CD2-CG	-7.66	101.17	107.30
1	L	368	TRP	CD1-CG-CD2	7.66	112.43	106.30
2	S	100	ARG	NE-CZ-NH1	7.62	124.11	120.30
2	S	4	TRP	CD1-CG-CD2	7.56	112.35	106.30
1	L	285	ARG	NE-CZ-NH2	-7.55	116.53	120.30
1	L	239	TYR	CB-CG-CD1	7.53	125.52	121.00
1	L	368	TRP	CE2-CD2-CG	-7.50	101.30	107.30
2	S	98	TRP	CE2-CD2-CG	-7.46	101.33	107.30
1	L	103	TYR	CB-CG-CD1	-7.43	116.54	121.00
2	S	65	ARG	NE-CZ-NH1	7.42	124.01	120.30
1	L	451	TRP	CD1-CG-CD2	7.32	112.15	106.30
2	S	38	TRP	CE2-CD2-CG	-7.32	101.45	107.30
1	L	214	TRP	CE2-CD2-CG	-7.30	101.46	107.30
1	L	451	TRP	CE2-CD2-CG	-7.18	101.56	107.30
1	L	139	ARG	NE-CZ-NH2	7.05	123.83	120.30
1	L	190	TYR	CB-CG-CD2	-7.03	116.78	121.00
2	S	17	TYR	CB-CG-CD2	-7.03	116.78	121.00
2	S	4	TRP	CE2-CD2-CG	-6.98	101.71	107.30
1	L	411	TRP	CE2-CD2-CG	-6.91	101.77	107.30
1	L	385	TRP	CE2-CD2-CG	-6.84	101.83	107.30
1	L	41	ARG	NE-CZ-NH2	-6.80	116.90	120.30
2	S	67	TRP	CE2-CD2-CG	-6.79	101.86	107.30
2	S	109	GLN	CA-CB-CG	6.79	128.34	113.40
1	L	97	TYR	CB-CG-CD2	-6.71	116.97	121.00
2	S	67	TRP	CB-CG-CD1	-6.70	118.29	127.00
2	S	67	TRP	CD1-CG-CD2	6.66	111.63	106.30
2	S	51	VAL	CG1-CB-CG2	-6.56	100.40	110.90
1	L	258	ARG	NE-CZ-NH2	-6.55	117.02	120.30
2	S	30	VAL	CA-CB-CG1	-6.55	101.08	110.90
1	L	167	ARG	NE-CZ-NH2	-6.47	117.06	120.30
1	L	185	TYR	CB-CG-CD1	-6.47	117.12	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	L	401	GLN	CA-CB-CG	6.46	127.62	113.40
1	L	70	TRP	CE2-CD2-CG	-6.45	102.14	107.30
1	L	462	TRP	CE2-CD2-CG	-6.41	102.17	107.30
1	L	219	LEU	CB-CG-CD1	-6.36	100.20	111.00
1	L	70	TRP	CD1-CG-CD2	6.33	111.36	106.30
1	L	411	TRP	CG-CD1-NE1	-6.28	103.83	110.10
1	L	69	VAL	CA-C-N	6.23	130.91	117.20
1	L	317	ALA	O-C-N	-6.12	112.91	122.70
1	L	239	TYR	CB-CG-CD2	-6.11	117.33	121.00
1	L	435	ARG	CB-CG-CD	-6.11	95.70	111.60
2	S	70	TRP	CG-CD1-NE1	-6.11	104.00	110.10
1	L	103	TYR	CD1-CG-CD2	6.08	124.59	117.90
1	L	447	GLU	CA-CB-CG	6.08	126.78	113.40
1	L	368	TRP	CG-CD2-CE3	6.05	139.35	133.90
1	L	313	VAL	CG1-CB-CG2	-6.05	101.22	110.90
2	S	52	TYR	CB-CG-CD2	-5.97	117.42	121.00
1	L	269	TYR	CA-CB-CG	5.86	124.53	113.40
2	S	98	TRP	N-CA-C	-5.86	95.19	111.00
1	L	215	ARG	CA-CB-CG	5.85	126.27	113.40
1	L	111	GLY	CA-C-N	5.80	129.95	117.20
1	L	368	TRP	CB-CG-CD1	-5.77	119.50	127.00
1	L	358	ARG	NE-CZ-NH2	-5.75	117.42	120.30
1	L	70	TRP	CA-C-N	5.75	129.85	117.20
1	L	250	MET	CG-SD-CE	5.69	109.31	100.20
1	L	318	LEU	CA-C-N	5.66	129.65	117.20
1	L	107	LEU	CA-CB-CG	5.63	128.24	115.30
1	L	462	TRP	CG-CD1-NE1	-5.61	104.49	110.10
1	L	253	ARG	NE-CZ-NH1	5.60	123.10	120.30
2	S	4	TRP	CB-CG-CD1	-5.58	119.74	127.00
1	L	198	ASP	CB-CG-OD2	-5.55	113.30	118.30
1	L	248	GLU	OE1-CD-OE2	-5.53	116.67	123.30
1	L	397	ASP	CB-CG-OD2	-5.50	113.35	118.30
1	L	397	ASP	CB-CG-OD1	5.49	123.24	118.30
1	L	131	ARG	NE-CZ-NH2	-5.45	117.58	120.30
1	L	462	TRP	CB-CG-CD1	-5.44	119.93	127.00
1	L	134	ARG	NE-CZ-NH2	-5.40	117.60	120.30
1	L	451	TRP	CB-CG-CD1	-5.35	120.05	127.00
1	L	193	LEU	CA-C-N	5.34	128.94	117.20
1	L	111	GLY	CA-C-O	-5.33	111.00	120.60
2	S	52	TYR	O-C-N	-5.32	114.19	122.70
1	L	170	LEU	CA-CB-CG	5.31	127.51	115.30
1	L	340	ASP	CA-C-N	5.31	128.88	117.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	L	202	ASP	CB-CG-OD1	5.30	123.07	118.30
1	L	147	THR	CA-C-N	5.29	128.84	117.20
1	L	69	VAL	CA-CB-CG2	-5.29	102.97	110.90
1	L	225	LEU	CA-CB-CG	5.29	127.46	115.30
1	L	139	ARG	NH1-CZ-NH2	-5.29	113.59	119.40
1	L	385	TRP	CG-CD1-NE1	-5.28	104.82	110.10
1	L	291	LEU	O-C-N	-5.28	114.25	122.70
1	L	200	THR	CA-C-N	5.27	128.79	117.20
1	L	194	ARG	NE-CZ-NH2	-5.26	117.67	120.30
1	L	462	TRP	CG-CD2-CE3	5.26	138.64	133.90
1	L	91	VAL	CA-C-N	-5.25	105.69	116.20
2	S	38	TRP	CG-CD1-NE1	-5.22	104.88	110.10
1	L	103	TYR	CG-CD1-CE1	-5.21	117.13	121.30
1	L	295	ARG	CB-CG-CD	-5.19	98.11	111.60
1	L	363	TYR	CB-CG-CD1	-5.18	117.89	121.00
1	L	31	THR	CA-CB-CG2	-5.15	105.19	112.40
1	L	89	ARG	NE-CZ-NH1	5.14	122.87	120.30
1	L	69	VAL	CA-CB-CG1	5.10	118.55	110.90
1	L	268	ASP	CB-CG-OD1	5.06	122.86	118.30
1	L	373	GLY	O-C-N	5.05	130.78	122.70
2	S	30	VAL	CA-CB-CG2	5.05	118.48	110.90

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	L	45	GLN	Peptide
2	S	65	ARG	Sidechain

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	L	3455	0	3394	87	0
2	S	1029	0	994	25	0
3	L	21	0	9	1	0
All	All	4505	0	4397	107	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 12.

All (107) 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:L:315:ALA:HB1	1:L:349:LEU:HD21	1.51	0.89
1:L:335:LEU:HD23	1:L:337:GLY:H	1.47	0.79
1:L:127:PHE:CE2	1:L:129:ALA:HB3	2.22	0.74
1:L:388:PRO:HG3	1:L:427:CYS:SG	2.31	0.70
1:L:195:GLY:HA3	1:L:417:ALA:HB3	1.74	0.70
1:L:316:LYS:NZ	1:L:366:GLN:HE21	1.92	0.67
1:L:170:LEU:HD21	1:L:424:LEU:HD22	1.75	0.67
2:S:10:LYS:HB3	2:S:50:PHE:CE1	2.35	0.61
1:L:402:PHE:HB2	1:L:406:THR:HG23	1.82	0.61
1:L:326:ILE:O	1:L:377:VAL:HG23	2.02	0.60
1:L:151:PRO:HD2	1:L:372:PRO:HD2	1.83	0.59
1:L:173:THR:HB	1:L:403:GLY:HA2	1.84	0.59
1:L:319:ARG:HG3	1:L:374:VAL:HG23	1.85	0.59
1:L:291:LEU:HG	1:L:293:ILE:HD11	1.85	0.58
2:S:51:VAL:HG13	2:S:62:TYR:HB3	1.85	0.58
2:S:33:LEU:HD12	2:S:38:TRP:HE3	1.68	0.58
1:L:235:ILE:HD11	2:S:51:VAL:HG21	1.87	0.57
1:L:140:ILE:H	1:L:366:GLN:HE22	1.52	0.56
1:L:446:ARG:HG3	1:L:467:PHE:CE2	2.40	0.56
1:L:428:VAL:HG13	1:L:431:ARG:HH21	1.69	0.56
1:L:134:ARG:HD3	1:L:136:GLU:OE2	2.06	0.56
1:L:86:ARG:HB3	1:L:100:TYR:CD1	2.41	0.56
1:L:127:PHE:HD2	1:L:130:LEU:HG	1.73	0.54
1:L:190:TYR:HB2	1:L:224:ALA:HB1	1.89	0.54
2:S:97:ALA:O	2:S:120:PRO:HD3	2.06	0.54
1:L:133:LEU:O	1:L:307:HIS:HA	2.08	0.54
2:S:87:VAL:O	2:S:91:LYS:HG3	2.08	0.53
1:L:446:ARG:HG3	1:L:467:PHE:CZ	2.44	0.53
1:L:142:PRO:HB3	1:L:369:VAL:HG11	1.91	0.53
2:S:10:LYS:HB3	2:S:50:PHE:CZ	2.44	0.52
2:S:89:GLU:O	2:S:92:LYS:HG3	2.10	0.52
1:L:429:LYS:HD3	2:S:18:LEU:HD13	1.93	0.51
1:L:411:TRP:HH2	1:L:453:PRO:HB2	1.75	0.51
1:L:319:ARG:NH1	1:L:368:TRP:CD1	2.78	0.51
2:S:33:LEU:HD12	2:S:38:TRP:HB2	1.93	0.51
1:L:284:CYS:HB3	1:L:289:LEU:O	2.12	0.50
1:L:312:ARG:HB3	1:L:345:PHE:HB3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:36:ILE:HD12	1:L:108:PHE:CE2	2.46	0.50
1:L:169:LEU:O	1:L:399:VAL:HA	2.12	0.50
1:L:194:ARG:HD2	1:L:231:GLU:OE2	2.11	0.50
2:S:11:LYS:HG3	2:S:17:TYR:CE1	2.48	0.49
1:L:25:TYR:HB2	1:L:55:ALA:HB2	1.94	0.49
1:L:29:TYR:OH	1:L:32:LYS:HE2	2.13	0.49
1:L:316:LYS:HZ1	1:L:366:GLN:HE21	1.56	0.49
2:S:94:TYR:HB3	2:S:97:ALA:HB2	1.94	0.48
1:L:451:TRP:HH2	2:S:18:LEU:HD23	1.78	0.48
1:L:428:VAL:HG13	1:L:431:ARG:NH2	2.29	0.48
1:L:168:PRO:HD2	1:L:424:LEU:HD11	1.97	0.47
1:L:110:GLU:HB3	1:L:147:THR:HB	1.96	0.47
1:L:295:ARG:HG2	1:L:327:HIS:HB2	1.97	0.47
2:S:42:LEU:HB3	2:S:70:TRP:HB3	1.97	0.47
1:L:414:ALA:O	1:L:418:VAL:HG23	2.15	0.47
1:L:199:PHE:HB3	1:L:239:TYR:CE1	2.50	0.46
1:L:292:HIS:HA	1:L:325:HIS:HB2	1.97	0.46
1:L:36:ILE:HD12	1:L:108:PHE:CD2	2.50	0.46
1:L:319:ARG:HG3	1:L:374:VAL:CG2	2.45	0.46
1:L:357:ASP:OD1	1:L:359:SER:HB3	2.15	0.46
1:L:170:LEU:CD2	1:L:424:LEU:HD22	2.45	0.46
2:S:38:TRP:CD2	2:S:112:CYS:SG	3.09	0.46
2:S:33:LEU:HD11	2:S:103:GLY:HA3	1.97	0.45
1:L:139:ARG:HH21	1:L:141:PRO:HB3	1.81	0.45
1:L:167:ARG:HG3	1:L:168:PRO:O	2.16	0.45
1:L:304:GLN:OE1	1:L:304:GLN:HA	2.17	0.45
2:S:73:PRO:HB2	2:S:75:PHE:CE2	2.52	0.45
1:L:344:GLY:HA2	1:L:362:ILE:HD11	1.98	0.45
1:L:201:LYS:HG2	1:L:202:ASP:O	2.17	0.45
2:S:60:GLY:O	2:S:65:ARG:NH2	2.50	0.45
1:L:136:GLU:O	1:L:312:ARG:NH1	2.51	0.44
1:L:169:LEU:HD12	1:L:375:LEU:HD21	1.98	0.44
2:S:106:ASN:OD1	2:S:107:VAL:HG13	2.17	0.44
1:L:422:VAL:HG23	1:L:455:LEU:HD22	1.99	0.44
1:L:140:ILE:HA	1:L:141:PRO:HD2	1.84	0.44
1:L:442:ASN:O	1:L:446:ARG:HB2	2.18	0.44
1:L:29:TYR:CE2	1:L:31:THR:HG22	2.53	0.43
1:L:175:LYS:HB2	1:L:407:LEU:HD12	2.00	0.43
1:L:159:ARG:NH2	1:L:396:ASP:O	2.52	0.43
1:L:169:LEU:HB2	1:L:399:VAL:HG22	2.00	0.43
2:S:27:LEU:HG	2:S:84:LEU:HD22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:331:VAL:HG21	1:L:393:ILE:HD13	2.01	0.43
1:L:72:ASP:HB3	1:L:77:LEU:N	2.33	0.43
1:L:461:VAL:O	1:L:464:GLU:HB2	2.18	0.42
1:L:155:ILE:HG23	1:L:375:LEU:HD13	2.00	0.42
1:L:329:GLY:O	1:L:378:ALA:HA	2.18	0.42
1:L:239:TYR:CE2	1:L:292:HIS:CE1	3.08	0.42
1:L:358:ARG:HD3	1:L:363:TYR:HA	1.99	0.42
1:L:155:ILE:HG12	1:L:375:LEU:HD13	2.01	0.42
1:L:140:ILE:HG22	1:L:145:VAL:HG23	2.00	0.42
1:L:293:ILE:HG23	1:L:293:ILE:HD12	1.79	0.42
1:L:132:ALA:HB1	1:L:306:ASN:O	2.19	0.42
1:L:414:ALA:O	1:L:417:ALA:HB3	2.19	0.41
1:L:411:TRP:CE2	2:S:1:MET:HA	2.55	0.41
1:L:379:SER:OG	3:L:490:CAP:H12	2.20	0.41
2:S:116:ILE:HG13	2:S:117:ALA:N	2.34	0.41
2:S:39:VAL:O	2:S:103:GLY:HA2	2.19	0.41
1:L:185:TYR:O	1:L:189:VAL:HG23	2.21	0.41
1:L:42:VAL:HG23	1:L:130:LEU:HD13	2.03	0.41
1:L:86:ARG:HB3	1:L:100:TYR:HD1	1.84	0.41
1:L:448:ALA:HA	1:L:451:TRP:CE2	2.56	0.41
1:L:130:LEU:HD12	1:L:133:LEU:HD22	2.02	0.41
1:L:140:ILE:HG22	1:L:145:VAL:CG2	2.50	0.41
1:L:90:VAL:HG22	1:L:98:ILE:HG12	2.02	0.41
1:L:264:ILE:HG21	1:L:264:ILE:HD13	1.75	0.41
1:L:411:TRP:CD2	2:S:1:MET:HA	2.56	0.41
1:L:193:LEU:HG	1:L:200:THR:HG23	2.02	0.40
1:L:168:PRO:HG3	1:L:396:ASP:HA	2.03	0.40
1:L:331:VAL:HG11	1:L:393:ILE:HG21	2.02	0.40
2:S:32:TYR:HA	2:S:35:LYS:HD3	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	L	439/477 (92%)	399 (91%)	34 (8%)	6 (1%)	14 35
2	S	121/123 (98%)	110 (91%)	9 (7%)	2 (2%)	11 29
All	All	560/600 (93%)	509 (91%)	43 (8%)	8 (1%)	14 35

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	46	PRO
1	L	69	VAL
1	L	94	LYS
1	L	380	GLY
1	L	62	SER
1	L	204	GLU
2	S	80	ALA
2	S	19	PRO

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	L	357/386 (92%)	328 (92%)	29 (8%)	15 33
2	S	110/110 (100%)	97 (88%)	13 (12%)	6 15
All	All	467/496 (94%)	425 (91%)	42 (9%)	12 27

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	22	LEU
1	L	23	THR
1	L	30	GLN
1	L	46	PRO
1	L	52	GLU
1	L	63	THR
1	L	69	VAL
1	L	70	TRP

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Mol	Chain	Res	Type
1	L	77	LEU
1	L	89	ARG
1	L	128	LYS
1	L	142	PRO
1	L	192	CYS
1	L	197	LEU
1	L	205	ASN
1	L	225	LEU
1	L	247	CYS
1	L	258	ARG
1	L	305	LYS
1	L	312	ARG
1	L	314	LEU
1	L	335	LEU
1	L	336	GLU
1	L	339	ARG
1	L	358	ARG
1	L	371	LEU
1	L	446	ARG
1	L	461	VAL
1	L	464	GLU
2	S	1	MET
2	S	7	ILE
2	S	19	PRO
2	S	24	GLU
2	S	42	LEU
2	S	96	GLN
2	S	105	ASP
2	S	106	ASN
2	S	109	GLN
2	S	110	VAL
2	S	113	ILE
2	S	116	ILE
2	S	119	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	156	GLN
1	L	207	ASN
1	L	327	HIS
1	L	366	GLN

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Mol	Chain	Res	Type
1	L	420	ASN
2	S	36	ASN
2	S	109	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)

1 ligand is 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	$\#\|Z\| > 2$
3	CAP	L	490	-	14,20,20	1.93	4 (28%)	15,31,31	3.21	10 (66%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	CAP	L	490	-	-	0/23/29/29	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	490	CAP	C5-C4	2.06	1.54	1.51
3	L	490	CAP	O5-C5	2.35	1.54	1.44
3	L	490	CAP	P2-O6P	2.83	1.64	1.54
3	L	490	CAP	P1-O1P	4.88	1.67	1.51

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	L	490	CAP	O2P-P1-O1P	-5.51	92.83	110.58
3	L	490	CAP	O1-P1-O1P	-3.59	97.99	107.14
3	L	490	CAP	O3-C3-C4	-3.03	102.42	108.91
3	L	490	CAP	O6P-P2-O5P	-2.94	96.18	107.38
3	L	490	CAP	O3P-P1-O2P	2.26	115.98	107.38
3	L	490	CAP	O2P-P1-O1	3.19	115.75	106.56
3	L	490	CAP	O4-C4-C5	3.54	117.89	110.19
3	L	490	CAP	O3P-P1-O1	4.28	118.88	106.56
3	L	490	CAP	O5-P2-O4P	4.33	118.17	107.14
3	L	490	CAP	O5P-P2-O5	4.67	120.02	106.56

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	L	490	CAP	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

6.1 Protein, DNA and RNA chains

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

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

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

6.4 Ligands

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

6.5 Other polymers

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