Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 07:10 PM GMT

PDB ID	:	1EE0
Title	:	2-PYRONE SYNTHASE COMPLEXED WITH ACETOACETYL-COA
Authors	:	Jez, J.M.; Austin, M.B.; Ferrer, J.; Bowmann, M.E.; Schroeder, J.; Noel, J.P.
Deposited on	:	2000-01-28
Resolution	:	2.05 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 (1)) were used in the production of this report:

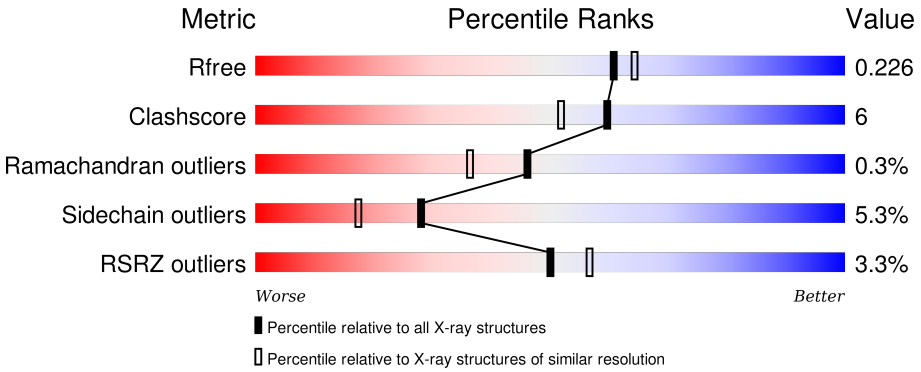
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.05 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	1192 (2.04-2.04)
Clashscore	102246	1269 (2.04-2.04)
Ramachandran outliers	100387	1258 (2.04-2.04)
Sidechain outliers	100360	1258 (2.04-2.04)
RSRZ outliers	91569	1194 (2.04-2.04)

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.

Mol	Chain	Length	Quality of chain
1	A	402	4% 75% 15% • 6%
1	B	402	2% 76% 14% •• 7%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 6451 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.

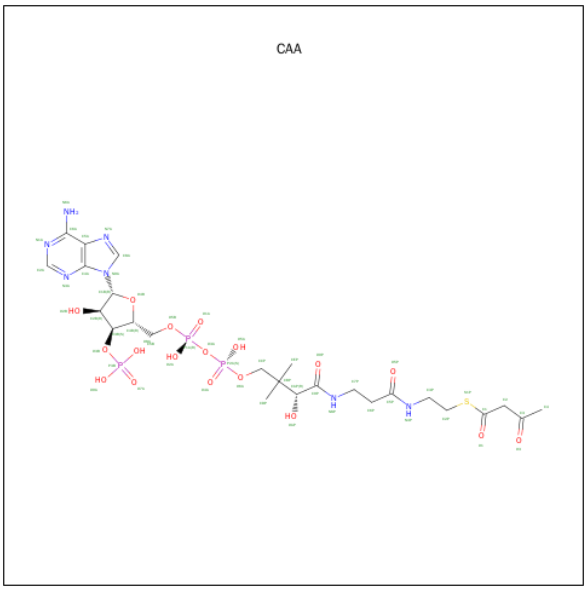
- Molecule 1 is a protein called 2-PYRONE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	376	Total	C	N	O	S	0	5	0
			2918	1851	502	546	19			
1	B	375	Total	C	N	O	S	0	2	0
			2884	1831	498	536	19			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	169	CSD	CYS	MODIFIED RESIDUE	UNP P48391
A	259	MET	VAL	CONFLICT	UNP P48391
B	169	CSD	CYS	MODIFIED RESIDUE	UNP P48391
B	259	MET	VAL	CONFLICT	UNP P48391

- Molecule 2 is ACETOACETYL-COENZYME A (three-letter code: CAA) (formula: $C_{25}H_{40}N_7O_{18}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			54	25	7	18	3	1		
2	B	1	Total	C	N	O	P	S	0	0
			54	25	7	18	3	1		

- Molecule 3 is water.

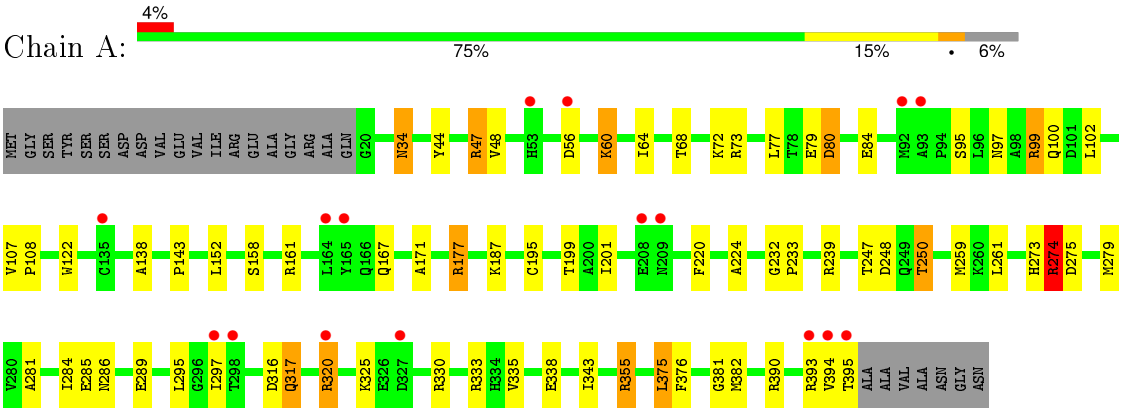
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	256	Total	O	0	0
			256	256		
3	B	285	Total	O	0	0
			285	285		

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.

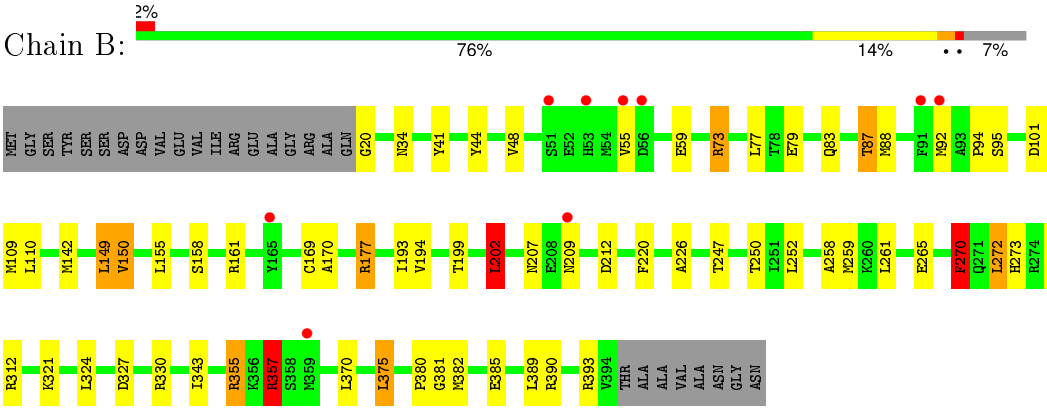
- Molecule 1: 2-PYRONE SYNTHASE

Chain A:



- Molecule 1: 2-PYRONE SYNTHASE

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	83.41Å 83.41Å 240.62Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.00 – 2.05 41.70 – 2.05	Depositor EDS
% Data completeness (in resolution range)	98.2 (42.00-2.05) 98.2 (41.70-2.05)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.13 (at 2.05Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.189 , 0.243 0.182 , 0.226	Depositor DCC
R_{free} test set	3085 reflections (5.35%)	DCC
Wilson B-factor (Å^2)	22.3	Xtriage
Anisotropy	0.385	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 60.8	EDS
Estimated twinning fraction	0.022 for -h,-k,l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Outliers	2 of 60757 reflections (0.003%)	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	6451	wwPDB-VP
Average B, all atoms (Å^2)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 61.62 % of the origin peak, indicating pseudo translational symmetry.The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.2749e-05.The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: CSD, CAA

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	0.77	0/2961	1.45	32/4010 (0.8%)
1	B	0.80	0/2927	1.46	25/3962 (0.6%)
All	All	0.79	0/5888	1.45	57/7972 (0.7%)

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	1
1	B	0	2
All	All	0	3

There are no bond length outliers.

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	390	ARG	NE-CZ-NH2	19.43	130.02	120.30
1	B	357	ARG	NE-CZ-NH1	-13.84	113.38	120.30
1	B	161	ARG	NE-CZ-NH1	13.61	127.10	120.30
1	A	330	ARG	NE-CZ-NH1	11.86	126.23	120.30
1	B	161	ARG	NE-CZ-NH2	-11.65	114.47	120.30
1	B	330	ARG	NE-CZ-NH1	-11.57	114.52	120.30
1	A	239	ARG	CD-NE-CZ	11.42	139.59	123.60
1	A	161	ARG	NE-CZ-NH2	-10.71	114.94	120.30
1	A	239	ARG	NE-CZ-NH1	10.62	125.61	120.30
1	A	381	GLY	O-C-N	-10.36	106.13	122.70
1	B	73	ARG	NE-CZ-NH1	10.20	125.40	120.30
1	A	390	ARG	NE-CZ-NH1	-9.90	115.35	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	390	ARG	CD-NE-CZ	9.22	136.51	123.60
1	B	355	ARG	NE-CZ-NH1	-8.82	115.89	120.30
1	B	375	LEU	CA-CB-CG	8.60	135.08	115.30
1	B	381	GLY	O-C-N	-8.20	109.59	122.70
1	A	47	ARG	NE-CZ-NH1	8.07	124.33	120.30
1	B	355	ARG	NE-CZ-NH2	8.03	124.31	120.30
1	A	73	ARG	NE-CZ-NH2	7.69	124.15	120.30
1	A	381	GLY	CA-C-N	7.63	133.99	117.20
1	B	381	GLY	CA-C-N	7.57	133.86	117.20
1	A	320	ARG	NE-CZ-NH1	7.41	124.00	120.30
1	B	357	ARG	NE-CZ-NH2	7.31	123.95	120.30
1	A	80	ASP	CB-CG-OD1	7.25	124.82	118.30
1	A	56	ASP	CB-CG-OD1	7.02	124.62	118.30
1	A	143	PRO	N-CA-CB	6.90	111.58	103.30
1	B	390	ARG	NE-CZ-NH2	-6.81	116.90	120.30
1	A	99	ARG	NE-CZ-NH2	-6.78	116.91	120.30
1	A	333	ARG	NE-CZ-NH2	6.54	123.57	120.30
1	A	375	LEU	CB-CG-CD1	6.43	121.94	111.00
1	B	177	ARG	NE-CZ-NH2	6.42	123.51	120.30
1	A	375	LEU	CA-CB-CG	6.42	130.07	115.30
1	B	390	ARG	CD-NE-CZ	6.40	132.56	123.60
1	B	92	MET	CG-SD-CE	6.12	109.99	100.20
1	B	101	ASP	CB-CG-OD1	6.06	123.76	118.30
1	B	88	MET	CG-SD-CE	5.99	109.78	100.20
1	B	327	ASP	CB-CG-OD1	5.93	123.63	118.30
1	B	212	ASP	CB-CG-OD1	5.88	123.59	118.30
1	B	312	ARG	NE-CZ-NH1	5.88	123.24	120.30
1	B	142	MET	CG-SD-CE	5.78	109.45	100.20
1	A	34	ASN	CA-CB-CG	5.77	126.09	113.40
1	B	202	LEU	CA-CB-CG	5.74	128.49	115.30
1	A	393	ARG	NE-CZ-NH2	-5.71	117.45	120.30
1	A	177	ARG	NE-CZ-NH2	5.66	123.13	120.30
1	A	382	MET	N-CA-CB	5.61	120.70	110.60
1	A	56	ASP	CB-CG-OD2	-5.60	113.26	118.30
1	A	355	ARG	NE-CZ-NH2	5.60	123.10	120.30
1	B	41	TYR	CB-CG-CD1	-5.56	117.66	121.00
1	A	316	ASP	CB-CG-OD2	5.46	123.21	118.30
1	A	161	ARG	NH1-CZ-NH2	5.41	125.35	119.40
1	B	202	LEU	CB-CG-CD2	5.35	120.10	111.00
1	A	274	ARG	NE-CZ-NH1	5.33	122.97	120.30
1	A	259	MET	CB-CA-C	-5.29	99.81	110.40
1	A	161	ARG	CD-NE-CZ	5.27	130.98	123.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	375	LEU	CB-CG-CD1	5.26	119.94	111.00
1	A	317	GLN	CG-CD-OE1	-5.15	111.30	121.60
1	A	259	MET	CG-SD-CE	5.12	108.40	100.20

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	122	TRP	Mainchain
1	B	270	PHE	Mainchain
1	B	380	PRO	Mainchain

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	2918	0	2970	34	0
1	B	2884	0	2946	38	0
2	A	54	0	35	3	0
2	B	54	0	34	3	0
3	A	256	0	0	1	0
3	B	285	0	0	7	0
All	All	6451	0	5985	68	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 6.

All (68) 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:A:273:HIS:HD2	1:A:275:ASP:H	1.19	0.88
1:B:321:LYS:HE2	3:B:850:HOH:O	1.84	0.77
1:A:97:ASN:HD21	1:B:265:GLU:H	1.34	0.74
1:B:73:ARG:HD3	3:B:894:HOH:O	1.89	0.72
1:A:261:LEU:HD11	2:A:600:CAA:H2'1	1.71	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:202:LEU:HG	1:B:270:PHE:HE1	1.52	0.72
1:B:273:HIS:HD2	1:B:275:ASP:H	1.38	0.71
1:A:273:HIS:CD2	1:A:275:ASP:H	2.07	0.70
1:A:248[B]:ASP:OD2	1:A:250[B]:THR:HG23	1.92	0.69
1:B:169:CSD:OD2	2:B:700:CAA:C3	2.43	0.67
1:B:87:THR:HG23	1:B:95:SER:HB3	1.78	0.65
1:B:87:THR:HB	3:B:721:HOH:O	1.94	0.65
1:B:87:THR:CG2	1:B:94:PRO:O	2.45	0.64
1:B:357:ARG:NH1	3:B:962:HOH:O	2.24	0.64
1:B:87:THR:HG21	1:B:94:PRO:O	2.00	0.62
1:B:149:LEU:HD11	1:B:194:VAL:HG21	1.82	0.61
1:A:250[B]:THR:HG22	1:B:158:SER:OG	2.02	0.60
1:A:47:ARG:NH2	1:A:79:GLU:OE2	2.36	0.58
1:A:60:LYS:HD3	2:A:600:CAA:O7A	2.05	0.57
1:A:281:ALA:HB2	1:A:317:GLN:HB3	1.86	0.57
1:B:34:ASN:HB3	3:B:744:HOH:O	2.05	0.56
1:B:273:HIS:CD2	1:B:275:ASP:H	2.22	0.55
1:A:138:ALA:HA	1:A:201[B]:ILE:HG12	1.89	0.55
1:B:79:GLU:O	1:B:83:GLN:HG3	2.08	0.53
1:B:87:THR:CG2	1:B:95:SER:HB3	2.40	0.52
1:B:207:ASN:ND2	1:B:209:ASN:H	2.08	0.51
1:B:321:LYS:HE3	3:B:754:HOH:O	2.10	0.51
1:A:97:ASN:ND2	1:B:265:GLU:H	2.07	0.51
1:B:20:GLY:N	3:B:967:HOH:O	2.44	0.51
1:B:87:THR:HG23	1:B:94:PRO:O	2.11	0.51
1:A:34:ASN:HB3	3:A:698:HOH:O	2.10	0.50
1:B:261:LEU:HD11	2:B:700:CAA:H2'1	1.92	0.50
1:A:177:ARG:HA	1:A:247:THR:HG21	1.93	0.50
1:B:177:ARG:HA	1:B:247:THR:HG21	1.95	0.48
1:A:80:ASP:O	1:A:84:GLU:HG3	2.14	0.47
1:A:167:GLN:HB3	1:A:171:ALA:HB2	1.96	0.47
1:A:44:TYR:CE1	1:A:48:VAL:HG11	2.50	0.47
1:A:107:VAL:N	1:A:108:PRO:HD2	2.29	0.47
1:B:199:THR:O	1:B:202:LEU:HD22	2.15	0.47
1:A:394:VAL:O	1:A:395:THR:HB	2.15	0.46
1:A:195:CYS:O	1:A:224:ALA:HA	2.15	0.46
1:A:285:GLU:OE2	1:A:289:GLU:OE1	2.33	0.46
1:A:72:LYS:HE3	1:A:338:GLU:OE1	2.16	0.46
1:A:284:ILE:HD11	1:A:376:PHE:CE1	2.50	0.46
1:A:64:ILE:O	1:A:68:THR:HG23	2.16	0.46
1:A:102:LEU:HD23	1:A:102:LEU:C	2.36	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:357:ARG:HH11	1:B:357:ARG:HD3	1.50	0.45
1:A:232:GLY:N	1:A:233:PRO:HD3	2.32	0.44
1:A:275:ASP:O	1:A:279:MET:HG3	2.17	0.44
1:A:199:THR:HG21	1:A:220:PHE:HB3	2.00	0.44
1:A:261:LEU:CD2	2:A:600:CAA:H4'3	2.48	0.43
1:A:47:ARG:NH2	1:A:79:GLU:CD	2.72	0.43
1:A:95:SER:O	1:A:99:ARG:HG3	2.18	0.43
1:A:47:ARG:HH22	1:A:79:GLU:CD	2.22	0.42
1:B:199:THR:HG21	1:B:220:PHE:HB3	2.00	0.42
1:B:48:VAL:HG11	1:B:79:GLU:HG2	2.00	0.42
1:A:100:GLN:HA	1:A:100:GLN:HE21	1.85	0.42
1:A:158:SER:OG	1:B:250:THR:HG23	2.19	0.42
1:B:170:ALA:HB1	1:B:385:GLU:HG3	2.02	0.42
1:B:258:ALA:HB3	1:B:382:MET:HG3	2.03	0.41
1:B:107:VAL:N	1:B:108:PRO:HD2	2.35	0.41
1:B:109:MET:HE3	1:B:109:MET:HA	2.01	0.41
1:A:274:ARG:O	1:A:274:ARG:HG3	2.20	0.41
1:B:55:VAL:O	1:B:59:GLU:HG3	2.20	0.41
1:B:193:ILE:O	1:B:226:ALA:HA	2.21	0.41
1:B:272:LEU:HD21	2:B:700:CAA:O5P	2.21	0.41
1:B:44:TYR:HE1	1:B:79:GLU:CG	2.34	0.41
1:B:150:VAL:HB	1:B:155:LEU:HB2	2.04	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	378/402 (94%)	372 (98%)	5 (1%)	1 (0%)	46	36
1	B	374/402 (93%)	367 (98%)	6 (2%)	1 (0%)	46	36
All	All	752/804 (94%)	739 (98%)	11 (2%)	2 (0%)	46	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	343	ILE
1	B	343	ILE

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	314/328 (96%)	299 (95%)	15 (5%)	31 22
1	B	310/328 (94%)	291 (94%)	19 (6%)	23 13
All	All	624/656 (95%)	590 (95%)	34 (5%)	28 17

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	LYS
1	A	77	LEU
1	A	152	LEU
1	A	187	LYS
1	A	250[A]	THR
1	A	250[B]	THR
1	A	274	ARG
1	A	286	ASN
1	A	295	LEU
1	A	297	ILE
1	A	320	ARG
1	A	325	LYS
1	A	335	VAL
1	A	355	ARG
1	A	375	LEU
1	B	77	LEU
1	B	87	THR
1	B	102	LEU
1	B	110	LEU
1	B	149	LEU
1	B	150	VAL

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Mol	Chain	Res	Type
1	B	202	LEU
1	B	252	LEU
1	B	259	MET
1	B	270	PHE
1	B	272	LEU
1	B	295	LEU
1	B	324	LEU
1	B	355	ARG
1	B	357	ARG
1	B	370	LEU
1	B	375	LEU
1	B	389	LEU
1	B	393	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	97	ASN
1	A	100	GLN
1	A	273	HIS
1	A	317	GLN
1	B	100	GLN
1	B	166	GLN
1	B	207	ASN
1	B	273	HIS
1	B	334	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

2 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	A	169	1	3,7,8	0.89	0	3,8,10	6.02	2 (66%)
1	CSD	B	169	1	3,7,8	1.13	0	3,8,10	8.07	2 (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
1	CSD	A	169	1	-	1/2/6/8	0/0/0/0
1	CSD	B	169	1	-	1/2/6/8	0/0/0/0

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	169	CSD	OD1-SG-CB	-13.34	83.17	105.40
1	A	169	CSD	OD1-SG-CB	-10.03	88.69	105.40
1	B	169	CSD	O-C-CA	-3.82	115.53	125.49
1	A	169	CSD	CB-CA-C	-2.28	105.20	111.46

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	169	CSD	CA-CB-SG-OD1
1	B	169	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	B	169	CSD	1	0

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

2 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	CAA	A	600	-	45,56,56	1.83	9 (20%)	58,83,83	2.87	23 (39%)
2	CAA	B	700	-	45,56,56	1.87	9 (20%)	58,83,83	3.41	24 (41%)

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	CAA	A	600	-	-	0/50/71/71	0/3/3/3
2	CAA	B	700	-	-	0/50/71/71	0/3/3/3

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	700	CAA	P1A-O5B	-5.43	1.34	1.59
2	A	600	CAA	P1A-O5B	-5.43	1.34	1.59
2	A	600	CAA	O2B-C2B	-2.76	1.36	1.43
2	B	700	CAA	O2B-C2B	-2.64	1.36	1.43
2	A	600	CAA	C4A-N3A	2.03	1.38	1.35
2	A	600	CAA	O3B-C3B	2.53	1.51	1.44
2	A	600	CAA	C3P-N4P	2.53	1.52	1.46
2	B	700	CAA	C2A-N1A	2.96	1.39	1.33
2	B	700	CAA	C3P-N4P	2.97	1.53	1.46
2	A	600	CAA	P3B-O7A	3.00	1.61	1.51
2	B	700	CAA	P3B-O7A	3.00	1.61	1.51
2	B	700	CAA	O5B-C5B	3.03	1.57	1.44
2	A	600	CAA	C7P-N8P	3.25	1.53	1.46
2	B	700	CAA	O4B-C4B	3.47	1.53	1.45
2	A	600	CAA	O4B-C4B	3.53	1.53	1.45
2	B	700	CAA	C7P-N8P	3.79	1.55	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	700	CAA	O4B-C1B	4.16	1.46	1.41
2	A	600	CAA	O4B-C1B	4.44	1.46	1.41

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	700	CAA	O1-C1-S1P	-11.10	114.03	122.83
2	B	700	CAA	O5P-C5P-C6P	-10.42	104.00	121.98
2	A	600	CAA	O5P-C5P-C6P	-7.59	108.88	121.98
2	A	600	CAA	O4B-C1B-N9A	-6.28	94.95	108.10
2	B	700	CAA	C7P-N8P-C9P	-5.56	111.53	122.53
2	B	700	CAA	O4B-C1B-N9A	-5.31	96.98	108.10
2	A	600	CAA	N3A-C2A-N1A	-5.16	124.95	128.89
2	B	700	CAA	N3A-C2A-N1A	-4.86	125.17	128.89
2	A	600	CAA	O1-C1-S1P	-4.76	119.06	122.83
2	B	700	CAA	C3P-N4P-C5P	-4.12	114.70	122.79
2	B	700	CAA	O3B-P3B-O7A	-3.72	97.82	107.11
2	A	600	CAA	O3B-P3B-O7A	-3.72	97.82	107.11
2	A	600	CAA	C7P-N8P-C9P	-3.65	115.30	122.53
2	B	700	CAA	OAP-CAP-C9P	-3.49	102.38	110.38
2	A	600	CAA	OAP-CAP-C9P	-3.44	102.50	110.38
2	A	600	CAA	O3B-C3B-C4B	-3.07	97.95	109.99
2	B	700	CAA	O3B-C3B-C4B	-2.88	98.71	109.99
2	B	700	CAA	C6P-C7P-N8P	-2.76	105.81	111.88
2	A	600	CAA	O4B-C4B-C3B	-2.76	98.50	104.86
2	B	700	CAA	C7P-C6P-C5P	-2.76	107.76	112.31
2	B	700	CAA	O4B-C4B-C3B	-2.53	99.02	104.86
2	A	600	CAA	C1B-N9A-C4A	-2.52	123.14	126.94
2	A	600	CAA	CDP-CBP-CCP	-2.50	105.26	108.50
2	B	700	CAA	C3P-C2P-S1P	-2.29	105.24	111.36
2	A	600	CAA	O6A-CCP-CBP	-2.18	107.04	110.55
2	B	700	CAA	O6A-CCP-CBP	-2.11	107.16	110.55
2	A	600	CAA	O5P-C5P-N4P	2.02	126.95	122.94
2	A	600	CAA	C4A-C5A-N7A	2.12	111.43	109.48
2	B	700	CAA	CDP-CBP-CAP	2.22	113.39	109.34
2	B	700	CAA	C4A-C5A-N7A	2.23	111.53	109.48
2	A	600	CAA	CEP-CBP-CCP	2.39	111.59	108.50
2	A	600	CAA	P2A-O3A-P1A	2.57	139.95	132.73
2	B	700	CAA	C6P-C5P-N4P	2.72	121.18	116.46
2	A	600	CAA	O3A-P2A-O6A	2.90	110.64	102.94
2	B	700	CAA	CEP-CBP-CCP	2.96	112.34	108.50
2	B	700	CAA	P2A-O3A-P1A	3.19	141.70	132.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	700	CAA	C2P-C3P-N4P	3.21	118.78	112.36
2	A	600	CAA	C2A-N1A-C6A	3.27	124.61	118.77
2	B	700	CAA	C2B-C3B-C4B	3.87	110.56	103.29
2	A	600	CAA	C2B-C3B-C4B	4.13	111.06	103.29
2	A	600	CAA	C6P-C5P-N4P	4.40	124.11	116.46
2	B	700	CAA	P3B-O3B-C3B	6.23	136.50	121.56
2	A	600	CAA	C2B-C1B-N9A	6.38	124.04	114.29
2	A	600	CAA	P3B-O3B-C3B	6.76	137.76	121.56
2	A	600	CAA	O2B-C2B-C3B	7.13	131.76	111.16
2	B	700	CAA	O2B-C2B-C3B	7.70	133.39	111.16
2	B	700	CAA	C2B-C1B-N9A	8.55	127.35	114.29

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	600	CAA	3	0
2	B	700	CAA	3	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

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, 95^{th} 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	375/402 (93%)	0.14	16 (4%) 39 44	11, 22, 39, 69	0
1	B	374/402 (93%)	0.11	9 (2%) 62 68	11, 21, 37, 51	0
All	All	749/804 (93%)	0.12	25 (3%) 50 57	11, 21, 38, 69	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	53	HIS	4.4
1	B	55	VAL	4.0
1	A	92	MET	3.8
1	A	395	THR	3.7
1	B	92	MET	3.6
1	B	56	ASP	3.5
1	A	56	ASP	3.3
1	A	394	VAL	3.2
1	A	164	LEU	3.2
1	A	208	GLU	3.1
1	B	53	HIS	2.9
1	B	359	MET	2.7
1	B	209	ASN	2.7
1	A	297	ILE	2.5
1	A	209	ASN	2.5
1	B	91	PHE	2.4
1	A	93	ALA	2.4
1	B	165	TYR	2.4
1	A	298	THR	2.3
1	A	165	TYR	2.2
1	B	51	SER	2.2
1	A	327	ASP	2.1
1	A	393	ARG	2.1
1	A	135	CYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	320	ARG	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. 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, 95^{th} 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
1	CSD	B	169	8/9	0.97	0.13	-	14,16,29,29	0
1	CSD	A	169	8/9	0.92	0.15	-	16,17,25,26	0

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, 95^{th} 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	CAA	B	700	54/54	0.89	0.14	0.22	27,32,36,37	0
2	CAA	A	600	54/54	0.92	0.13	0.15	27,32,37,38	0

6.5 Other polymers (i)

There are no such residues in this entry.