Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 05:33 AM GMT

PDB ID	:	2R98
Title	:	Crystal Structure of N-acetylglutamate synthase (selenoMet substituted) from Neisseria gonorrhoeae
Authors	:	Shi, D.; Sagar, V.; Jin, Z.; Yu, X.; Caldovic, L.; Morizono, H.; Allewell, N.M.; Tuchman, M.
Deposited on	:	2007-09-12
Resolution	:	2.40 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:

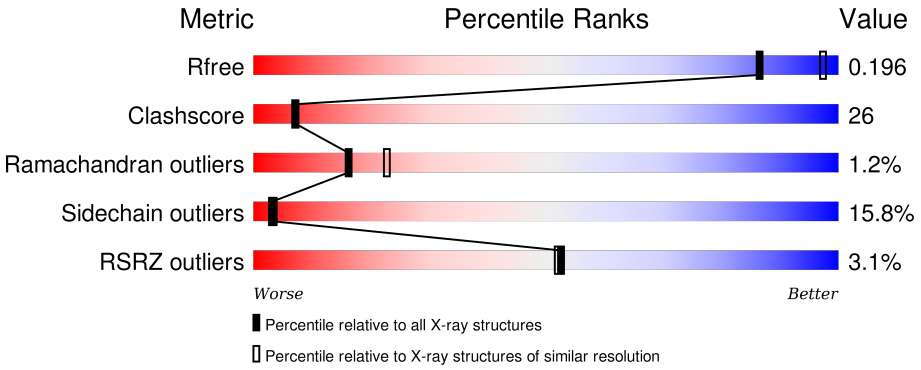
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance

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

The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	456	3%, 55%, 29%, 9%, 7%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ACO	A	1	X	-	-	-

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3370 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 Putative acetylglutamate synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	424	Total	C	N	O	S	Se	0	0	0
			3227	2012	595	611	5	4			

There are 23 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MSE	-	EXPRESSION TAG	UNP Q5FAK7
A	-18	GLY	-	EXPRESSION TAG	UNP Q5FAK7
A	-17	SER	-	EXPRESSION TAG	UNP Q5FAK7
A	-16	SER	-	EXPRESSION TAG	UNP Q5FAK7
A	-15	HIS	-	EXPRESSION TAG	UNP Q5FAK7
A	-14	HIS	-	EXPRESSION TAG	UNP Q5FAK7
A	-13	HIS	-	EXPRESSION TAG	UNP Q5FAK7
A	-12	HIS	-	EXPRESSION TAG	UNP Q5FAK7
A	-11	HIS	-	EXPRESSION TAG	UNP Q5FAK7
A	-10	HIS	-	EXPRESSION TAG	UNP Q5FAK7
A	-9	SER	-	EXPRESSION TAG	UNP Q5FAK7
A	-8	SER	-	EXPRESSION TAG	UNP Q5FAK7
A	-7	GLY	-	EXPRESSION TAG	UNP Q5FAK7
A	-6	LEU	-	EXPRESSION TAG	UNP Q5FAK7
A	-5	VAL	-	EXPRESSION TAG	UNP Q5FAK7
A	-4	PRO	-	EXPRESSION TAG	UNP Q5FAK7
A	-3	ARG	-	EXPRESSION TAG	UNP Q5FAK7
A	-2	GLY	-	EXPRESSION TAG	UNP Q5FAK7
A	-1	SER	-	EXPRESSION TAG	UNP Q5FAK7
A	0	HIS	-	EXPRESSION TAG	UNP Q5FAK7
A	312	ILE	VAL	ENGINEERED	UNP Q5FAK7
A	336	ASN	ASP	ENGINEERED	UNP Q5FAK7
A	427	SER	PRO	ENGINEERED	UNP Q5FAK7

- Molecule 2 is ACETYL COENZYME *A (three-letter code: ACO) (formula: $C_{23}H_{38}N_7O_{17}P_3S$).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			51	23	7	17	3	1		

- Molecule 3 is water.

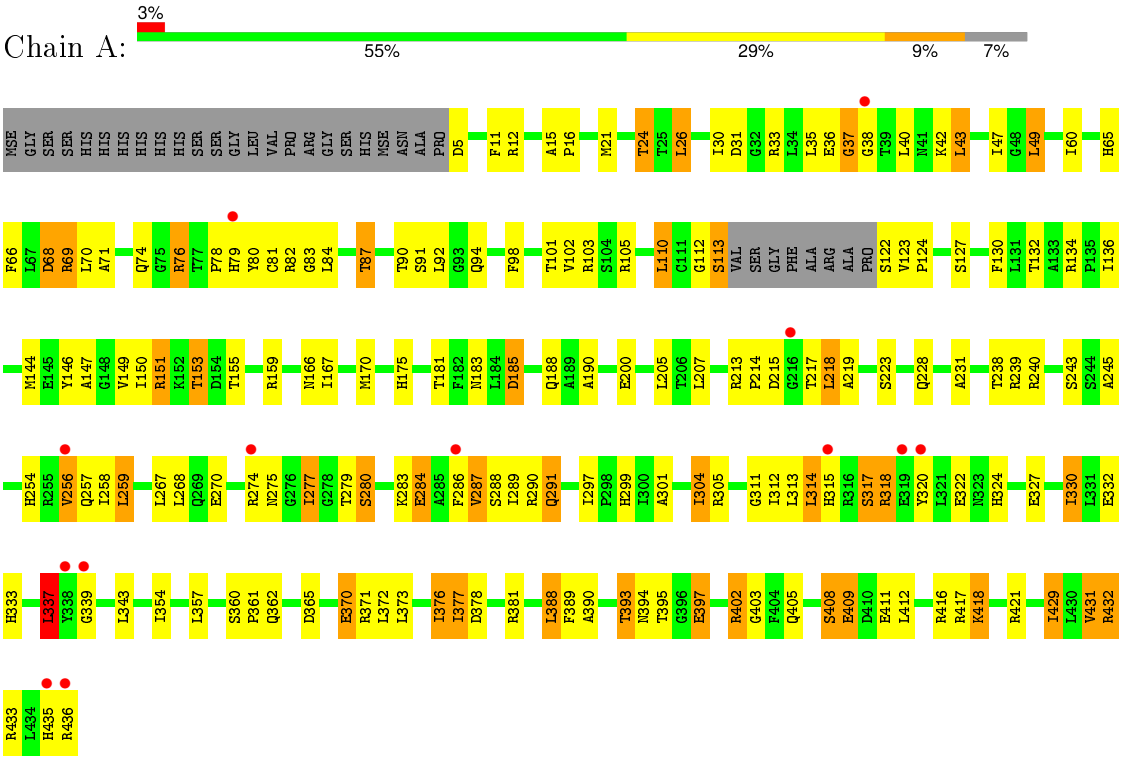
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	92	Total	O	0	0
			92	92		

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: Putative acetylglutamate synthase

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 3 1 2	Depositor
Cell constants a, b, c, α, β, γ	98.97Å 98.97Å 89.29Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.40 19.80 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.5 (20.00-2.40) 99.7 (19.80-2.40)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.06 (at 2.41Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.194 , 0.244 0.194 , 0.196	Depositor DCC
R_{free} test set	953 reflections (5.09%)	DCC
Wilson B-factor (Å^2)	57.0	Xtriage
Anisotropy	0.267	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 48.1	EDS
Estimated twinning fraction	0.041 for -h,-k,l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 19676 reflections	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	3370	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% of the height of the origin peak. No significant pseudotranslation is detected.*

[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: ACO

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	0.52	0/3276	1.01	9/4421 (0.2%)

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	5

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	12	ARG	NE-CZ-NH1	-8.33	116.13	120.30
1	A	337	LEU	CA-CB-CG	6.66	130.62	115.30
1	A	110	LEU	CA-CB-CG	6.58	130.44	115.30
1	A	49	LEU	CB-CG-CD2	5.97	121.15	111.00
1	A	431	VAL	CB-CA-C	-5.89	100.21	111.40
1	A	68	ASP	CB-CG-OD2	5.71	123.44	118.30
1	A	259	LEU	CB-CG-CD2	-5.50	101.64	111.00
1	A	26	LEU	CB-CG-CD1	-5.22	102.12	111.00
1	A	432	ARG	NE-CZ-NH2	-5.04	117.78	120.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	122	SER	Peptide

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Mol	Chain	Res	Type	Group
1	A	283	LYS	Peptide
1	A	37	GLY	Peptide
1	A	390	ALA	Mainchain
1	A	431	VAL	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	3227	0	3208	167	0
2	A	51	0	34	7	0
3	A	92	0	0	14	0
All	All	3370	0	3242	170	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 26.

All (170) 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:377:ILE:CD1	1:A:381:ARG:HH21	1.53	1.21
1:A:435:HIS:HB2	3:A:514:HOH:O	1.46	1.13
1:A:377:ILE:HD11	1:A:381:ARG:HH21	1.10	1.12
1:A:112:GLY:O	1:A:113:SER:HB2	1.47	1.09
1:A:291:GLN:HG2	3:A:480:HOH:O	1.54	1.07
1:A:377:ILE:HD11	1:A:381:ARG:NH2	1.70	1.06
1:A:239:ARG:HH21	1:A:243:SER:HB2	1.18	1.05
1:A:24:THR:HG23	3:A:491:HOH:O	1.53	1.05
1:A:393:THR:HG23	1:A:394:ASN:ND2	1.72	1.05
1:A:370:GLU:OE2	1:A:402:ARG:HD3	1.58	1.04
1:A:103:ARG:NH2	1:A:127:SER:HB3	1.71	1.04
1:A:103:ARG:HH21	1:A:127:SER:HB3	0.90	1.03
1:A:91:SER:HA	1:A:94:GLN:HE21	1.26	1.00
1:A:277:ILE:HD12	1:A:277:ILE:N	1.76	0.98
1:A:151:ARG:HG3	1:A:151:ARG:HH11	1.25	0.98
1:A:231:ALA:HB1	1:A:239:ARG:HD2	1.46	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:393:THR:CG2	1:A:394:ASN:ND2	2.29	0.95
1:A:98:PHE:O	1:A:101:THR:HG22	1.67	0.94
1:A:405:GLN:HE22	1:A:433:ARG:NH1	1.65	0.94
1:A:103:ARG:HH21	1:A:127:SER:CB	1.80	0.94
1:A:287:VAL:HG21	1:A:371:ARG:HD3	1.49	0.94
1:A:277:ILE:HD12	1:A:277:ILE:H	1.27	0.93
1:A:245:ALA:HB1	1:A:256:VAL:HG21	1.51	0.92
1:A:71:ALA:O	1:A:76:ARG:O	1.92	0.87
1:A:151:ARG:HG3	1:A:151:ARG:NH1	1.86	0.86
1:A:397:GLU:OE2	2:A:1:ACO:C2A	2.24	0.85
1:A:397:GLU:OE2	2:A:1:ACO:N1A	2.09	0.85
1:A:377:ILE:CD1	1:A:381:ARG:NH2	2.35	0.84
1:A:123:VAL:CG1	1:A:167:ILE:HD12	2.09	0.82
1:A:151:ARG:CG	1:A:151:ARG:HH11	1.92	0.81
1:A:218:LEU:HD23	1:A:219:ALA:N	1.96	0.80
1:A:84:LEU:HD13	1:A:136:ILE:CD1	2.11	0.80
1:A:228:GLN:NE2	1:A:239:ARG:HH12	1.79	0.79
1:A:123:VAL:HG12	1:A:167:ILE:HD12	1.63	0.79
1:A:397:GLU:OE2	2:A:1:ACO:H2A	1.83	0.77
1:A:33:ARG:CZ	1:A:207:LEU:HD22	2.15	0.77
1:A:239:ARG:NH2	1:A:243:SER:HB2	1.98	0.76
1:A:36:GLU:HG2	1:A:105:ARG:HH22	1.50	0.75
1:A:376:ILE:HD11	1:A:388:LEU:HD11	1.68	0.75
1:A:21:MSE:O	1:A:24:THR:HB	1.87	0.74
1:A:377:ILE:HD12	1:A:381:ARG:HH21	1.49	0.74
1:A:123:VAL:HG11	1:A:167:ILE:CD1	2.18	0.74
1:A:123:VAL:CG1	1:A:167:ILE:CD1	2.66	0.74
1:A:31:ASP:OD2	1:A:33:ARG:HD3	1.89	0.72
1:A:36:GLU:HG2	1:A:105:ARG:NH2	2.05	0.72
1:A:218:LEU:C	1:A:218:LEU:HD23	2.08	0.72
1:A:270:GLU:HG2	1:A:277:ILE:CD1	2.20	0.71
1:A:213:ARG:HB3	1:A:214:PRO:HD2	1.71	0.71
1:A:231:ALA:CB	1:A:239:ARG:HH11	2.03	0.71
1:A:200:GLU:HA	1:A:254:HIS:CD2	2.26	0.71
1:A:405:GLN:HE22	1:A:433:ARG:HH12	1.38	0.70
1:A:84:LEU:HD13	1:A:136:ILE:HD13	1.74	0.70
1:A:76:ARG:HH22	1:A:90:THR:HB	1.57	0.69
1:A:393:THR:HG23	1:A:394:ASN:CG	2.13	0.68
1:A:123:VAL:HG11	1:A:167:ILE:HD11	1.74	0.68
1:A:228:GLN:HE22	1:A:239:ARG:HH12	1.40	0.68
1:A:287:VAL:HG12	1:A:289:ILE:HG13	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:CYS:O	1:A:82:ARG:HB3	1.93	0.66
1:A:35:LEU:HD13	1:A:102:VAL:CG1	2.26	0.66
1:A:90:THR:O	1:A:94:GLN:HG3	1.96	0.65
1:A:105:ARG:NH1	3:A:495:HOH:O	2.30	0.64
1:A:259:LEU:HD21	1:A:270:GLU:HG3	1.79	0.64
1:A:218:LEU:C	1:A:218:LEU:CD2	2.67	0.63
1:A:408:SER:HB2	1:A:411:GLU:OE1	1.98	0.63
1:A:101:THR:HG23	1:A:102:VAL:N	2.13	0.63
1:A:354:ILE:HD12	1:A:376:ILE:HG12	1.80	0.63
1:A:270:GLU:HG2	1:A:277:ILE:HD11	1.81	0.62
1:A:377:ILE:HG13	1:A:378:ASP:N	2.13	0.62
1:A:231:ALA:HB3	1:A:239:ARG:HH11	1.63	0.62
1:A:223:SER:HB2	1:A:284:GLU:OE1	2.00	0.62
1:A:66:PHE:HD1	1:A:69:ARG:HH21	1.47	0.61
1:A:258:ILE:C	1:A:259:LEU:HD23	2.20	0.61
1:A:277:ILE:CD1	1:A:277:ILE:N	2.51	0.61
1:A:284:GLU:O	1:A:284:GLU:HG2	2.00	0.61
1:A:409:GLU:O	1:A:417:ARG:HD3	2.00	0.60
1:A:245:ALA:CB	1:A:256:VAL:HG21	2.28	0.60
1:A:21:MSE:HE1	1:A:26:LEU:HD12	1.83	0.59
1:A:405:GLN:NE2	1:A:433:ARG:NH1	2.45	0.59
1:A:82:ARG:HG3	1:A:82:ARG:O	2.01	0.59
1:A:37:GLY:O	3:A:474:HOH:O	2.17	0.59
1:A:24:THR:CG2	3:A:491:HOH:O	2.27	0.58
1:A:112:GLY:O	1:A:113:SER:CB	2.31	0.58
1:A:35:LEU:HD13	1:A:102:VAL:HG11	1.86	0.58
1:A:82:ARG:CG	3:A:492:HOH:O	2.52	0.58
1:A:33:ARG:NH2	1:A:207:LEU:HD22	2.20	0.57
1:A:33:ARG:HD2	1:A:65:HIS:CG	2.40	0.57
1:A:38:GLY:HA2	3:A:474:HOH:O	2.05	0.56
1:A:113:SER:C	3:A:456:HOH:O	2.43	0.56
1:A:91:SER:HA	1:A:94:GLN:NE2	2.09	0.56
1:A:327:GLU:HB2	1:A:343:LEU:O	2.06	0.56
1:A:395:THR:HG22	1:A:395:THR:O	2.06	0.56
1:A:254:HIS:HB3	3:A:498:HOH:O	2.05	0.56
1:A:405:GLN:NE2	1:A:433:ARG:HH12	2.01	0.55
1:A:389:PHE:HD1	1:A:429:ILE:CD1	2.19	0.55
1:A:304:ILE:HG13	1:A:313:LEU:CD1	2.35	0.55
1:A:412:LEU:HD21	1:A:429:ILE:HD13	1.87	0.55
1:A:317:SER:HB3	1:A:320:TYR:CD1	2.42	0.54
1:A:259:LEU:CD2	1:A:270:GLU:HG3	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:257:GLN:OE1	1:A:280:SER:HB2	2.07	0.54
1:A:15:ALA:HB3	1:A:16:PRO:HD3	1.90	0.53
1:A:324:HIS:O	1:A:327:GLU:HG2	2.08	0.53
1:A:101:THR:CG2	1:A:102:VAL:N	2.72	0.53
1:A:82:ARG:HG2	3:A:492:HOH:O	2.07	0.52
1:A:239:ARG:HH21	1:A:243:SER:CB	2.07	0.52
1:A:83:GLY:HA3	1:A:240:ARG:NE	2.24	0.52
1:A:318:ARG:O	1:A:322:GLU:HG3	2.10	0.52
1:A:30:ILE:HG12	1:A:205:LEU:HD12	1.91	0.52
1:A:35:LEU:HD13	1:A:102:VAL:HG13	1.93	0.51
1:A:228:GLN:HE22	1:A:239:ARG:NH1	2.08	0.51
1:A:317:SER:HB3	1:A:320:TYR:CE1	2.46	0.51
1:A:286:PHE:O	1:A:287:VAL:HG23	2.12	0.51
1:A:35:LEU:O	1:A:36:GLU:HG3	2.12	0.50
2:A:1:ACO:OAP	2:A:1:ACO:O6A	2.30	0.50
1:A:185:ASP:OD1	1:A:188:GLN:HG2	2.11	0.50
1:A:144:MSE:HB3	1:A:147:ALA:HB3	1.95	0.49
1:A:403:GLY:O	1:A:432:ARG:HG3	2.12	0.49
1:A:320:TYR:O	1:A:324:HIS:CD2	2.65	0.49
1:A:66:PHE:HD1	1:A:69:ARG:NH2	2.09	0.49
1:A:103:ARG:NH2	1:A:127:SER:CB	2.54	0.48
1:A:70:LEU:HD23	1:A:70:LEU:HA	1.57	0.48
1:A:123:VAL:HG13	1:A:124:PRO:HD2	1.96	0.47
1:A:286:PHE:O	1:A:287:VAL:CG2	2.62	0.47
1:A:330:ILE:HB	1:A:337:LEU:HD22	1.97	0.47
1:A:297:ILE:HD13	1:A:322:GLU:HG2	1.96	0.47
1:A:155:THR:HG22	1:A:159:ARG:HH21	1.78	0.47
1:A:207:LEU:HD12	1:A:207:LEU:N	2.29	0.47
1:A:228:GLN:NE2	1:A:239:ARG:NH1	2.56	0.47
1:A:132:THR:HG23	3:A:477:HOH:O	2.15	0.46
1:A:213:ARG:HB3	1:A:214:PRO:CD	2.41	0.46
1:A:299:HIS:HB3	1:A:337:LEU:HG	1.96	0.46
1:A:409:GLU:OE2	1:A:421:ARG:NE	2.48	0.46
1:A:98:PHE:O	1:A:101:THR:CG2	2.50	0.45
1:A:259:LEU:HD21	1:A:270:GLU:CG	2.46	0.45
1:A:82:ARG:HG3	3:A:492:HOH:O	2.14	0.45
1:A:36:GLU:HA	1:A:105:ARG:NH2	2.31	0.45
1:A:130:PHE:HB2	1:A:170:MSE:CE	2.46	0.45
1:A:301:ALA:O	1:A:305:ARG:HG3	2.17	0.45
1:A:66:PHE:CD1	1:A:69:ARG:NH2	2.84	0.45
1:A:60:ILE:HD12	1:A:190:ALA:HA	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:THR:O	1:A:146:TYR:HB2	2.17	0.44
1:A:136:ILE:HG22	1:A:136:ILE:O	2.17	0.44
1:A:155:THR:HG22	1:A:159:ARG:NH2	2.33	0.44
1:A:360:SER:HA	1:A:361:PRO:HD3	1.79	0.44
1:A:259:LEU:CD2	1:A:270:GLU:CG	2.96	0.43
1:A:357:LEU:O	2:A:1:ACO:N4P	2.42	0.43
1:A:312:ILE:HA	1:A:394:ASN:HD22	1.83	0.43
1:A:98:PHE:HA	1:A:101:THR:HG22	2.00	0.43
2:A:1:ACO:H8A	3:A:487:HOH:O	2.17	0.43
1:A:311:GLY:O	1:A:393:THR:HG21	2.19	0.43
1:A:314:LEU:O	1:A:315:HIS:C	2.57	0.43
1:A:416:ARG:HA	1:A:416:ARG:HD2	1.78	0.43
1:A:74:GLN:HE21	1:A:76:ARG:HD3	1.84	0.42
1:A:343:LEU:HD13	1:A:376:ILE:HD13	2.01	0.42
1:A:43:LEU:HD22	1:A:47:ILE:HD11	2.01	0.42
1:A:339:GLY:HA3	1:A:372:LEU:HD11	2.01	0.42
1:A:175:HIS:CE1	1:A:181:THR:HG23	2.55	0.42
1:A:150:ILE:HG21	1:A:153:THR:HG22	2.01	0.42
1:A:43:LEU:HD22	1:A:47:ILE:CD1	2.49	0.42
1:A:124:PRO:HB2	1:A:166:ASN:HD22	1.85	0.41
1:A:318:ARG:HE	1:A:318:ARG:HB2	1.43	0.41
1:A:78:PRO:HB2	1:A:80:TYR:HE1	1.85	0.41
1:A:418:LYS:HD2	1:A:418:LYS:HA	1.69	0.41
1:A:231:ALA:CB	1:A:239:ARG:NH1	2.79	0.41
1:A:219:ALA:O	1:A:279:THR:HG23	2.21	0.41
1:A:134:ARG:O	1:A:134:ARG:HG3	2.20	0.41
1:A:68:ASP:N	1:A:68:ASP:OD1	2.54	0.41
1:A:218:LEU:HD23	1:A:219:ALA:C	2.41	0.41
2:A:1:ACO:O9P	2:A:1:ACO:H62	2.21	0.40
1:A:373:LEU:O	1:A:377:ILE:HG23	2.21	0.40
1:A:218:LEU:CD2	1:A:219:ALA:C	2.89	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	420/456 (92%)	390 (93%)	25 (6%)	5 (1%)	16 23

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	275	ASN
1	A	288	SER
1	A	284	GLU
1	A	333	HIS
1	A	287	VAL

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	330/349 (95%)	278 (84%)	52 (16%)	3 3

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ASP
1	A	11	PHE
1	A	24	THR
1	A	40	LEU
1	A	42	LYS
1	A	43	LEU
1	A	49	LEU
1	A	69	ARG
1	A	76	ARG
1	A	79	HIS
1	A	87	THR
1	A	92	LEU
1	A	110	LEU
1	A	113	SER

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Mol	Chain	Res	Type
1	A	149	VAL
1	A	151	ARG
1	A	153	THR
1	A	183	ASN
1	A	185	ASP
1	A	215	ASP
1	A	217	THR
1	A	218	LEU
1	A	238	THR
1	A	256	VAL
1	A	267	LEU
1	A	268	LEU
1	A	274	ARG
1	A	277	ILE
1	A	280	SER
1	A	290	ARG
1	A	291	GLN
1	A	304	ILE
1	A	314	LEU
1	A	317	SER
1	A	318	ARG
1	A	330	ILE
1	A	332	GLU
1	A	337	LEU
1	A	362	GLN
1	A	365	ASP
1	A	370	GLU
1	A	376	ILE
1	A	377	ILE
1	A	388	LEU
1	A	393	THR
1	A	397	GLU
1	A	402	ARG
1	A	408	SER
1	A	409	GLU
1	A	418	LYS
1	A	429	ILE
1	A	436	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	GLN
1	A	94	GLN
1	A	161	GLN
1	A	166	ASN
1	A	198	GLN
1	A	225	GLN
1	A	228	GLN
1	A	254	HIS
1	A	257	GLN
1	A	275	ASN
1	A	324	HIS
1	A	362	GLN
1	A	394	ASN
1	A	405	GLN
1	A	423	ASN

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

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	$\#\|Z\|>2$
2	ACO	A	1	-	43,53,53	1.25	3 (6%)	55,79,79	2.06	11 (20%)

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	ACO	A	1	-	1/1/12/14	0/47/67/67	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	ACO	CH3-C	-2.30	1.40	1.50
2	A	1	ACO	P3B-O7A	3.86	1.63	1.51
2	A	1	ACO	O4B-C1B	4.22	1.46	1.41

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1	ACO	N3A-C2A-N1A	-9.45	121.66	128.89
2	A	1	ACO	C2B-C1B-N9A	-5.96	105.19	114.29
2	A	1	ACO	C6P-C5P-N4P	-3.56	110.27	116.46
2	A	1	ACO	O3A-P2A-O6A	-2.94	95.12	102.94
2	A	1	ACO	C7P-N8P-C9P	-2.72	117.15	122.53
2	A	1	ACO	O3A-P1A-O5B	-2.69	95.80	102.94
2	A	1	ACO	C4A-C5A-N7A	-2.31	107.35	109.48
2	A	1	ACO	C4B-O4B-C1B	-2.21	107.29	109.72
2	A	1	ACO	O8A-P3B-O7A	2.15	117.51	110.58
2	A	1	ACO	O5P-C5P-C6P	2.29	125.93	121.98
2	A	1	ACO	O6A-CCP-CBP	2.90	115.22	110.55

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	1	ACO	CAP

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1	ACO	7	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

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	420/456 (92%)	-0.04	13 (3%) 52 52	27, 43, 61, 78	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	436	ARG	4.8
1	A	286	PHE	4.2
1	A	320	TYR	3.3
1	A	319	GLU	3.1
1	A	79	HIS	3.1
1	A	315	HIS	2.8
1	A	274	ARG	2.8
1	A	339	GLY	2.7
1	A	338	TYR	2.5
1	A	256	VAL	2.3
1	A	38	GLY	2.3
1	A	216	GLY	2.2
1	A	435	HIS	2.2

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

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	ACO	A	1	51/51	0.97	0.08	-1.63	38,53,64,67	0

6.5 Other polymers

There are no such residues in this entry.