Full wwPDB X-ray Structure Validation Report (i)

Jun 18, 2024 – 11:24 PM EDT

PDB ID	:	4AMA
Title	:	Crystal Structure of N-acetylneuraminic acid lyase from Staphylococcus aureus with the chemical modification thia-lysine at position 165 in complex with pyruvate
Authors	:	Timms, N.; Polyakova, A.; Windle, C.L.; Trinh, C.H.; Nelson, A.; Pearson, A.R.; Berry, A.
Deposited on	:	2012-03-08
Resolution	:	2.35 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

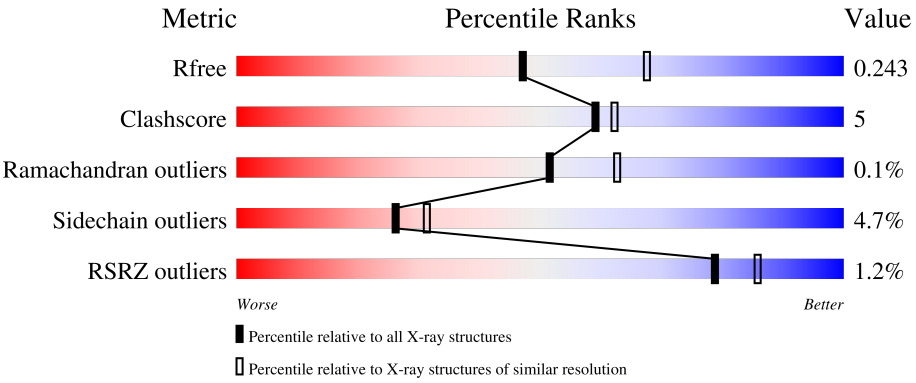
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 2.35 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}	130704	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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	298	87% 10%
1	B	298	% 85% 9%
1	C	298	2% 86% 12%
1	D	298	2% 88% 8%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 9639 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 N-ACETYLNEURAMINATE LYASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	293	Total	C	N	O	S	0	0	0
			2324	1488	387	445	4			
1	B	294	Total	C	N	O	S	0	0	0
			2315	1482	386	443	4			
1	C	298	Total	C	N	O	S	0	0	0
			2344	1501	393	446	4			
1	D	291	Total	C	N	O	S	0	0	0
			2300	1475	381	440	4			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	HIS	-	expression tag	UNP Q2G160
A	-3	HIS	-	expression tag	UNP Q2G160
A	-2	HIS	-	expression tag	UNP Q2G160
A	-1	HIS	-	expression tag	UNP Q2G160
A	0	HIS	-	expression tag	UNP Q2G160
A	1	HIS	-	expression tag	UNP Q2G160
B	-4	HIS	-	expression tag	UNP Q2G160
B	-3	HIS	-	expression tag	UNP Q2G160
B	-2	HIS	-	expression tag	UNP Q2G160
B	-1	HIS	-	expression tag	UNP Q2G160
B	0	HIS	-	expression tag	UNP Q2G160
B	1	HIS	-	expression tag	UNP Q2G160
C	-4	HIS	-	expression tag	UNP Q2G160
C	-3	HIS	-	expression tag	UNP Q2G160
C	-2	HIS	-	expression tag	UNP Q2G160
C	-1	HIS	-	expression tag	UNP Q2G160
C	0	HIS	-	expression tag	UNP Q2G160
C	1	HIS	-	expression tag	UNP Q2G160
D	-4	HIS	-	expression tag	UNP Q2G160
D	-3	HIS	-	expression tag	UNP Q2G160
D	-2	HIS	-	expression tag	UNP Q2G160

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	HIS	-	expression tag	UNP Q2G160
D	0	HIS	-	expression tag	UNP Q2G160
D	1	HIS	-	expression tag	UNP Q2G160

- Molecule 2 is water.

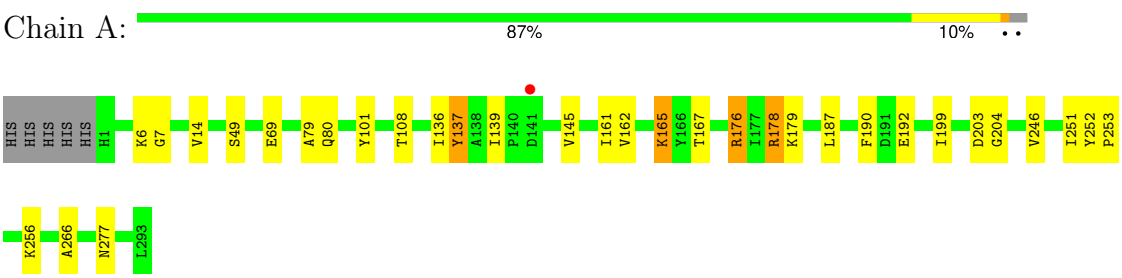
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	100	Total O 100 100	0	0
2	B	96	Total O 96 96	0	0
2	C	56	Total O 56 56	0	0
2	D	104	Total O 104 104	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

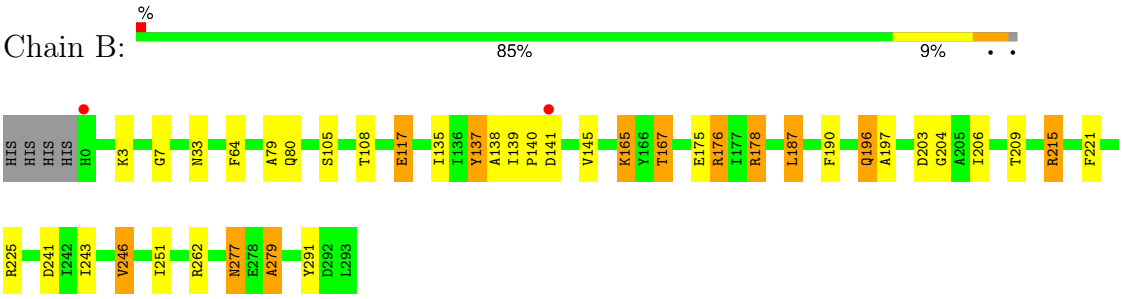
• Molecule 1: N-ACETYLNEURAMINATE LYASE

Chain A:



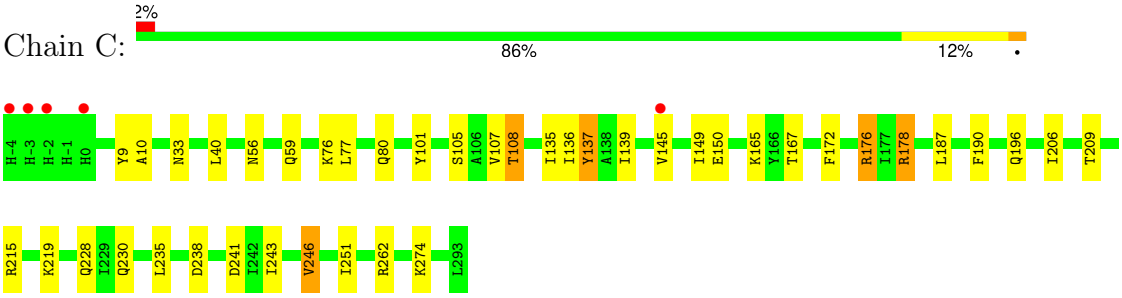
• Molecule 1: N-ACETYLNEURAMINATE LYASE

Chain B:



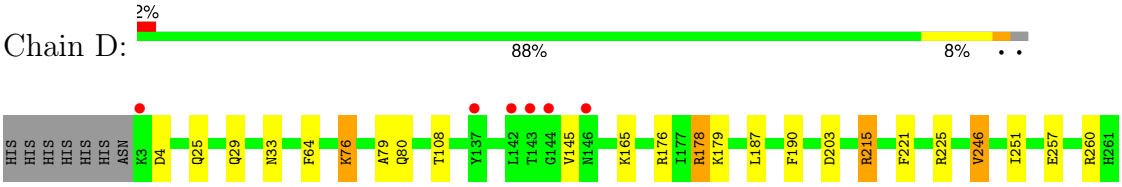
• Molecule 1: N-ACETYLNEURAMINATE LYASE

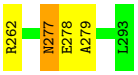
Chain C:



• Molecule 1: N-ACETYLNEURAMINATE LYASE

Chain D:




R262
N277
E278
A279
L293

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	82.94Å 110.28Å 132.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.63 – 2.35 29.63 – 2.35	Depositor EDS
% Data completeness (in resolution range)	98.2 (29.63-2.35) 98.4 (29.63-2.35)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.92 (at 2.36Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.189 , 0.244 0.189 , 0.243	Depositor DCC
R_{free} test set	2506 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å^2)	25.7	Xtriage
Anisotropy	0.423	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 28.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	9639	wwPDB-VP
Average B, all atoms (Å^2)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.76% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: KPY

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	0.70	0/2353	0.77	0/3182
1	B	0.66	0/2344	0.74	0/3172
1	C	0.69	0/2376	0.75	1/3219 (0.0%)
1	D	0.70	0/2328	0.77	0/3150
All	All	0.69	0/9401	0.76	1/12723 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	136	ILE	CB-CA-C	-6.12	99.36	111.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2324	0	2292	19	0
1	B	2315	0	2270	29	0
1	C	2344	0	2282	23	0
1	D	2300	0	2267	20	0
2	A	100	0	0	0	0
2	B	96	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	56	0	0	1	0
2	D	104	0	0	0	0
All	All	9639	0	9111	87	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 5.

All (87) 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:D:215:ARG:HH11	1:D:215:ARG:HG2	1.28	0.98
1:B:33:ASN:HD21	1:B:262:ARG:HH11	1.23	0.87
1:B:139:ILE:HG13	1:B:167:THR:HG21	1.64	0.79
1:A:6:LYS:HG2	1:A:203:ASP:HB3	1.65	0.78
1:D:277:ASN:ND2	1:D:279:ALA:H	1.88	0.71
1:C:108:THR:HG21	1:C:145:VAL:HB	1.73	0.69
1:D:33:ASN:HD21	1:D:262:ARG:HH11	1.38	0.68
1:D:178:ARG:O	1:D:178:ARG:HD3	1.95	0.66
1:D:215:ARG:HH11	1:D:215:ARG:CG	2.06	0.64
1:C:33:ASN:HD21	1:C:262:ARG:HH11	1.45	0.63
1:B:138:ALA:O	1:B:167:THR:HB	1.99	0.62
1:B:209:THR:HG21	1:B:243:ILE:HG12	1.81	0.62
1:C:139:ILE:HA	1:C:167:THR:OG1	2.02	0.60
1:B:33:ASN:HD21	1:B:262:ARG:NH1	1.98	0.60
1:C:107:VAL:HA	1:C:137:TYR:HB3	1.85	0.59
1:D:277:ASN:HD22	1:D:279:ALA:H	1.48	0.59
1:A:192:GLU:HG3	1:C:172:PHE:CD2	2.39	0.58
1:C:209:THR:HG21	1:C:243:ILE:HG12	1.84	0.58
1:D:215:ARG:HG2	1:D:215:ARG:NH1	2.08	0.57
1:D:277:ASN:HD22	1:D:277:ASN:C	2.08	0.56
1:B:105:SER:HB2	1:B:135:ILE:O	2.05	0.56
1:D:257:GLU:OE1	1:D:260:ARG:NH1	2.39	0.56
1:B:277:ASN:ND2	1:B:279:ALA:HB3	2.21	0.56
1:B:196:GLN:NE2	1:B:196:GLN:H	2.05	0.55
1:B:221:PHE:O	1:B:225:ARG:HG3	2.06	0.55
1:A:199:ILE:HD11	1:C:196:GLN:HG3	1.87	0.55
1:C:219:LYS:HB3	1:C:235:LEU:CD1	2.37	0.55
1:D:246:VAL:CG1	1:D:251:ILE:HG13	2.37	0.55
1:B:187:LEU:HD23	1:B:206:ILE:HD11	1.88	0.54
1:C:56:ASN:OD1	1:C:59:GLN:HG3	2.08	0.54
1:B:175:GLU:OE1	1:B:176:ARG:HD3	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:178:ARG:HD3	1:C:178:ARG:O	2.09	0.52
1:B:215:ARG:HG2	1:B:215:ARG:HH11	1.74	0.52
1:D:277:ASN:HD22	1:D:278:GLU:N	2.07	0.52
1:D:246:VAL:HG13	1:D:251:ILE:HG13	1.91	0.52
1:B:196:GLN:H	1:B:196:GLN:HE21	1.58	0.51
1:A:136:ILE:HD11	1:A:161:ILE:HD13	1.93	0.51
1:B:246:VAL:HG13	1:B:251:ILE:HG13	1.93	0.50
1:D:4:ASP:O	1:D:76:LYS:NZ	2.43	0.50
1:D:108:THR:HG21	1:D:145:VAL:HG11	1.94	0.49
1:D:25:GLN:O	1:D:29:GLN:HG3	2.13	0.49
1:C:178:ARG:HD3	1:C:178:ARG:C	2.27	0.49
1:B:176:ARG:HA	1:B:176:ARG:HD2	1.59	0.48
1:B:139:ILE:HA	1:B:167:THR:CG2	2.44	0.48
1:B:139:ILE:HA	1:B:167:THR:HG21	1.96	0.48
1:C:33:ASN:HD21	1:C:262:ARG:NH1	2.10	0.48
1:D:178:ARG:NH2	1:D:203:ASP:OD2	2.45	0.47
1:B:190:PHE:HE1	1:B:197:ALA:HB2	1.78	0.47
1:A:256:LYS:HB3	1:A:266:ALA:HB1	1.96	0.47
1:D:221:PHE:O	1:D:225:ARG:HG3	2.13	0.47
1:A:14:VAL:O	1:A:256:LYS:HE3	2.15	0.47
1:A:252:TYR:N	1:A:253:PRO:HD2	2.29	0.47
1:D:64:PHE:CE1	1:D:79:ALA:HB1	2.49	0.47
1:A:176:ARG:HA	1:A:176:ARG:HD2	1.55	0.46
1:C:246:VAL:CG1	1:C:251:ILE:HG13	2.45	0.46
1:B:117:GLU:HG3	1:C:274:LYS:HE3	1.98	0.46
1:A:246:VAL:CG1	1:A:251:ILE:HG13	2.45	0.46
1:B:178:ARG:O	1:B:178:ARG:HD3	2.15	0.46
1:D:176:ARG:HA	1:D:176:ARG:HD2	1.69	0.45
1:D:178:ARG:O	1:D:178:ARG:CD	2.62	0.45
1:A:192:GLU:HG3	1:C:172:PHE:CG	2.51	0.45
1:A:108:THR:HG21	1:A:145:VAL:HG11	1.98	0.45
1:B:108:THR:HG21	1:B:145:VAL:HB	1.99	0.45
1:B:7:GLY:O	1:B:204:GLY:HA3	2.17	0.44
1:A:178:ARG:HD3	1:A:178:ARG:HA	1.50	0.44
1:A:49:SER:OG	1:A:165:KPY:OL2	2.23	0.44
1:A:139:ILE:HA	1:A:167:THR:OG1	2.18	0.44
1:C:105:SER:HB3	1:C:135:ILE:HB	2.00	0.44
1:C:176:ARG:HA	1:C:176:ARG:HD2	1.65	0.44
1:C:9:TYR:HB2	1:C:206:ILE:HG12	1.99	0.43
1:C:77:LEU:HB3	1:C:101:TYR:CD2	2.53	0.43
1:B:241:ASP:OD2	1:B:291:TYR:OH	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:178:ARG:HD3	1:B:178:ARG:HA	1.59	0.43
1:B:137:TYR:CE1	1:B:165:KPY:HD2C	2.54	0.43
1:A:178:ARG:NH2	1:A:203:ASP:OD2	2.51	0.42
1:C:238:ASP:O	1:C:241:ASP:HB2	2.19	0.42
1:A:79:ALA:HB2	1:A:101:TYR:CD2	2.54	0.42
1:A:108:THR:HG21	1:A:145:VAL:CB	2.50	0.42
1:B:277:ASN:ND2	1:B:279:ALA:H	2.16	0.42
1:C:10:ALA:HB2	1:C:40:LEU:CD1	2.50	0.42
1:B:140:PRO:HD2	1:B:167:THR:HG22	2.02	0.42
1:A:137:TYR:C	1:A:137:TYR:CD1	2.94	0.41
1:C:219:LYS:HB3	1:C:235:LEU:HD13	2.02	0.41
1:C:215:ARG:HB2	2:C:2012:HOH:O	2.20	0.40
1:B:178:ARG:NH2	1:B:203:ASP:OD2	2.55	0.40
1:A:7:GLY:O	1:A:204:GLY:HA3	2.22	0.40
1:B:64:PHE:CE1	1:B:79:ALA:HB1	2.57	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	290/298 (97%)	284 (98%)	6 (2%)	0	100	100
1	B	291/298 (98%)	285 (98%)	5 (2%)	1 (0%)	41	47
1	C	295/298 (99%)	288 (98%)	7 (2%)	0	100	100
1	D	288/298 (97%)	282 (98%)	6 (2%)	0	100	100
All	All	1164/1192 (98%)	1139 (98%)	24 (2%)	1 (0%)	51	63

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	279	ALA

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	243/253 (96%)	233 (96%)	10 (4%)	30	37
1	B	240/253 (95%)	227 (95%)	13 (5%)	22	25
1	C	242/253 (96%)	229 (95%)	13 (5%)	22	25
1	D	240/253 (95%)	231 (96%)	9 (4%)	33	41
All	All	965/1012 (95%)	920 (95%)	45 (5%)	26	31

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	69	GLU
1	A	80	GLN
1	A	137	TYR
1	A	162	VAL
1	A	176	ARG
1	A	178	ARG
1	A	179	LYS
1	A	187	LEU
1	A	190	PHE
1	A	277	ASN
1	B	3	LYS
1	B	80	GLN
1	B	117	GLU
1	B	137	TYR
1	B	141	ASP
1	B	167	THR
1	B	176	ARG
1	B	178	ARG
1	B	187	LEU
1	B	196	GLN
1	B	215	ARG

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Mol	Chain	Res	Type
1	B	246	VAL
1	B	277	ASN
1	C	76	LYS
1	C	80	GLN
1	C	108	THR
1	C	137	TYR
1	C	149	ILE
1	C	150	GLU
1	C	176	ARG
1	C	178	ARG
1	C	187	LEU
1	C	190	PHE
1	C	228	GLN
1	C	230	GLN
1	C	246	VAL
1	D	76	LYS
1	D	80	GLN
1	D	178	ARG
1	D	179	LYS
1	D	187	LEU
1	D	190	PHE
1	D	215	ARG
1	D	246	VAL
1	D	277	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	ASN
1	A	277	ASN
1	A	280	HIS
1	B	33	ASN
1	B	196	GLN
1	B	226	GLN
1	B	277	ASN
1	C	33	ASN
1	D	33	ASN
1	D	277	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

4 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	KPY	A	165	1	11,13,14	1.70	3 (27%)	11,15,17	2.06	4 (36%)
1	KPY	D	165	1	11,13,14	2.05	3 (27%)	11,15,17	2.43	4 (36%)
1	KPY	B	165	1	11,13,14	2.38	3 (27%)	11,15,17	1.79	4 (36%)
1	KPY	C	165	1	11,13,14	1.63	3 (27%)	11,15,17	2.05	1 (9%)

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	KPY	A	165	1	-	0/12/14/16	-
1	KPY	D	165	1	-	0/12/14/16	-
1	KPY	B	165	1	-	1/12/14/16	-
1	KPY	C	165	1	-	0/12/14/16	-

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	165	KPY	CL-CI	6.30	1.57	1.49
1	D	165	KPY	CL-CI	4.10	1.54	1.49
1	D	165	KPY	CK-CI	3.98	1.57	1.49
1	A	165	KPY	CK-CI	3.71	1.57	1.49
1	C	165	KPY	CK-CI	3.50	1.56	1.49
1	B	165	KPY	CK-CI	3.26	1.56	1.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	165	KPY	OL1-CL	-2.89	1.22	1.30
1	A	165	KPY	CL-CI	2.87	1.53	1.49
1	C	165	KPY	CL-CI	2.78	1.53	1.49
1	B	165	KPY	OL1-CL	-2.45	1.23	1.30
1	C	165	KPY	OL1-CL	-2.33	1.24	1.30
1	A	165	KPY	CB-SG	2.04	1.88	1.81

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	165	KPY	CD-CE-NZ	-5.85	105.56	110.75
1	D	165	KPY	CK-CI-CL	5.68	123.47	118.11
1	A	165	KPY	CD-CE-NZ	-4.83	106.47	110.75
1	D	165	KPY	OL2-CL-CI	3.47	125.58	121.35
1	B	165	KPY	CD-CE-NZ	-3.44	107.70	110.75
1	D	165	KPY	OL1-CL-OL2	-2.64	117.62	123.90
1	D	165	KPY	CD-CE-NZ	-2.58	108.46	110.75
1	A	165	KPY	OL1-CL-CI	2.53	121.89	116.50
1	B	165	KPY	OL1-CL-OL2	-2.46	118.03	123.90
1	A	165	KPY	CK-CI-CL	2.37	120.35	118.11
1	B	165	KPY	OL1-CL-CI	2.30	121.41	116.50
1	B	165	KPY	CL-CI-NZ	2.29	120.28	114.88
1	A	165	KPY	OL2-CL-CI	-2.00	118.90	121.35

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	165	KPY	CA-CB-SG-CD

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	165	KPY	1	0
1	B	165	KPY	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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	292/298 (97%)	-0.40	1 (0%)	94	97	16, 25, 45, 65	0
1	B	293/298 (98%)	-0.37	2 (0%)	87	92	17, 26, 43, 71	0
1	C	297/298 (99%)	-0.27	5 (1%)	70	78	18, 26, 45, 79	0
1	D	290/298 (97%)	-0.41	6 (2%)	63	74	14, 21, 43, 65	0
All	All	1172/1192 (98%)	-0.36	14 (1%)	79	86	14, 24, 45, 79	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	-3	HIS	3.3
1	D	137	TYR	3.3
1	C	-4	HIS	3.2
1	C	-2	HIS	3.2
1	C	0	HIS	3.0
1	A	141	ASP	2.9
1	B	0	HIS	2.7
1	D	142	LEU	2.4
1	D	144	GLY	2.2
1	B	141	ASP	2.2
1	D	146	ASN	2.1
1	C	145	VAL	2.1
1	D	3	LYS	2.0
1	D	143	THR	2.0

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

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. 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	B-factors(Å^2)	Q<0.9
1	KPY	B	165	14/15	0.95	0.20	21,27,29,30	0
1	KPY	C	165	14/15	0.95	0.18	22,29,33,33	0
1	KPY	A	165	14/15	0.96	0.17	24,29,31,32	0
1	KPY	D	165	14/15	0.97	0.17	19,23,25,26	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.