



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 01:53 am BST

PDB ID : 6ARH
Title : Crystal structure of Human NAL at a resolution of 1.6 Angstrom
Authors : Pearce, F.G.; Bundela, R.; Keown, J.R.
Deposited on : 2017-08-22
Resolution : 1.60 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

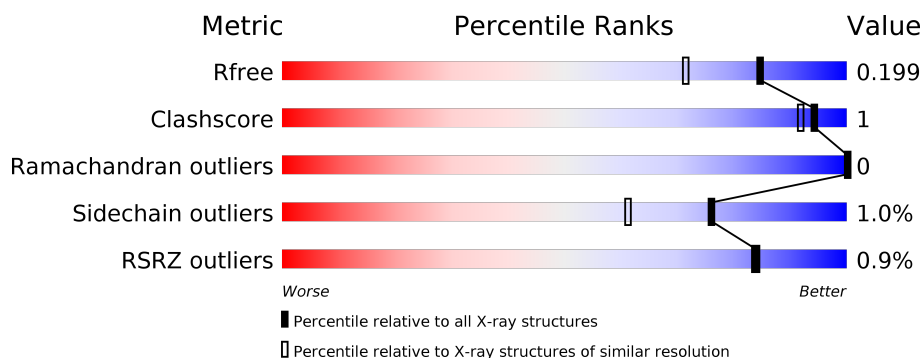
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	338	<div> <div>%</div> <div> <div></div> <div>85%</div> <div>11%</div> </div> </div>
1	B	338	<div> <div>%</div> <div> <div></div> <div>86%</div> <div>10%</div> </div> </div>
1	C	338	<div> <div>%</div> <div> <div></div> <div>86%</div> <div>11%</div> </div> </div>
1	D	338	<div> <div>%</div> <div> <div></div> <div>85%</div> <div>11%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10729 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-acetylneuraminase lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	300	Total	C	N	O	S	0	7	0
			2368	1529	389	442	8			
1	B	303	Total	C	N	O	S	0	4	0
			2359	1526	382	443	8			
1	C	300	Total	C	N	O	S	0	4	0
			2347	1516	382	441	8			
1	D	301	Total	C	N	O	S	0	4	0
			2352	1519	383	442	8			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	MET	-	expression tag	UNP Q9BXD5
A	-14	GLY	-	expression tag	UNP Q9BXD5
A	-13	HIS	-	expression tag	UNP Q9BXD5
A	-12	HIS	-	expression tag	UNP Q9BXD5
A	-11	HIS	-	expression tag	UNP Q9BXD5
A	-10	HIS	-	expression tag	UNP Q9BXD5
A	-9	HIS	-	expression tag	UNP Q9BXD5
A	-8	HIS	-	expression tag	UNP Q9BXD5
A	-7	GLY	-	expression tag	UNP Q9BXD5
A	-6	GLU	-	expression tag	UNP Q9BXD5
A	-5	ASN	-	expression tag	UNP Q9BXD5
A	-4	LEU	-	expression tag	UNP Q9BXD5
A	-3	TYR	-	expression tag	UNP Q9BXD5
A	-2	PHE	-	expression tag	UNP Q9BXD5
A	-1	GLN	-	expression tag	UNP Q9BXD5
A	0	GLY	-	expression tag	UNP Q9BXD5
A	1	GLY	-	expression tag	UNP Q9BXD5
A	2	SER	-	expression tag	UNP Q9BXD5
B	-15	MET	-	expression tag	UNP Q9BXD5
B	-14	GLY	-	expression tag	UNP Q9BXD5
B	-13	HIS	-	expression tag	UNP Q9BXD5

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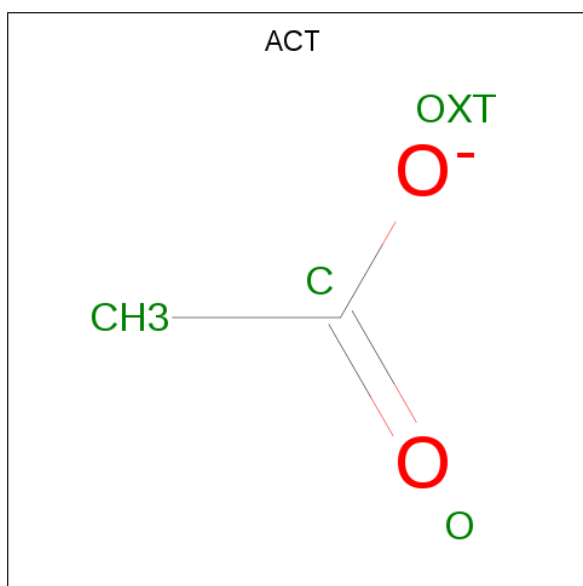
Chain	Residue	Modelled	Actual	Comment	Reference
B	-12	HIS	-	expression tag	UNP Q9BXD5
B	-11	HIS	-	expression tag	UNP Q9BXD5
B	-10	HIS	-	expression tag	UNP Q9BXD5
B	-9	HIS	-	expression tag	UNP Q9BXD5
B	-8	HIS	-	expression tag	UNP Q9BXD5
B	-7	GLY	-	expression tag	UNP Q9BXD5
B	-6	GLU	-	expression tag	UNP Q9BXD5
B	-5	ASN	-	expression tag	UNP Q9BXD5
B	-4	LEU	-	expression tag	UNP Q9BXD5
B	-3	TYR	-	expression tag	UNP Q9BXD5
B	-2	PHE	-	expression tag	UNP Q9BXD5
B	-1	GLN	-	expression tag	UNP Q9BXD5
B	0	GLY	-	expression tag	UNP Q9BXD5
B	1	GLY	-	expression tag	UNP Q9BXD5
B	2	SER	-	expression tag	UNP Q9BXD5
C	-15	MET	-	expression tag	UNP Q9BXD5
C	-14	GLY	-	expression tag	UNP Q9BXD5
C	-13	HIS	-	expression tag	UNP Q9BXD5
C	-12	HIS	-	expression tag	UNP Q9BXD5
C	-11	HIS	-	expression tag	UNP Q9BXD5
C	-10	HIS	-	expression tag	UNP Q9BXD5
C	-9	HIS	-	expression tag	UNP Q9BXD5
C	-8	HIS	-	expression tag	UNP Q9BXD5
C	-7	GLY	-	expression tag	UNP Q9BXD5
C	-6	GLU	-	expression tag	UNP Q9BXD5
C	-5	ASN	-	expression tag	UNP Q9BXD5
C	-4	LEU	-	expression tag	UNP Q9BXD5
C	-3	TYR	-	expression tag	UNP Q9BXD5
C	-2	PHE	-	expression tag	UNP Q9BXD5
C	-1	GLN	-	expression tag	UNP Q9BXD5
C	0	GLY	-	expression tag	UNP Q9BXD5
C	1	GLY	-	expression tag	UNP Q9BXD5
C	2	SER	-	expression tag	UNP Q9BXD5
D	-15	MET	-	expression tag	UNP Q9BXD5
D	-14	GLY	-	expression tag	UNP Q9BXD5
D	-13	HIS	-	expression tag	UNP Q9BXD5
D	-12	HIS	-	expression tag	UNP Q9BXD5
D	-11	HIS	-	expression tag	UNP Q9BXD5
D	-10	HIS	-	expression tag	UNP Q9BXD5
D	-9	HIS	-	expression tag	UNP Q9BXD5
D	-8	HIS	-	expression tag	UNP Q9BXD5
D	-7	GLY	-	expression tag	UNP Q9BXD5

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-6	GLU	-	expression tag	UNP Q9BXD5
D	-5	ASN	-	expression tag	UNP Q9BXD5
D	-4	LEU	-	expression tag	UNP Q9BXD5
D	-3	TYR	-	expression tag	UNP Q9BXD5
D	-2	PHE	-	expression tag	UNP Q9BXD5
D	-1	GLN	-	expression tag	UNP Q9BXD5
D	0	GLY	-	expression tag	UNP Q9BXD5
D	1	GLY	-	expression tag	UNP Q9BXD5
D	2	SER	-	expression tag	UNP Q9BXD5

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	319	Total	O	0	0
			319	319		
3	B	318	Total	O	0	0
			318	318		
3	C	347	Total	O	0	0
			347	347		
3	D	287	Total	O	0	0
			287	287		

- Molecule 1: N-acetylneuraminate lyase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	96.99Å 106.90Å 127.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.27 – 1.60 49.27 – 1.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.27-1.60) 100.0 (49.27-1.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 1.60Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.164 , 0.189 0.176 , 0.199	Depositor DCC
R_{free} test set	8739 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	21.4	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 35.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10729	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.49% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.43	0/2419	0.64	1/3272 (0.0%)
1	B	0.43	0/2408	0.66	1/3260 (0.0%)
1	C	0.45	0/2395	0.67	1/3241 (0.0%)
1	D	0.42	0/2400	0.66	1/3248 (0.0%)
All	All	0.43	0/9622	0.66	4/13021 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	178	ASP	CB-CG-OD1	6.73	124.36	118.30
1	C	178	ASP	CB-CG-OD1	6.14	123.83	118.30
1	A	178	ASP	CB-CG-OD1	5.45	123.20	118.30
1	B	178	ASP	CB-CG-OD1	5.02	122.82	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2368	0	2390	6	0
1	B	2359	0	2369	7	0
1	C	2347	0	2368	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2352	0	2370	8	0
2	A	8	0	6	0	0
2	B	4	0	3	0	0
2	C	12	0	9	0	0
2	D	8	0	6	1	0
3	A	319	0	0	3	0
3	B	318	0	0	0	0
3	C	347	0	0	0	0
3	D	287	0	0	0	0
All	All	10729	0	9521	24	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 1.

All (24) 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:B:66:ARG:NH2	1:B:97:GLU:OE2	2.02	0.90
1:A:22:THR:OG1	1:A:24:ASN:OD1	2.01	0.79
1:B:127:ILE:HD11	1:B:160:LEU:HG	1.67	0.77
3:A:615:HOH:O	1:D:284:ARG:HD3	1.92	0.69
1:A:158:GLU:OE1	3:A:501:HOH:O	2.11	0.69
1:D:127:ILE:HD12	1:D:164:ILE:CD1	2.29	0.62
1:B:22:THR:OG1	1:B:24:ASN:OD1	2.16	0.59
1:C:22:THR:OG1	1:C:24:ASN:OD1	2.16	0.58
1:B:127:ILE:HD12	1:B:164:ILE:CD1	2.41	0.50
1:C:127:ILE:HD11	1:C:160:LEU:HG	1.93	0.50
1:D:110:ILE:HG23	1:D:140:LEU:HD23	1.95	0.49
1:D:14:VAL:HB	1:D:218:VAL:HG22	1.95	0.48
1:D:54:THR:OG1	2:D:401:ACT:H1	2.14	0.48
1:A:228[B]:LYS:HG3	1:A:251:ILE:HD11	1.99	0.45
1:A:176:PHE:CE2	1:A:178:ASP:HB3	2.52	0.45
1:D:127:ILE:HD12	1:D:164:ILE:HD13	1.99	0.44
1:A:160:LEU:HD23	1:A:164:ILE:HD12	2.02	0.42
1:C:160:LEU:HD23	1:C:164:ILE:HD12	2.01	0.42
1:B:14:VAL:HB	1:B:218:VAL:HG22	2.02	0.42
1:A:253[B]:ARG:NH2	3:A:503:HOH:O	2.49	0.41
1:D:204:GLU:HB3	1:D:255:ILE:HG21	2.02	0.41
1:B:49:PHE:HE1	1:B:111:ALA:HB3	1.86	0.40
1:D:49:PHE:HE1	1:D:111:ALA:HB3	1.85	0.40
1:B:127:ILE:HD12	1:B:164:ILE:HD11	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	305/338 (90%)	300 (98%)	5 (2%)	0	100	100
1	B	305/338 (90%)	298 (98%)	7 (2%)	0	100	100
1	C	302/338 (89%)	295 (98%)	7 (2%)	0	100	100
1	D	303/338 (90%)	296 (98%)	7 (2%)	0	100	100
All	All	1215/1352 (90%)	1189 (98%)	26 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	254/279 (91%)	251 (99%)	3 (1%)	71	54
1	B	252/279 (90%)	250 (99%)	2 (1%)	81	70
1	C	252/279 (90%)	249 (99%)	3 (1%)	71	54
1	D	252/279 (90%)	250 (99%)	2 (1%)	81	70
All	All	1010/1116 (90%)	1000 (99%)	10 (1%)	76	61

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	TYR
1	A	145	TYR
1	A	203	ASP
1	B	35	TYR
1	B	203	ASP
1	C	35	TYR
1	C	145	TYR
1	C	203	ASP
1	D	35	TYR
1	D	203	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	GLN
1	B	67	GLN
1	D	67	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](#)

8 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	ACT	A	402	-	1,3,3	0.85	0	0,3,3	0.00	-
2	ACT	A	401	-	1,3,3	1.31	0	0,3,3	0.00	-
2	ACT	B	401	-	1,3,3	1.47	0	0,3,3	0.00	-
2	ACT	C	401	-	1,3,3	1.41	0	0,3,3	0.00	-
2	ACT	D	401	-	1,3,3	0.93	0	0,3,3	0.00	-
2	ACT	C	402	-	1,3,3	1.07	0	0,3,3	0.00	-
2	ACT	D	402	-	1,3,3	0.95	0	0,3,3	0.00	-
2	ACT	C	403	-	1,3,3	0.20	0	0,3,3	0.00	-

There are no bond length outliers.

There are no bond angle outliers.

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
2	D	401	ACT	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	300/338 (88%)	-0.42	2 (0%) 87 87	16, 21, 35, 47	0
1	B	303/338 (89%)	-0.35	2 (0%) 87 87	16, 21, 32, 50	0
1	C	300/338 (88%)	-0.34	2 (0%) 87 87	15, 19, 32, 55	0
1	D	301/338 (89%)	-0.40	5 (1%) 70 69	16, 23, 36, 62	0
All	All	1204/1352 (89%)	-0.38	11 (0%) 84 84	15, 21, 34, 62	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	117	PHE	4.6
1	D	308	LEU	3.9
1	B	309	SER	3.6
1	D	195	GLN	3.6
1	B	310	PHE	3.0
1	D	193	GLN	3.0
1	A	121	TRP	3.0
1	D	307	PHE	2.8
1	D	8	LYS	2.7
1	C	193	GLN	2.6
1	C	195	GLN	2.2

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	ACT	B	401	4/4	0.77	0.17	39,39,40,44	0
2	ACT	D	401	4/4	0.87	0.18	44,45,46,49	0
2	ACT	C	401	4/4	0.88	0.11	25,26,27,27	0
2	ACT	C	403	4/4	0.89	0.12	36,38,38,38	0
2	ACT	A	401	4/4	0.92	0.08	28,29,31,32	0
2	ACT	D	402	4/4	0.95	0.13	42,42,44,44	0
2	ACT	C	402	4/4	0.95	0.11	26,26,27,27	0
2	ACT	A	402	4/4	0.96	0.07	29,29,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.