



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:03 PM GMT

PDB ID : 4H0Y
Title : Crystal structure of NAD⁺-Ia(E380S)-actin complex
Authors : Tsurumura, T.; Oda, M.; Nagahama, M.; Tsuge, H.
Deposited on : 2012-09-10
Resolution : 1.94 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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) 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

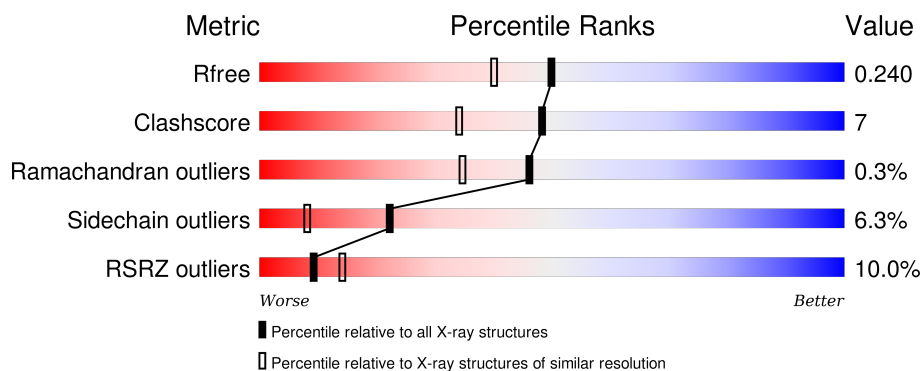
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.94 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	2910 (1.96-1.92)
Clashscore	102246	3095 (1.96-1.92)
Ramachandran outliers	100387	3062 (1.96-1.92)
Sidechain outliers	100360	3062 (1.96-1.92)
RSRZ outliers	91569	2915 (1.96-1.92)

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 $\leq 5\%$. 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	418	
2	B	375	

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
5	EDO	A	507	-	-	-	X
5	EDO	A	508	-	-	-	X
5	EDO	A	511	-	-	-	X
5	EDO	A	513	-	-	-	X
5	EDO	A	514	-	-	-	X
5	EDO	A	517	-	-	-	X
5	EDO	A	518	-	-	-	X
5	EDO	A	519	-	-	-	X
5	EDO	A	521	-	-	-	X
5	EDO	A	522	-	-	-	X
5	EDO	A	523	-	-	-	X
5	EDO	A	527	-	-	-	X
5	EDO	B	404	-	-	-	X
5	EDO	B	406	-	-	-	X
5	EDO	B	409	-	-	-	X
5	EDO	B	410	-	-	-	X
5	EDO	B	411	-	-	-	X
5	EDO	B	412	-	-	-	X
5	EDO	B	413	-	-	-	X
5	EDO	B	416	-	-	-	X
5	EDO	B	418	-	-	-	X
5	EDO	B	422	-	-	-	X
5	EDO	B	423	-	-	-	X
5	EDO	B	424	-	-	-	X
5	EDO	B	425	-	-	-	X
5	EDO	B	426	-	-	-	X
5	EDO	B	428	-	-	-	X
5	EDO	B	429	-	-	-	X

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 6739 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 Iota toxin component Ia.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	413	Total	C	N	O	S	0	0	0
			3361	2144	553	661	3			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	ARG	-	EXPRESSION TAG	UNP Q46220
A	-3	GLY	-	EXPRESSION TAG	UNP Q46220
A	-2	SER	-	EXPRESSION TAG	UNP Q46220
A	-1	HIS	-	EXPRESSION TAG	UNP Q46220
A	0	MET	-	EXPRESSION TAG	UNP Q46220
A	380	SER	GLU	ENGINEERED MUTATION	UNP Q46220

- Molecule 2 is a protein called Actin, alpha skeletal muscle.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	358	Total	C	N	O	S	0	0	0
			2799	1774	469	537	19			

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	
			44	21	7	14	2	

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	O P		
			5	4 1	0	0

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

Continued on next page...

Continued from previous page...

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

Continued on next page...

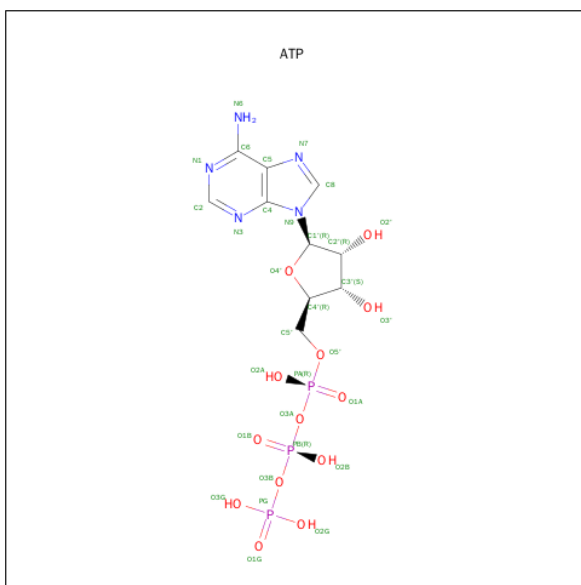
Continued from previous page...

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

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

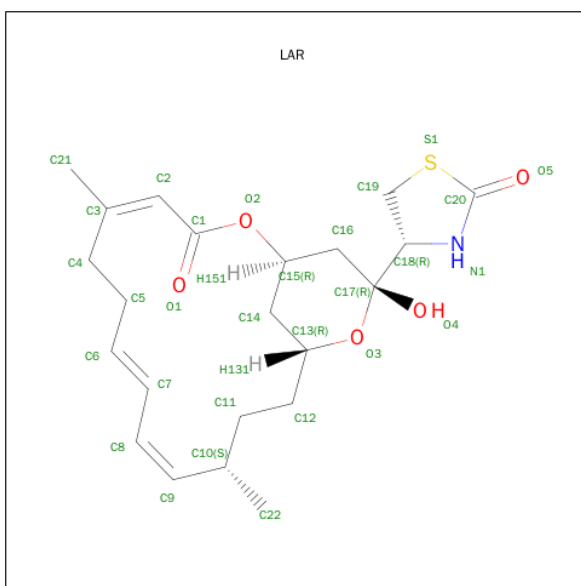
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Ca	0	0
			1	1		

- Molecule 7 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 8 is LATRUNCULIN A (three-letter code: LAR) (formula: $C_{22}H_{31}NO_5S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	B	1	Total	C	N	O	S	0	0
			29	22	1	5	1		

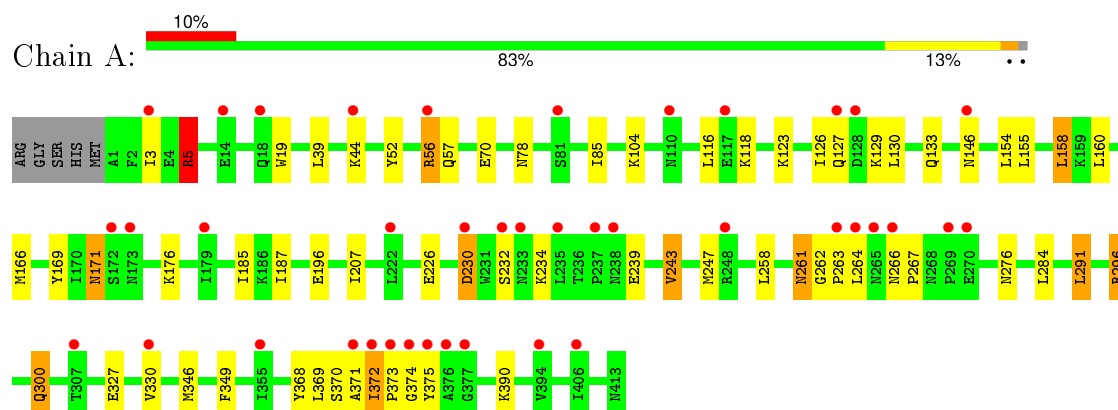
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	149	Total 149	O 149	0	0
9	B	112	Total 112	O 112	0	0

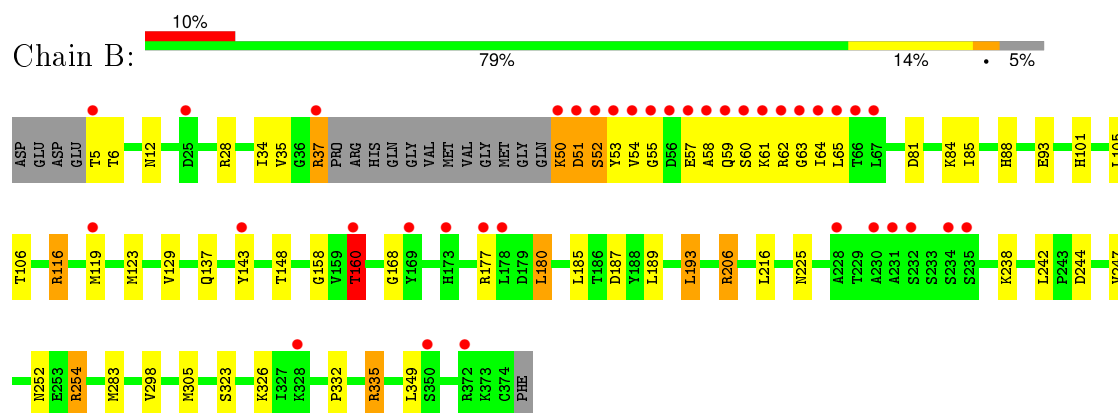
3 Residue-property plots [i](#)

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.

- Molecule 1: Iota toxin component Ia



- Molecule 2: Actin, alpha skeletal muscle



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	53.87Å 134.94Å 154.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.91 – 1.94 30.91 – 1.94	Depositor EDS
% Data completeness (in resolution range)	99.6 (30.91-1.94) 99.6 (30.91-1.94)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.93 (at 1.94Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.220 , 0.233 0.228 , 0.240	Depositor DCC
R_{free} test set	4182 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	25.9	Xtriage
Anisotropy	0.127	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 51.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 83784 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6739	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% 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, 0.375 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: NAD, PO4, EDO, ATP, CA, LAR

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	0.35	0/3430	0.81	8/4633 (0.2%)
2	B	0.42	1/2858 (0.0%)	0.65	3/3873 (0.1%)
All	All	0.38	1/6288 (0.0%)	0.74	11/8506 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	160	THR	CB-OG1	12.83	1.69	1.43

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	5	ARG	NE-CZ-NH2	-16.57	112.01	120.30
1	A	296	ARG	NE-CZ-NH2	-15.27	112.67	120.30
1	A	296	ARG	NE-CZ-NH1	14.11	127.36	120.30
1	A	5	ARG	NE-CZ-NH1	13.01	126.80	120.30
2	B	254	ARG	NE-CZ-NH1	-8.66	115.97	120.30
1	A	296	ARG	CG-CD-NE	-7.82	95.38	111.80
2	B	254	ARG	NE-CZ-NH2	7.50	124.05	120.30
1	A	296	ARG	CD-NE-CZ	7.06	133.49	123.60
1	A	291	LEU	CA-CB-CG	5.73	128.49	115.30
2	B	206	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	A	5	ARG	CD-NE-CZ	5.38	131.13	123.60

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	3361	0	3343	40	0
2	B	2799	0	2773	54	1
3	A	44	0	26	0	0
4	A	5	0	0	1	0
5	A	104	0	155	4	0
5	B	104	0	156	10	0
6	B	1	0	0	0	0
7	B	31	0	12	0	0
8	B	29	0	31	1	0
9	A	149	0	0	1	0
9	B	112	0	0	1	0
All	All	6739	0	6496	95	1

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 7.

All (95) 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)
2:B:160:THR:CB	2:B:160:THR:OG1	1.68	1.41
1:A:104:LYS:NZ	5:A:519:EDO:O1	1.67	1.24
2:B:51:ASP:O	2:B:52:SER:HB3	1.51	1.08
2:B:116:ARG:NH1	9:B:586:HOH:O	1.98	0.96
2:B:216:LEU:O	2:B:254:ARG:HD2	1.69	0.92
2:B:160:THR:HG23	2:B:180:LEU:O	1.72	0.89
2:B:332:PRO:O	2:B:335:ARG:HG2	1.74	0.88
2:B:51:ASP:O	2:B:52:SER:CB	2.29	0.81
2:B:283:MET:CE	5:B:409:EDO:O2	2.30	0.80
1:A:146:ASN:HA	1:A:196:GLU:HG3	1.65	0.78
2:B:61:LYS:HD3	2:B:64:ILE:HG22	1.67	0.77
2:B:283:MET:HE1	5:B:409:EDO:O2	1.85	0.76
2:B:298:VAL:CG1	5:B:408:EDO:H21	2.17	0.75
1:A:56:ARG:NH2	1:A:57:GLN:HB3	2.10	0.67
1:A:372:ILE:HG22	1:A:372:ILE:O	1.95	0.67
1:A:129:LYS:HE2	1:A:390:LYS:HE2	1.76	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:ILE:HD13	1:A:166:MET:HE3	1.79	0.64
1:A:261:ASN:HD21	1:A:264:LEU:HD12	1.64	0.62
2:B:216:LEU:HD22	2:B:238:LYS:HG2	1.80	0.62
2:B:216:LEU:O	2:B:254:ARG:CD	2.46	0.62
1:A:5:ARG:HD3	4:A:502:PO4:O3	1.99	0.62
1:A:230:ASP:N	1:A:230:ASP:OD2	2.19	0.61
1:A:247:MET:HE1	1:A:372:ILE:HG21	1.82	0.61
1:A:56:ARG:HH21	1:A:57:GLN:HB3	1.63	0.61
1:A:5:ARG:HD2	1:A:19:TRP:CZ2	2.37	0.60
2:B:54:VAL:HG13	2:B:85:ILE:HD13	1.83	0.60
1:A:158:LEU:HD21	1:A:185:ILE:HD12	1.84	0.59
2:B:283:MET:HE3	5:B:409:EDO:O2	2.03	0.59
1:A:239:GLU:O	1:A:243:VAL:HG13	2.02	0.58
1:A:371:ALA:O	1:A:373:PRO:HD3	2.02	0.58
2:B:52:SER:HB2	2:B:84:LYS:HG3	1.85	0.57
2:B:143:TYR:OH	2:B:349:LEU:HD11	2.04	0.57
2:B:53:TYR:CE2	2:B:60:SER:CB	2.88	0.57
2:B:298:VAL:HG12	5:B:408:EDO:H21	1.87	0.56
2:B:6:THR:O	2:B:101:HIS:HD2	1.88	0.56
2:B:148:THR:H	5:B:413:EDO:H12	1.71	0.56
1:A:78:ASN:ND2	9:A:609:HOH:O	2.38	0.56
1:A:146:ASN:HA	1:A:196:GLU:CG	2.37	0.55
2:B:323:SER:HA	5:B:410:EDO:H22	1.89	0.55
2:B:160:THR:CG2	2:B:180:LEU:O	2.52	0.54
2:B:50:LYS:O	2:B:51:ASP:C	2.45	0.54
1:A:346:MET:HG3	1:A:349:PHE:CE1	2.44	0.53
2:B:148:THR:OG1	2:B:168:GLY:N	2.41	0.53
2:B:51:ASP:O	2:B:84:LYS:HE3	2.10	0.52
2:B:37:ARG:NH1	2:B:81:ASP:OD1	2.43	0.51
2:B:158:GLY:HA2	5:B:404:EDO:H11	1.92	0.51
1:A:123:LYS:O	1:A:127:GLN:HB2	2.10	0.51
2:B:53:TYR:CD2	2:B:60:SER:HB2	2.45	0.51
2:B:53:TYR:HB3	2:B:58:ALA:HA	1.92	0.51
1:A:126:ILE:HD12	1:A:187:ILE:HD13	1.93	0.51
1:A:263:PRO:O	1:A:267:PRO:HG3	2.12	0.50
1:A:296:ARG:HD3	1:A:349:PHE:CE1	2.46	0.50
2:B:50:LYS:O	2:B:51:ASP:O	2.29	0.49
2:B:35:VAL:HA	2:B:53:TYR:O	2.12	0.49
2:B:106:THR:HB	2:B:137:GLN:HG3	1.95	0.49
1:A:133:GLN:HG3	1:A:185:ILE:HG12	1.95	0.48
2:B:53:TYR:CE2	2:B:60:SER:HB2	2.48	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:346:MET:HG3	1:A:349:PHE:CZ	2.49	0.48
2:B:53:TYR:CD2	2:B:60:SER:CB	2.96	0.48
2:B:52:SER:O	2:B:53:TYR:CD1	2.67	0.48
1:A:171:ASN:HD22	1:A:171:ASN:C	2.17	0.47
1:A:160:LEU:HD13	1:A:166:MET:HE2	1.97	0.47
1:A:52:TYR:CZ	1:A:56:ARG:HB3	2.49	0.47
1:A:327:GLU:OE1	5:A:507:EDO:H21	2.15	0.47
2:B:187:ASP:OD1	2:B:206:ARG:HD2	2.15	0.46
1:A:5:ARG:HD2	1:A:19:TRP:CH2	2.50	0.46
1:A:3:ILE:O	1:A:3:ILE:HG22	2.14	0.46
1:A:374:GLY:O	1:A:375:TYR:CD1	2.69	0.46
2:B:50:LYS:C	2:B:51:ASP:O	2.53	0.46
2:B:63:GLY:N	2:B:65:LEU:O	2.50	0.45
1:A:185:ILE:CD1	1:A:207:ILE:HG22	2.47	0.45
2:B:61:LYS:HD3	2:B:64:ILE:CG2	2.43	0.44
2:B:298:VAL:HG11	5:B:408:EDO:H21	1.98	0.44
1:A:368:TYR:CZ	1:A:370:SER:HB2	2.53	0.44
2:B:53:TYR:HD2	2:B:57:GLU:O	2.00	0.43
1:A:5:ARG:CD	1:A:19:TRP:CZ2	3.00	0.43
1:A:300:GLN:NE2	1:A:300:GLN:H	2.16	0.43
1:A:104:LYS:NZ	5:A:519:EDO:C1	2.75	0.43
2:B:305:MET:SD	2:B:335:ARG:HG3	2.58	0.43
2:B:34:ILE:HD11	2:B:55:GLY:O	2.19	0.42
2:B:105:LEU:CD1	2:B:119:MET:HG2	2.48	0.42
2:B:206:ARG:HG2	8:B:403:LAR:S1	2.59	0.42
2:B:54:VAL:HG13	2:B:85:ILE:CD1	2.46	0.42
1:A:169:TYR:CE2	1:A:176:LYS:HB2	2.55	0.42
1:A:262:GLY:N	1:A:263:PRO:CD	2.82	0.42
5:B:411:EDO:H21	5:B:412:EDO:O2	2.19	0.42
1:A:70:GLU:OE2	5:A:525:EDO:O1	2.37	0.41
2:B:160:THR:CB	2:B:160:THR:HG1	2.12	0.41
2:B:88:HIS:HE1	2:B:93:GLU:OE2	2.03	0.41
2:B:119:MET:HB2	2:B:119:MET:HE2	1.93	0.41
2:B:189:LEU:HG	2:B:193:LEU:HD22	2.03	0.41
1:A:146:ASN:HD22	1:A:196:GLU:HG3	1.84	0.41
2:B:54:VAL:HA	2:B:58:ALA:HB2	2.03	0.41
2:B:123:MET:CE	2:B:129:VAL:HG11	2.51	0.40
2:B:62:ARG:C	2:B:65:LEU:O	2.60	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:244:ASP:OD1	2:B:326:LYS:NZ[1_655]	2.10	0.10

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	411/418 (98%)	396 (96%)	15 (4%)	0	100	100
2	B	354/375 (94%)	347 (98%)	5 (1%)	2 (1%)	30	16
All	All	765/793 (96%)	743 (97%)	20 (3%)	2 (0%)	46	35

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	51	ASP
2	B	52	SER

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	376/380 (99%)	350 (93%)	26 (7%)	19	6
2	B	304/318 (96%)	287 (94%)	17 (6%)	26	11
All	All	680/698 (97%)	637 (94%)	43 (6%)	22	8

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ARG
1	A	39	LEU
1	A	44	LYS
1	A	56	ARG
1	A	116	LEU
1	A	118	LYS
1	A	130	LEU
1	A	154	LEU
1	A	155	LEU
1	A	158	LEU
1	A	171	ASN
1	A	226	GLU
1	A	230	ASP
1	A	232	SER
1	A	234	LYS
1	A	243	VAL
1	A	258	LEU
1	A	261	ASN
1	A	266	ASN
1	A	276	ASN
1	A	284	LEU
1	A	291	LEU
1	A	300	GLN
1	A	330	VAL
1	A	369	LEU
1	A	372	ILE
2	B	5	THR
2	B	12	ASN
2	B	28	ARG
2	B	37	ARG
2	B	50	LYS
2	B	59	GLN
2	B	116	ARG
2	B	160	THR
2	B	177	ARG
2	B	180	LEU
2	B	185	LEU
2	B	193	LEU
2	B	225	ASN
2	B	242	LEU
2	B	247	VAL
2	B	252	ASN
2	B	335	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	ASN
1	A	51	ASN
1	A	54	GLN
1	A	78	ASN
1	A	90	ASN
1	A	146	ASN
1	A	157	HIS
1	A	171	ASN
1	A	181	GLN
1	A	209	ASN
1	A	244	ASN
1	A	256	ASN
1	A	261	ASN
1	A	266	ASN
1	A	300	GLN
1	A	392	ASN
2	B	73	HIS
2	B	88	HIS
2	B	101	HIS
2	B	128	ASN
2	B	162	ASN
2	B	252	ASN
2	B	263	GLN

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 ⓘ

Of 57 ligands modelled in this entry, 1 is monoatomic - leaving 56 for Mogul analysis.

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	NAD	A	501	-	38,48,48	0.93	3 (7%)	47,73,73	1.88	9 (19%)
4	PO4	A	502	-	4,4,4	0.35	0	6,6,6	0.26	0
5	EDO	A	503	-	3,3,3	0.50	0	2,2,2	0.40	0
5	EDO	A	504	-	3,3,3	0.48	0	2,2,2	0.36	0
5	EDO	A	505	-	3,3,3	0.49	0	2,2,2	0.30	0
5	EDO	A	506	-	3,3,3	0.45	0	2,2,2	0.44	0
5	EDO	A	507	-	3,3,3	0.47	0	2,2,2	0.31	0
5	EDO	A	508	-	3,3,3	0.49	0	2,2,2	0.27	0
5	EDO	A	509	-	3,3,3	0.51	0	2,2,2	0.28	0
5	EDO	A	510	-	3,3,3	0.47	0	2,2,2	0.34	0
5	EDO	A	511	-	3,3,3	0.47	0	2,2,2	0.35	0
5	EDO	A	512	-	3,3,3	0.47	0	2,2,2	0.24	0
5	EDO	A	513	-	3,3,3	0.47	0	2,2,2	0.38	0
5	EDO	A	514	-	3,3,3	0.49	0	2,2,2	0.29	0
5	EDO	A	515	-	3,3,3	0.50	0	2,2,2	0.23	0
5	EDO	A	516	-	3,3,3	0.47	0	2,2,2	0.22	0
5	EDO	A	517	-	3,3,3	0.48	0	2,2,2	0.35	0
5	EDO	A	518	-	3,3,3	0.48	0	2,2,2	0.31	0
5	EDO	A	519	-	3,3,3	0.40	0	2,2,2	0.69	0
5	EDO	A	520	-	3,3,3	0.45	0	2,2,2	0.31	0
5	EDO	A	521	-	3,3,3	0.49	0	2,2,2	0.23	0
5	EDO	A	522	-	3,3,3	0.49	0	2,2,2	0.32	0
5	EDO	A	523	-	3,3,3	0.48	0	2,2,2	0.32	0
5	EDO	A	524	-	3,3,3	0.45	0	2,2,2	0.36	0
5	EDO	A	525	-	3,3,3	0.69	0	2,2,2	0.26	0
5	EDO	A	526	-	3,3,3	0.46	0	2,2,2	0.31	0
5	EDO	A	527	-	3,3,3	0.50	0	2,2,2	0.20	0
5	EDO	A	528	-	3,3,3	0.49	0	2,2,2	0.33	0
7	ATP	B	402	-	24,33,33	0.94	1 (4%)	31,52,52	1.77	5 (16%)
8	LAR	B	403	-	29,31,31	1.40	3 (10%)	29,43,43	2.31	9 (31%)
5	EDO	B	404	-	3,3,3	0.52	0	2,2,2	0.12	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EDO	B	405	-	3,3,3	0.48	0	2,2,2	0.37	0
5	EDO	B	406	-	3,3,3	0.50	0	2,2,2	0.22	0
5	EDO	B	407	-	3,3,3	0.46	0	2,2,2	0.40	0
5	EDO	B	408	-	3,3,3	0.47	0	2,2,2	0.34	0
5	EDO	B	409	-	3,3,3	0.43	0	2,2,2	0.55	0
5	EDO	B	410	-	3,3,3	0.39	0	2,2,2	0.41	0
5	EDO	B	411	-	3,3,3	0.49	0	2,2,2	0.20	0
5	EDO	B	412	-	3,3,3	0.44	0	2,2,2	0.47	0
5	EDO	B	413	-	3,3,3	0.43	0	2,2,2	0.30	0
5	EDO	B	414	-	3,3,3	0.47	0	2,2,2	0.31	0
5	EDO	B	415	-	3,3,3	0.47	0	2,2,2	0.33	0
5	EDO	B	416	-	3,3,3	0.52	0	2,2,2	0.34	0
5	EDO	B	417	-	3,3,3	0.46	0	2,2,2	0.42	0
5	EDO	B	418	-	3,3,3	0.47	0	2,2,2	0.32	0
5	EDO	B	419	-	3,3,3	0.49	0	2,2,2	0.29	0
5	EDO	B	420	-	3,3,3	0.50	0	2,2,2	0.22	0
5	EDO	B	421	-	3,3,3	0.50	0	2,2,2	0.31	0
5	EDO	B	422	-	3,3,3	0.48	0	2,2,2	0.32	0
5	EDO	B	423	-	3,3,3	0.49	0	2,2,2	0.29	0
5	EDO	B	424	-	3,3,3	0.45	0	2,2,2	0.30	0
5	EDO	B	425	-	3,3,3	0.44	0	2,2,2	0.45	0
5	EDO	B	426	-	3,3,3	0.49	0	2,2,2	0.27	0
5	EDO	B	427	-	3,3,3	0.52	0	2,2,2	0.35	0
5	EDO	B	428	-	3,3,3	0.46	0	2,2,2	0.31	0
5	EDO	B	429	-	3,3,3	0.51	0	2,2,2	0.27	0

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	NAD	A	501	-	-	0/22/62/62	0/5/5/5
4	PO4	A	502	-	-	0/0/0/0	0/0/0/0
5	EDO	A	503	-	-	0/1/1/1	0/0/0/0
5	EDO	A	504	-	-	0/1/1/1	0/0/0/0
5	EDO	A	505	-	-	0/1/1/1	0/0/0/0
5	EDO	A	506	-	-	0/1/1/1	0/0/0/0
5	EDO	A	507	-	-	0/1/1/1	0/0/0/0
5	EDO	A	508	-	-	0/1/1/1	0/0/0/0
5	EDO	A	509	-	-	0/1/1/1	0/0/0/0
5	EDO	A	510	-	-	0/1/1/1	0/0/0/0
5	EDO	A	511	-	-	0/1/1/1	0/0/0/0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	A	512	-	-	0/1/1/1	0/0/0/0
5	EDO	A	513	-	-	0/1/1/1	0/0/0/0
5	EDO	A	514	-	-	0/1/1/1	0/0/0/0
5	EDO	A	515	-	-	0/1/1/1	0/0/0/0
5	EDO	A	516	-	-	0/1/1/1	0/0/0/0
5	EDO	A	517	-	-	0/1/1/1	0/0/0/0
5	EDO	A	518	-	-	0/1/1/1	0/0/0/0
5	EDO	A	519	-	-	0/1/1/1	0/0/0/0
5	EDO	A	520	-	-	0/1/1/1	0/0/0/0
5	EDO	A	521	-	-	0/1/1/1	0/0/0/0
5	EDO	A	522	-	-	0/1/1/1	0/0/0/0
5	EDO	A	523	-	-	0/1/1/1	0/0/0/0
5	EDO	A	524	-	-	0/1/1/1	0/0/0/0
5	EDO	A	525	-	-	0/1/1/1	0/0/0/0
5	EDO	A	526	-	-	0/1/1/1	0/0/0/0
5	EDO	A	527	-	-	0/1/1/1	0/0/0/0
5	EDO	A	528	-	-	0/1/1/1	0/0/0/0
7	ATP	B	402	-	-	0/18/38/38	0/3/3/3
8	LAR	B	403	-	-	0/23/51/51	0/1/3/3
5	EDO	B	404	-	-	0/1/1/1	0/0/0/0
5	EDO	B	405	-	-	0/1/1/1	0/0/0/0
5	EDO	B	406	-	-	0/1/1/1	0/0/0/0
5	EDO	B	407	-	-	0/1/1/1	0/0/0/0
5	EDO	B	408	-	-	0/1/1/1	0/0/0/0
5	EDO	B	409	-	-	0/1/1/1	0/0/0/0
5	EDO	B	410	-	-	0/1/1/1	0/0/0/0
5	EDO	B	411	-	-	0/1/1/1	0/0/0/0
5	EDO	B	412	-	-	0/1/1/1	0/0/0/0
5	EDO	B	413	-	-	0/1/1/1	0/0/0/0
5	EDO	B	414	-	-	0/1/1/1	0/0/0/0
5	EDO	B	415	-	-	0/1/1/1	0/0/0/0
5	EDO	B	416	-	-	0/1/1/1	0/0/0/0
5	EDO	B	417	-	-	0/1/1/1	0/0/0/0
5	EDO	B	418	-	-	0/1/1/1	0/0/0/0
5	EDO	B	419	-	-	0/1/1/1	0/0/0/0
5	EDO	B	420	-	-	0/1/1/1	0/0/0/0
5	EDO	B	421	-	-	0/1/1/1	0/0/0/0
5	EDO	B	422	-	-	0/1/1/1	0/0/0/0
5	EDO	B	423	-	-	0/1/1/1	0/0/0/0
5	EDO	B	424	-	-	0/1/1/1	0/0/0/0
5	EDO	B	425	-	-	0/1/1/1	0/0/0/0
5	EDO	B	426	-	-	0/1/1/1	0/0/0/0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	B	427	-	-	0/1/1/1	0/0/0/0
5	EDO	B	428	-	-	0/1/1/1	0/0/0/0
5	EDO	B	429	-	-	0/1/1/1	0/0/0/0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	403	LAR	C20-N1	-2.83	1.32	1.36
8	B	403	LAR	C19-S1	-2.15	1.77	1.81
3	A	501	NAD	O4B-C1B	2.25	1.44	1.41
3	A	501	NAD	O4D-C1D	2.26	1.44	1.41
3	A	501	NAD	C5A-C4A	2.89	1.47	1.40
7	B	402	ATP	C5-C4	2.98	1.47	1.40
8	B	403	LAR	O2-C1	5.50	1.46	1.34

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	NAD	N3A-C2A-N1A	-8.24	122.58	128.89
8	B	403	LAR	C4-C5-C6	-7.32	94.60	112.79
7	B	402	ATP	N3-C2-N1	-7.28	123.32	128.89
8	B	403	LAR	C8-C7-C6	-5.22	100.25	124.44
8	B	403	LAR	O1-C1-C2	-2.88	118.63	126.20
8	B	403	LAR	O2-C1-O1	-2.71	119.07	123.30
7	B	402	ATP	PA-O3A-PB	-2.67	125.23	132.73
3	A	501	NAD	C1B-N9A-C4A	-2.66	122.94	126.94
3	A	501	NAD	C5D-C4D-C3D	-2.63	104.76	115.21
3	A	501	NAD	C4A-C5A-N7A	-2.58	107.10	109.48
7	B	402	ATP	C1'-N9-C4	-2.34	123.41	126.94
3	A	501	NAD	PN-O3-PA	-2.32	126.20	132.73
7	B	402	ATP	C4-C5-N7	-2.17	107.48	109.48
3	A	501	NAD	O7N-C7N-N7N	-2.01	119.77	122.59
3	A	501	NAD	C2A-N1A-C6A	2.02	122.37	118.77
8	B	403	LAR	O2-C15-C16	2.07	112.71	107.61
8	B	403	LAR	C19-C18-N1	2.12	106.06	100.73
8	B	403	LAR	C5-C6-C7	2.19	138.34	126.01
7	B	402	ATP	C2-N1-C6	2.22	122.73	118.77
8	B	403	LAR	C18-N1-C20	2.74	116.63	113.04
3	A	501	NAD	C3N-C7N-N7N	3.41	121.55	117.82
8	B	403	LAR	O2-C1-C2	4.38	122.30	111.51
3	A	501	NAD	O4D-C1D-N1N	5.01	113.64	108.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

12 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	502	PO4	1	0
5	A	507	EDO	1	0
5	A	519	EDO	2	0
5	A	525	EDO	1	0
8	B	403	LAR	1	0
5	B	404	EDO	1	0
5	B	408	EDO	3	0
5	B	409	EDO	3	0
5	B	410	EDO	1	0
5	B	411	EDO	1	0
5	B	412	EDO	1	0
5	B	413	EDO	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 ⓘ

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	413/418 (98%)	0.78	40 (9%) 10 15	23, 33, 63, 98	0
2	B	358/375 (95%)	0.77	37 (10%) 9 13	15, 23, 59, 119	0
All	All	771/793 (97%)	0.77	77 (9%) 9 14	15, 29, 61, 119	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	63	GLY	13.1
2	B	60	SER	13.0
1	A	376	ALA	9.7
2	B	52	SER	9.6
2	B	57	GLU	9.5
2	B	51	ASP	9.2
2	B	56	ASP	8.2
2	B	62	ARG	8.0
2	B	53	TYR	8.0
2	B	59	GLN	7.2
2	B	64	ILE	7.2
1	A	375	TYR	6.7
1	A	372	ILE	6.3
2	B	54	VAL	5.9
2	B	61	LYS	5.8
2	B	50	LYS	5.1
2	B	58	ALA	5.0
1	A	373	PRO	4.7
2	B	65	LEU	4.7
2	B	66	THR	4.6
2	B	37	ARG	4.5
1	A	237	PRO	4.5
1	A	238	ASN	4.4
1	A	264	LEU	4.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	B	232	SER	4.1
2	B	55	GLY	3.8
1	A	44	LYS	3.7
1	A	371	ALA	3.7
2	B	231	ALA	3.7
1	A	173	ASN	3.6
1	A	117	GLU	3.6
1	A	222	LEU	3.4
1	A	377	GLY	3.3
1	A	110	ASN	3.2
1	A	128	ASP	3.2
1	A	270	GLU	3.1
1	A	233	ASN	3.1
1	A	266	ASN	3.1
1	A	330	VAL	3.0
1	A	179	ILE	2.9
1	A	374	GLY	2.9
1	A	307	THR	2.8
2	B	143	TYR	2.8
2	B	119	MET	2.7
2	B	67	LEU	2.7
1	A	235	LEU	2.6
1	A	56	ARG	2.6
1	A	230	ASP	2.6
2	B	169	TYR	2.6
1	A	3	ILE	2.5
1	A	127	GLN	2.5
1	A	265	ASN	2.5
2	B	372	ARG	2.5
2	B	160	THR	2.4
2	B	350	SER	2.4
2	B	328	LYS	2.4
1	A	146	ASN	2.3
1	A	355	ILE	2.3
1	A	269	PRO	2.3
2	B	228	ALA	2.3
2	B	235	SER	2.3
1	A	406	ILE	2.3
1	A	263	PRO	2.2
1	A	81	SER	2.2
1	A	232	SER	2.2
1	A	248	ARG	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	B	178	LEU	2.1
1	A	394	VAL	2.1
1	A	14	GLU	2.1
1	A	18	GLN	2.1
2	B	177	ARG	2.1
2	B	5	THR	2.1
1	A	172	SER	2.0
2	B	230	ALA	2.0
2	B	173	HIS	2.0
2	B	25	ASP	2.0
2	B	234	SER	2.0

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. 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, 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	LLDF	B-factors(Å ²)	Q<0.9
5	EDO	B	426	4/4	0.77	0.51	19.21	43,43,43,43	0
5	EDO	A	514	4/4	0.75	0.39	15.19	36,36,36,36	0
5	EDO	B	424	4/4	0.65	0.32	13.92	37,37,37,37	0
5	EDO	B	425	4/4	0.78	0.34	13.11	31,31,32,32	0
5	EDO	A	518	4/4	0.72	0.28	10.14	37,38,38,38	0
5	EDO	B	410	4/4	0.92	0.36	9.38	32,32,32,32	0
5	EDO	B	423	4/4	0.74	0.25	8.92	45,45,45,46	0
5	EDO	B	406	4/4	0.62	0.37	7.34	45,45,45,45	0
5	EDO	A	522	4/4	0.74	0.30	6.80	34,34,34,34	0
5	EDO	A	521	4/4	0.76	0.41	6.01	47,47,47,47	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
5	EDO	A	507	4/4	0.67	0.28	5.85	38,38,38,38	0
5	EDO	B	418	4/4	0.78	0.27	5.79	39,39,39,39	0
5	EDO	B	422	4/4	0.97	0.22	5.75	23,24,24,24	0
5	EDO	A	523	4/4	0.85	0.24	5.62	36,36,36,36	0
5	EDO	B	413	4/4	0.73	0.30	5.17	32,32,32,32	0
5	EDO	B	428	4/4	0.75	0.22	4.73	38,38,38,38	0
5	EDO	B	416	4/4	0.83	0.33	4.23	37,38,38,38	0
5	EDO	A	508	4/4	0.81	0.21	3.85	30,30,30,30	0
5	EDO	B	429	4/4	0.83	0.20	3.66	39,39,39,39	0
5	EDO	B	404	4/4	0.88	0.22	3.43	33,33,34,34	0
5	EDO	A	527	4/4	0.77	0.29	3.43	46,46,46,46	0
5	EDO	B	412	4/4	0.84	0.21	3.30	48,48,48,48	0
5	EDO	B	411	4/4	0.77	0.20	3.17	42,42,42,42	0
5	EDO	A	511	4/4	0.81	0.24	2.92	42,42,42,42	0
5	EDO	B	409	4/4	0.90	0.19	2.69	34,35,35,35	0
5	EDO	A	513	4/4	0.81	0.17	2.24	38,38,38,38	0
5	EDO	A	519	4/4	0.68	0.24	2.21	46,46,47,47	0
5	EDO	A	517	4/4	0.90	0.17	2.15	34,34,34,34	0
5	EDO	A	520	4/4	0.86	0.21	1.81	37,37,37,37	0
5	EDO	B	419	4/4	0.89	0.15	1.77	31,32,32,32	0
5	EDO	B	427	4/4	0.78	0.23	1.75	46,46,46,46	0
5	EDO	B	415	4/4	0.85	0.24	1.69	40,40,40,40	0
5	EDO	A	516	4/4	0.86	0.20	1.39	37,37,37,37	0
4	PO4	A	502	5/5	0.94	0.21	1.37	42,42,42,42	0
5	EDO	A	503	4/4	0.88	0.19	1.30	27,27,27,27	0
5	EDO	B	408	4/4	0.80	0.20	1.10	22,22,22,23	0
5	EDO	B	421	4/4	0.78	0.19	0.79	38,38,38,39	0
5	EDO	A	505	4/4	0.90	0.15	0.60	30,30,30,30	0
5	EDO	B	417	4/4	0.95	0.15	0.52	26,26,26,26	0
3	NAD	A	501	44/44	0.93	0.13	-0.16	29,34,38,38	0
5	EDO	B	414	4/4	0.94	0.14	-0.20	27,27,27,27	0
5	EDO	A	509	4/4	0.90	0.14	-0.30	27,27,27,27	0
5	EDO	A	504	4/4	0.89	0.15	-0.33	31,31,31,31	0
5	EDO	B	407	4/4	0.93	0.11	-0.38	27,27,27,27	0
8	LAR	B	403	29/29	0.93	0.12	-0.48	23,25,30,30	0
7	ATP	B	402	31/31	0.98	0.12	-0.64	17,17,17,17	0
5	EDO	A	510	4/4	0.94	0.12	-0.77	43,43,43,43	0
6	CA	B	401	1/1	1.00	0.09	-	17,17,17,17	0
5	EDO	A	506	4/4	0.73	0.29	-	43,43,43,44	0
5	EDO	A	515	4/4	0.84	0.31	-	48,48,48,48	0
5	EDO	A	525	4/4	0.70	0.32	-	40,41,41,41	0
5	EDO	A	528	4/4	0.70	0.28	-	35,35,35,36	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	EDO	A	526	4/4	0.72	0.39	-	43,43,44,44	0
5	EDO	A	512	4/4	0.80	0.29	-	40,40,40,40	0
5	EDO	A	524	4/4	0.85	0.45	-	41,41,41,41	0
5	EDO	B	420	4/4	0.76	0.35	-	40,40,40,40	0
5	EDO	B	405	4/4	0.87	0.26	-	43,43,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.