



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:55 AM GMT

PDB ID : 3NKP  
Title : Crystal structure of mouse autotaxin in complex with 18:1-LPA  
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Deposited on : 2010-06-20  
Resolution : 1.75 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](mailto: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.

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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

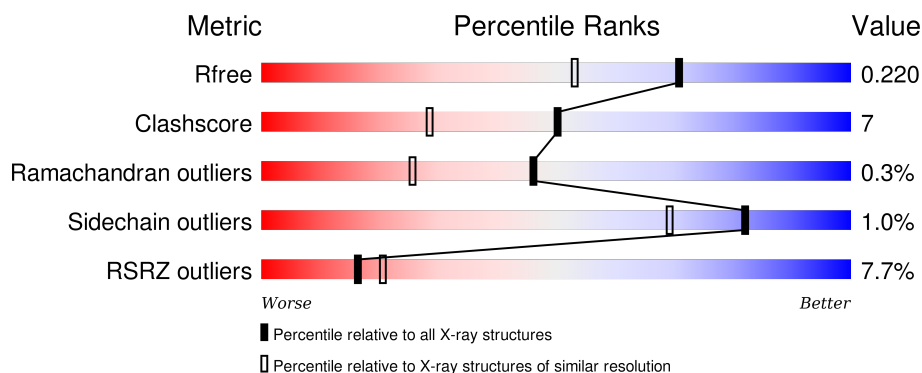
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*


The reported resolution of this entry is 1.75 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	1609 (1.76-1.76)
Clashscore	102246	1730 (1.76-1.76)
Ramachandran outliers	100387	1711 (1.76-1.76)
Sidechain outliers	100360	1711 (1.76-1.76)
RSRZ outliers	91569	1610 (1.76-1.76)

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  $\geq 3$ , 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	831	

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
10	EDO	A	1010	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
10	EDO	A	1012	-	-	X	-
10	EDO	A	1013	-	-	-	X
10	EDO	A	1016	-	-	-	X
10	EDO	A	1017	-	-	-	X
10	EDO	A	1018	-	-	-	X
10	EDO	A	1021	-	-	-	X
10	EDO	A	1024	-	-	X	X
10	EDO	A	1026	-	-	-	X
10	EDO	A	1027	-	-	-	X
10	EDO	A	1028	-	-	-	X
10	EDO	A	1030	-	-	-	X
3	MAN	A	904	-	-	-	X
8	NKP	A	1006	-	-	-	X
8	NKP	A	1007	-	-	-	X
9	SCN	A	1008	-	-	-	X

## 2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 7182 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 Ectonucleotide pyrophosphatase/phosphodiesterase family member 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	788	Total	C	N	O	S	0	4	0
			6312	4007	1078	1177	50			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	LYS	SEE REMARK 999	UNP Q9R1E6
A	?	-	VAL	SEE REMARK 999	UNP Q9R1E6
A	?	-	GLU	SEE REMARK 999	UNP Q9R1E6
A	?	-	PRO	SEE REMARK 999	UNP Q9R1E6
A	859	SER	-	EXPRESSION TAG	UNP Q9R1E6
A	860	ARG	-	EXPRESSION TAG	UNP Q9R1E6
A	861	GLU	-	EXPRESSION TAG	UNP Q9R1E6
A	862	ASN	-	EXPRESSION TAG	UNP Q9R1E6
A	863	LEU	-	EXPRESSION TAG	UNP Q9R1E6
A	864	TYR	-	EXPRESSION TAG	UNP Q9R1E6
A	865	PHE	-	EXPRESSION TAG	UNP Q9R1E6
A	866	GLN	-	EXPRESSION TAG	UNP Q9R1E6

- Molecule 2 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	2	Total	C	N	O	0	0
			28	16	2	10		
2	A	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 3 is a polymer of unknown type called SUGAR (6-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	6	Total	C	N	O	0	0
			72	40	2	30		

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Zn	0	0
			2	2		

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Ca	0	0
			1	1		

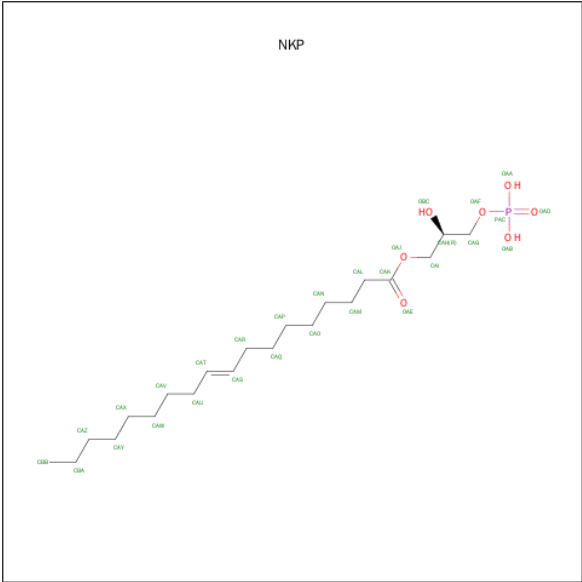
- Molecule 6 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Na	0	0
			1	1		

- Molecule 7 is POTASSIUM ION (three-letter code: K) (formula: K).

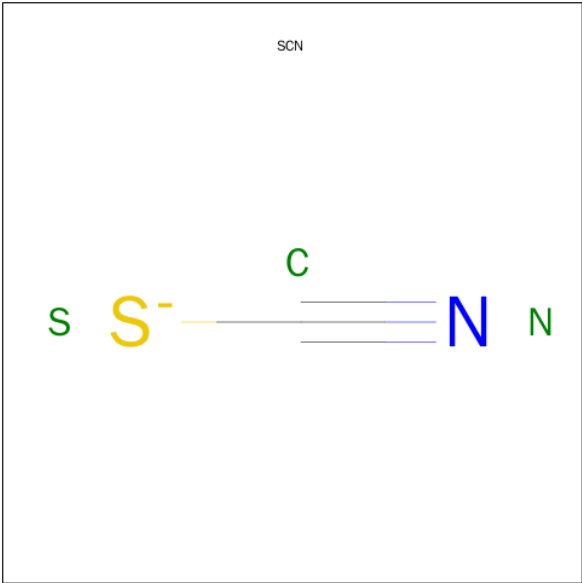
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	K	0	0
			1	1		

- Molecule 8 is (2R)-2-HYDROXY-3-(PHOSPHONOOXY)PROPYL (9E)-OCTADEC-9-EN OATE (three-letter code: NKP) (formula: C<sub>21</sub>H<sub>41</sub>O<sub>7</sub>P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	O	P	0	0
			29	21	7	1		
8	A	1	Total	C			0	0
			6	6				

- Molecule 9 is THIOCYANATE ION (three-letter code: SCN) (formula: CNS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	A	1	Total	C	N	S	0	0
			3	1	1	1		
9	A	1	Total	C	N	S	0	0
			3	1	1	1		

- Molecule 10 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



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

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

- Molecule 11 is water.

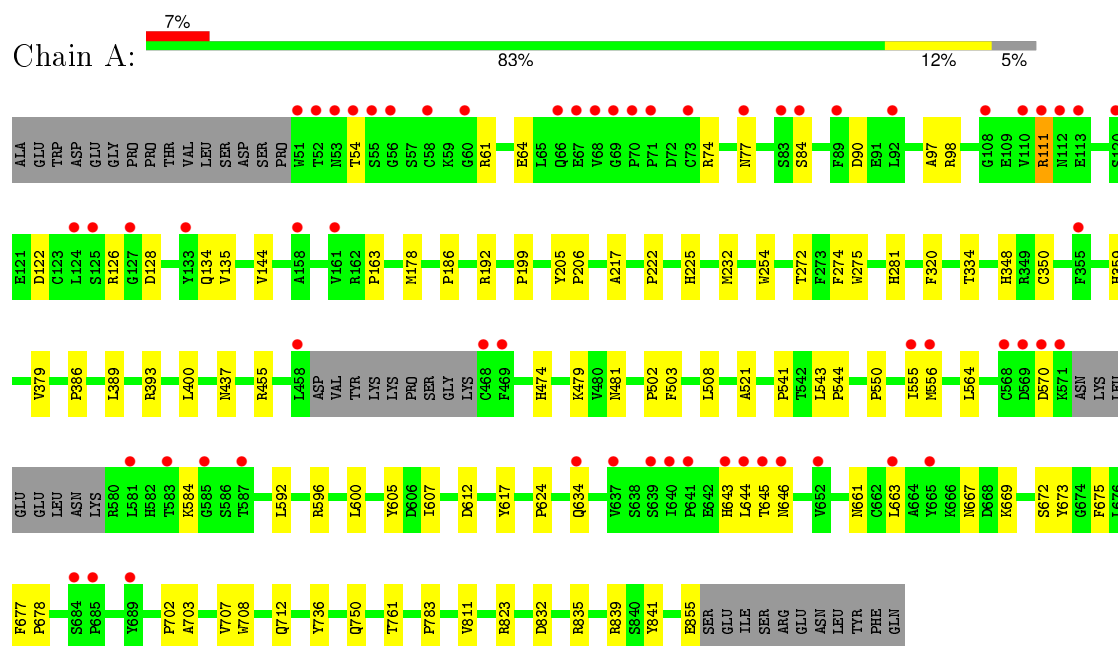
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
11	A	612	Total O 612 612	0	0



### 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: Ectonucleotide pyrophosphatase/phosphodiesterase family member 2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	61.52Å 94.52Å 75.50Å 90.00° 94.55° 90.00°	Depositor
Resolution (Å)	26.53 – 1.75 49.51 – 1.75	Depositor EDS
% Data completeness (in resolution range)	97.7 (26.53-1.75) 97.6 (49.51-1.75)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.06 (at 1.75Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6 _289)	Depositor
R, $R_{free}$	0.189 , 0.222 0.188 , 0.220	Depositor DCC
$R_{free}$ test set	4242 reflections (5.29%)	DCC
Wilson B-factor (Å <sup>2</sup> )	25.3	Xtriage
Anisotropy	0.175	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 50.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 84509 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	7182	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.83% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, BMA, NAG, NA, K, EDO, NKP, SCN, CA, MAN

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	0.37	0/6506	0.53	0/8844

There are no bond length outliers.

There are no bond angle outliers.

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	6312	0	5991	85	0
2	A	56	0	50	0	0
3	A	72	0	61	0	0
4	A	2	0	0	0	0
5	A	1	0	0	0	0
6	A	1	0	0	0	0
7	A	1	0	0	0	0
8	A	35	0	50	12	0
9	A	6	0	0	0	0
10	A	84	0	126	25	0
11	A	612	0	0	6	0
All	All	7182	0	6278	90	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 7.

All (90) 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:77:ASN:HD21	1:A:272:THR:HG21	1.12	1.11
1:A:77:ASN:ND2	1:A:272:THR:HG21	1.66	1.09
8:A:1006:NKP:HAT	8:A:1006:NKP:CAO	1.84	1.07
1:A:675:PHE:H	10:A:1012:EDO:H11	0.98	1.07
8:A:1006:NKP:HAOA	8:A:1006:NKP:HAT	1.40	1.03
1:A:835:ARG:HE	10:A:1024:EDO:H22	1.18	1.03
1:A:77:ASN:HD21	1:A:272:THR:CG2	1.77	0.97
1:A:503:PHE:HE2	1:A:508:LEU:HD21	1.28	0.97
1:A:675:PHE:N	10:A:1012:EDO:H11	1.82	0.94
8:A:1006:NKP:CAT	8:A:1006:NKP:CAO	2.54	0.85
1:A:334:THR:OG1	10:A:1023:EDO:H22	1.76	0.85
1:A:281:HIS:ND1	10:A:1023:EDO:H21	1.91	0.84
8:A:1006:NKP:HAO	8:A:1006:NKP:CAT	2.10	0.81
1:A:275:TRP:H	8:A:1006:NKP:CAQ	1.95	0.80
1:A:503:PHE:CE2	1:A:508:LEU:HD21	2.18	0.78
1:A:111:ARG:CZ	1:A:126:ARG:HD3	2.16	0.76
1:A:503:PHE:HE2	1:A:508:LEU:CD2	2.00	0.74
1:A:275:TRP:H	8:A:1006:NKP:HAQA	1.52	0.73
1:A:77:ASN:HD21	1:A:272:THR:CB	2.02	0.72
1:A:111:ARG:NH1	1:A:122:ASP:OD1	2.23	0.71
1:A:702:PRO:HB2	10:A:1025:EDO:H22	1.72	0.69
1:A:570:ASP:HB2	1:A:643:HIS:HD2	1.58	0.67
1:A:178:MET:HE2	1:A:192:ARG:HD3	1.76	0.66
1:A:835:ARG:NE	10:A:1024:EDO:H22	2.01	0.65
1:A:634:GLN:HG3	11:A:1392:HOH:O	1.97	0.63
1:A:543:LEU:HD12	1:A:544:PRO:HD2	1.82	0.62
1:A:678:PRO:HB3	1:A:712:GLN:HB3	1.82	0.61
1:A:835:ARG:HH21	10:A:1024:EDO:H11	1.65	0.60
1:A:281:HIS:ND1	10:A:1023:EDO:C2	2.64	0.60
1:A:600:LEU:HD11	1:A:832:ASP:HB2	1.82	0.60
1:A:275:TRP:N	8:A:1006:NKP:HAR	2.19	0.57
1:A:550:PRO:HB2	1:A:607:ILE:HG12	1.86	0.56
1:A:275:TRP:H	8:A:1006:NKP:CAR	2.19	0.55
1:A:199:PRO:HG2	1:A:502:PRO:HG3	1.89	0.55
1:A:111:ARG:NH2	1:A:126:ARG:HD3	2.22	0.54
1:A:222:PRO:HA	1:A:225:HIS:CE1	2.42	0.54
1:A:74:ARG:HD3	1:A:84:SER:OG	2.08	0.54
1:A:550:PRO:HB3	1:A:605:TYR:CE1	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:ARG:NH2	1:A:128:ASP:OD1	2.41	0.53
1:A:708:TRP:CD1	10:A:1012:EDO:H12	2.43	0.53
10:A:1030:EDO:H21	11:A:24:HOH:O	2.08	0.53
1:A:570:ASP:HB2	1:A:643:HIS:CD2	2.42	0.53
1:A:835:ARG:HE	10:A:1024:EDO:C2	2.06	0.53
1:A:592:LEU:HD21	1:A:596:ARG:HB3	1.91	0.53
1:A:275:TRP:H	8:A:1006:NKP:HAR	1.74	0.52
1:A:541:PRO:HG2	1:A:841:TYR:CD2	2.44	0.52
1:A:393:ARG:HA	10:A:1017:EDO:H21	1.92	0.51
1:A:134:GLN:NE2	11:A:1431:HOH:O	2.42	0.51
1:A:643:HIS:CD2	1:A:644:LEU:HG	2.45	0.51
1:A:661:ASN:OD1	1:A:663:LEU:HB3	2.11	0.51
1:A:61:ARG:HD2	1:A:64:GLU:OE1	2.10	0.50
1:A:783:PRO:HB3	10:A:1026:EDO:H21	1.92	0.50
1:A:437:ASN:HD21	10:A:1030:EDO:H22	1.76	0.50
1:A:555:ILE:O	1:A:556:MET:HG3	2.13	0.49
1:A:274:PHE:HA	8:A:1006:NKP:HAQ	1.95	0.49
1:A:703:ALA:HB2	10:A:1025:EDO:H21	1.95	0.49
1:A:379:VAL:HB	1:A:455:ARG:HG2	1.96	0.48
1:A:612:ASP:HB3	1:A:677:PHE:CZ	2.49	0.48
1:A:437:ASN:HD21	10:A:1030:EDO:C2	2.27	0.48
1:A:675:PHE:H	10:A:1012:EDO:C1	1.93	0.47
1:A:144:VAL:HG21	1:A:186:PRO:HB2	1.97	0.47
1:A:521:ALA:O	10:A:1018:EDO:H22	2.15	0.47
1:A:761:THR:HA	10:A:1015:EDO:H12	1.97	0.47
1:A:217:ALA:HA	8:A:1006:NKP:HBBB	1.97	0.47
1:A:672:SER:OG	1:A:673:TYR:N	2.47	0.47
1:A:479:LYS:O	1:A:855:GLU:HG3	2.15	0.46
1:A:272:THR:O	1:A:272:THR:HG23	2.14	0.46
1:A:564:LEU:O	1:A:661:ASN:HB2	2.16	0.46
1:A:205:TYR:CD2	1:A:206:PRO:HA	2.51	0.46
1:A:320:PHE:HE2	11:A:1464:HOH:O	1.99	0.46
1:A:359:HIS:CE1	1:A:474:HIS:CE1	3.04	0.45
1:A:541:PRO:HG2	1:A:841:TYR:HD2	1.81	0.45
1:A:839:ARG:HE	10:A:1026:EDO:H22	1.82	0.45
1:A:163:PRO:HB3	1:A:350:CYS:O	2.17	0.44
1:A:839:ARG:NE	10:A:1026:EDO:H22	2.32	0.44
1:A:90:ASP:HB3	11:A:1477:HOH:O	2.19	0.43
1:A:707:VAL:HG13	1:A:811:VAL:HG13	2.00	0.42
1:A:645:THR:HG22	1:A:646:ASN:ND2	2.35	0.42
1:A:675:PHE:HB2	10:A:1012:EDO:H21	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:736:TYR:OH	1:A:750:GLN:HB3	2.18	0.42
1:A:570:ASP:OD1	1:A:643:HIS:NE2	2.52	0.42
1:A:348:HIS:CD2	1:A:348:HIS:H	2.38	0.42
1:A:135:VAL:HG23	11:A:1308:HOH:O	2.19	0.42
1:A:97:ALA:O	1:A:98:ARG:HB2	2.20	0.41
1:A:254:TRP:HB2	8:A:1007:NKP:HAYA	2.01	0.41
1:A:823:ARG:CZ	10:A:1010:EDO:H22	2.50	0.41
1:A:617:TYR:HA	1:A:624:PRO:HA	2.02	0.41
1:A:272:THR:O	1:A:272:THR:CG2	2.69	0.41
1:A:675:PHE:HB2	10:A:1012:EDO:C2	2.51	0.40
1:A:206:PRO:HB3	1:A:389:LEU:HD13	2.04	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	786/831 (95%)	762 (97%)	22 (3%)	2 (0%)	46 25

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	584	LYS
1	A	386	PRO

### 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	703/756 (93%)	696 (99%)	7 (1%)	82	69

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	THR
1	A	111	ARG
1	A	232	MET
1	A	400	LEU
1	A	481	ASN
1	A	667	ASN
1	A	669	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	GLN
1	A	77	ASN
1	A	118	HIS
1	A	134	GLN
1	A	337	GLN
1	A	348	HIS
1	A	374	ASN
1	A	398	ASN
1	A	643	HIS
1	A	646	ASN
1	A	667	ASN
1	A	827	HIS

### 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 ⓘ

10 carbohydrates 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	NAG	A	898	1,2	14,14,15	0.43	0	15,19,21	1.56	1 (6%)
2	NAG	A	899	2	14,14,15	0.54	0	15,19,21	0.99	1 (6%)
3	NAG	A	900	1,3	14,14,15	0.60	0	15,19,21	1.22	1 (6%)
3	NAG	A	901	3	14,14,15	0.53	0	15,19,21	0.84	0
3	BMA	A	902	3	11,11,12	0.68	0	14,15,17	0.83	1 (7%)
3	MAN	A	903	3	11,11,12	0.53	0	14,15,17	0.77	0
3	MAN	A	904	3	11,11,12	0.59	0	14,15,17	0.97	0
3	MAN	A	905	3	11,11,12	0.64	0	14,15,17	0.92	0
2	NAG	A	907	1,2	14,14,15	0.68	0	15,19,21	1.02	1 (6%)
2	NAG	A	908	2	14,14,15	0.62	0	15,19,21	0.80	1 (6%)

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	NAG	A	898	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	899	2	-	0/6/23/26	0/1/1/1
3	NAG	A	900	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	901	3	-	0/6/23/26	0/1/1/1
3	BMA	A	902	3	-	0/2/19/22	0/1/1/1
3	MAN	A	903	3	-	0/2/19/22	0/1/1/1
3	MAN	A	904	3	-	0/2/19/22	0/1/1/1
3	MAN	A	905	3	-	0/2/19/22	0/1/1/1
2	NAG	A	907	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	908	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	902	BMA	O5-C1-C2	-2.19	107.31	110.86
2	A	907	NAG	O4-C4-C5	-2.10	103.69	109.24
2	A	899	NAG	C1-O5-C5	2.12	114.94	112.25
3	A	900	NAG	C2-N2-C7	2.14	125.79	123.04
2	A	908	NAG	C1-O5-C5	2.17	115.00	112.25
2	A	898	NAG	C1-O5-C5	5.21	118.86	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.6 Ligand geometry

Of 30 ligands modelled in this entry, 5 are monoatomic - leaving 25 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	$\# Z  > 2$
8	NKP	A	1006	4	28,28,28	0.79	1 (3%)	30,32,32	1.60	8 (26%)
8	NKP	A	1007	-	5,5,28	0.27	0	4,4,32	0.55	0
9	SCN	A	1008	-	2,2,2	1.64	1 (50%)	1,1,1	0.31	0
9	SCN	A	1009	-	2,2,2	1.81	1 (50%)	1,1,1	0.35	0
10	EDO	A	1010	-	3,3,3	0.50	0	2,2,2	0.60	0
10	EDO	A	1011	-	3,3,3	0.65	0	2,2,2	0.08	0
10	EDO	A	1012	-	3,3,3	0.33	0	2,2,2	0.71	0
10	EDO	A	1013	-	3,3,3	0.51	0	2,2,2	0.38	0
10	EDO	A	1014	-	3,3,3	0.53	0	2,2,2	0.14	0
10	EDO	A	1015	-	3,3,3	0.47	0	2,2,2	0.35	0
10	EDO	A	1016	-	3,3,3	0.53	0	2,2,2	0.39	0
10	EDO	A	1017	-	3,3,3	0.49	0	2,2,2	0.36	0
10	EDO	A	1018	-	3,3,3	0.55	0	2,2,2	0.26	0
10	EDO	A	1019	-	3,3,3	0.44	0	2,2,2	0.39	0
10	EDO	A	1020	-	3,3,3	0.47	0	2,2,2	0.55	0
10	EDO	A	1021	-	3,3,3	0.49	0	2,2,2	0.40	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
10	EDO	A	1022	-	3,3,3	0.49	0	2,2,2	0.40	0
10	EDO	A	1023	-	3,3,3	0.50	0	2,2,2	0.19	0
10	EDO	A	1024	-	3,3,3	0.51	0	2,2,2	0.35	0
10	EDO	A	1025	-	3,3,3	0.48	0	2,2,2	0.37	0
10	EDO	A	1026	-	3,3,3	0.43	0	2,2,2	0.43	0
10	EDO	A	1027	-	3,3,3	0.52	0	2,2,2	0.38	0
10	EDO	A	1028	-	3,3,3	0.51	0	2,2,2	0.45	0
10	EDO	A	1029	-	3,3,3	0.54	0	2,2,2	0.39	0
10	EDO	A	1030	-	3,3,3	0.52	0	2,2,2	0.11	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
8	NKP	A	1006	4	-	0/28/28/28	0/0/0/0
8	NKP	A	1007	-	-	0/3/3/28	0/0/0/0
9	SCN	A	1008	-	-	0/0/0/0	0/0/0/0
9	SCN	A	1009	-	-	0/0/0/0	0/0/0/0
10	EDO	A	1010	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1011	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1012	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1013	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1014	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1015	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1016	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1017	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1018	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1019	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1020	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1021	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1022	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1023	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1024	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1025	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1026	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1027	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1028	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1029	-	-	0/1/1/1	0/0/0/0
10	EDO	A	1030	-	-	0/1/1/1	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	1006	NKP	CAR-CAS	-2.61	1.38	1.50
9	A	1008	SCN	C-S	2.26	1.78	1.63
9	A	1009	SCN	C-S	2.51	1.79	1.63

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	1006	NKP	CAI-OAJ-CAK	-3.44	107.22	116.85
8	A	1006	NKP	CAW-CAV-CAU	-2.33	104.76	113.86
8	A	1006	NKP	OAA-PAC-OAF	-2.28	99.99	106.56
8	A	1006	NKP	OAJ-CAK-OAE	-2.03	118.26	123.49
8	A	1006	NKP	CAN-CAM-CAL	-2.02	105.90	113.29
8	A	1006	NKP	OAJ-CAK-CAL	2.40	119.20	111.90
8	A	1006	NKP	CAP-CAQ-CAR	2.69	124.35	113.86
8	A	1006	NKP	OBC-CAH-CAG	3.09	121.13	109.35

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

12 monomers are involved in 37 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	1006	NKP	11	0
8	A	1007	NKP	1	0
10	A	1010	EDO	1	0
10	A	1012	EDO	6	0
10	A	1015	EDO	1	0
10	A	1017	EDO	1	0
10	A	1018	EDO	1	0
10	A	1023	EDO	3	0
10	A	1024	EDO	4	0
10	A	1025	EDO	2	0
10	A	1026	EDO	3	0
10	A	1030	EDO	3	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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	788/831 (94%)	0.42	61 (7%) 16 20	14, 30, 56, 73	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	51	TRP	7.3
1	A	110	VAL	7.2
1	A	83	SER	6.4
1	A	585	GLY	5.1
1	A	569	ASP	4.6
1	A	583	THR	4.3
1	A	52	THR	4.2
1	A	124	LEU	4.1
1	A	637	VAL	4.0
1	A	73	CYS	4.0
1	A	581	LEU	4.0
1	A	71	PRO	3.9
1	A	69	GLY	3.9
1	A	55	SER	3.8
1	A	68	VAL	3.6
1	A	640	ILE	3.6
1	A	641	PRO	3.6
1	A	644	LEU	3.5
1	A	70	PRO	3.4
1	A	469	PHE	3.4
1	A	645	THR	3.3
1	A	571	LYS	3.2
1	A	92	LEU	3.2
1	A	685	PRO	3.2
1	A	66	GLN	3.2
1	A	458	LEU	3.1
1	A	125	SER	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	56	GLY	3.0
1	A	108	GLY	2.9
1	A	684	SER	2.8
1	A	67	GLU	2.8
1	A	58	CYS	2.8
1	A	556	MET	2.7
1	A	643	HIS	2.7
1	A	53	ASN	2.7
1	A	639	SER	2.7
1	A	587	THR	2.6
1	A	665	TYR	2.6
1	A	111	ARG	2.6
1	A	158	ALA	2.6
1	A	120	SER	2.5
1	A	54	THR	2.5
1	A	161	VAL	2.5
1	A	646	ASN	2.5
1	A	77	ASN	2.4
1	A	570	ASP	2.4
1	A	634	GLN	2.4
1	A	652	VAL	2.4
1	A	112	ASN	2.4
1	A	568	CYS	2.3
1	A	84	SER	2.3
1	A	663	LEU	2.3
1	A	468	CYS	2.3
1	A	555	ILE	2.2
1	A	127	GLY	2.2
1	A	113	GLU	2.2
1	A	60	GLY	2.2
1	A	89	PHE	2.1
1	A	689	TYR	2.0
1	A	355	PHE	2.0
1	A	133	TYR	2.0

## 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 ⓘ

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	MAN	A	904	11/12	0.94	0.18	4.80	32,35,41,51	0
2	NAG	A	907	14/15	0.93	0.11	1.72	26,31,39,47	0
2	NAG	A	898	14/15	0.77	0.27	0.52	66,68,71,72	0
3	NAG	A	900	14/15	0.96	0.09	-1.91	16,20,22,23	0
3	MAN	A	905	11/12	0.94	0.10	-	25,31,37,38	0
3	MAN	A	903	11/12	0.80	0.13	-	39,41,46,47	0
2	NAG	A	899	14/15	0.83	0.21	-	65,72,77,77	0
3	NAG	A	901	14/15	0.95	0.08	-	23,31,39,42	0
3	BMA	A	902	11/12	0.84	0.13	-	40,43,51,51	0
2	NAG	A	908	14/15	0.82	0.22	-	50,56,58,58	0

### 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
10	EDO	A	1030	4/4	0.92	0.18	7.97	18,27,33,37	0
10	EDO	A	1021	4/4	0.94	0.18	7.48	34,37,39,43	0
10	EDO	A	1018	4/4	0.83	0.18	7.46	25,31,32,40	0
10	EDO	A	1013	4/4	0.93	0.15	6.35	33,33,33,33	0
8	NKP	A	1007	6/29	0.92	0.16	5.82	34,37,38,40	0
10	EDO	A	1017	4/4	0.95	0.16	4.41	22,25,29,29	0
9	SCN	A	1008	3/3	0.81	0.16	3.86	24,24,36,41	0
8	NKP	A	1006	29/29	0.93	0.19	3.73	21,37,46,51	0
10	EDO	A	1026	4/4	0.92	0.16	3.70	21,28,34,39	0
10	EDO	A	1024	4/4	0.72	0.20	3.46	46,47,49,50	0
10	EDO	A	1016	4/4	0.94	0.12	3.18	27,29,30,31	0
10	EDO	A	1010	4/4	0.92	0.14	2.92	33,33,35,36	0
10	EDO	A	1027	4/4	0.93	0.12	2.12	31,33,41,43	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
10	EDO	A	1028	4/4	0.97	0.09	2.04	19,21,23,24	0
10	EDO	A	1011	4/4	0.86	0.14	1.91	26,30,32,34	0
10	EDO	A	1029	4/4	0.95	0.09	1.68	25,26,32,34	0
10	EDO	A	1022	4/4	0.88	0.11	1.62	34,35,38,46	0
9	SCN	A	1009	3/3	0.98	0.12	0.77	26,26,36,50	0
10	EDO	A	1020	4/4	0.96	0.10	0.51	25,27,33,37	0
10	EDO	A	1023	4/4	0.95	0.14	0.46	25,30,32,35	0
4	ZN	A	1002	1/1	1.00	0.11	0.27	18,18,18,18	0
10	EDO	A	1015	4/4	0.95	0.09	-0.25	28,30,32,35	0
10	EDO	A	1012	4/4	0.94	0.11	-0.37	30,33,33,35	0
10	EDO	A	1014	4/4	0.93	0.10	-0.50	28,29,29,33	0
5	CA	A	1003	1/1	1.00	0.09	-0.53	23,23,23,23	0
10	EDO	A	1019	4/4	0.95	0.08	-0.94	32,32,36,42	0
4	ZN	A	1001	1/1	0.99	0.10	-1.46	21,21,21,21	0
7	K	A	1005	1/1	0.93	0.09	-1.55	38,38,38,38	0
10	EDO	A	1025	4/4	0.96	0.29	-	37,39,41,44	0
6	NA	A	1004	1/1	0.98	0.07	-	25,25,25,25	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.