



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:38 PM GMT

PDB ID : 1GKA  
Title : THE MOLECULAR BASIS OF THE COLORATION MECHANISM IN LOBSTER SHELL. BETA-CRUSTACYANIN AT 3.2 Å RESOLUTION  
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Deposited on : 2001-08-10  
Resolution : 3.23 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

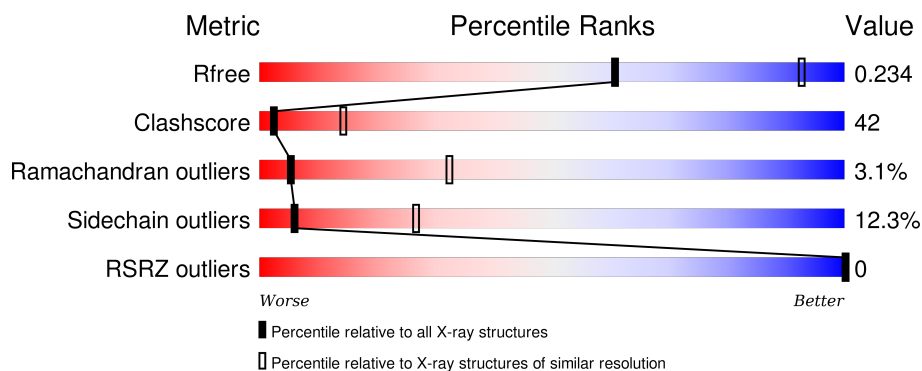
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.23 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	1092 (3.28-3.20)
Clashscore	102246	1227 (3.28-3.20)
Ramachandran outliers	100387	1204 (3.28-3.20)
Sidechain outliers	100360	1203 (3.28-3.20)
RSRZ outliers	91569	1097 (3.28-3.20)

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	180	<div> <div>28%</div> <div>59%</div> <div>12%</div> <div>.</div> </div>
2	B	174	<div> <div>49%</div> <div>40%</div> <div>10%</div> <div>.</div> </div>

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
3	AXT	A	1182	-	-	-	X
3	AXT	B	1175	-	-	-	X
4	D12	A	1183	-	-	X	X
5	TRS	B	1176	-	X	-	X
6	EPE	B	1177	-	-	-	X

## 2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 2996 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.

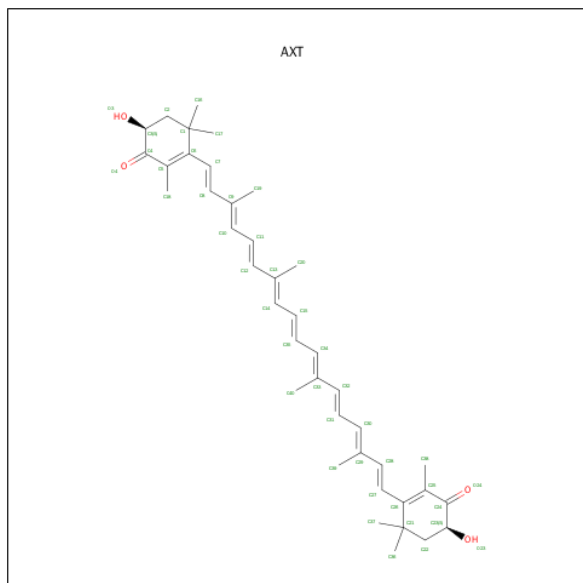
- Molecule 1 is a protein called CRUSTACYANIN A1 SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	180	Total	C	N	O	S	0	0	0
			1452	931	240	275	6			

- Molecule 2 is a protein called CRUSTACYANIN A2 SUBUNIT.

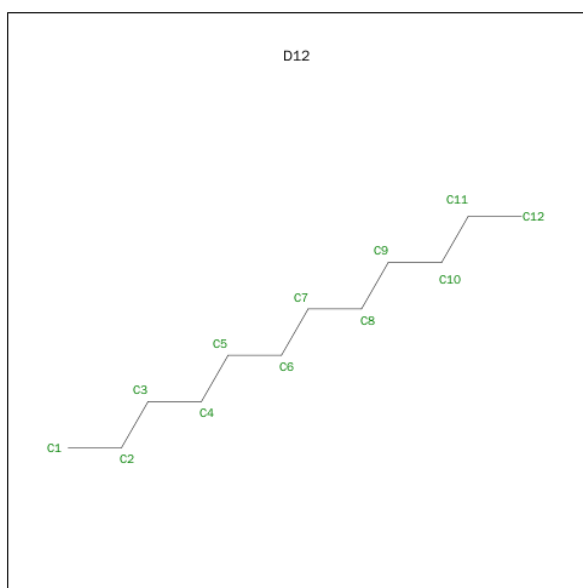
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	174	Total	C	N	O	S	0	0	0
			1391	891	224	270	6			

- Molecule 3 is ASTAXANTHIN (three-letter code: AXT) (formula:  $C_{40}H_{52}O_4$ ).



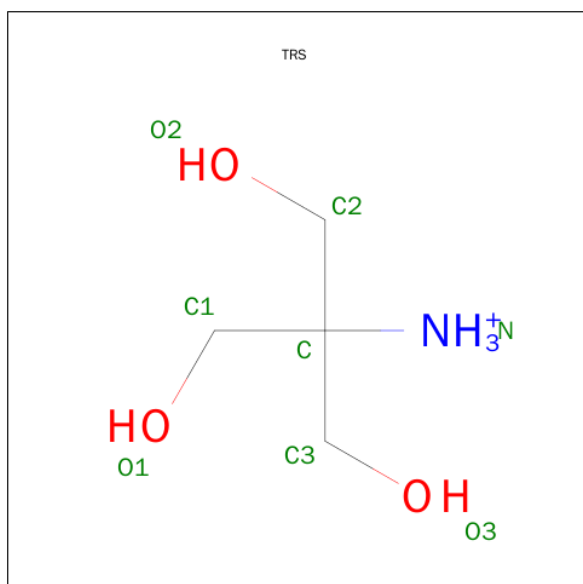
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			44	40	4		
3	B	1	Total	C	O	0	0
			44	40	4		

- Molecule 4 is DODECANE (three-letter code: D12) (formula:  $C_{12}H_{26}$ ).



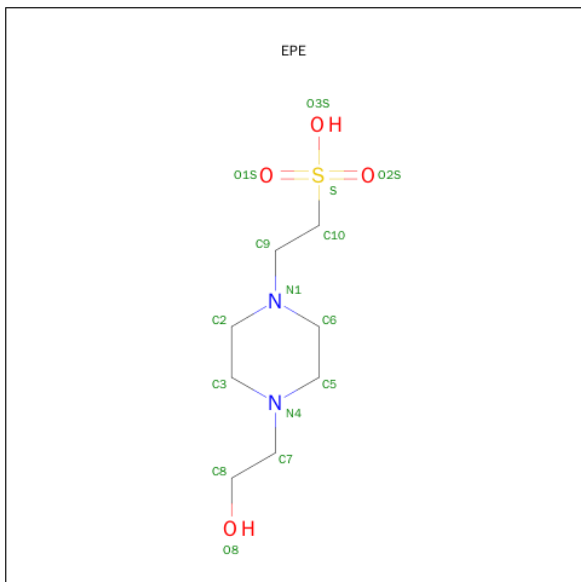
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C		0	0
			12	12			

- Molecule 5 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula:  $C_4H_{12}NO_3$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 6 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula:  $C_8H_{18}N_2O_4S$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

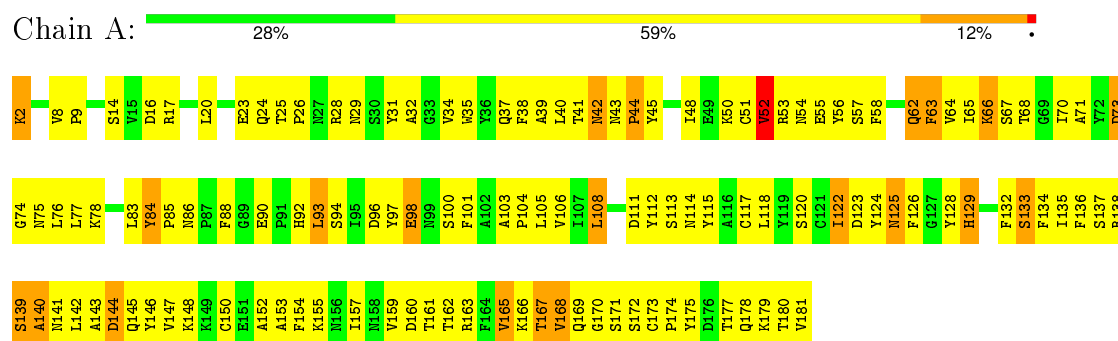
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	11	Total	O	0	0
			11	11		
7	B	19	Total	O	0	0
			19	19		

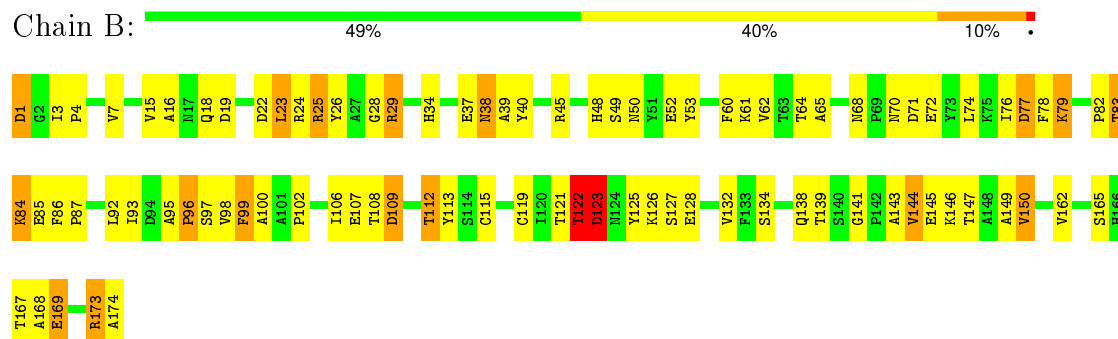
### 3 Residue-property plots

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: CRUSTACYANIN A1 SUBUNIT



#### • Molecule 2: CRUSTACYANIN A2 SUBUNIT



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	155.47Å 155.47Å 168.50Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	100.00 – 3.23 77.73 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.7 (100.00-3.23) 99.6 (77.73-3.00)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.37 (at 3.01Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.213 , 0.252 0.207 , 0.234	Depositor DCC
$R_{free}$ test set	1012 reflections (5.10%)	DCC
Wilson B-factor (Å <sup>2</sup> )	71.2	Xtriage
Anisotropy	0.174	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 46.0	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 24516 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	2996	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.88% 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: TRS, EPE, D12, AXT

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	1.45	12/1492 (0.8%)	1.25	7/2025 (0.3%)
2	B	1.55	13/1433 (0.9%)	1.33	14/1953 (0.7%)
All	All	1.50	25/2925 (0.9%)	1.29	21/3978 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	2

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	98	GLU	CG-CD	8.88	1.65	1.51
1	A	63	PHE	CB-CG	-7.25	1.39	1.51
1	A	98	GLU	CB-CG	7.18	1.65	1.52
2	B	85	GLU	CG-CD	7.09	1.62	1.51
1	A	66	LYS	CD-CE	7.05	1.68	1.51
1	A	64	VAL	CA-CB	-6.71	1.40	1.54
2	B	145	GLU	CG-CD	6.60	1.61	1.51
2	B	145	GLU	CD-OE1	6.41	1.32	1.25
2	B	99	PHE	CB-CG	-6.09	1.41	1.51
2	B	150	VAL	CB-CG2	-6.07	1.40	1.52
1	A	23	GLU	CG-CD	5.96	1.60	1.51
1	A	106	VAL	CB-CG2	-5.77	1.40	1.52
2	B	145	GLU	CD-OE2	5.75	1.31	1.25
2	B	122	THR	CB-CG2	-5.72	1.33	1.52
1	A	55	GLU	CG-CD	5.62	1.60	1.51
2	B	162	VAL	CB-CG2	-5.58	1.41	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	79	LYS	CE-NZ	5.37	1.62	1.49
1	A	155	LYS	CD-CE	5.30	1.64	1.51
2	B	19	ASP	CB-CG	5.20	1.62	1.51
2	B	52	GLU	CD-OE1	5.19	1.31	1.25
2	B	169	GLU	CG-CD	5.18	1.59	1.51
2	B	144	VAL	CB-CG1	-5.18	1.42	1.52
1	A	55	GLU	CD-OE2	5.13	1.31	1.25
1	A	135	ILE	CA-CB	-5.13	1.43	1.54
1	A	52	VAL	CB-CG1	-5.12	1.42	1.52

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	70	ASN	N-CA-C	-6.60	93.18	111.00
1	A	122	ILE	CG1-CB-CG2	-6.57	96.95	111.40
1	A	138	ARG	NE-CZ-NH2	-6.10	117.25	120.30
2	B	25	ARG	NE-CZ-NH1	-6.01	117.29	120.30
1	A	96	ASP	N-CA-CB	-5.95	99.90	110.60
2	B	23	LEU	CA-CB-CG	-5.74	102.11	115.30
2	B	123	ASP	N-CA-C	-5.65	95.75	111.00
1	A	66	LYS	CD-CE-NZ	5.64	124.68	111.70
2	B	19	ASP	C-N-CA	-5.54	107.86	121.70
2	B	92	LEU	CB-CG-CD1	-5.41	101.80	111.00
2	B	92	LEU	CA-CB-CG	5.39	127.69	115.30
2	B	127	SER	N-CA-C	-5.36	96.53	111.00
2	B	24	ARG	NE-CZ-NH2	-5.36	117.62	120.30
2	B	61	LYS	N-CA-C	-5.30	96.68	111.00
1	A	16	ASP	CB-CG-OD2	5.25	123.03	118.30
2	B	61	LYS	CD-CE-NZ	5.19	123.63	111.70
2	B	71	ASP	N-CA-C	5.15	124.90	111.00
1	A	63	PHE	CB-CA-C	-5.13	100.13	110.40
2	B	79	LYS	CD-CE-NZ	5.11	123.45	111.70
1	A	138	ARG	NE-CZ-NH1	5.05	122.83	120.30
2	B	25	ARG	NE-CZ-NH2	5.03	122.81	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	26	TYR	Sidechain
2	B	53	TYR	Sidechain

## 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	1452	0	1383	134	1
2	B	1391	0	1299	103	0
3	A	44	0	52	16	0
3	B	44	0	52	8	0
4	A	12	0	26	10	0
5	B	8	0	12	3	0
6	B	15	0	18	0	0
7	A	11	0	0	2	0
7	B	19	0	0	7	1
All	All	2996	0	2842	245	2

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 42.

All (245) 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:38:ASN:C	2:B:38:ASN:HD22	1.55	1.10
1:A:142:LEU:HD23	1:A:147:VAL:HG23	1.09	1.07
2:B:115:CYS:SG	2:B:147:THR:OG1	2.06	1.05
1:A:40:LEU:O	1:A:40:LEU:HD12	1.58	1.03
2:B:122:THR:HG22	7:B:2014:HOH:O	1.57	1.03
2:B:112:THR:HG22	2:B:113:TYR:HD2	1.22	1.02
1:A:2:LYS:HB3	2:B:100:ALA:HB2	1.47	0.95
1:A:142:LEU:CD2	1:A:147:VAL:HG23	1.99	0.93
1:A:37:GLN:NE2	1:A:40:LEU:HD23	1.89	0.88
2:B:23:LEU:HD12	2:B:23:LEU:H	1.36	0.88
3:B:1175:AXT:H8	3:B:1175:AXT:H161	1.56	0.87
1:A:48:ILE:HG22	1:A:52:VAL:HG22	1.56	0.87
2:B:23:LEU:HD12	2:B:23:LEU:N	1.90	0.86
3:A:1182:AXT:H28	3:A:1182:AXT:H361	1.61	0.83
2:B:45:ARG:HH22	2:B:169:GLU:HG2	1.43	0.83
2:B:45:ARG:HH12	2:B:169:GLU:HB3	1.41	0.82
2:B:95:ALA:HB1	2:B:96:PRO:HD2	1.63	0.81
2:B:45:ARG:HH22	2:B:169:GLU:CG	1.94	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:173:ARG:HD3	7:B:2018:HOH:O	1.80	0.80
2:B:38:ASN:C	2:B:38:ASN:ND2	2.32	0.80
1:A:108:LEU:HD11	1:A:157:ILE:HD13	1.65	0.79
3:A:1182:AXT:H8	3:A:1182:AXT:H161	1.66	0.77
3:B:1175:AXT:H361	3:B:1175:AXT:H28	1.64	0.77
1:A:160:ASP:OD2	1:A:162:THR:HG23	1.84	0.76
1:A:66:LYS:HE3	1:A:78:LYS:HD2	1.69	0.73
2:B:112:THR:HG22	2:B:113:TYR:CD2	2.15	0.73
1:A:28:ARG:HG2	1:A:58:PHE:CE2	2.23	0.73
2:B:132:VAL:HG22	2:B:147:THR:HG21	1.70	0.72
1:A:86:ASN:HD22	1:A:90:GLU:H	1.35	0.72
1:A:143:ALA:O	1:A:145:GLN:N	2.21	0.72
1:A:93:LEU:H	1:A:93:LEU:HD12	1.56	0.70
2:B:45:ARG:NE	7:B:2008:HOH:O	2.18	0.69
2:B:76:ILE:HG22	2:B:77:ASP:N	2.07	0.69
1:A:40:LEU:HD12	1:A:40:LEU:C	2.13	0.69
1:A:62:GLN:HB2	1:A:84:TYR:HA	1.75	0.69
1:A:48:ILE:CG2	1:A:52:VAL:HG22	2.23	0.69
2:B:109:ASP:OD2	2:B:112:THR:HB	1.92	0.68
4:A:1183:D12:H122	2:B:99:PHE:CD2	2.27	0.68
1:A:180:THR:O	1:A:181:VAL:HB	1.92	0.68
1:A:78:LYS:HD3	1:A:181:VAL:HG22	1.76	0.68
1:A:133:SER:HB3	1:A:154:PHE:HE2	1.58	0.68
1:A:98:GLU:OE2	2:B:1:ASP:HB3	1.94	0.68
1:A:97:TYR:CE1	3:A:1182:AXT:H173	2.29	0.67
1:A:142:LEU:HD23	1:A:147:VAL:CG2	2.04	0.67
2:B:18:GLN:HG2	2:B:107:GLU:HA	1.76	0.67
3:B:1175:AXT:H361	3:B:1175:AXT:C28	2.24	0.67
2:B:115:CYS:SG	2:B:147:THR:CB	2.83	0.67
1:A:101:PHE:HB2	4:A:1183:D12:H61	1.78	0.65
1:A:84:TYR:HE1	1:A:94:SER:HB2	1.61	0.65
1:A:100:SER:O	2:B:1:ASP:HA	1.96	0.64
2:B:132:VAL:HG22	2:B:147:THR:CG2	2.27	0.64
1:A:84:TYR:CE1	1:A:94:SER:HB2	2.33	0.64
1:A:8:VAL:HG13	1:A:9:PRO:HD2	1.80	0.63
1:A:25:THR:N	1:A:26:PRO:CD	2.61	0.63
2:B:115:CYS:SG	2:B:147:THR:HG23	2.39	0.63
1:A:8:VAL:O	1:A:129:HIS:HD2	1.82	0.63
2:B:68:ASN:HD22	2:B:72:GLU:HG3	1.62	0.63
1:A:133:SER:HB3	1:A:154:PHE:CE2	2.33	0.62
2:B:126:LYS:HE2	7:B:2003:HOH:O	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:25:ARG:HH12	5:B:1176:TRS:C3	2.12	0.62
2:B:74:LEU:N	2:B:74:LEU:HD12	2.15	0.62
1:A:108:LEU:HD11	1:A:157:ILE:CD1	2.29	0.62
3:B:1175:AXT:C16	3:B:1175:AXT:H8	2.28	0.61
1:A:86:ASN:ND2	1:A:88:PHE:H	1.98	0.61
1:A:84:TYR:O	1:A:93:LEU:HB3	2.01	0.61
1:A:86:ASN:ND2	1:A:90:GLU:H	1.97	0.60
2:B:122:THR:HG22	2:B:123:ASP:N	2.16	0.60
1:A:125:ASN:OD1	1:A:125:ASN:N	2.33	0.60
1:A:56:TYR:CZ	1:A:65:ILE:HD13	2.37	0.60
2:B:86:PHE:HB3	2:B:87:PRO:HD2	1.83	0.60
1:A:51:CYS:HA	1:A:169:GLN:OE1	2.02	0.59
2:B:96:PRO:HG2	2:B:97:SER:H	1.67	0.59
2:B:4:PRO:HG2	3:B:1175:AXT:H403	1.84	0.58
1:A:50:LYS:HB2	1:A:168:VAL:HG12	1.86	0.58
2:B:76:ILE:CG2	2:B:77:ASP:N	2.67	0.58
3:A:1182:AXT:O24	2:B:102:PRO:HB3	2.03	0.58
2:B:45:ARG:HH12	2:B:169:GLU:CB	2.14	0.57
1:A:124:TYR:OH	3:A:1182:AXT:H34	2.04	0.57
2:B:15:VAL:HG11	2:B:106:ILE:HD11	1.87	0.57
1:A:114:ASN:ND2	1:A:146:TYR:OH	2.38	0.57
2:B:25:ARG:HH12	5:B:1176:TRS:H32	1.70	0.57
1:A:143:ALA:C	1:A:145:GLN:H	2.06	0.57
1:A:160:ASP:O	1:A:163:ARG:HB2	2.05	0.57
1:A:93:LEU:HD12	7:A:2008:HOH:O	2.06	0.56
2:B:109:ASP:OD1	2:B:112:THR:N	2.31	0.56
2:B:99:PHE:O	2:B:100:ALA:C	2.43	0.56
1:A:34:VAL:O	1:A:34:VAL:HG23	2.03	0.56
1:A:25:THR:N	1:A:26:PRO:HD3	2.21	0.56
1:A:39:ALA:HB2	1:A:166:LYS:HA	1.88	0.56
2:B:123:ASP:N	7:B:2014:HOH:O	2.11	0.55
1:A:174:PRO:O	1:A:177:THR:N	2.38	0.55
3:A:1182:AXT:H361	3:A:1182:AXT:C28	2.35	0.55
4:A:1183:D12:H81	4:A:1183:D12:H41	1.88	0.55
3:A:1182:AXT:H402	2:B:99:PHE:HE2	1.71	0.55
3:A:1182:AXT:C8	3:A:1182:AXT:H161	2.36	0.55
1:A:45:TYR:HE2	4:A:1183:D12:H32	1.72	0.55
1:A:2:LYS:CB	2:B:100:ALA:HB2	2.31	0.54
2:B:115:CYS:SG	2:B:147:THR:CG2	2.96	0.54
1:A:168:VAL:O	1:A:168:VAL:CG1	2.56	0.54
2:B:23:LEU:N	2:B:23:LEU:CD1	2.66	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:62:VAL:HG11	2:B:78:PHE:CZ	2.43	0.54
2:B:23:LEU:H	2:B:23:LEU:CD1	2.13	0.54
1:A:68:THR:HB	1:A:178:GLN:HE21	1.73	0.54
1:A:45:TYR:CE2	4:A:1183:D12:H32	2.43	0.53
1:A:28:ARG:HG2	1:A:58:PHE:CZ	2.43	0.53
2:B:38:ASN:HD22	2:B:39:ALA:N	2.05	0.53
2:B:125:TYR:N	7:B:2014:HOH:O	2.27	0.53
2:B:64:THR:CG2	2:B:65:ALA:N	2.72	0.53
1:A:115:TYR:N	1:A:115:TYR:CD2	2.76	0.53
1:A:97:TYR:CD1	3:A:1182:AXT:H173	2.43	0.52
1:A:56:TYR:CZ	1:A:65:ILE:CD1	2.92	0.52
2:B:68:ASN:HD22	2:B:72:GLU:CG	2.22	0.52
2:B:29:ARG:NH2	2:B:174:ALA:O	2.30	0.51
1:A:58:PHE:HD2	1:A:63:PHE:CD2	2.28	0.51
2:B:106:ILE:HG21	2:B:150:VAL:HG12	1.92	0.51
1:A:38:PHE:CZ	1:A:147:VAL:HG21	2.46	0.51
1:A:118:LEU:HB2	1:A:134:PHE:HB2	1.91	0.51
1:A:54:ASN:OD1	1:A:67:SER:OG	2.20	0.51
2:B:68:ASN:ND2	2:B:74:LEU:HD11	2.26	0.51
1:A:29:ASN:C	1:A:31:TYR:H	2.13	0.51
1:A:117:CYS:C	1:A:118:LEU:HD23	2.31	0.51
1:A:77:LEU:HD12	1:A:78:LYS:N	2.26	0.51
1:A:93:LEU:N	1:A:93:LEU:HD12	2.25	0.50
1:A:40:LEU:CD1	1:A:40:LEU:C	2.80	0.50
1:A:37:GLN:HE22	1:A:40:LEU:HD23	1.73	0.50
2:B:119:CYS:SG	2:B:128:GLU:HB2	2.52	0.50
1:A:78:LYS:HD3	1:A:181:VAL:HG13	1.93	0.50
1:A:115:TYR:HA	1:A:136:PHE:O	2.10	0.50
1:A:137:SER:CB	1:A:142:LEU:HD13	2.40	0.50
2:B:23:LEU:HD23	2:B:60:PHE:CE1	2.47	0.49
1:A:165:VAL:O	1:A:165:VAL:HG12	2.11	0.49
1:A:37:GLN:HE21	1:A:167:THR:HG21	1.76	0.49
2:B:122:THR:CG2	2:B:123:ASP:N	2.73	0.49
2:B:78:PHE:N	2:B:78:PHE:CD2	2.80	0.49
1:A:146:TYR:O	1:A:148:LYS:N	2.46	0.49
1:A:146:TYR:O	1:A:147:VAL:C	2.52	0.48
1:A:48:ILE:O	1:A:48:ILE:HG22	2.13	0.48
1:A:93:LEU:HD11	7:A:2003:HOH:O	2.13	0.48
1:A:53:ARG:NH1	1:A:53:ARG:HG2	2.29	0.48
1:A:139:SER:O	1:A:141:ASN:N	2.46	0.48
1:A:168:VAL:HG12	1:A:168:VAL:O	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:121:THR:HG22	2:B:122:THR:N	2.29	0.48
1:A:84:TYR:HB2	1:A:85:PRO:HD2	1.96	0.48
4:A:1183:D12:C12	2:B:99:PHE:CD2	2.98	0.47
2:B:134:SER:HB2	2:B:143:ALA:CB	2.44	0.47
1:A:43:ASN:O	1:A:45:TYR:N	2.47	0.47
1:A:45:TYR:N	1:A:45:TYR:CD1	2.82	0.47
1:A:139:SER:O	1:A:140:ALA:C	2.53	0.47
1:A:146:TYR:C	1:A:148:LYS:N	2.68	0.47
3:B:1175:AXT:H7	3:B:1175:AXT:H181	1.61	0.47
1:A:78:LYS:HD3	1:A:181:VAL:CG2	2.45	0.47
1:A:181:VAL:OXT	1:A:181:VAL:HG12	2.15	0.47
2:B:83:THR:HG23	2:B:84:LYS:N	2.30	0.47
2:B:45:ARG:HH22	2:B:169:GLU:HG3	1.76	0.46
1:A:62:GLN:HB2	1:A:84:TYR:CA	2.45	0.46
1:A:118:LEU:HD23	1:A:118:LEU:N	2.28	0.46
1:A:103:ALA:HB1	1:A:104:PRO:HD2	1.95	0.46
1:A:173:CYS:HB3	1:A:175:TYR:CE2	2.50	0.46
3:A:1182:AXT:H7	3:A:1182:AXT:H181	1.49	0.46
1:A:45:TYR:HE2	4:A:1183:D12:C3	2.28	0.46
2:B:22:ASP:O	2:B:23:LEU:C	2.53	0.46
1:A:65:ILE:HG22	1:A:66:LYS:N	2.31	0.46
1:A:56:TYR:CE1	1:A:65:ILE:HD13	2.51	0.46
2:B:125:TYR:HB2	7:B:2014:HOH:O	2.15	0.46
1:A:174:PRO:O	1:A:175:TYR:C	2.54	0.46
2:B:122:THR:HG22	2:B:123:ASP:H	1.80	0.46
1:A:157:ILE:HG13	1:A:159:VAL:HG23	1.96	0.46
1:A:52:VAL:N	1:A:169:GLN:HE22	2.14	0.45
2:B:3:ILE:O	2:B:4:PRO:C	2.54	0.45
2:B:109:ASP:HB3	2:B:113:TYR:CE2	2.50	0.45
1:A:128:TYR:N	1:A:128:TYR:CD1	2.83	0.45
1:A:35:TRP:CE3	1:A:136:PHE:HB3	2.51	0.45
1:A:123:ASP:O	2:B:122:THR:O	2.35	0.45
3:A:1182:AXT:H27	3:A:1182:AXT:H381	1.60	0.45
3:A:1182:AXT:H31	3:A:1182:AXT:H391	1.85	0.45
1:A:29:ASN:C	1:A:31:TYR:N	2.70	0.45
1:A:41:THR:O	1:A:42:ASN:C	2.54	0.45
2:B:121:THR:CG2	2:B:122:THR:N	2.80	0.45
2:B:29:ARG:HA	2:B:49:SER:O	2.17	0.45
1:A:150:CYS:O	1:A:153:ALA:N	2.45	0.45
2:B:113:TYR:CE2	2:B:146:LYS:HG2	2.52	0.45
1:A:58:PHE:CD2	1:A:63:PHE:CE2	3.05	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:ASP:O	1:A:113:SER:N	2.50	0.44
2:B:64:THR:HG22	2:B:65:ALA:N	2.31	0.44
2:B:16:ALA:O	2:B:106:ILE:HA	2.17	0.44
1:A:122:ILE:HG21	1:A:122:ILE:HD13	1.59	0.44
4:A:1183:D12:C8	4:A:1183:D12:H41	2.47	0.44
1:A:68:THR:CB	1:A:178:GLN:HE21	2.30	0.44
2:B:113:TYR:N	2:B:113:TYR:CD2	2.85	0.44
1:A:73:ASP:OD1	1:A:73:ASP:C	2.56	0.44
2:B:45:ARG:NH1	2:B:169:GLU:HB3	2.22	0.44
1:A:100:SER:O	2:B:1:ASP:N	2.50	0.43
2:B:34:HIS:CD2	2:B:144:VAL:HG22	2.53	0.43
2:B:49:SER:HB3	2:B:64:THR:HA	1.99	0.43
1:A:97:TYR:CZ	3:A:1182:AXT:H3	2.54	0.43
1:A:78:LYS:CD	1:A:181:VAL:HG22	2.46	0.43
2:B:146:LYS:O	2:B:149:ALA:HB3	2.19	0.43
2:B:49:SER:HB3	2:B:64:THR:OG1	2.19	0.43
2:B:167:THR:HB	2:B:168:ALA:H	1.56	0.43
1:A:170:GLY:O	1:A:173:CYS:HB2	2.19	0.43
1:A:100:SER:O	2:B:1:ASP:CA	2.66	0.43
2:B:40:TYR:HB3	2:B:98:VAL:HG22	2.00	0.43
1:A:83:LEU:HD11	1:A:105:LEU:HD23	2.01	0.43
1:A:143:ALA:HB3	1:A:146:TYR:CD1	2.54	0.43
1:A:17:ARG:NH1	1:A:92:HIS:ND1	2.66	0.43
2:B:68:ASN:ND2	2:B:72:GLU:HG3	2.32	0.42
1:A:53:ARG:HH11	1:A:53:ARG:CG	2.32	0.42
1:A:98:GLU:CD	2:B:1:ASP:HB3	2.39	0.42
1:A:40:LEU:CD1	1:A:40:LEU:O	2.48	0.42
1:A:132:PHE:CD2	3:A:1182:AXT:H203	2.54	0.42
2:B:93:ILE:HG21	2:B:93:ILE:HD13	1.78	0.42
2:B:23:LEU:HD21	2:B:82:PRO:HD3	2.00	0.42
2:B:15:VAL:HB	2:B:106:ILE:HG12	2.00	0.42
1:A:50:LYS:O	1:A:51:CYS:HB2	2.20	0.42
1:A:68:THR:HB	1:A:178:GLN:HB3	2.01	0.42
2:B:106:ILE:HG21	2:B:150:VAL:CG1	2.50	0.42
1:A:37:GLN:NE2	1:A:40:LEU:CD2	2.72	0.42
4:A:1183:D12:H122	2:B:99:PHE:HD2	1.83	0.42
2:B:3:ILE:HG22	2:B:4:PRO:O	2.19	0.42
1:A:62:GLN:HB2	1:A:84:TYR:HB3	2.01	0.42
2:B:18:GLN:HE21	2:B:108:THR:H	1.67	0.42
2:B:86:PHE:CB	2:B:87:PRO:HD2	2.47	0.41
1:A:20:LEU:HA	1:A:20:LEU:HD23	1.80	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:1182:AXT:C36	3:A:1182:AXT:C28	2.99	0.41
3:B:1175:AXT:H373	3:B:1175:AXT:H23	1.84	0.41
2:B:48:HIS:CE1	2:B:65:ALA:HB3	2.55	0.41
1:A:43:ASN:HA	1:A:44:PRO:HD2	1.85	0.41
1:A:20:LEU:HB3	1:A:24:GLN:NE2	2.35	0.41
1:A:111:ASP:OD1	1:A:114:ASN:HB2	2.20	0.41
1:A:126:PHE:CD1	3:A:1182:AXT:H23	2.56	0.41
2:B:25:ARG:HH12	5:B:1176:TRS:H31	1.84	0.41
1:A:150:CYS:C	1:A:152:ALA:N	2.73	0.41
1:A:25:THR:HB	1:A:26:PRO:HD3	2.03	0.41
1:A:43:ASN:C	1:A:43:ASN:OD1	2.60	0.40
2:B:3:ILE:HG23	2:B:4:PRO:HD2	2.04	0.40
2:B:7:VAL:HA	2:B:126:LYS:O	2.20	0.40
1:A:73:ASP:OD1	1:A:74:GLY:N	2.54	0.40
2:B:83:THR:CG2	2:B:84:LYS:N	2.81	0.40
2:B:173:ARG:O	2:B:174:ALA:C	2.59	0.40
1:A:178:GLN:O	1:A:180:THR:N	2.54	0.40
1:A:58:PHE:CD2	1:A:63:PHE:CD2	3.09	0.40
2:B:28:GLY:O	2:B:50:ASN:HA	2.21	0.40
1:A:37:GLN:HB3	1:A:52:VAL:HB	2.03	0.40
4:A:1183:D12:H122	2:B:99:PHE:CE2	2.55	0.40
2:B:78:PHE:CE2	3:B:1175:AXT:H3	2.57	0.40

All (2) 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)
7:B:2011:HOH:O	7:B:2011:HOH:O[2_665]	0.76	1.44
1:A:75:ASN:ND2	1:A:75:ASN:ND2[10_665]	2.14	0.06

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	178/180 (99%)	148 (83%)	21 (12%)	9 (5%)	2	19
2	B	172/174 (99%)	155 (90%)	15 (9%)	2 (1%)	16	58
All	All	350/354 (99%)	303 (87%)	36 (10%)	11 (3%)	5	32

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	140	ALA
1	A	71	ALA
1	A	144	ASP
2	B	141	GLY
1	A	179	LYS
1	A	44	PRO
1	A	32	ALA
1	A	42	ASN
1	A	112	TYR
2	B	96	PRO
1	A	52	VAL

### 5.3.2 Protein sidechains ⓘ

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	159/159 (100%)	137 (86%)	22 (14%)	4	20
2	B	150/150 (100%)	134 (89%)	16 (11%)	8	33
All	All	309/309 (100%)	271 (88%)	38 (12%)	6	27

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	14	SER
1	A	57	SER
1	A	62	GLN
1	A	70	ILE

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Mol	Chain	Res	Type
1	A	73	ASP
1	A	76	LEU
1	A	84	TYR
1	A	93	LEU
1	A	108	LEU
1	A	120	SER
1	A	125	ASN
1	A	129	HIS
1	A	133	SER
1	A	139	SER
1	A	144	ASP
1	A	161	THR
1	A	165	VAL
1	A	167	THR
1	A	168	VAL
1	A	171	SER
1	A	172	SER
2	B	1	ASP
2	B	29	ARG
2	B	37	GLU
2	B	38	ASN
2	B	77	ASP
2	B	79	LYS
2	B	83	THR
2	B	84	LYS
2	B	109	ASP
2	B	112	THR
2	B	122	THR
2	B	123	ASP
2	B	138	GLN
2	B	139	THR
2	B	165	SER
2	B	173	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	86	ASN
1	A	114	ASN
1	A	129	HIS
1	A	178	GLN
2	B	38	ASN

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Mol	Chain	Res	Type
2	B	68	ASN
2	B	70	ASN
2	B	90	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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

5 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
3	AXT	A	1182	-	43,45,45	2.05	9 (20%)	48,64,64	1.99	14 (29%)
4	D12	A	1183	-	11,11,11	0.48	0	10,10,10	0.72	0
3	AXT	B	1175	-	43,45,45	2.42	16 (37%)	48,64,64	1.73	14 (29%)
5	TRS	B	1176	-	7,7,7	5.92	7 (100%)	9,9,9	3.36	6 (66%)
6	EPE	B	1177	-	14,15,15	2.84	8 (57%)	18,20,20	3.72	8 (44%)

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	AXT	A	1182	-	-	0/29/75/75	0/2/2/2
4	D12	A	1183	-	-	0/9/9/9	0/0/0/0
3	AXT	B	1175	-	-	0/29/75/75	0/2/2/2
5	TRS	B	1176	-	-	0/9/9/9	0/0/0/0
6	EPE	B	1177	-	-	0/9/19/19	0/1/1/1

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1182	AXT	C32-C33	-4.14	1.36	1.45
3	B	1175	AXT	C4-C5	-3.94	1.40	1.47
3	B	1175	AXT	C16-C1	-3.86	1.45	1.53
3	A	1182	AXT	C8-C9	-3.77	1.37	1.45
3	B	1175	AXT	C22-C23	-3.38	1.46	1.52
3	A	1182	AXT	C28-C29	-3.29	1.38	1.45
3	B	1175	AXT	C8-C9	-3.01	1.39	1.45
3	A	1182	AXT	C21-C26	-2.99	1.49	1.53
3	A	1182	AXT	C24-C25	-2.94	1.42	1.47
3	B	1175	AXT	C12-C13	-2.61	1.40	1.45
3	B	1175	AXT	C28-C29	-2.21	1.41	1.45
3	B	1175	AXT	C36-C21	-2.08	1.49	1.53
3	B	1175	AXT	C31-C32	2.32	1.40	1.34
6	B	1177	EPE	C7-C8	2.61	1.62	1.51
3	B	1175	AXT	C1-C6	2.70	1.57	1.53
6	B	1177	EPE	C9-C10	2.87	1.62	1.52
6	B	1177	EPE	C2-N1	2.95	1.55	1.46
5	B	1176	TRS	O1-C1	2.99	1.52	1.42
3	B	1175	AXT	C28-C27	3.09	1.42	1.33
6	B	1177	EPE	O2S-S	3.11	1.55	1.45
3	A	1182	AXT	C10-C9	3.13	1.39	1.35
6	B	1177	EPE	C6-N1	3.38	1.56	1.46
3	B	1175	AXT	C14-C13	3.68	1.40	1.35
6	B	1177	EPE	C9-N1	3.81	1.56	1.47
3	B	1175	AXT	C30-C29	4.10	1.41	1.35
5	B	1176	TRS	O3-C3	4.32	1.56	1.42
3	A	1182	AXT	C26-C25	4.48	1.42	1.35
6	B	1177	EPE	O1S-S	4.50	1.59	1.45
3	A	1182	AXT	C14-C13	4.50	1.41	1.35
5	B	1176	TRS	C-N	4.56	1.57	1.50
3	B	1175	AXT	C26-C25	4.56	1.42	1.35
3	B	1175	AXT	C34-C33	4.76	1.42	1.35
6	B	1177	EPE	C7-N4	4.93	1.59	1.47
5	B	1176	TRS	O2-C2	5.00	1.58	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1182	AXT	C34-C33	5.01	1.42	1.35
3	B	1175	AXT	C6-C5	5.08	1.42	1.35
5	B	1176	TRS	C1-C	5.41	1.63	1.53
5	B	1176	TRS	C3-C	5.45	1.63	1.53
3	B	1175	AXT	C10-C9	5.83	1.43	1.35
5	B	1176	TRS	C2-C	10.63	1.73	1.53

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1176	TRS	O1-C1-C	-5.75	99.54	111.18
3	A	1182	AXT	C18-C5-C6	-5.05	116.94	123.87
3	A	1182	AXT	C27-C28-C29	-4.78	118.93	126.22
3	A	1182	AXT	C30-C31-C32	-3.64	112.02	123.13
5	B	1176	TRS	C1-C-N	-3.37	101.95	108.09
3	A	1182	AXT	C31-C30-C29	-3.37	122.33	127.20
3	B	1175	AXT	C38-C25-C26	-3.17	119.51	123.87
6	B	1177	EPE	C3-C2-N1	-3.12	105.05	110.63
5	B	1176	TRS	C2-C-C1	-2.95	104.40	110.78
3	A	1182	AXT	C10-C11-C12	-2.84	114.48	123.13
3	A	1182	AXT	C38-C25-C26	-2.80	120.02	123.87
6	B	1177	EPE	C7-N4-C5	-2.70	104.34	111.27
3	B	1175	AXT	C15-C35-C34	-2.67	117.50	123.39
3	B	1175	AXT	C18-C5-C6	-2.63	120.26	123.87
3	B	1175	AXT	C15-C14-C13	-2.62	123.41	127.20
3	A	1182	AXT	C32-C33-C34	-2.52	114.92	118.98
3	B	1175	AXT	C35-C34-C33	-2.47	123.63	127.20
3	A	1182	AXT	C15-C35-C34	-2.47	117.93	123.39
3	B	1175	AXT	O3-C3-C2	-2.45	105.72	110.33
6	B	1177	EPE	C9-N1-C6	-2.43	105.03	111.27
3	A	1182	AXT	C7-C8-C9	-2.40	122.56	126.22
3	B	1175	AXT	C10-C11-C12	-2.26	116.23	123.13
3	B	1175	AXT	C31-C30-C29	-2.26	123.94	127.20
3	B	1175	AXT	C8-C9-C10	-2.24	115.38	118.98
6	B	1177	EPE	O2S-S-O1S	-2.16	105.59	113.48
6	B	1177	EPE	C7-N4-C3	2.03	116.48	111.27
6	B	1177	EPE	C9-N1-C2	2.03	116.48	111.27
3	B	1175	AXT	C20-C13-C12	2.11	121.61	118.10
3	B	1175	AXT	C22-C21-C26	2.40	114.31	110.49
5	B	1176	TRS	O3-C3-C	2.42	116.08	111.18
3	B	1175	AXT	C16-C1-C6	2.51	114.24	110.30
3	A	1182	AXT	C36-C21-C26	2.73	114.59	110.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1182	AXT	C21-C26-C27	2.75	123.52	115.82
3	B	1175	AXT	C38-C25-C24	2.85	120.78	115.69
3	A	1182	AXT	C18-C5-C4	2.92	120.90	115.69
3	A	1182	AXT	C20-C13-C12	2.92	122.96	118.10
3	A	1182	AXT	C1-C6-C7	3.05	124.36	115.82
6	B	1177	EPE	O2S-S-C10	3.22	109.66	106.91
3	B	1175	AXT	C1-C6-C7	3.67	126.10	115.82
5	B	1176	TRS	C3-C-N	4.39	116.07	108.09
5	B	1176	TRS	C2-C-N	4.52	116.31	108.09
6	B	1177	EPE	O1S-S-C10	13.59	118.50	106.91

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 37 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1182	AXT	16	0
4	A	1183	D12	10	0
3	B	1175	AXT	8	0
5	B	1176	TRS	3	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, 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	180/180 (100%)	-0.04	0 100 100	20, 44, 70, 78	0
2	B	174/174 (100%)	-0.06	0 100 100	13, 31, 64, 74	0
All	All	354/354 (100%)	-0.05	0 100 100	13, 39, 67, 78	0

There are no RSRZ outliers to report.

### 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, 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
5	TRS	B	1176	8/8	0.69	0.61	7.98	69,70,71,71	0
6	EPE	B	1177	15/15	0.83	0.36	5.32	87,89,97,97	0
4	D12	A	1183	12/12	0.81	0.46	4.18	78,80,82,82	0
3	AXT	B	1175	44/44	0.97	0.31	2.11	14,22,26,29	0
3	AXT	A	1182	44/44	0.97	0.31	2.01	25,31,38,39	0



## 6.5 Other polymers [i](#)

There are no such residues in this entry.