



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 08:22 PM UTC

PDB ID : 4KC7 / pdb_00004kc7
Title : Crystal Structure of Endo-1,5-alpha-L-arabinanase from Thermotoga petrophila RKU-1
Authors : Nascimento, A.F.Z.; Polo, C.C.; Santos, C.R.; Costa, M.C.M.F.; Mesa, A.N.; Prade, R.A.; Ruller, R.; Squina, F.M.; Murakami, M.T.
Deposited on : 2013-04-24
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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

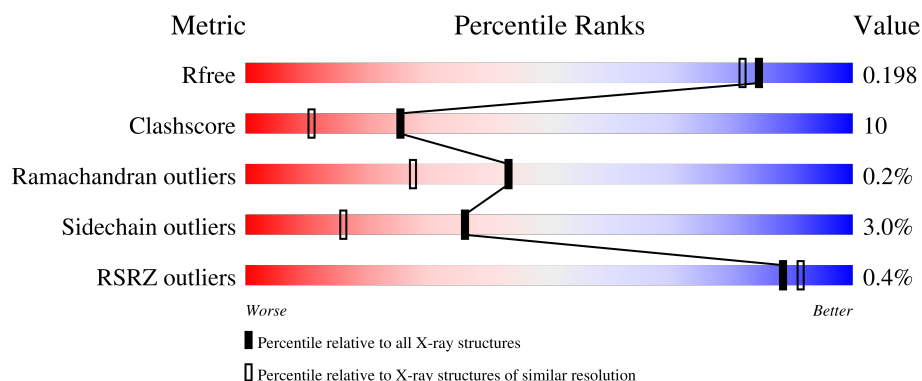
1 Overall quality at a glance

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}	180053	3183 (1.76-1.76)
Clashscore	190562	3299 (1.76-1.76)
Ramachandran outliers	187476	3274 (1.76-1.76)
Sidechain outliers	187428	3274 (1.76-1.76)
RSRZ outliers	180081	3183 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	474	
1	B	474	
1	C	474	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11493 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 Glycoside hydrolase, family 43.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	449	Total	C	N	O	S	0	6	0
			3623	2337	602	674	10			
1	B	449	Total	C	N	O	S	0	7	0
			3632	2345	605	672	10			
1	C	433	Total	C	N	O	S	0	3	0
			3476	2244	582	640	10			

There are 69 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	MET	-	initiating methionine	UNP A5IKD4
A	-1	GLY	-	expression tag	UNP A5IKD4
A	0	SER	-	expression tag	UNP A5IKD4
A	1	SER	-	expression tag	UNP A5IKD4
A	2	HIS	-	expression tag	UNP A5IKD4
A	3	HIS	-	expression tag	UNP A5IKD4
A	4	HIS	-	expression tag	UNP A5IKD4
A	5	HIS	-	expression tag	UNP A5IKD4
A	6	HIS	-	expression tag	UNP A5IKD4
A	7	HIS	-	expression tag	UNP A5IKD4
A	8	SER	-	expression tag	UNP A5IKD4
A	9	SER	-	expression tag	UNP A5IKD4
A	10	GLY	-	expression tag	UNP A5IKD4
A	11	LEU	-	expression tag	UNP A5IKD4
A	12	VAL	-	expression tag	UNP A5IKD4
A	13	PRO	-	expression tag	UNP A5IKD4
A	14	ARG	-	expression tag	UNP A5IKD4
A	15	GLY	-	expression tag	UNP A5IKD4
A	16	SER	-	expression tag	UNP A5IKD4
A	17	HIS	-	expression tag	UNP A5IKD4
A	18	MET	-	expression tag	UNP A5IKD4
A	19	ALA	-	expression tag	UNP A5IKD4
A	20	SER	-	expression tag	UNP A5IKD4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	MET	-	initiating methionine	UNP A5IKD4
B	-1	GLY	-	expression tag	UNP A5IKD4
B	0	SER	-	expression tag	UNP A5IKD4
B	1	SER	-	expression tag	UNP A5IKD4
B	2	HIS	-	expression tag	UNP A5IKD4
B	3	HIS	-	expression tag	UNP A5IKD4
B	4	HIS	-	expression tag	UNP A5IKD4
B	5	HIS	-	expression tag	UNP A5IKD4
B	6	HIS	-	expression tag	UNP A5IKD4
B	7	HIS	-	expression tag	UNP A5IKD4
B	8	SER	-	expression tag	UNP A5IKD4
B	9	SER	-	expression tag	UNP A5IKD4
B	10	GLY	-	expression tag	UNP A5IKD4
B	11	LEU	-	expression tag	UNP A5IKD4
B	12	VAL	-	expression tag	UNP A5IKD4
B	13	PRO	-	expression tag	UNP A5IKD4
B	14	ARG	-	expression tag	UNP A5IKD4
B	15	GLY	-	expression tag	UNP A5IKD4
B	16	SER	-	expression tag	UNP A5IKD4
B	17	HIS	-	expression tag	UNP A5IKD4
B	18	MET	-	expression tag	UNP A5IKD4
B	19	ALA	-	expression tag	UNP A5IKD4
B	20	SER	-	expression tag	UNP A5IKD4
C	-2	MET	-	initiating methionine	UNP A5IKD4
C	-1	GLY	-	expression tag	UNP A5IKD4
C	0	SER	-	expression tag	UNP A5IKD4
C	1	SER	-	expression tag	UNP A5IKD4
C	2	HIS	-	expression tag	UNP A5IKD4
C	3	HIS	-	expression tag	UNP A5IKD4
C	4	HIS	-	expression tag	UNP A5IKD4
C	5	HIS	-	expression tag	UNP A5IKD4
C	6	HIS	-	expression tag	UNP A5IKD4
C	7	HIS	-	expression tag	UNP A5IKD4
C	8	SER	-	expression tag	UNP A5IKD4
C	9	SER	-	expression tag	UNP A5IKD4
C	10	GLY	-	expression tag	UNP A5IKD4
C	11	LEU	-	expression tag	UNP A5IKD4
C	12	VAL	-	expression tag	UNP A5IKD4
C	13	PRO	-	expression tag	UNP A5IKD4
C	14	ARG	-	expression tag	UNP A5IKD4
C	15	GLY	-	expression tag	UNP A5IKD4
C	16	SER	-	expression tag	UNP A5IKD4

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Chain	Residue	Modelled	Actual	Comment	Reference
C	17	HIS	-	expression tag	UNP A5IKD4
C	18	MET	-	expression tag	UNP A5IKD4
C	19	ALA	-	expression tag	UNP A5IKD4
C	20	SER	-	expression tag	UNP A5IKD4

- Molecule 2 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			7	4	3		
2	B	1	Total	C	O	0	0
			7	4	3		
2	B	1	Total	C	O	0	0
			7	4	3		
2	C	1	Total	C	O	0	0
			7	4	3		

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		
3	B	1	Total	Ca	0	0
			1	1		
3	C	1	Total	Ca	0	0
			1	1		

- Molecule 4 is water.

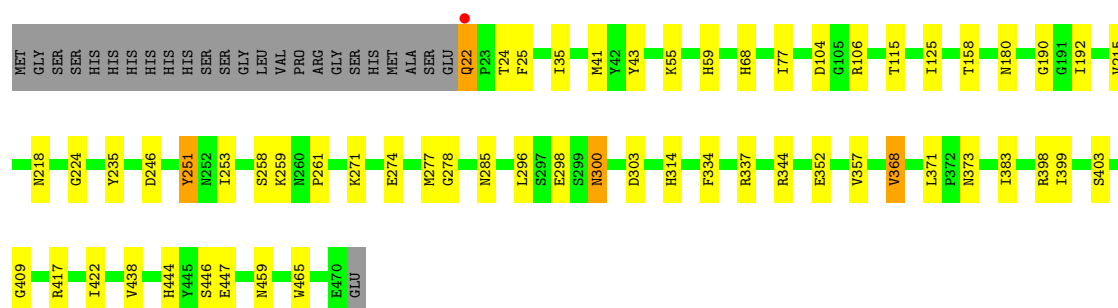
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	355	Total 355	O 355	0	0
4	B	176	Total 176	O 176	0	0
4	C	200	Total 200	O 200	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

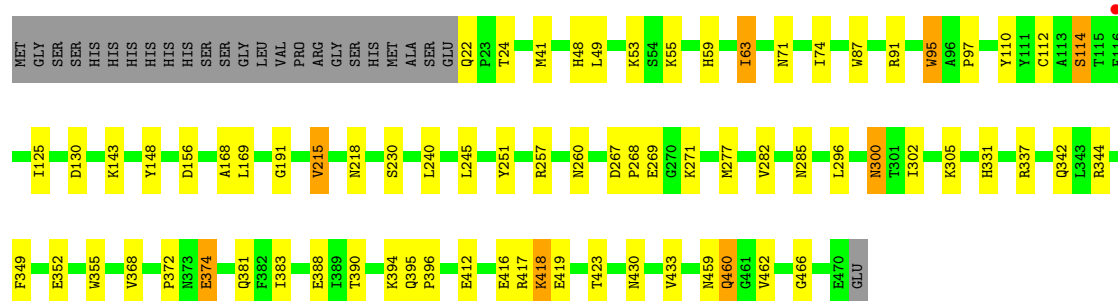
• Molecule 1: Glycoside hydrolase, family 43

Chain A: 



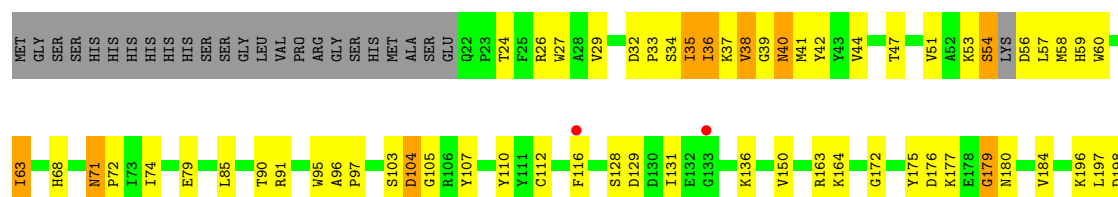
• Molecule 1: Glycoside hydrolase, family 43

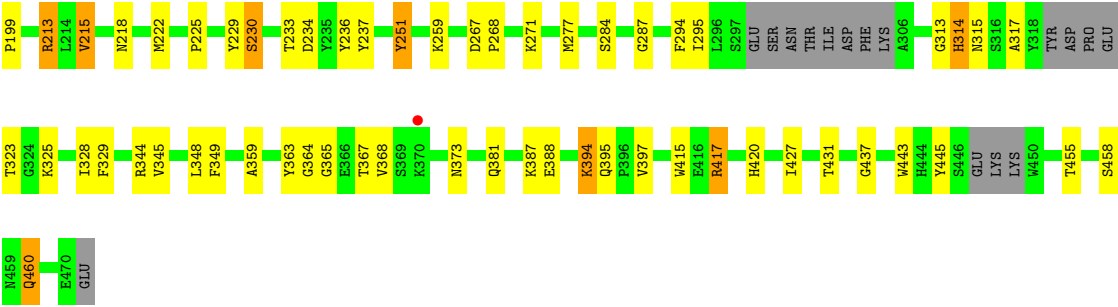
Chain B: 



• Molecule 1: Glycoside hydrolase, family 43

Chain C: 





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	41.67Å 86.91Å 194.85Å 90.00° 90.15° 90.00°	Depositor
Resolution (Å)	40.77 – 1.75 40.77 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.1 (40.77-1.75) 99.1 (40.77-1.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.49 (at 1.75Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.160 , 0.197 0.168 , 0.198	Depositor DCC
R_{free} test set	7010 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	20.1	Xtriage
Anisotropy	0.534	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 34.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.398 for h,-k,-l	Xtriage
Reported twinning fraction	0.612 for H, K, L 0.388 for -h,-k,l	Depositor
Outliers	0 of 139453 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	11493	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.27% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	1.40	17/3766 (0.5%)	1.30	11/5117 (0.2%)
1	B	1.22	8/3780 (0.2%)	1.22	17/5134 (0.3%)
1	C	1.11	1/3598 (0.0%)	1.19	10/4885 (0.2%)
All	All	1.25	26/11144 (0.2%)	1.24	38/15136 (0.3%)

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	303	ASP	C-O	7.58	1.31	1.23
1	A	278	GLY	C-O	-6.81	1.19	1.24
1	A	409	GLY	N-CA	6.79	1.52	1.45
1	A	115	THR	C-O	-6.77	1.15	1.23
1	A	251	TYR	C-O	6.67	1.31	1.23
1	B	423	THR	C-O	6.67	1.31	1.24
1	B	230	SER	C-O	-6.52	1.17	1.24
1	A	158	THR	C-O	6.22	1.30	1.24
1	A	344	ARG	NE-CZ	-6.08	1.26	1.33
1	A	125	ILE	C-O	-6.00	1.17	1.24
1	A	344	ARG	CZ-NH2	5.61	1.40	1.33
1	B	74	ILE	CA-C	5.46	1.58	1.52
1	B	396	PRO	N-CA	5.44	1.53	1.46
1	A	253	ILE	C-O	-5.34	1.18	1.24
1	A	224	GLY	N-CA	-5.31	1.37	1.44
1	C	458	SER	N-CA	5.28	1.52	1.45
1	B	95	TRP	CA-C	5.27	1.58	1.52
1	B	412	GLU	CA-C	-5.19	1.46	1.52
1	B	168	ALA	CA-C	-5.19	1.47	1.53
1	A	258	SER	N-CA	5.18	1.52	1.45
1	B	125	ILE	CA-CB	-5.17	1.48	1.54
1	A	261	PRO	C-O	5.12	1.30	1.24
1	A	192	ILE	C-O	5.10	1.29	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	334	PHE	CA-CB	-5.05	1.47	1.54
1	A	422	ILE	CA-C	-5.03	1.46	1.52
1	A	465	TRP	C-O	-5.01	1.17	1.24

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	35	ILE	CB-CA-C	-10.21	97.22	111.21
1	A	371	LEU	CA-C-N	-9.85	110.28	120.03
1	A	371	LEU	C-N-CA	-9.85	110.28	120.03
1	C	71	ASN	CA-C-N	9.06	129.25	119.28
1	C	71	ASN	C-N-CA	9.06	129.25	119.28
1	B	395	GLN	CA-C-N	-8.01	112.46	120.31
1	B	395	GLN	C-N-CA	-8.01	112.46	120.31
1	B	230	SER	CA-C-N	-7.75	111.73	119.56
1	B	230	SER	C-N-CA	-7.75	111.73	119.56
1	B	230	SER	CB-CA-C	-7.68	101.64	110.17
1	A	22	GLN	CA-C-N	-7.36	112.42	119.85
1	A	22	GLN	C-N-CA	-7.36	112.42	119.85
1	B	22	GLN	CA-C-N	7.35	127.40	119.90
1	B	22	GLN	C-N-CA	7.35	127.40	119.90
1	C	96	ALA	CA-C-N	-7.05	112.97	120.66
1	C	96	ALA	C-N-CA	-7.05	112.97	120.66
1	C	388	GLU	N-CA-C	6.41	119.20	110.55
1	A	77	ILE	N-CA-C	6.20	117.69	110.62
1	C	315	ASN	N-CA-C	6.09	119.09	110.14
1	A	403	SER	N-CA-C	5.92	119.32	111.75
1	A	373	ASN	N-CA-C	5.74	118.27	111.33
1	B	71	ASN	CA-C-N	5.70	125.40	119.64
1	B	71	ASN	C-N-CA	5.70	125.40	119.64
1	B	267	ASP	CA-C-N	-5.60	113.91	119.56
1	B	267	ASP	C-N-CA	-5.60	113.91	119.56
1	C	36	ILE	CB-CA-C	-5.53	100.91	110.65
1	C	397	VAL	N-CA-C	-5.51	100.67	108.65
1	B	245	LEU	N-CA-C	5.41	119.13	112.54
1	A	25	PHE	N-CA-C	5.29	117.43	109.23
1	A	368	VAL	CB-CA-C	-5.27	102.73	111.32
1	B	114	SER	N-CA-C	5.26	115.48	108.38
1	B	257	ARG	N-CA-C	5.26	117.06	109.07
1	B	282	VAL	N-CA-C	5.23	115.45	110.53
1	A	190	GLY	N-CA-C	5.22	120.42	114.67
1	B	260	ASN	N-CA-C	5.22	117.17	110.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	445	TYR	N-CA-C	5.15	116.58	111.07
1	A	180	ASN	N-CA-C	-5.09	103.68	110.55
1	B	419	GLU	CB-CA-C	-5.02	110.80	116.63

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	3623	0	3479	30	0
1	B	3632	0	3501	46	0
1	C	3476	0	3335	126	2
2	A	7	0	10	0	0
2	B	14	0	20	3	0
2	C	7	0	10	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	A	355	0	0	12	2
4	B	176	0	0	9	0
4	C	200	0	0	46	0
All	All	11493	0	10355	203	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 10.

All (203) 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:C:58:MET:HB3	4:C:779:HOH:O	1.29	1.25
1:B:459:ASN:HB3	4:B:707:HOH:O	1.39	1.23
1:C:136:LYS:HB3	4:C:621:HOH:O	1.51	1.08
1:C:38:VAL:HG21	1:C:107:TYR:OH	1.55	1.05
1:C:175:TYR:CD1	4:C:789:HOH:O	2.12	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:39:GLY:O	1:C:41:MET:N	1.96	0.98
1:A:459:ASN:ND2	4:A:866:HOH:O	1.97	0.96
1:C:373[A]:ASN:OD1	1:C:417:ARG:NH2	2.01	0.93
1:C:175:TYR:HD1	4:C:789:HOH:O	1.46	0.92
1:A:274:GLU:OE2	4:A:788:HOH:O	1.88	0.91
1:C:229:TYR:OH	1:C:234:ASP:OD1	1.87	0.91
1:C:345:VAL:HG21	4:C:677:HOH:O	1.72	0.90
1:C:38:VAL:HG13	1:C:39:GLY:CA	2.03	0.88
1:C:51:VAL:HB	1:C:63:ILE:HG22	1.55	0.88
1:B:433:VAL:HG11	2:B:502:PEG:H41	1.58	0.85
1:A:300:ASN:HB3	4:A:863:HOH:O	1.76	0.85
1:C:230:SER:OG	1:C:233:THR:HG22	1.79	0.83
1:B:24[A]:THR:HG23	1:B:59:HIS:NE2	1.96	0.81
1:B:416:GLU:OE2	4:B:732:HOH:O	1.99	0.79
1:B:352:GLU:OE2	1:B:418:LYS:NZ	2.15	0.78
1:C:38:VAL:HG13	1:C:39:GLY:HA3	1.65	0.78
1:B:296:LEU:CD2	1:B:383:ILE:HD11	2.16	0.75
1:C:395:GLN:HA	4:C:759:HOH:O	1.85	0.75
1:B:218:ASN:OD1	1:B:277:MET:HE1	1.86	0.75
1:C:236:TYR:CE2	4:C:790:HOH:O	2.41	0.73
1:C:417:ARG:HD3	1:C:417:ARG:C	2.12	0.73
1:C:131:ILE:HG22	4:C:714:HOH:O	1.88	0.73
1:C:215:VAL:HG22	1:C:222:MET:SD	2.29	0.71
1:C:27:TRP:CD1	4:C:666:HOH:O	2.42	0.71
1:A:296:LEU:CD2	1:A:383:ILE:HD11	2.21	0.71
1:C:179:GLY:O	4:C:789:HOH:O	2.08	0.71
1:C:36:ILE:HG22	4:C:672:HOH:O	1.91	0.70
1:C:363:TYR:CD2	4:C:780:HOH:O	2.43	0.70
1:A:235:TYR:CE1	1:A:259:LYS:HE2	2.27	0.70
1:C:51:VAL:HB	1:C:63:ILE:CG2	2.21	0.70
1:C:415:TRP:HZ3	1:C:417:ARG:HG3	1.56	0.70
1:C:177:LYS:HG2	4:C:788:HOH:O	1.91	0.70
1:C:36:ILE:HD11	4:C:778:HOH:O	1.92	0.69
1:C:38:VAL:HG13	1:C:39:GLY:HA2	1.74	0.69
1:B:459:ASN:CB	4:B:707:HOH:O	2.15	0.68
1:C:27:TRP:NE1	4:C:666:HOH:O	2.27	0.68
1:B:271:LYS:HE3	1:B:285:ASN:O	1.93	0.67
1:C:230:SER:HG	1:C:233:THR:HG22	1.58	0.67
1:C:54:SER:HG	1:C:56:ASP:N	1.92	0.67
1:C:105:GLY:O	4:C:799:HOH:O	2.12	0.67
1:C:417:ARG:HD3	1:C:417:ARG:O	1.93	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:39:GLY:C	1:C:41:MET:N	2.52	0.66
1:C:38:VAL:CG1	1:C:39:GLY:HA3	2.25	0.66
1:B:342:GLN:OE1	1:B:344:ARG:NH1	2.22	0.64
1:C:60:TRP:HZ2	4:C:677:HOH:O	1.79	0.63
1:C:236:TYR:CD2	4:C:790:HOH:O	2.50	0.63
1:C:387:LYS:NZ	1:C:460:GLN:O	2.31	0.63
1:A:24[B]:THR:HG21	4:A:888:HOH:O	1.98	0.63
1:C:233:THR:HG21	1:C:237:TYR:OH	1.99	0.63
1:B:352:GLU:CD	1:B:418:LYS:HZ3	2.07	0.61
1:C:323:THR:HB	1:C:325:LYS:HE3	1.82	0.61
1:C:40:ASN:HB2	4:C:800:HOH:O	1.99	0.61
1:C:90:THR:O	1:C:91:ARG:HG2	2.01	0.61
1:B:296:LEU:HD23	1:B:383:ILE:HD11	1.82	0.60
1:A:235:TYR:CZ	1:A:259:LYS:HE2	2.36	0.60
1:A:246:ASP:OD1	4:A:859:HOH:O	2.17	0.59
1:C:90:THR:C	1:C:91:ARG:HG2	2.28	0.59
1:C:39:GLY:C	1:C:41:MET:H	2.11	0.58
1:C:287:GLY:O	1:C:387:LYS:HE3	2.03	0.58
1:C:344:ARG:NH2	4:C:666:HOH:O	2.36	0.58
1:C:58:MET:CB	4:C:779:HOH:O	2.10	0.57
1:A:41:MET:SD	1:A:55:LYS:HG2	2.44	0.57
1:A:68:HIS:HD2	4:A:813:HOH:O	1.88	0.56
1:B:191:GLY:HA2	4:B:691:HOH:O	2.06	0.55
1:C:218:ASN:OD1	1:C:277:MET:HE1	2.07	0.55
1:C:437:GLY:HA3	1:C:455:THR:O	2.05	0.55
1:C:349:PHE:CE1	1:C:359:ALA:HB2	2.42	0.55
1:C:349:PHE:HE1	1:C:359:ALA:HB2	1.72	0.55
1:B:388:GLU:HG2	1:B:390:THR:HG22	1.88	0.55
1:B:53:LYS:HE2	1:B:63:ILE:HD11	1.88	0.55
1:C:365:GLY:HA2	4:C:782:HOH:O	2.07	0.55
1:C:395:GLN:CA	4:C:759:HOH:O	2.47	0.55
1:C:294:PHE:HA	1:C:394:LYS:O	2.07	0.55
1:C:97:PRO:HA	1:C:110:TYR:O	2.07	0.54
1:A:296:LEU:HD23	1:A:383:ILE:HD11	1.89	0.54
1:C:417:ARG:C	1:C:417:ARG:CD	2.80	0.54
1:C:26:ARG:HB3	1:C:26:ARG:CZ	2.36	0.54
1:C:35:ILE:HG12	1:C:329:PHE:CD1	2.43	0.53
1:C:229:TYR:CZ	1:C:234:ASP:HA	2.43	0.53
1:C:417:ARG:O	1:C:417:ARG:CD	2.56	0.53
1:C:68:HIS:O	1:C:71:ASN:ND2	2.42	0.52
1:C:363:TYR:CG	4:C:780:HOH:O	2.61	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:175:TYR:CE1	4:C:789:HOH:O	2.53	0.52
1:C:236:TYR:HE2	4:C:790:HOH:O	1.85	0.52
1:C:95:TRP:HB2	1:C:112:CYS:SG	2.49	0.52
1:B:112:CYS:SG	1:B:169:LEU:HD22	2.50	0.52
1:C:345:VAL:HG11	4:C:677:HOH:O	2.10	0.52
1:C:24:THR:OG1	1:C:59:HIS:NE2	2.42	0.52
1:C:35:ILE:HD11	1:C:329:PHE:CD1	2.44	0.52
1:B:374[A]:GLU:H	1:B:374[A]:GLU:CD	2.17	0.52
1:C:116:PHE:HB3	4:C:675:HOH:O	2.09	0.52
1:C:196:LYS:NZ	4:C:731:HOH:O	2.42	0.52
1:A:444:HIS:CD2	1:A:447:GLU:H	2.28	0.52
1:C:36:ILE:CD1	4:C:778:HOH:O	2.55	0.51
1:C:42:TYR:O	1:C:53:LYS:HA	2.10	0.51
1:B:41:MET:SD	1:B:55:LYS:HG3	2.50	0.51
1:C:325:LYS:HE2	4:C:644:HOH:O	2.08	0.51
1:A:24[B]:THR:HG22	1:A:59:HIS:NE2	2.26	0.51
1:A:218:ASN:OD1	1:A:277:MET:HE1	2.10	0.51
1:A:337:ARG:HD2	4:A:707:HOH:O	2.10	0.51
1:A:337:ARG:NH1	4:A:877:HOH:O	2.44	0.51
1:C:34:SER:O	1:C:44:VAL:HA	2.10	0.50
1:C:176:ASP:CG	4:C:788:HOH:O	2.54	0.50
1:A:271:LYS:HE3	1:A:285:ASN:O	2.11	0.50
1:C:85:LEU:HD22	1:C:91:ARG:HA	1.94	0.50
1:C:443:TRP:HB3	4:C:783:HOH:O	2.11	0.50
1:A:444:HIS:HD2	1:A:446:SER:H	1.59	0.50
1:C:40:ASN:C	4:C:800:HOH:O	2.55	0.50
1:C:328:ILE:HB	1:C:348:LEU:HD11	1.94	0.49
1:B:296:LEU:HD21	1:B:383:ILE:HD11	1.92	0.49
1:C:42:TYR:CD2	1:C:57:LEU:HD21	2.47	0.49
1:C:180:ASN:C	4:C:789:HOH:O	2.56	0.49
1:C:136:LYS:CB	4:C:621:HOH:O	2.32	0.49
1:C:177:LYS:HD3	1:C:229:TYR:CE2	2.48	0.49
1:B:95:TRP:CZ3	1:B:114:SER:HB3	2.48	0.49
1:C:79:GLU:HG3	4:C:687:HOH:O	2.13	0.48
1:C:163:ARG:HE	1:C:218:ASN:ND2	2.10	0.48
1:B:388:GLU:CD	1:B:394:LYS:HZ1	2.22	0.47
1:C:427:ILE:HD12	1:C:427:ILE:N	2.28	0.47
1:B:148:TYR:C	1:B:148:TYR:CD1	2.93	0.47
1:C:284:SER:O	1:C:387:LYS:HG3	2.15	0.47
1:A:104:ASP:OD2	1:A:106:ARG:HD3	2.15	0.47
1:C:35:ILE:HG12	1:C:329:PHE:CE1	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:352:GLU:HG3	4:B:641:HOH:O	2.12	0.47
1:B:352:GLU:HG3	4:B:605:HOH:O	2.13	0.47
1:A:104:ASP:OD2	1:A:106:ARG:CD	2.63	0.47
1:B:48:HIS:NE2	2:B:501:PEG:H22	2.30	0.47
1:A:271:LYS:NZ	4:A:873:HOH:O	2.48	0.47
1:B:349:PHE:CE1	1:B:368:VAL:CG2	2.97	0.46
1:C:26:ARG:NH1	1:C:60:TRP:O	2.47	0.46
1:C:329:PHE:CD1	1:C:329:PHE:N	2.83	0.46
1:B:269:GLU:OE2	1:B:460:GLN:NE2	2.49	0.46
1:B:372:PRO:HB2	1:B:374[A]:GLU:HG2	1.98	0.46
1:C:42:TYR:CE2	1:C:57:LEU:HD21	2.50	0.46
1:C:236:TYR:HE1	1:C:259:LYS:O	1.99	0.46
1:C:234:ASP:O	1:C:234:ASP:CG	2.58	0.46
1:B:337[A]:ARG:HD2	4:B:710:HOH:O	2.15	0.46
1:C:128:SER:OG	1:C:129:ASP:N	2.50	0.45
1:A:298[B]:GLU:OE2	1:A:398:ARG:NH2	2.49	0.45
1:B:130:ASP:CG	4:B:760:HOH:O	2.59	0.45
1:C:198:ASP:O	1:C:199:PRO:C	2.58	0.45
1:C:367:THR:N	4:C:616:HOH:O	2.48	0.45
1:B:215:VAL:HG21	1:B:240:LEU:HD11	1.97	0.45
1:C:32:ASP:N	1:C:33:PRO:CD	2.80	0.45
1:C:40:ASN:CB	4:C:800:HOH:O	2.58	0.45
1:C:314:HIS:HD2	4:C:797:HOH:O	1.98	0.45
1:C:29:VAL:CG1	4:C:668:HOH:O	2.65	0.45
1:C:35:ILE:HG22	1:C:36:ILE:N	2.32	0.44
1:C:39:GLY:O	1:C:41:MET:CA	2.64	0.44
1:A:24[B]:THR:CG2	4:A:889:HOH:O	2.65	0.44
1:C:104:ASP:OD1	1:C:104:ASP:N	2.49	0.44
1:C:38:VAL:CG2	1:C:107:TYR:OH	2.45	0.44
1:C:229:TYR:HB2	4:C:790:HOH:O	2.17	0.44
1:C:32:ASP:OD2	1:C:47:THR:OG1	2.27	0.43
1:C:58:MET:CA	4:C:779:HOH:O	2.55	0.43
1:C:267:ASP:HB2	1:C:268:PRO:HD2	2.00	0.43
1:C:295:ILE:HG23	1:C:395:GLN:HG2	2.00	0.43
1:C:184:VAL:CG2	1:C:225:PRO:HB2	2.49	0.43
1:B:462:VAL:HB	2:B:502:PEG:H42	1.99	0.43
1:B:87:TRP:CD2	1:B:143:LYS:HD2	2.53	0.43
1:B:331:HIS:CD2	1:B:331:HIS:C	2.96	0.43
1:B:388:GLU:HG2	1:B:390:THR:CG2	2.48	0.43
1:A:314:HIS:CE1	4:A:840:HOH:O	2.72	0.42
1:C:176:ASP:OD1	1:C:176:ASP:C	2.62	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:97:PRO:HA	1:B:110:TYR:O	2.19	0.42
1:C:381:GLN:HG2	4:C:795:HOH:O	2.20	0.42
1:A:444:HIS:HE1	4:A:897:HOH:O	2.03	0.42
1:C:35:ILE:HD13	1:C:317:ALA:C	2.45	0.42
1:A:417:ARG:HD2	1:A:417:ARG:C	2.44	0.42
1:B:114:SER:HB2	1:B:169:LEU:HD21	2.02	0.42
1:C:71:ASN:HA	1:C:72:PRO:HD3	1.84	0.42
1:A:41:MET:SD	1:A:55:LYS:CG	3.08	0.41
1:B:268:PRO:HG3	1:B:355:TRP:CE2	2.55	0.41
1:C:35:ILE:CG1	1:C:329:PHE:CD1	3.04	0.41
1:B:300:ASN:OD1	1:B:302:ILE:O	2.37	0.41
1:C:365:GLY:CA	4:C:782:HOH:O	2.66	0.41
1:C:197:LEU:O	1:C:199:PRO:HD3	2.21	0.41
1:C:35:ILE:CD1	1:C:329:PHE:CD1	3.03	0.41
1:C:39:GLY:O	1:C:41:MET:CB	2.69	0.41
1:C:54:SER:CB	4:C:678:HOH:O	2.69	0.41
1:B:271:LYS:CE	1:B:285:ASN:O	2.65	0.41
1:C:38:VAL:H	1:C:39:GLY:HA3	1.85	0.41
1:C:172:GLY:N	1:C:184:VAL:O	2.45	0.41
1:C:251:TYR:CD2	1:C:313:GLY:HA3	2.56	0.41
1:A:35:ILE:HD12	1:A:43:TYR:O	2.21	0.41
1:C:26:ARG:HH11	1:C:26:ARG:CG	2.34	0.41
1:C:267:ASP:HB2	1:C:268:PRO:CD	2.51	0.41
1:C:58:MET:HE3	1:C:364:GLY:HA2	2.02	0.40
1:B:49:LEU:HD23	1:B:49:LEU:HA	1.93	0.40
1:B:156:ASP:OD1	1:B:156:ASP:C	2.63	0.40
1:A:357:VAL:HG21	1:A:438[B]:VAL:HG12	2.03	0.40
1:C:368:VAL:HG11	1:C:420:HIS:CG	2.56	0.40
1:A:399:ILE:HD13	1:A:399:ILE:HG21	1.91	0.40
1:B:381:GLN:O	1:B:466:GLY:HA2	2.21	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)
1:C:213:ARG:NH2	4:A:725:HOH:O[2_555]	1.23	0.97
1:C:213:ARG:CZ	4:A:725:HOH:O[2_555]	1.92	0.28

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	453/474 (96%)	436 (96%)	17 (4%)	0	100	100
1	B	454/474 (96%)	441 (97%)	13 (3%)	0	100	100
1	C	426/474 (90%)	407 (96%)	16 (4%)	3 (1%)	18	7
All	All	1333/1422 (94%)	1284 (96%)	46 (4%)	3 (0%)	43	27

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	40	ASN
1	C	314	HIS
1	C	179	GLY

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	391/406 (96%)	385 (98%)	6 (2%)	57	41
1	B	392/406 (97%)	381 (97%)	11 (3%)	38	18
1	C	372/406 (92%)	354 (95%)	18 (5%)	23	6
All	All	1155/1218 (95%)	1120 (97%)	35 (3%)	36	16

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	215	VAL
1	A	251	TYR
1	A	300	ASN
1	A	352	GLU
1	A	368	VAL
1	B	63	ILE
1	B	91	ARG
1	B	215	VAL
1	B	251	TYR
1	B	300	ASN
1	B	305	LYS
1	B	374[A]	GLU
1	B	374[B]	GLU
1	B	418	LYS
1	B	430	ASN
1	B	460	GLN
1	C	37	LYS
1	C	38	VAL
1	C	54	SER
1	C	63	ILE
1	C	74	ILE
1	C	103	SER
1	C	104	ASP
1	C	150	VAL
1	C	164	LYS
1	C	213	ARG
1	C	215	VAL
1	C	230	SER
1	C	251	TYR
1	C	271	LYS
1	C	394	LYS
1	C	417	ARG
1	C	431	THR
1	C	460	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	HIS
1	A	300	ASN
1	A	381	GLN
1	A	444	HIS

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Mol	Chain	Res	Type
1	A	459	ASN
1	B	68	HIS
1	B	76	ASN
1	B	167	ASN
1	B	275	ASN
1	B	381	GLN
1	B	395	GLN
1	B	402	ASN
1	B	420	HIS
1	B	444	HIS
1	C	76	ASN
1	C	167	ASN
1	C	200	ASN
1	C	275	ASN
1	C	381	GLN
1	C	430	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 3 are monoatomic - leaving 4 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
2	PEG	B	501	-	6,6,6	0.44	0	5,5,5	1.46	1 (20%)
2	PEG	B	502	-	6,6,6	0.61	0	5,5,5	0.72	0
2	PEG	A	501	-	6,6,6	0.58	0	5,5,5	0.81	0
2	PEG	C	501	-	6,6,6	0.48	0	5,5,5	1.00	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
2	PEG	B	501	-	-	1/4/4/4	-
2	PEG	B	502	-	-	3/4/4/4	-
2	PEG	A	501	-	-	0/4/4/4	-
2	PEG	C	501	-	-	2/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	PEG	O2-C2-C1	2.40	120.71	110.11

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	502	PEG	O1-C1-C2-O2
2	C	501	PEG	C1-C2-O2-C3
2	B	502	PEG	C1-C2-O2-C3
2	B	501	PEG	C1-C2-O2-C3
2	B	502	PEG	O2-C3-C4-O4
2	C	501	PEG	C4-C3-O2-C2

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	PEG	1	0
2	B	502	PEG	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	449/474 (94%)	-1.14	1 (0%) 91 93	8, 16, 28, 54	8 (1%)
1	B	449/474 (94%)	-0.80	1 (0%) 91 93	14, 25, 38, 54	8 (1%)
1	C	433/474 (91%)	-0.25	3 (0%) 84 88	18, 34, 51, 63	8 (1%)
All	All	1331/1422 (93%)	-0.73	5 (0%) 88 91	8, 25, 46, 63	24 (1%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	116	PHE	3.7
1	C	133	GLY	3.1
1	A	22	GLN	3.0
1	C	116	PHE	2.9
1	C	370	LYS	2.8

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PEG	B	502	7/7	0.96	0.07	31,36,40,46	0
2	PEG	B	501	7/7	0.97	0.05	33,35,43,48	0
2	PEG	C	501	7/7	0.97	0.06	40,41,42,42	0
3	CA	A	502	1/1	0.97	0.10	39,39,39,39	0
3	CA	B	503	1/1	0.97	0.13	39,39,39,39	0
2	PEG	A	501	7/7	0.98	0.05	28,29,34,39	0
3	CA	C	502	1/1	0.98	0.06	41,41,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.