



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 12:42 AM UTC

PDB ID : 5WMS / pdb_00005wms
Title : Phosphotriesterase variant S7
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Deposited on : 2017-07-31
Resolution : 1.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The 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

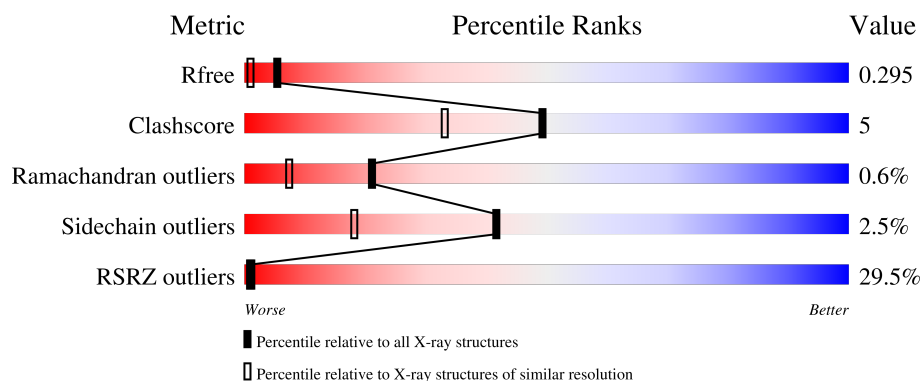
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (1.60-1.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	333	18% 83% 11% 5%
1	G	333	25% 80% 13% 5%
1	Q	333	26% 81% 13% 5%
1	S	333	42% 82% 11% 6%

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	CAC	A	2403	-	X	-	-
3	CAC	G	2403	-	X	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10306 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 Phosphotriesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	315	Total	C	N	O	S	0	5	0
			2434	1537	436	455	6			
1	G	315	Total	C	N	O	S	0	6	0
			2444	1545	435	458	6			
1	Q	316	Total	C	N	O	S	0	4	0
			2439	1539	442	452	6			
1	S	314	Total	C	N	O	S	0	4	0
			2423	1528	435	454	6			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	216	LEU	PHE	conflict	UNP A0A060GYS1
A	233	ALA	ASP	conflict	UNP A0A060GYS1
A	234	ALA	THR	conflict	UNP A0A060GYS1
A	254	SER	HIS	conflict	UNP A0A060GYS1
A	260	VAL	ILE	conflict	UNP A0A060GYS1
A	263	GLY	GLU	conflict	UNP A0A060GYS1
A	271	HIS	LEU	conflict	UNP A0A060GYS1
A	293	THR	MET	conflict	UNP A0A060GYS1
A	306	ILE	PHE	conflict	UNP A0A060GYS1
A	320	GLY	VAL	conflict	UNP A0A060GYS1
G	216	LEU	PHE	conflict	UNP A0A060GYS1
G	233	ALA	ASP	conflict	UNP A0A060GYS1
G	234	ALA	THR	conflict	UNP A0A060GYS1
G	254	SER	HIS	conflict	UNP A0A060GYS1
G	260	VAL	ILE	conflict	UNP A0A060GYS1
G	263	GLY	GLU	conflict	UNP A0A060GYS1
G	271	HIS	LEU	conflict	UNP A0A060GYS1
G	293	THR	MET	conflict	UNP A0A060GYS1
G	306	ILE	PHE	conflict	UNP A0A060GYS1
G	320	GLY	VAL	conflict	UNP A0A060GYS1
Q	216	LEU	PHE	conflict	UNP A0A060GYS1

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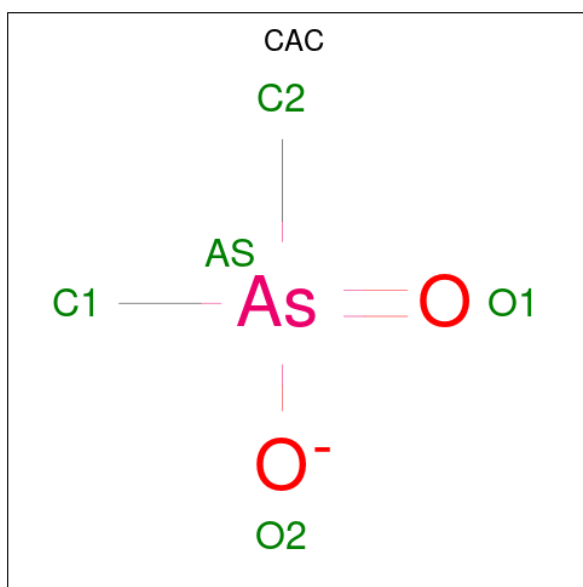
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Chain	Residue	Modelled	Actual	Comment	Reference
Q	233	ALA	ASP	conflict	UNP A0A060GYS1
Q	234	ALA	THR	conflict	UNP A0A060GYS1
Q	254	SER	HIS	conflict	UNP A0A060GYS1
Q	260	VAL	ILE	conflict	UNP A0A060GYS1
Q	263	GLY	GLU	conflict	UNP A0A060GYS1
Q	271	HIS	LEU	conflict	UNP A0A060GYS1
Q	293	THR	MET	conflict	UNP A0A060GYS1
Q	306	ILE	PHE	conflict	UNP A0A060GYS1
Q	320	GLY	VAL	conflict	UNP A0A060GYS1
S	216	LEU	PHE	conflict	UNP A0A060GYS1
S	233	ALA	ASP	conflict	UNP A0A060GYS1
S	234	ALA	THR	conflict	UNP A0A060GYS1
S	254	SER	HIS	conflict	UNP A0A060GYS1
S	260	VAL	ILE	conflict	UNP A0A060GYS1
S	263	GLY	GLU	conflict	UNP A0A060GYS1
S	271	HIS	LEU	conflict	UNP A0A060GYS1
S	293	THR	MET	conflict	UNP A0A060GYS1
S	306	ILE	PHE	conflict	UNP A0A060GYS1
S	320	GLY	VAL	conflict	UNP A0A060GYS1

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

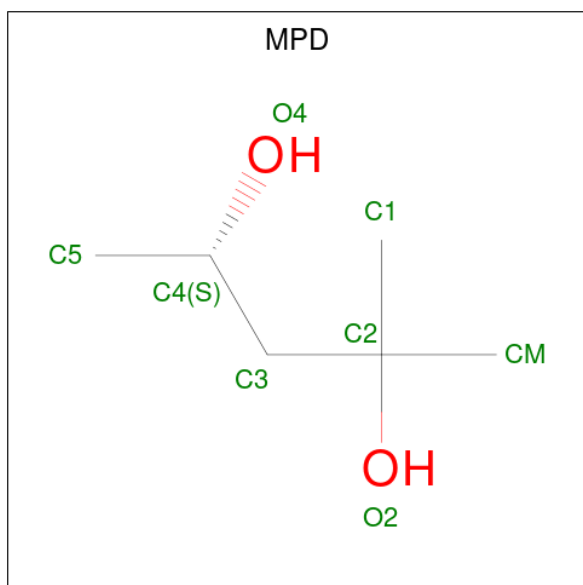
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Zn 2 2	0	0
2	G	2	Total Zn 2 2	0	0
2	Q	2	Total Zn 2 2	0	0
2	S	2	Total Zn 2 2	0	0

- Molecule 3 is CACODYLATE ION (CCD ID: CAC) (formula: C₂H₆AsO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	As	C	O	0	0
			5	1	2	2		
3	G	1	Total	As	C	O	0	0
			5	1	2	2		
3	Q	1	Total	As	C	O	0	0
			5	1	2	2		
3	S	1	Total	As	C	O	0	0
			5	1	2	2		

- Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 8 6 2	0	0
4	A	1	Total C O 8 6 2	0	0
4	G	1	Total C O 8 6 2	0	0
4	G	1	Total C O 8 6 2	0	0
4	Q	1	Total C O 8 6 2	0	0
4	Q	1	Total C O 8 6 2	0	0
4	Q	1	Total C O 8 6 2	0	0
4	S	1	Total C O 8 6 2	0	0
4	S	1	Total C O 8 6 2	0	0

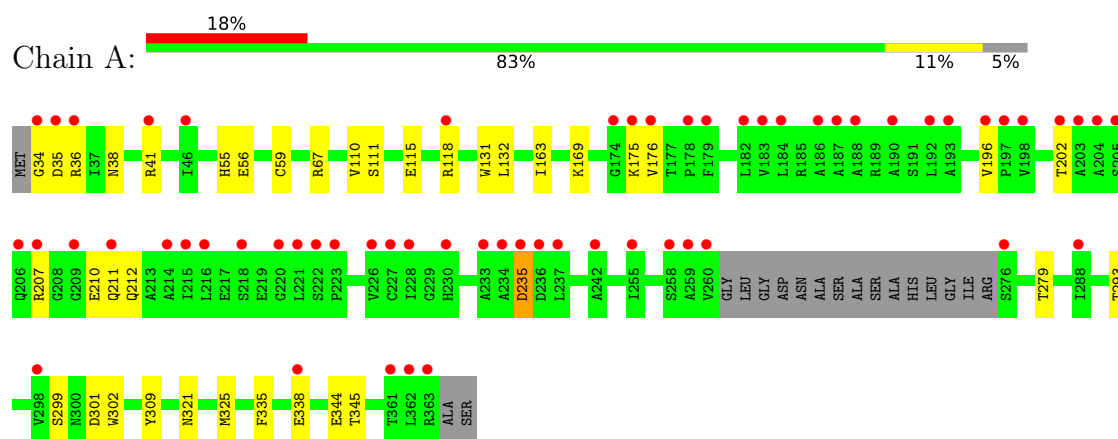
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	131	Total O 131 131	0	0
5	G	107	Total O 107 107	0	0
5	Q	131	Total O 131 131	0	0
5	S	97	Total O 97 97	0	0

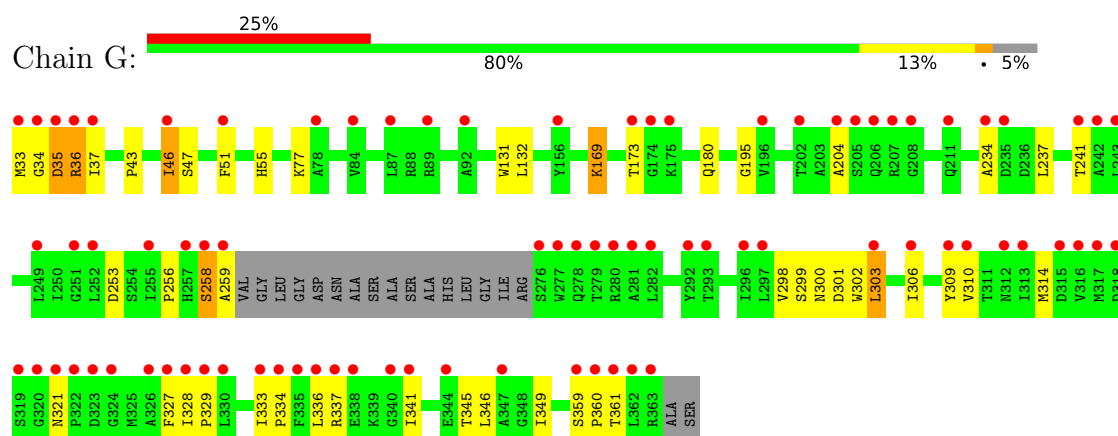
3 Residue-property plots

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.

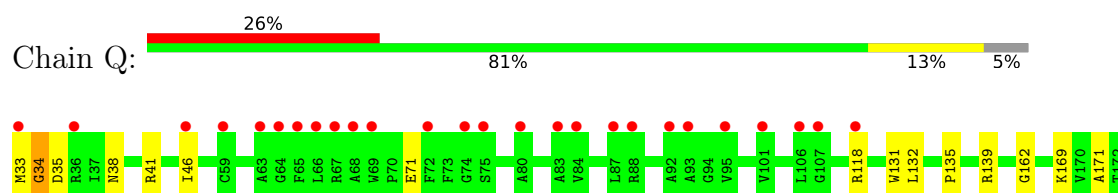
• Molecule 1: Phosphotriesterase

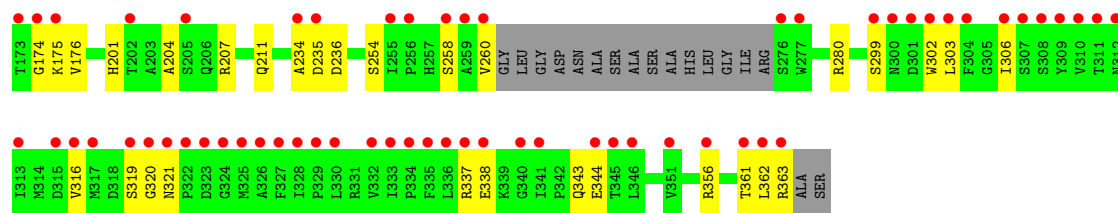


• Molecule 1: Phosphotriesterase

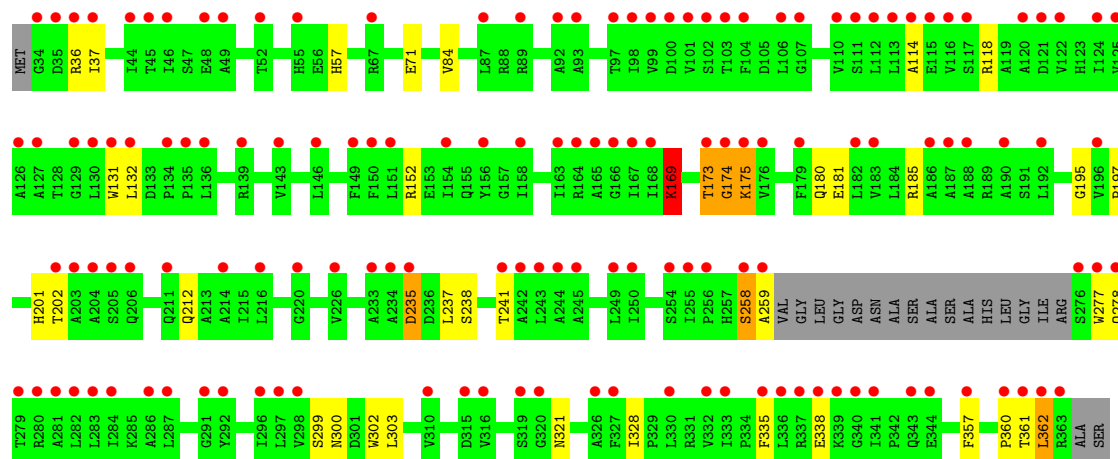
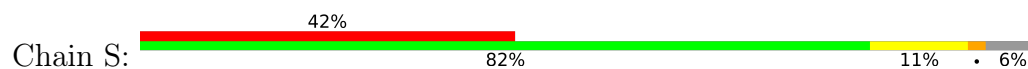


• Molecule 1: Phosphotriesterase





● Molecule 1: Phosphotriesterase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	85.66Å 86.12Å 176.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.06 – 1.60 43.06 – 1.60	Depositor EDS
% Data completeness (in resolution range)	99.8 (43.06-1.60) 99.7 (43.06-1.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.02 (at 1.60Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.264 , 0.292 0.266 , 0.295	Depositor DCC
R_{free} test set	8654 reflections (2.96%)	wwPDB-VP
Wilson B-factor (Å ²)	19.2	Xtriage
Anisotropy	0.530	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 48.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10306	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 89.64 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.1526e-08. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: CAC, MPD, ZN, KCX

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	0.29	0/2464	0.51	0/3346
1	G	0.27	0/2475	0.52	2/3362 (0.1%)
1	Q	0.27	0/2469	0.50	0/3351
1	S	0.26	0/2453	0.52	2/3330 (0.1%)
All	All	0.27	0/9861	0.51	4/13389 (0.0%)

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
1	S	0	1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	G	258	SER	CA-C-N	6.04	132.57	121.70
1	G	258	SER	C-N-CA	6.04	132.57	121.70
1	S	258	SER	CA-C-N	5.96	132.42	121.70
1	S	258	SER	C-N-CA	5.96	132.42	121.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	S	169	KCX	Mainchain

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	2434	0	2465	24	0
1	G	2444	0	2466	33	0
1	Q	2439	0	2474	25	0
1	S	2423	0	2449	25	0
2	A	2	0	0	0	0
2	G	2	0	0	0	0
2	Q	2	0	0	0	0
2	S	2	0	0	0	0
3	A	5	0	0	0	0
3	G	5	0	0	1	0
3	Q	5	0	0	0	0
3	S	5	0	0	2	0
4	A	16	0	28	2	0
4	G	16	0	28	0	0
4	Q	24	0	42	2	0
4	S	16	0	28	1	0
5	A	131	0	0	5	0
5	G	107	0	0	0	1
5	Q	131	0	0	4	0
5	S	97	0	0	0	0
All	All	10306	0	9980	106	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (106) 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:Q:33:MET:O	1:Q:35:ASP:N	2.04	0.90
1:A:34:GLY:O	1:A:36:ARG:N	2.16	0.78
1:S:173:THR:HG22	1:S:180:GLN:HE22	1.50	0.75
1:Q:207:ARG:NH1	1:Q:236:ASP:OD2	2.19	0.74
1:S:258:SER:H	1:S:259:ALA:HB3	1.55	0.71
1:G:35:ASP:N	1:G:35:ASP:OD1	2.24	0.71
1:Q:41:ARG:HD3	1:Q:118:ARG:HG2	1.70	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:235:ASP:HB2	4:Q:2405:MPD:H31	1.74	0.70
1:Q:38:ASN:ND2	5:Q:2503:HOH:O	2.25	0.68
1:Q:260:VAL:HA	1:Q:320:GLY:HA3	1.76	0.68
1:G:55:HIS:CD2	1:G:301:ASP:OD1	2.49	0.65
1:G:258:SER:HB2	1:G:259:ALA:HA	1.79	0.65
1:S:71[B]:GLU:H	1:S:71[B]:GLU:CD	2.05	0.64
1:G:237:LEU:O	1:G:241:THR:HG23	1.99	0.63
1:S:131:TRP:HE1	3:S:2403:CAC:C1	2.15	0.58
1:A:202:THR:HB	1:A:212:GLN:NE2	2.19	0.58
1:S:131:TRP:CG	1:S:132:LEU:H	2.21	0.58
1:G:33:MET:HA	1:G:360:PRO:HG2	1.86	0.58
1:A:309:TYR:HB2	4:A:2405:MPD:C1	2.34	0.57
1:S:278:GLN:HG2	1:S:335:PHE:CE1	2.40	0.57
1:A:41[B]:ARG:NH1	1:A:118:ARG:HD3	2.20	0.57
1:Q:131:TRP:CG	1:Q:132:LEU:H	2.23	0.56
1:S:84:VAL:HG11	4:S:2404:MPD:H53	1.87	0.56
1:S:197:PRO:HB2	1:S:357:PHE:CZ	2.41	0.55
1:S:235:ASP:N	1:S:235:ASP:OD1	2.37	0.55
1:G:298:VAL:HG23	1:G:349[B]:ILE:HD11	1.89	0.55
1:Q:135:PRO:O	1:Q:139[A]:ARG:HG3	2.08	0.54
1:Q:302:TRP:CH2	1:Q:321:ASN:HB3	2.43	0.54
1:A:110[B]:VAL:HG23	5:A:2568:HOH:O	2.08	0.53
1:Q:174:GLY:O	5:Q:2501:HOH:O	2.18	0.53
1:A:38:ASN:OD1	5:A:2502:HOH:O	2.18	0.52
1:Q:361:THR:HG22	1:Q:363:ARG:H	1.74	0.52
1:A:344:GLU:CD	1:A:344:GLU:H	2.18	0.52
1:G:33:MET:O	1:G:35:ASP:N	2.42	0.52
1:G:33:MET:N	1:G:36:ARG:O	2.42	0.52
1:S:237:LEU:O	1:S:241:THR:HG23	2.09	0.52
1:A:202:THR:HG23	5:A:2505:HOH:O	2.09	0.51
1:G:337:ARG:HG3	1:G:346:LEU:HD12	1.93	0.51
1:A:131:TRP:CG	1:A:132:LEU:H	2.29	0.50
1:G:258:SER:N	1:G:259:ALA:HB2	2.27	0.50
1:Q:34:GLY:N	5:Q:2507:HOH:O	2.43	0.50
1:Q:204:ALA:HB1	1:Q:234:ALA:HB2	1.93	0.50
1:S:302:TRP:CH2	1:S:321:ASN:HB3	2.46	0.50
1:S:174:GLY:HA2	1:S:175:LYS:C	2.36	0.50
1:G:51:PHE:CE2	1:G:329:PRO:HB3	2.47	0.49
1:G:302:TRP:CH2	1:G:321:ASN:HB3	2.47	0.49
1:S:361:THR:O	1:S:362:LEU:HB2	2.10	0.49
1:G:37:ILE:HG22	1:G:360:PRO:HD3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:202[B]:THR:HG23	1:S:212:GLN:NE2	2.26	0.49
1:G:131:TRP:CG	1:G:132:LEU:H	2.30	0.49
1:A:175:LYS:HA	1:A:211:GLN:HE21	1.77	0.49
1:Q:171:ALA:HA	1:Q:201:HIS:HB3	1.95	0.49
1:A:110[B]:VAL:HG11	1:A:163:ILE:HG21	1.94	0.48
1:G:314:MET:HA	1:G:314:MET:HE3	1.95	0.48
1:A:293:THR:HG21	1:A:345:THR:HG23	1.95	0.48
1:Q:337:ARG:NH1	1:Q:343:GLN:OE1	2.45	0.48
1:Q:316:VAL:O	1:Q:319:SER:OG	2.31	0.47
1:G:336:LEU:HB3	1:G:341:ILE:HD12	1.96	0.47
1:A:309:TYR:HB2	4:A:2405:MPD:H11	1.96	0.47
1:Q:344:GLU:CD	1:Q:344:GLU:H	2.22	0.47
1:S:201:HIS:CE1	3:S:2403:CAC:C1	2.98	0.46
1:G:204:ALA:HB1	1:G:234:ALA:HB2	1.97	0.46
1:S:259:ALA:HB2	1:S:277:TRP:HB3	1.96	0.46
1:S:57:HIS:CE1	1:S:169:KCX:OQ2	2.69	0.45
1:Q:303:LEU:HD22	1:Q:306:ILE:HG12	1.98	0.45
1:Q:235:ASP:OD1	5:Q:2502:HOH:O	2.21	0.45
1:Q:254:SER:O	1:Q:280:ARG:NH1	2.50	0.45
1:Q:303:LEU:HD13	1:Q:306:ILE:HD11	1.98	0.44
1:A:56:GLU:HB3	1:A:325:MET:HE2	2.00	0.44
1:G:258:SER:HB2	1:G:259:ALA:CA	2.46	0.44
1:A:67:ARG:HG3	5:A:2612:HOH:O	2.18	0.43
1:G:359[B]:SER:O	1:G:361:THR:OG1	2.31	0.43
1:Q:258:SER:OG	1:Q:280:ARG:NE	2.51	0.43
1:S:258:SER:N	1:S:259:ALA:HB3	2.28	0.43
1:A:111:SER:O	1:A:115:GLU:HG3	2.18	0.43
1:S:195:GLY:O	1:S:360:PRO:HA	2.19	0.42
1:A:59:CYS:HB3	5:A:2563:HOH:O	2.19	0.42
1:G:169:KCX:OQ1	3:G:2403:CAC:O1	2.37	0.42
1:G:173:THR:HG23	1:G:180:GLN:HE22	1.85	0.42
1:G:256:PRO:HB3	1:G:327:PHE:CD2	2.55	0.42
1:G:333:ILE:HB	1:G:334:PRO:HD3	2.02	0.42
1:A:335:PHE:O	1:A:338:GLU:HG2	2.20	0.42
1:G:298:VAL:CG2	1:G:349[B]:ILE:HD11	2.49	0.42
1:Q:71:GLU:CD	1:S:152:ARG:HE	2.28	0.42
1:Q:175:LYS:HA	1:Q:211:GLN:HE21	1.84	0.42
1:S:175:LYS:H	1:S:175:LYS:HG3	1.50	0.42
1:Q:162:GLY:O	4:Q:2405:MPD:O4	2.38	0.41
1:S:131:TRP:CG	1:S:132:LEU:N	2.87	0.41
1:S:300:ASN:OD1	1:S:328:ILE:HG12	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:131:TRP:CG	1:Q:132:LEU:N	2.88	0.41
1:A:302:TRP:CH2	1:A:321:ASN:HB3	2.56	0.41
1:G:55:HIS:HB3	1:G:253:ASP:HB2	2.01	0.41
1:G:195:GLY:O	1:G:360:PRO:HA	2.20	0.41
1:S:114:ALA:O	1:S:118:ARG:HG3	2.21	0.41
1:G:258:SER:H	1:G:259:ALA:HB2	1.86	0.41
1:G:46:ILE:HD12	1:G:46:ILE:HA	1.89	0.41
1:G:345:THR:O	1:G:349[B]:ILE:HG22	2.20	0.41
1:G:303:LEU:HD13	1:G:306:ILE:HD11	2.03	0.41
1:S:181:GLU:O	1:S:185:ARG:HG3	2.21	0.41
1:A:55:HIS:CD2	1:A:301:ASP:OD1	2.74	0.41
1:A:207:ARG:NH2	1:A:210:GLU:OE2	2.44	0.41
1:G:36:ARG:HE	1:G:43:PRO:HB2	1.86	0.40
1:G:309:TYR:CD2	1:G:310:VAL:HG22	2.57	0.40
1:A:110[B]:VAL:HG12	1:A:163:ILE:HD13	2.03	0.40
1:A:110[B]:VAL:CG1	1:A:163:ILE:HG21	2.51	0.40
1:G:300:ASN:OD1	1:G:328:ILE:HG12	2.21	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:G:2560:HOH:O	5:G:2597:HOH:O[4_556]	2.13	0.07

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	315/333 (95%)	306 (97%)	7 (2%)	2 (1%)	21	7
1	G	316/333 (95%)	305 (96%)	10 (3%)	1 (0%)	36	20
1	Q	315/333 (95%)	304 (96%)	9 (3%)	2 (1%)	21	7

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	S	313/333 (94%)	305 (97%)	6 (2%)	2 (1%)	21	7
All	All	1259/1332 (94%)	1220 (97%)	32 (2%)	7 (1%)	21	7

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	35	ASP
1	Q	34	GLY
1	A	176	VAL
1	G	34	GLY
1	S	174	GLY
1	Q	176	VAL
1	S	362	LEU

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	256/262 (98%)	252 (98%)	4 (2%)	55	33
1	G	256/262 (98%)	249 (97%)	7 (3%)	39	16
1	Q	255/262 (97%)	249 (98%)	6 (2%)	43	19
1	S	254/262 (97%)	245 (96%)	9 (4%)	32	11
All	All	1021/1048 (97%)	995 (98%)	26 (2%)	42	18

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	196	VAL
1	A	235	ASP
1	A	279	THR
1	A	299	SER
1	G	35	ASP
1	G	36	ARG
1	G	46	ILE

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Mol	Chain	Res	Type
1	G	47	SER
1	G	77	LYS
1	G	299	SER
1	G	303	LEU
1	Q	46	ILE
1	Q	299	SER
1	Q	338	GLU
1	Q	356[A]	ARG
1	Q	356[B]	ARG
1	Q	362	LEU
1	S	36	ARG
1	S	37	ILE
1	S	173	THR
1	S	175	LYS
1	S	235	ASP
1	S	238	SER
1	S	299	SER
1	S	303	LEU
1	S	338	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	G	148	GLN
1	Q	38	ASN
1	S	180	GLN
1	S	206	GLN
1	S	278	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues 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
1	KCX	S	169	2,1	10,11,12	2.25	2 (20%)	6,12,14	2.35	1 (16%)
1	KCX	G	169	2,1	10,11,12	2.39	2 (20%)	6,12,14	2.31	2 (33%)
1	KCX	Q	169	2,1	10,11,12	2.00	2 (20%)	6,12,14	1.19	1 (16%)
1	KCX	A	169	2,1	10,11,12	1.77	2 (20%)	6,12,14	2.65	2 (33%)

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
1	KCX	S	169	2,1	-	1/9/10/12	-
1	KCX	G	169	2,1	-	1/9/10/12	-
1	KCX	Q	169	2,1	-	1/9/10/12	-
1	KCX	A	169	2,1	-	1/9/10/12	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	169	KCX	CX-NZ	6.57	1.46	1.35
1	S	169	KCX	CX-NZ	6.27	1.46	1.35
1	Q	169	KCX	CX-NZ	5.35	1.44	1.35
1	A	169	KCX	CX-NZ	4.66	1.43	1.35
1	G	169	KCX	OQ1-CX	3.10	1.27	1.21
1	Q	169	KCX	OQ1-CX	2.71	1.26	1.21
1	S	169	KCX	OQ1-CX	2.62	1.26	1.21
1	A	169	KCX	OQ1-CX	2.46	1.26	1.21

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	169	KCX	OQ1-CX-NZ	-5.56	116.48	124.92
1	S	169	KCX	OQ1-CX-NZ	-5.33	116.83	124.92
1	G	169	KCX	OQ1-CX-NZ	-5.08	117.20	124.92
1	Q	169	KCX	OQ1-CX-NZ	-2.46	121.18	124.92
1	A	169	KCX	CE-NZ-CX	-2.40	117.90	121.98
1	G	169	KCX	CE-NZ-CX	-2.12	118.39	121.98

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	169	KCX	C-CA-CB-CG
1	S	169	KCX	C-CA-CB-CG
1	G	169	KCX	C-CA-CB-CG
1	Q	169	KCX	C-CA-CB-CG

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	S	169	KCX	1	0
1	G	169	KCX	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 21 ligands modelled in this entry, 8 are monoatomic - leaving 13 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	$\# Z > 2$
4	MPD	Q	2404	-	7,7,7	0.70	0	9,10,10	0.68	0
3	CAC	Q	2403	2	2,4,4	2.71	2 (100%)	4,6,6	2.11	1 (25%)
4	MPD	Q	2405	-	7,7,7	0.68	0	9,10,10	0.54	0
4	MPD	S	2405	-	7,7,7	0.78	0	9,10,10	0.37	0
3	CAC	G	2403	2	2,4,4	2.63	2 (100%)	4,6,6	9.00	3 (75%)
4	MPD	G	2405	-	7,7,7	0.70	0	9,10,10	0.42	0
4	MPD	A	2404	-	7,7,7	0.75	0	9,10,10	0.40	0
3	CAC	S	2403	2	2,4,4	1.93	1 (50%)	4,6,6	3.98	1 (25%)
4	MPD	Q	2406	-	7,7,7	0.74	0	9,10,10	0.37	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	CAC	A	2403	2	2,4,4	2.76	2 (100%)	4,6,6	1.95	2 (50%)
4	MPD	A	2405	-	7,7,7	0.68	0	9,10,10	0.56	0
4	MPD	G	2404	-	7,7,7	0.64	0	9,10,10	0.70	0
4	MPD	S	2404	-	7,7,7	0.69	0	9,10,10	0.36	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
4	MPD	Q	2404	-	-	0/5/5/5	-
4	MPD	Q	2405	-	-	3/5/5/5	-
4	MPD	S	2405	-	-	3/5/5/5	-
4	MPD	G	2405	-	-	3/5/5/5	-
4	MPD	A	2404	-	-	5/5/5/5	-
4	MPD	Q	2406	-	-	2/5/5/5	-
4	MPD	A	2405	-	-	0/5/5/5	-
4	MPD	G	2404	-	-	1/5/5/5	-
4	MPD	S	2404	-	-	3/5/5/5	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	2403	CAC	AS-C2	3.10	1.97	1.90
3	A	2403	CAC	AS-C2	2.93	1.97	1.90
3	Q	2403	CAC	AS-C2	2.88	1.97	1.90
3	A	2403	CAC	AS-C1	2.58	1.96	1.90
3	Q	2403	CAC	AS-C1	2.53	1.96	1.90
3	S	2403	CAC	AS-C2	2.35	1.95	1.90
3	G	2403	CAC	AS-C1	2.05	1.95	1.90

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	2403	CAC	O1-AS-C2	-17.31	89.98	111.50
3	S	2403	CAC	O1-AS-C2	-7.61	102.03	111.50
3	G	2403	CAC	O2-AS-C1	3.79	115.02	105.84
3	Q	2403	CAC	O1-AS-C2	-3.52	107.12	111.50
3	G	2403	CAC	O2-AS-C2	3.17	113.51	105.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	2403	CAC	O1-AS-C2	-2.51	108.38	111.50
3	A	2403	CAC	O2-AS-C2	2.05	110.81	105.84

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	2404	MPD	C2-C3-C4-O4
4	A	2404	MPD	C2-C3-C4-C5
4	Q	2405	MPD	C1-C2-C3-C4
4	Q	2405	MPD	O2-C2-C3-C4
4	A	2404	MPD	C1-C2-C3-C4
4	G	2405	MPD	C1-C2-C3-C4
4	G	2405	MPD	CM-C2-C3-C4
4	Q	2405	MPD	CM-C2-C3-C4
4	Q	2406	MPD	C1-C2-C3-C4
4	S	2405	MPD	C1-C2-C3-C4
4	A	2404	MPD	O2-C2-C3-C4
4	G	2404	MPD	O2-C2-C3-C4
4	G	2405	MPD	O2-C2-C3-C4
4	S	2404	MPD	O2-C2-C3-C4
4	S	2405	MPD	O2-C2-C3-C4
4	A	2404	MPD	CM-C2-C3-C4
4	Q	2406	MPD	CM-C2-C3-C4
4	S	2404	MPD	C1-C2-C3-C4
4	S	2404	MPD	CM-C2-C3-C4
4	S	2405	MPD	CM-C2-C3-C4

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	Q	2405	MPD	2	0
3	G	2403	CAC	1	0
3	S	2403	CAC	2	0
4	A	2405	MPD	2	0
4	S	2404	MPD	1	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, 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	314/333 (94%)	1.03	60 (19%) 3 3	9, 25, 42, 58	7 (2%)
1	G	314/333 (94%)	1.43	83 (26%) 1 1	14, 33, 48, 59	7 (2%)
1	Q	315/333 (94%)	1.28	86 (27%) 1 1	9, 29, 48, 61	6 (1%)
1	S	313/333 (93%)	2.01	141 (45%) 0 0	11, 39, 49, 58	4 (1%)
All	All	1256/1332 (94%)	1.44	370 (29%) 1 1	9, 33, 47, 61	24 (1%)

All (370) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	S	362	LEU	9.1
1	G	362	LEU	7.6
1	S	259	ALA	7.4
1	A	34	GLY	7.1
1	S	37	ILE	6.8
1	S	174	GLY	6.4
1	G	33	MET	5.7
1	A	204	ALA	5.0
1	A	258[A]	SER	5.0
1	G	363	ARG	4.9
1	G	34	GLY	4.7
1	G	258	SER	4.7
1	Q	327	PHE	4.7
1	A	362	LEU	4.7
1	S	46	ILE	4.6
1	G	259	ALA	4.5
1	S	175	LYS	4.5
1	S	204	ALA	4.4
1	S	234	ALA	4.4
1	Q	302	TRP	4.3
1	Q	260	VAL	4.3

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Mol	Chain	Res	Type	RSRZ
1	A	215	ILE	4.3
1	A	363	ARG	4.3
1	S	156	TYR	4.2
1	S	101	VAL	4.2
1	Q	330	LEU	4.2
1	S	255	ILE	4.2
1	S	363	ARG	4.2
1	S	124	ILE	4.2
1	A	203	ALA	4.1
1	Q	309	TYR	4.1
1	A	175	LYS	4.1
1	A	233	ALA	4.0
1	S	173	THR	4.0
1	S	34	GLY	4.0
1	Q	362	LEU	4.0
1	S	192	LEU	4.0
1	Q	316	VAL	4.0
1	Q	310	VAL	3.9
1	A	182	LEU	3.9
1	Q	304	PHE	3.9
1	S	150	PHE	3.9
1	S	125	VAL	3.8
1	Q	363	ARG	3.8
1	Q	319	SER	3.8
1	Q	326	ALA	3.8
1	Q	306	ILE	3.7
1	S	106	LEU	3.7
1	A	235	ASP	3.7
1	A	205	SER	3.7
1	G	173	THR	3.7
1	Q	361	THR	3.7
1	S	283	LEU	3.7
1	G	255	ILE	3.6
1	G	329	PRO	3.6
1	S	203	ALA	3.6
1	Q	106	LEU	3.6
1	Q	259	ALA	3.5
1	S	130	LEU	3.5
1	S	154	ILE	3.5
1	A	178	PRO	3.5
1	A	46	ILE	3.5
1	A	228	ILE	3.5

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Mol	Chain	Res	Type	RSRZ
1	S	35	ASP	3.5
1	A	183	VAL	3.5
1	S	103	THR	3.5
1	A	190	ALA	3.4
1	S	258	SER	3.4
1	S	298	VAL	3.4
1	S	235	ASP	3.4
1	S	44	ILE	3.4
1	G	326	ALA	3.4
1	G	46	ILE	3.4
1	S	344	GLU	3.4
1	A	361	THR	3.4
1	Q	299	SER	3.4
1	Q	33	MET	3.4
1	S	281	ALA	3.3
1	Q	338	GLU	3.3
1	A	179	PHE	3.3
1	Q	258	SER	3.3
1	S	233	ALA	3.3
1	S	244	ALA	3.3
1	S	99	VAL	3.3
1	G	324	GLY	3.3
1	S	277	TRP	3.3
1	S	151	LEU	3.3
1	S	186	ALA	3.3
1	S	205	SER	3.2
1	S	360	PRO	3.2
1	A	260	VAL	3.2
1	Q	235	ASP	3.2
1	S	168	ILE	3.2
1	Q	69	TRP	3.2
1	A	259	ALA	3.2
1	G	35	ASP	3.2
1	S	113	LEU	3.2
1	S	282	LEU	3.2
1	S	276	SER	3.2
1	Q	335	PHE	3.2
1	S	335	PHE	3.2
1	A	202	THR	3.2
1	G	205	SER	3.1
1	S	183	VAL	3.1
1	G	360	PRO	3.1

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Mol	Chain	Res	Type	RSRZ
1	G	328	ILE	3.1
1	G	330	LEU	3.1
1	Q	303	LEU	3.1
1	A	214	ALA	3.1
1	A	226	VAL	3.1
1	G	310	VAL	3.1
1	S	319	SER	3.1
1	Q	312	ASN	3.1
1	S	149	PHE	3.1
1	A	35	ASP	3.1
1	A	196	VAL	3.0
1	Q	255	ILE	3.0
1	Q	313	ILE	3.0
1	S	336	LEU	3.0
1	A	198	VAL	3.0
1	Q	332	VAL	3.0
1	G	313	ILE	3.0
1	S	165	ALA	3.0
1	S	211[A]	GLN	3.0
1	S	361	THR	3.0
1	S	187	ALA	3.0
1	Q	328	ILE	3.0
1	S	139	ARG	3.0
1	Q	344	GLU	3.0
1	Q	329	PRO	3.0
1	S	320	GLY	3.0
1	G	309	TYR	2.9
1	A	221	LEU	2.9
1	G	87	LEU	2.9
1	Q	324	GLY	2.9
1	S	45	THR	2.9
1	Q	66	LEU	2.9
1	S	132	LEU	2.9
1	Q	59	CYS	2.9
1	Q	341	ILE	2.9
1	S	167	ILE	2.9
1	S	100	ASP	2.9
1	S	357	PHE	2.9
1	S	110	VAL	2.9
1	A	234	ALA	2.9
1	Q	80	ALA	2.9
1	S	36	ARG	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	184	LEU	2.9
1	S	327	PHE	2.8
1	S	216	LEU	2.8
1	G	327	PHE	2.8
1	G	320	GLY	2.8
1	S	286	ALA	2.8
1	Q	307	SER	2.8
1	A	41[A]	ARG	2.8
1	G	37	ILE	2.8
1	S	163	ILE	2.8
1	S	284	ILE	2.8
1	G	323	ASP	2.8
1	Q	320	GLY	2.8
1	G	204	ALA	2.8
1	G	319	SER	2.8
1	S	120	ALA	2.8
1	S	188	ALA	2.8
1	G	84	VAL	2.8
1	G	316	VAL	2.8
1	G	337	ARG	2.8
1	A	192	LEU	2.8
1	G	175	LYS	2.8
1	G	293	THR	2.8
1	G	359[A]	SER	2.8
1	S	190	ALA	2.7
1	A	211	GLN	2.7
1	Q	72	PHE	2.7
1	Q	277	TRP	2.7
1	S	131	TRP	2.7
1	S	206	GLN	2.7
1	S	136	LEU	2.7
1	Q	46	ILE	2.7
1	S	92	ALA	2.7
1	S	242	ALA	2.7
1	A	236	ASP	2.7
1	G	156[A]	TYR	2.7
1	S	112	LEU	2.7
1	S	330	LEU	2.7
1	S	97	THR	2.7
1	Q	92	ALA	2.7
1	Q	323	ASP	2.7
1	S	337	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
1	S	338	GLU	2.7
1	A	276	SER	2.7
1	A	176	VAL	2.7
1	G	202	THR	2.7
1	A	174	GLY	2.7
1	Q	234	ALA	2.7
1	S	89	ARG	2.6
1	S	135	PRO	2.6
1	A	216	LEU	2.6
1	S	127	ALA	2.6
1	G	51	PHE	2.6
1	S	179	PHE	2.6
1	G	292	TYR	2.6
1	S	310	VAL	2.6
1	S	98	ILE	2.6
1	S	158	ILE	2.6
1	S	339	LYS	2.6
1	A	220	GLY	2.6
1	G	36	ARG	2.6
1	G	280	ARG	2.6
1	S	166	GLY	2.6
1	Q	315	ASP	2.6
1	Q	93	ALA	2.6
1	G	277	TRP	2.6
1	G	322	PRO	2.6
1	A	36	ARG	2.6
1	A	338	GLU	2.6
1	G	89	ARG	2.6
1	Q	356[A]	ARG	2.6
1	G	235	ASP	2.5
1	Q	74	GLY	2.5
1	G	278[A]	GLN	2.5
1	A	222	SER	2.5
1	S	249	LEU	2.5
1	S	116	VAL	2.5
1	S	143	VAL	2.5
1	S	114	ALA	2.5
1	S	241	THR	2.5
1	Q	65	PHE	2.5
1	S	104	PHE	2.5
1	S	243	LEU	2.5
1	S	196	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	G	341	ILE	2.5
1	Q	333	ILE	2.5
1	S	129	GLY	2.5
1	S	340	GLY	2.5
1	S	111	SER	2.5
1	S	278	GLN	2.5
1	S	182	LEU	2.5
1	S	287	LEU	2.5
1	G	281	ALA	2.5
1	Q	322	PRO	2.5
1	S	245	ALA	2.5
1	G	196	VAL	2.5
1	G	306	ILE	2.5
1	S	250	ILE	2.5
1	G	338	GLU	2.5
1	A	218	SER	2.5
1	G	315	ASP	2.4
1	A	223	PRO	2.4
1	A	186	ALA	2.4
1	G	234	ALA	2.4
1	Q	63	ALA	2.4
1	A	255	ILE	2.4
1	G	333	ILE	2.4
1	Q	205	SER	2.4
1	Q	345	THR	2.4
1	S	176	VAL	2.4
1	G	206	GLN	2.4
1	G	251	GLY	2.4
1	S	291	GLY	2.4
1	Q	321	ASN	2.4
1	Q	101	VAL	2.4
1	Q	351	VAL	2.4
1	Q	301	ASP	2.4
1	Q	107	GLY	2.4
1	Q	83	ALA	2.4
1	S	279	THR	2.4
1	A	288	ILE	2.4
1	S	341	ILE	2.4
1	G	321	ASN	2.3
1	G	303	LEU	2.3
1	G	335	PHE	2.3
1	S	122	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	S	117	SER	2.3
1	G	344	GLU	2.3
1	S	164	ARG	2.3
1	G	318	ASP	2.3
1	A	227	CYS	2.3
1	Q	311	THR	2.3
1	S	126	ALA	2.3
1	G	282	LEU	2.3
1	S	292	TYR	2.3
1	S	333	ILE	2.3
1	G	174	GLY	2.3
1	G	334	PRO	2.3
1	Q	202	THR	2.3
1	S	115	GLU	2.3
1	G	276	SER	2.3
1	S	102	SER	2.3
1	G	296	ILE	2.3
1	S	296	ILE	2.3
1	A	298[A]	VAL	2.3
1	Q	95	VAL	2.3
1	G	257	HIS	2.2
1	G	92	ALA	2.2
1	G	317	MET	2.2
1	S	332	VAL	2.2
1	S	256	PRO	2.2
1	Q	88[A]	ARG	2.2
1	S	280	ARG	2.2
1	G	78	ALA	2.2
1	G	347	ALA	2.2
1	G	252	LEU	2.2
1	G	336	LEU	2.2
1	Q	87	LEU	2.2
1	Q	337	ARG	2.2
1	Q	256	PRO	2.2
1	S	134	PRO	2.2
1	Q	84	VAL	2.2
1	S	93	ALA	2.2
1	G	211	GLN	2.2
1	S	48	GLU	2.2
1	A	209	GLY	2.2
1	G	249	LEU	2.2
1	G	297	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	Q	336	LEU	2.2
1	S	146	LEU	2.2
1	Q	118	ARG	2.2
1	G	279	THR	2.2
1	Q	308	SER	2.2
1	Q	317	MET	2.1
1	Q	325	MET	2.1
1	A	207	ARG	2.1
1	Q	67	ARG	2.1
1	S	107	GLY	2.1
1	G	243	LEU	2.1
1	Q	346	LEU	2.1
1	Q	276	SER	2.1
1	S	49	ALA	2.1
1	Q	64	GLY	2.1
1	Q	174	GLY	2.1
1	S	220	GLY	2.1
1	S	297	LEU	2.1
1	G	241	THR	2.1
1	G	361	THR	2.1
1	Q	334	PRO	2.1
1	A	193	ALA	2.1
1	S	214	ALA	2.1
1	G	312	ASN	2.1
1	Q	75	SER	2.1
1	S	121	ASP	2.1
1	A	118	ARG	2.1
1	Q	36	ARG	2.1
1	S	87	LEU	2.1
1	S	52	THR	2.1
1	S	202[A]	THR	2.1
1	G	208	GLY	2.0
1	G	340	GLY	2.0
1	A	187	ALA	2.0
1	A	188	ALA	2.0
1	A	242	ALA	2.0
1	Q	300	ASN	2.0
1	S	326	ALA	2.0
1	A	206	GLN	2.0
1	S	226	VAL	2.0
1	S	315	ASP	2.0
1	G	207	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
1	S	254	SER	2.0
1	A	237	LEU	2.0
1	Q	173	THR	2.0
1	A	197	PRO	2.0
1	Q	175	LYS	2.0
1	Q	340	GLY	2.0
1	S	343	GLN	2.0
1	S	67[A]	ARG	2.0
1	G	242	ALA	2.0
1	Q	68	ALA	2.0
1	S	316	VAL	2.0
1	A	230	HIS	2.0
1	S	55	HIS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	KCX	S	169	12/13	0.81	0.20	39,40,40,40	0
1	KCX	G	169	12/13	0.87	0.11	18,22,25,26	0
1	KCX	A	169	12/13	0.92	0.09	15,17,19,20	0
1	KCX	Q	169	12/13	0.95	0.07	17,18,21,22	0

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(Å ²)	Q<0.9
4	MPD	G	2404	8/8	0.47	0.25	35,41,44,45	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MPD	S	2405	8/8	0.60	0.20	42,43,46,48	0
4	MPD	S	2404	8/8	0.65	0.22	35,39,45,45	0
4	MPD	G	2405	8/8	0.69	0.18	41,43,44,47	0
4	MPD	A	2405	8/8	0.72	0.22	18,25,28,32	8
4	MPD	Q	2405	8/8	0.74	0.16	34,37,41,44	0
4	MPD	Q	2404	8/8	0.76	0.15	39,43,44,44	0
4	MPD	Q	2406	8/8	0.76	0.17	33,38,40,46	0
4	MPD	A	2404	8/8	0.83	0.14	32,34,40,44	0
2	ZN	S	2401	1/1	0.89	0.09	20,20,20,20	1
2	ZN	G	2401	1/1	0.90	0.08	20,20,20,20	0
2	ZN	S	2402	1/1	0.91	0.07	21,21,21,21	1
2	ZN	G	2402	1/1	0.91	0.09	21,21,21,21	1
3	CAC	G	2403	5/5	0.93	0.20	13,23,25,33	5
3	CAC	S	2403	5/5	0.96	0.11	10,15,20,22	5
2	ZN	Q	2401	1/1	0.98	0.03	15,15,15,15	1
3	CAC	Q	2403	5/5	0.98	0.08	15,16,20,22	5
2	ZN	Q	2402	1/1	0.99	0.03	17,17,17,17	1
2	ZN	A	2402	1/1	0.99	0.03	14,14,14,14	1
2	ZN	A	2401	1/1	0.99	0.04	14,14,14,14	0
3	CAC	A	2403	5/5	0.99	0.09	12,13,15,18	5

6.5 Other polymers

There are no such residues in this entry.