



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:03 AM UTC

PDB ID : 3D9S / pdb_00003d9s
Title : Human Aquaporin 5 (AQP5) - High Resolution X-ray Structure
Authors : Horsefield, R.; Norden, K.; Fellert, M.; Backmark, A.; Tornroth-Horsefield, S.;
Terwisscha Van Scheltinga, A.C.; Kvassman, J.; Kjellbom, P.; Johanson, U.;
Neutze, R.
Deposited on : 2008-05-27
Resolution : 2.00 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

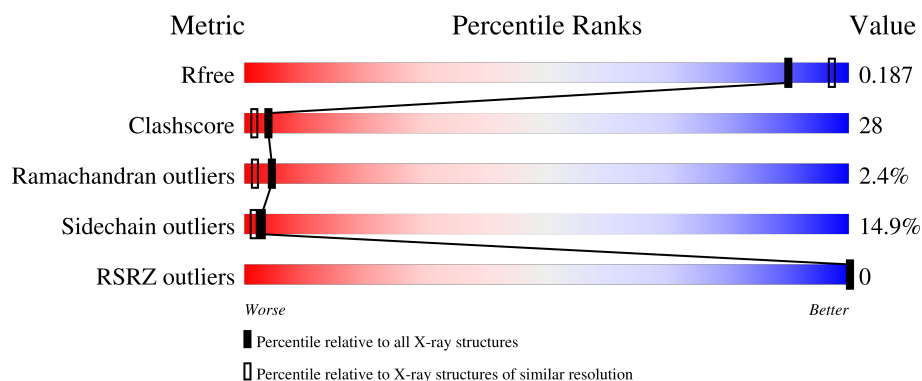
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

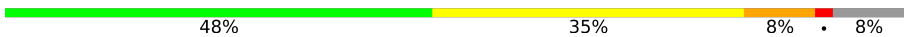
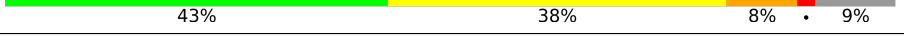
The reported resolution of this entry is 2.00 Å.

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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	266	
1	B	266	
1	C	266	
1	D	266	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7390 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.

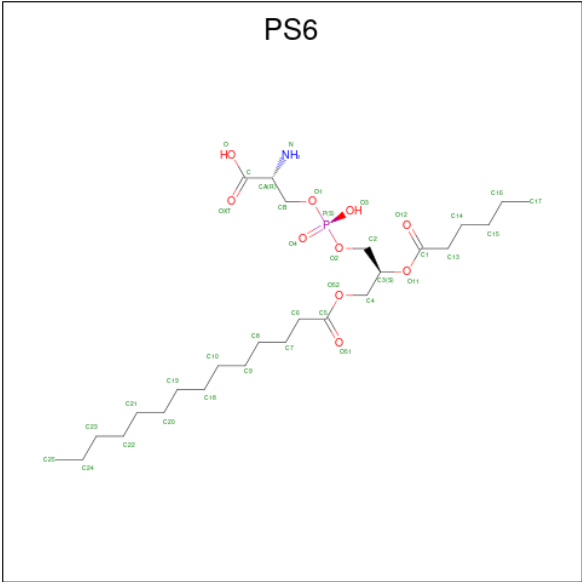
- Molecule 1 is a protein called Aquaporin-5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	245	Total	C	N	O	S	0	0	0
			1814	1199	298	311	6			
1	B	245	Total	C	N	O	S	0	0	0
			1814	1199	298	311	6			
1	C	243	Total	C	N	O	S	0	0	0
			1799	1190	295	308	6			
1	D	245	Total	C	N	O	S	0	0	0
			1814	1199	298	311	6			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP P55064
A	1	SER	-	expression tag	UNP P55064
B	0	MET	-	initiating methionine	UNP P55064
B	1	SER	-	expression tag	UNP P55064
C	0	MET	-	initiating methionine	UNP P55064
C	1	SER	-	expression tag	UNP P55064
D	0	MET	-	initiating methionine	UNP P55064
D	1	SER	-	expression tag	UNP P55064

- Molecule 2 is O-[(S)-{[(2S)-2-(hexanoyloxy)-3-(tetradecanoyloxy)propyl]oxy} (hydroxy)phosphoryl]-D-serine (CCD ID: PS6) (formula: C₂₆H₅₀NO₁₀P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	D	1	Total	C	N	O	P	0	0
			38	26	1	10	1		

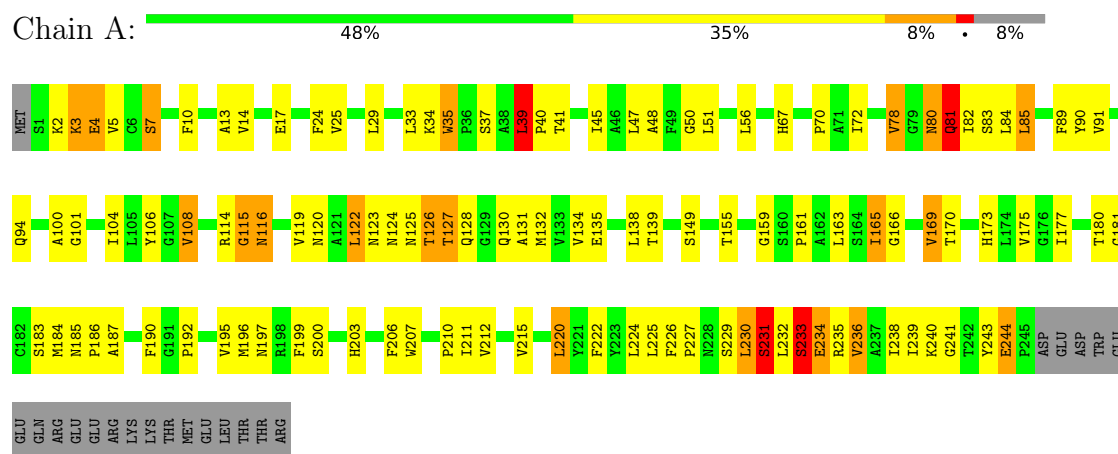
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	26	Total	O	0	0
			26	26		
3	B	27	Total	O	0	0
			27	27		
3	C	32	Total	O	0	0
			32	32		
3	D	26	Total	O	0	0
			26	26		

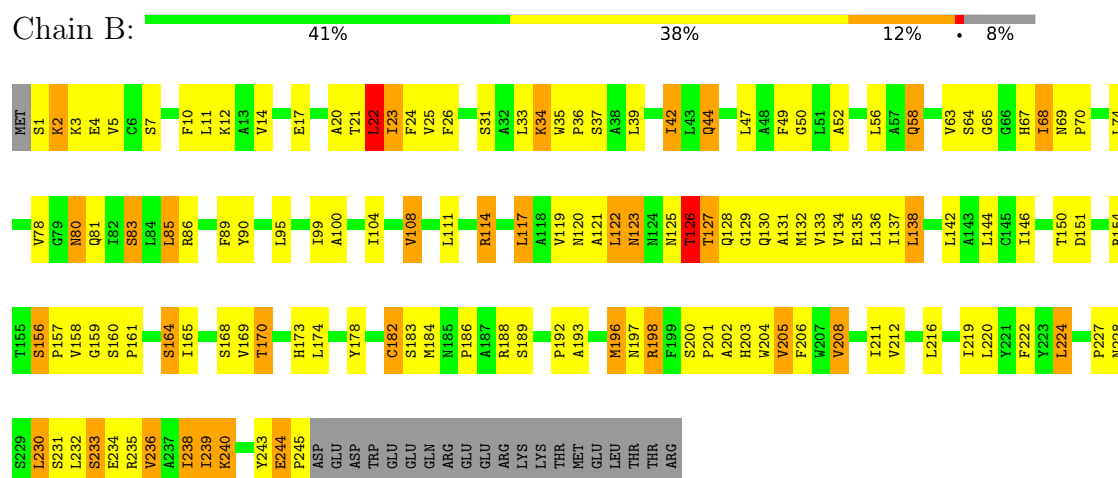
3 Residue-property plots

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.

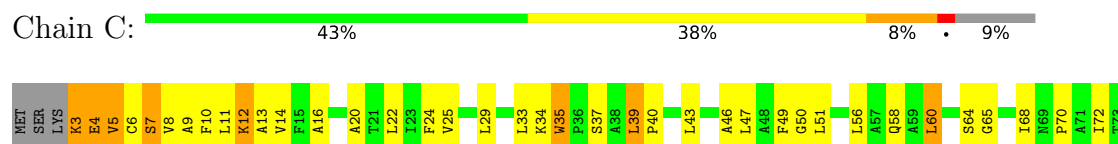
• Molecule 1: Aquaporin-5

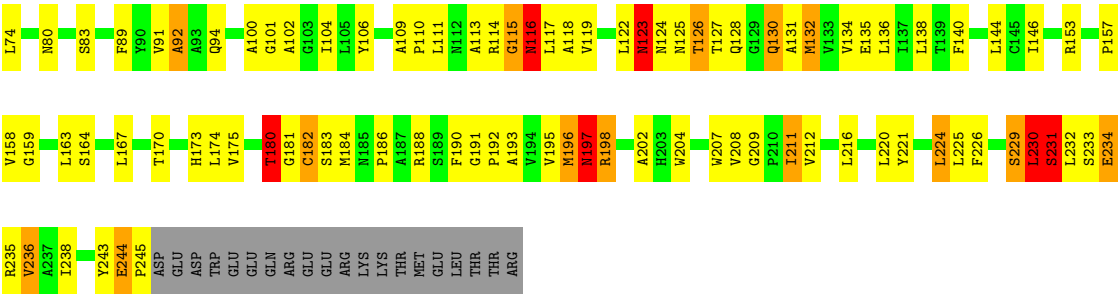


• Molecule 1: Aquaporin-5

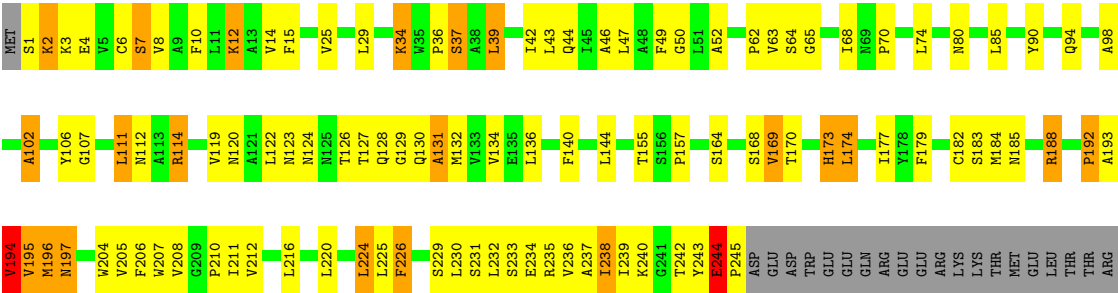


• Molecule 1: Aquaporin-5





● Molecule 1: Aquaporin-5



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	90.48Å 90.64Å 184.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.00 10.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	78.0 (10.00-2.00) 81.8 (10.00-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 2.00Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.162 , 0.193 0.160 , 0.187	Depositor DCC
R_{free} test set	4222 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	29.2	Xtriage
Anisotropy	0.074	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 51.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.21$	Xtriage
Estimated twinning fraction	0.447 for k,h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7390	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.68% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 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: PS6

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	0.47	0/1858	1.66	18/2539 (0.7%)
1	B	0.45	0/1858	1.56	27/2539 (1.1%)
1	C	0.47	0/1843	1.66	25/2520 (1.0%)
1	D	0.47	0/1858	1.60	34/2539 (1.3%)
All	All	0.47	0/7417	1.62	104/10137 (1.0%)

There are no bond length outliers.

All (104) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	115	GLY	CA-C-N	15.46	149.53	121.70
1	A	115	GLY	C-N-CA	15.46	149.53	121.70
1	C	116	ASN	CA-C-N	14.07	147.04	121.70
1	C	116	ASN	C-N-CA	14.07	147.04	121.70
1	A	39	LEU	O-C-N	12.43	132.12	121.31
1	C	115	GLY	CA-C-N	10.39	141.39	121.54
1	C	115	GLY	C-N-CA	10.39	141.39	121.54
1	C	229	SER	CA-C-N	9.83	140.31	121.54
1	C	229	SER	C-N-CA	9.83	140.31	121.54
1	C	49	PHE	CA-C-N	8.43	129.29	119.94
1	C	49	PHE	C-N-CA	8.43	129.29	119.94
1	C	175	VAL	CA-C-N	8.26	129.15	119.98
1	C	175	VAL	C-N-CA	8.26	129.15	119.98
1	C	35	TRP	CA-C-O	8.19	127.44	119.75
1	D	111	LEU	CA-C-N	8.03	136.88	121.54
1	D	111	LEU	C-N-CA	8.03	136.88	121.54
1	C	130	GLN	CA-C-N	7.89	130.86	120.28
1	C	130	GLN	C-N-CA	7.89	130.86	120.28
1	A	230	LEU	CA-C-N	7.73	135.61	121.70
1	A	230	LEU	C-N-CA	7.73	135.61	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	205	VAL	O-C-N	7.52	129.16	121.87
1	B	86	ARG	CA-C-N	7.50	130.94	120.29
1	B	86	ARG	C-N-CA	7.50	130.94	120.29
1	D	212	VAL	CA-C-N	7.29	127.88	119.94
1	D	212	VAL	C-N-CA	7.29	127.88	119.94
1	C	181	GLY	N-CA-C	-6.82	106.35	115.21
1	A	175	VAL	CA-C-N	6.80	127.53	119.98
1	A	175	VAL	C-N-CA	6.80	127.53	119.98
1	B	22	LEU	CA-CB-CG	6.76	139.98	116.30
1	D	194	VAL	O-C-N	-6.69	114.21	122.57
1	B	151	ASP	CA-CB-CG	6.68	119.28	112.60
1	D	65	GLY	O-C-N	6.60	127.99	122.71
1	D	107	GLY	CA-C-N	-6.57	114.84	122.14
1	D	107	GLY	C-N-CA	-6.57	114.84	122.14
1	D	226	PHE	CA-CB-CG	-6.49	107.31	113.80
1	B	65	GLY	N-CA-C	-6.47	107.39	115.08
1	B	174	LEU	CA-C-N	-6.41	116.60	123.08
1	B	174	LEU	C-N-CA	-6.41	116.60	123.08
1	B	44	GLN	CA-C-O	6.29	127.43	120.82
1	B	108	VAL	O-C-N	-6.27	116.08	122.23
1	C	180	THR	CA-C-O	6.18	126.15	118.43
1	A	126	THR	CA-C-N	6.15	130.30	121.50
1	A	126	THR	C-N-CA	6.15	130.30	121.50
1	A	212	VAL	CA-C-N	6.08	126.84	120.03
1	A	212	VAL	C-N-CA	6.08	126.84	120.03
1	B	170	THR	CA-C-N	5.98	129.79	120.82
1	B	170	THR	C-N-CA	5.98	129.79	120.82
1	D	196	MET	N-CA-C	-5.92	102.71	110.53
1	D	188	ARG	CD-NE-CZ	5.89	132.65	124.40
1	D	197	ASN	CA-CB-CG	5.89	118.49	112.60
1	B	108	VAL	CA-C-O	5.88	124.84	119.20
1	C	83	SER	CA-C-N	5.88	128.64	120.29
1	C	83	SER	C-N-CA	5.88	128.64	120.29
1	B	236	VAL	CA-C-N	5.87	128.15	120.28
1	B	236	VAL	C-N-CA	5.87	128.15	120.28
1	B	58	GLN	CB-CG-CD	5.87	122.58	112.60
1	C	92	ALA	CA-C-N	5.77	127.94	120.44
1	C	92	ALA	C-N-CA	5.77	127.94	120.44
1	D	44	GLN	CA-C-N	5.74	127.80	120.56
1	D	44	GLN	C-N-CA	5.74	127.80	120.56
1	D	192	PRO	CA-C-N	5.71	127.86	120.44
1	D	192	PRO	C-N-CA	5.71	127.86	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	64	SER	CA-C-O	5.59	124.84	118.47
1	A	181	GLY	CA-C-N	5.55	132.14	121.54
1	A	181	GLY	C-N-CA	5.55	132.14	121.54
1	D	12	LYS	CA-C-N	5.55	127.71	120.28
1	D	12	LYS	C-N-CA	5.55	127.71	120.28
1	A	234	GLU	N-CA-C	-5.53	106.30	113.16
1	D	102	ALA	CA-C-N	5.46	126.00	119.94
1	D	102	ALA	C-N-CA	5.46	126.00	119.94
1	B	23	ILE	O-C-N	5.45	129.03	121.84
1	C	197	ASN	CA-C-N	5.45	131.95	121.54
1	C	197	ASN	C-N-CA	5.45	131.95	121.54
1	B	126	THR	CA-C-O	5.45	126.92	120.66
1	B	49	PHE	CA-C-N	5.40	125.94	119.94
1	B	49	PHE	C-N-CA	5.40	125.94	119.94
1	A	78	VAL	O-C-N	-5.39	116.63	121.91
1	C	58	GLN	O-C-N	5.38	127.83	122.12
1	D	131	ALA	CA-C-N	5.37	127.79	120.54
1	D	131	ALA	C-N-CA	5.37	127.79	120.54
1	D	185	ASN	CA-C-N	5.37	125.70	119.47
1	D	185	ASN	C-N-CA	5.37	125.70	119.47
1	D	168	SER	CA-C-N	-5.33	111.55	120.30
1	D	168	SER	C-N-CA	-5.33	111.55	120.30
1	D	173	HIS	CA-CB-CG	-5.29	108.51	113.80
1	D	196	MET	CA-C-N	-5.28	111.45	121.54
1	D	196	MET	C-N-CA	-5.28	111.45	121.54
1	B	58	GLN	O-C-N	5.26	127.49	122.07
1	B	151	ASP	CA-C-N	5.26	131.59	121.54
1	B	151	ASP	C-N-CA	5.26	131.59	121.54
1	B	36	PRO	CA-C-N	5.22	127.53	120.38
1	B	36	PRO	C-N-CA	5.22	127.53	120.38
1	B	219	ILE	CA-C-N	5.22	127.70	120.29
1	B	219	ILE	C-N-CA	5.22	127.70	120.29
1	D	164	SER	CA-C-N	5.21	127.23	120.56
1	D	164	SER	C-N-CA	5.21	127.23	120.56
1	C	65	GLY	N-CA-C	-5.21	107.50	115.00
1	C	116	ASN	CA-CB-CG	5.20	117.80	112.60
1	A	199	PHE	CA-CB-CG	-5.15	108.65	113.80
1	D	64	SER	CA-C-N	-5.13	117.86	123.30
1	D	64	SER	C-N-CA	-5.13	117.86	123.30
1	C	234	GLU	CA-CB-CG	5.12	124.34	114.10
1	A	80	ASN	CA-C-N	5.08	131.24	121.54
1	A	80	ASN	C-N-CA	5.08	131.24	121.54

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	1814	0	1887	93	0
1	B	1814	0	1888	133	0
1	C	1799	0	1867	107	0
1	D	1814	0	1888	96	0
2	D	38	0	48	16	0
3	A	26	0	0	6	0
3	B	27	0	0	6	0
3	C	32	0	0	2	0
3	D	26	0	0	1	0
All	All	7390	0	7578	418	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 28.

All (418) 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:123:ASN:ND2	1:C:123:ASN:O	1.87	1.07
1:B:228:ASN:CG	1:B:230:LEU:HD23	1.81	1.05
1:A:127:THR:HG22	1:A:130:GLN:H	1.23	1.04
1:D:236:VAL:HG13	1:D:240:LYS:HE3	1.40	1.02
1:B:228:ASN:ND2	1:B:230:LEU:HD21	1.75	1.00
1:B:228:ASN:CG	1:B:230:LEU:CD2	2.36	0.98
1:A:236:VAL:HA	1:A:239:ILE:HD12	1.45	0.97
1:C:157:PRO:HB3	2:D:266:PS6:H15A	1.47	0.96
1:B:123:ASN:HD21	1:B:125:ASN:HB2	1.38	0.88
1:C:157:PRO:HA	2:D:266:PS6:H17B	1.54	0.87
1:B:228:ASN:ND2	1:B:230:LEU:CD2	2.38	0.86
1:D:132:MET:HE1	1:D:208:VAL:HG12	1.55	0.86
1:C:102:ALA:HB3	1:C:195:VAL:HG21	1.59	0.84
1:B:5:VAL:HG21	1:B:85:LEU:HD22	1.62	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122:LEU:HD12	1:B:131:ALA:HB2	1.64	0.80
1:A:195:VAL:HG12	1:A:196:MET:HE3	1.62	0.79
1:A:230:LEU:HB2	1:A:234:GLU:OE1	1.82	0.79
1:B:228:ASN:HB2	1:B:230:LEU:CD2	2.12	0.79
1:A:220:LEU:HA	1:A:224:LEU:HD13	1.64	0.78
1:D:236:VAL:O	1:D:240:LYS:HG2	1.84	0.78
1:C:231:SER:OG	1:C:234:GLU:HG3	1.83	0.78
1:A:135:GLU:O	1:A:139:THR:HG23	1.84	0.77
1:D:127:THR:HG23	1:D:130:GLN:H	1.50	0.77
1:B:138:LEU:O	1:B:184:MET:HE3	1.85	0.77
1:C:111:LEU:HD12	1:C:111:LEU:H	1.49	0.76
1:B:233:SER:O	1:B:236:VAL:HG22	1.85	0.76
1:B:212:VAL:O	1:B:216:LEU:HD13	1.85	0.76
1:B:122:LEU:HD11	1:B:126:THR:HG22	1.68	0.76
1:D:1:SER:HB2	1:D:240:LYS:O	1.86	0.75
1:C:127:THR:OG1	1:C:130:GLN:HG3	1.86	0.75
1:C:221:TYR:HA	1:C:225:LEU:HB2	1.70	0.74
1:B:228:ASN:CB	1:B:230:LEU:CD2	2.66	0.73
1:D:206:PHE:O	1:D:210:PRO:HG2	1.89	0.72
1:A:41:THR:O	1:A:45:ILE:HD12	1.90	0.71
1:B:161:PRO:O	1:B:165:ILE:HG12	1.91	0.71
1:D:50:GLY:HA2	1:D:169:VAL:HG22	1.73	0.71
1:A:230:LEU:HD12	1:A:230:LEU:O	1.91	0.71
1:C:135:GLU:OE1	1:C:180:THR:HG21	1.91	0.71
1:B:95:LEU:O	1:B:99:ILE:HG13	1.90	0.71
1:C:232:LEU:O	1:C:236:VAL:HG13	1.91	0.70
1:D:231:SER:OG	1:D:234:GLU:HG3	1.91	0.70
1:C:134:VAL:O	1:C:138:LEU:HG	1.92	0.70
1:D:234:GLU:O	1:D:238:ILE:HG12	1.91	0.69
1:A:119:VAL:HG12	1:A:120:ASN:O	1.92	0.69
1:B:35:TRP:O	1:B:39:LEU:HD23	1.93	0.69
1:B:123:ASN:ND2	1:B:125:ASN:HB2	2.07	0.69
1:D:157:PRO:HA	2:D:266:PS6:O	1.92	0.68
1:B:138:LEU:HD23	3:B:276:HOH:O	1.92	0.68
1:D:34:LYS:HB3	1:D:39:LEU:HD21	1.74	0.68
1:A:78:VAL:HG11	1:A:215:VAL:HG13	1.76	0.68
1:A:139:THR:HG22	1:A:184:MET:HA	1.75	0.68
1:D:122:LEU:HD21	1:D:131:ALA:HB2	1.75	0.68
1:D:220:LEU:O	1:D:224:LEU:HB2	1.93	0.68
1:C:191:GLY:O	1:C:195:VAL:HG23	1.93	0.67
1:C:233:SER:O	1:C:236:VAL:HG22	1.95	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:THR:HG21	1:A:210:PRO:HA	1.75	0.67
1:B:220:LEU:HA	1:B:224:LEU:HB2	1.76	0.67
1:D:128:GLN:HE22	1:D:204:TRP:HE1	1.43	0.67
1:B:228:ASN:CG	1:B:230:LEU:HD21	2.13	0.67
1:C:122:LEU:HD21	1:C:131:ALA:HB2	1.77	0.66
1:A:231:SER:HB2	1:A:235:ARG:HH21	1.58	0.66
1:B:21:THR:O	1:B:25:VAL:HG23	1.94	0.66
1:B:120:ASN:HD21	1:B:188:ARG:HD3	1.60	0.66
1:B:182:CYS:HA	3:B:276:HOH:O	1.95	0.66
1:B:228:ASN:CB	1:B:230:LEU:HD23	2.25	0.66
1:D:173:HIS:O	1:D:177:ILE:HG13	1.96	0.66
1:B:100:ALA:O	1:B:104:ILE:HD12	1.96	0.66
1:C:109:ALA:HB3	3:C:284:HOH:O	1.96	0.65
1:D:123:ASN:HB3	1:D:126:THR:HB	1.77	0.65
1:C:140:PHE:O	1:C:144:LEU:HD23	1.97	0.64
1:D:119:VAL:HG12	1:D:192:PRO:HB2	1.80	0.64
1:B:204:TRP:O	1:B:208:VAL:HG13	1.97	0.64
1:A:4:GLU:HG3	3:A:278:HOH:O	1.96	0.64
1:C:116:ASN:HB2	1:C:118:ALA:H	1.63	0.64
1:B:231:SER:N	1:B:234:GLU:HB2	2.13	0.63
1:A:48:ALA:HB2	1:C:174:LEU:HD23	1.80	0.63
1:A:159:GLY:O	2:D:266:PS6:H14	1.99	0.63
1:C:159:GLY:H	2:D:266:PS6:H16	1.63	0.63
1:C:235:ARG:O	1:C:238:ILE:HB	1.99	0.62
1:C:180:THR:HG22	1:C:182:CYS:H	1.63	0.62
1:A:127:THR:HB	1:A:130:GLN:OE1	1.99	0.62
1:D:8:VAL:HG22	1:D:12:LYS:HE2	1.81	0.62
1:D:129:GLY:O	1:D:132:MET:HB3	1.99	0.62
1:A:123:ASN:HB3	1:A:126:THR:OG1	1.98	0.62
1:B:135:GLU:OE1	1:B:183:SER:HB3	2.00	0.62
1:C:207:TRP:O	1:C:211:ILE:HD12	2.00	0.62
1:D:128:GLN:NE2	1:D:204:TRP:HE1	1.96	0.62
1:A:5:VAL:HG23	3:A:278:HOH:O	2.00	0.61
1:B:127:THR:OG1	1:B:130:GLN:HG3	2.00	0.61
1:D:80:ASN:HD22	1:D:235:ARG:HH11	1.48	0.61
1:D:194:VAL:HG12	1:D:195:VAL:H	1.64	0.61
1:C:70:PRO:HG3	1:C:94:GLN:HE21	1.65	0.61
1:A:206:PHE:O	1:A:210:PRO:HG2	2.00	0.61
1:A:243:TYR:CZ	1:A:244:GLU:HG2	2.36	0.61
1:B:126:THR:HG23	1:B:130:GLN:HB2	1.83	0.61
1:B:47:LEU:HD12	1:D:43:LEU:HD21	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:134:VAL:O	1:B:138:LEU:HD13	2.00	0.61
1:A:196:MET:HE2	1:A:196:MET:HA	1.83	0.60
1:B:236:VAL:O	1:B:240:LYS:HD3	2.02	0.60
1:D:207:TRP:O	1:D:211:ILE:HD12	2.01	0.60
1:C:119:VAL:HG12	1:C:192:PRO:HB2	1.84	0.60
1:D:50:GLY:HA3	1:D:170:THR:HG22	1.83	0.60
1:C:91:VAL:HA	1:C:94:GLN:NE2	2.17	0.60
1:C:231:SER:HB3	1:C:234:GLU:OE1	2.01	0.60
1:C:123:ASN:HB3	1:C:126:THR:HB	1.83	0.59
1:B:127:THR:O	1:B:130:GLN:HB2	2.01	0.59
1:C:89:PHE:O	1:C:92:ALA:HB3	2.02	0.59
1:B:235:ARG:O	1:B:239:ILE:HG13	2.01	0.59
1:C:158:VAL:H	2:D:266:PS6:C17	2.16	0.59
1:D:239:ILE:HB	1:D:240:LYS:HE2	1.84	0.59
1:B:33:LEU:HB2	1:B:35:TRP:NE1	2.18	0.59
1:A:122:LEU:CD2	1:A:180:THR:HG22	2.33	0.59
1:C:230:LEU:O	1:C:234:GLU:HB2	2.02	0.59
1:D:119:VAL:CG1	1:D:192:PRO:HB2	2.33	0.58
1:D:132:MET:CE	1:D:208:VAL:HG12	2.31	0.58
1:A:50:GLY:HA2	1:A:169:VAL:HG13	1.85	0.58
1:C:116:ASN:HB2	1:C:118:ALA:N	2.18	0.58
1:C:22:LEU:HD21	1:C:104:ILE:HD13	1.85	0.58
1:A:10:PHE:O	1:A:14:VAL:HG23	2.02	0.58
1:B:228:ASN:HD22	1:B:230:LEU:HD21	1.62	0.58
2:D:266:PS6:H2A	2:D:266:PS6:H14A	1.85	0.58
1:A:134:VAL:O	1:A:138:LEU:HG	2.04	0.58
1:A:195:VAL:HG12	1:A:196:MET:CE	2.33	0.58
1:D:7:SER:HB2	1:D:10:PHE:H	1.69	0.57
1:B:243:TYR:OH	1:B:245:PRO:HB3	2.05	0.57
1:D:50:GLY:HA3	1:D:170:THR:CG2	2.35	0.57
1:C:33:LEU:O	1:C:40:PRO:HG2	2.04	0.57
1:C:122:LEU:HD21	1:C:131:ALA:CB	2.34	0.57
1:A:173:HIS:O	1:A:177:ILE:HG13	2.05	0.56
1:B:202:ALA:HA	1:B:204:TRP:NE1	2.20	0.56
1:B:70:PRO:O	1:B:74:LEU:HG	2.05	0.56
1:D:123:ASN:HB3	1:D:126:THR:CB	2.34	0.56
1:D:127:THR:CG2	1:D:130:GLN:H	2.18	0.56
1:B:47:LEU:HD11	2:D:266:PS6:C25	2.34	0.56
1:C:114:ARG:HD3	1:C:196:MET:HE1	1.86	0.56
1:A:33:LEU:O	1:A:40:PRO:HG2	2.05	0.56
1:B:126:THR:HG23	1:B:130:GLN:CB	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:PHE:O	1:A:227:PRO:HA	2.06	0.56
1:B:63:VAL:HG22	1:D:226:PHE:CZ	2.41	0.56
1:C:24:PHE:HB2	1:C:56:LEU:HD12	1.87	0.55
1:A:135:GLU:OE2	1:A:183:SER:HB3	2.07	0.55
1:C:5:VAL:HA	1:C:10:PHE:CD2	2.41	0.55
1:B:231:SER:H	1:B:234:GLU:HB2	1.71	0.54
1:B:10:PHE:O	1:B:14:VAL:HG23	2.07	0.54
1:C:230:LEU:HD11	1:C:238:ILE:HD11	1.89	0.54
1:D:123:ASN:O	1:D:124:ASN:HB2	2.07	0.54
1:B:208:VAL:O	1:B:212:VAL:HG23	2.08	0.54
1:C:13:ALA:O	1:C:16:ALA:HB3	2.06	0.54
1:C:198:ARG:O	1:C:198:ARG:HG2	2.07	0.54
1:B:42:ILE:HD13	1:B:178:TYR:CZ	2.42	0.54
1:D:136:LEU:HD21	1:D:216:LEU:HD12	1.89	0.54
1:D:122:LEU:HD21	1:D:131:ALA:CB	2.36	0.53
1:B:169:VAL:O	1:B:173:HIS:HD2	1.91	0.53
1:A:100:ALA:O	1:A:104:ILE:HD12	2.07	0.53
1:B:1:SER:HB3	1:B:4:GLU:HB3	1.90	0.53
1:B:160:SER:O	1:B:164:SER:HB2	2.09	0.53
1:B:74:LEU:HD12	1:B:211:ILE:HG12	1.90	0.53
1:A:234:GLU:O	1:A:238:ILE:HG12	2.08	0.53
1:D:25:VAL:HG21	1:D:98:ALA:HA	1.90	0.53
1:D:234:GLU:HB3	1:D:243:TYR:CE1	2.42	0.53
1:A:10:PHE:CE2	1:A:14:VAL:HG21	2.44	0.53
1:B:146:ILE:HA	1:B:165:ILE:HD11	1.91	0.53
1:B:236:VAL:C	1:B:240:LYS:HD3	2.33	0.53
1:C:50:GLY:HA3	1:C:170:THR:OG1	2.09	0.53
1:D:2:LYS:HD2	1:D:3:LYS:HE2	1.89	0.53
1:A:17:GLU:OE1	1:A:67:HIS:HB2	2.09	0.52
1:A:4:GLU:O	1:A:7:SER:OG	2.28	0.52
1:A:91:VAL:HA	1:A:94:GLN:NE2	2.24	0.52
1:B:146:ILE:O	1:B:150:THR:HG23	2.09	0.52
1:C:102:ALA:CB	1:C:195:VAL:HG21	2.37	0.52
1:B:238:ILE:HG23	1:B:243:TYR:CG	2.44	0.52
1:A:13:ALA:HB1	1:A:90:TYR:OH	2.10	0.52
1:B:126:THR:HG23	1:B:130:GLN:HE21	1.75	0.52
1:C:20:ALA:O	1:C:56:LEU:HD13	2.09	0.52
1:C:180:THR:HG22	1:C:182:CYS:N	2.24	0.52
1:D:2:LYS:O	1:D:6:CYS:HB2	2.10	0.52
1:A:51:LEU:HD22	1:C:167:LEU:HB3	1.90	0.52
1:D:182:CYS:SG	1:D:184:MET:HE2	2.48	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:195:VAL:O	1:A:196:MET:HE2	2.10	0.52
1:C:159:GLY:N	2:D:266:PS6:H16	2.25	0.52
1:C:110:PRO:HG2	1:C:113:ALA:CB	2.40	0.51
1:A:131:ALA:HB1	1:A:180:THR:CG2	2.40	0.51
1:C:10:PHE:O	1:C:14:VAL:HG23	2.10	0.51
1:A:70:PRO:HA	1:A:94:GLN:HE22	1.74	0.51
1:C:4:GLU:HB3	1:C:10:PHE:CD1	2.44	0.51
1:C:47:LEU:O	1:C:51:LEU:HG	2.10	0.51
1:D:204:TRP:HE3	1:D:208:VAL:HG21	1.74	0.51
1:D:244:GLU:HG3	1:D:245:PRO:HD2	1.93	0.51
1:B:122:LEU:HD11	1:B:126:THR:O	2.10	0.51
1:C:193:ALA:O	1:C:197:ASN:O	2.29	0.51
1:B:193:ALA:O	1:B:196:MET:O	2.29	0.51
1:C:106:TYR:HE1	1:C:114:ARG:NH1	2.09	0.51
1:D:70:PRO:O	1:D:74:LEU:HG	2.10	0.51
1:B:127:THR:H	1:B:130:GLN:NE2	2.09	0.51
1:C:80:ASN:HD21	1:C:235:ARG:HB3	1.76	0.51
1:D:192:PRO:O	1:D:194:VAL:O	2.28	0.51
1:D:232:LEU:O	1:D:236:VAL:HG23	2.11	0.50
1:A:3:LYS:HG3	1:A:4:GLU:N	2.26	0.50
1:A:225:LEU:O	1:A:227:PRO:HD3	2.11	0.50
1:B:128:GLN:O	1:B:131:ALA:HB3	2.11	0.50
1:A:220:LEU:HD11	1:D:15:PHE:CE2	2.47	0.50
1:C:70:PRO:HG3	1:C:94:GLN:NE2	2.26	0.50
1:D:234:GLU:HB3	1:D:243:TYR:HE1	1.76	0.50
1:A:139:THR:CG2	1:A:184:MET:HA	2.39	0.50
1:C:230:LEU:O	1:C:231:SER:O	2.30	0.50
1:C:173:HIS:HE1	3:C:271:HOH:O	1.94	0.50
1:D:140:PHE:CE2	1:D:144:LEU:HD22	2.46	0.50
1:A:122:LEU:HD11	1:A:128:GLN:HA	1.94	0.50
1:A:185:ASN:HD21	1:A:187:ALA:HB3	1.77	0.50
1:A:25:VAL:O	1:A:29:LEU:HG	2.12	0.49
1:D:8:VAL:HG22	1:D:12:LYS:CE	2.42	0.49
1:D:68:ILE:O	1:D:68:ILE:HG22	2.13	0.49
1:D:193:ALA:O	1:D:196:MET:O	2.30	0.49
1:A:39:LEU:HD23	3:A:291:HOH:O	2.12	0.49
1:A:166:GLY:O	1:A:169:VAL:HG13	2.12	0.49
1:B:159:GLY:HA2	2:D:266:PS6:H4	1.94	0.49
1:C:100:ALA:O	1:C:104:ILE:HD12	2.12	0.49
1:D:243:TYR:O	1:D:244:GLU:HB2	2.12	0.49
1:C:183:SER:O	1:C:184:MET:HB2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:LEU:HD11	1:D:15:PHE:HE2	1.78	0.49
1:B:24:PHE:HB2	1:B:56:LEU:HD12	1.94	0.49
1:B:52:ALA:O	1:B:56:LEU:HG	2.13	0.49
1:B:136:LEU:HD21	1:B:216:LEU:HD22	1.95	0.49
1:D:127:THR:HG22	1:D:130:GLN:CD	2.37	0.49
1:D:120:ASN:HD21	1:D:188:ARG:HD3	1.78	0.49
1:A:211:ILE:O	1:A:215:VAL:HG23	2.12	0.49
1:B:119:VAL:CG1	1:B:192:PRO:HB2	2.42	0.49
1:B:198:ARG:NH2	1:B:200:SER:HA	2.28	0.49
1:C:7:SER:HG	1:C:10:PHE:H	1.61	0.48
1:C:153:ARG:NH2	1:C:226:PHE:HB3	2.28	0.48
1:D:85:LEU:HD22	1:D:85:LEU:H	1.77	0.48
1:B:63:VAL:HG22	1:D:226:PHE:CE2	2.48	0.48
1:B:159:GLY:C	2:D:266:PS6:H4	2.38	0.48
1:B:236:VAL:O	1:B:239:ILE:HB	2.13	0.48
1:A:83:SER:HB3	3:A:281:HOH:O	2.12	0.48
1:D:237:ALA:HB1	1:D:243:TYR:N	2.28	0.48
1:A:226:PHE:CZ	1:D:63:VAL:HG22	2.49	0.48
1:B:158:VAL:HG12	1:B:158:VAL:O	2.13	0.48
1:C:163:LEU:HD13	2:D:266:PS6:H9	1.95	0.48
1:D:127:THR:HG22	1:D:130:GLN:CG	2.43	0.48
1:B:12:LYS:HE3	1:D:225:LEU:HA	1.96	0.48
1:C:25:VAL:O	1:C:29:LEU:HG	2.14	0.48
1:A:47:LEU:HD21	2:D:266:PS6:H24A	1.96	0.48
1:D:194:VAL:HG12	1:D:195:VAL:N	2.29	0.48
1:C:231:SER:CB	1:C:234:GLU:HG3	2.43	0.48
1:A:106:TYR:CE1	1:A:196:MET:HE1	2.48	0.47
1:B:127:THR:HG23	1:B:130:GLN:CD	2.39	0.47
1:A:85:LEU:O	1:A:89:PHE:HD2	1.97	0.47
1:C:158:VAL:N	2:D:266:PS6:H16	2.30	0.47
1:B:146:ILE:HA	1:B:165:ILE:CD1	2.44	0.47
1:A:225:LEU:C	1:A:227:PRO:HD3	2.40	0.47
1:B:222:PHE:O	1:B:227:PRO:HA	2.13	0.47
1:C:138:LEU:O	1:C:184:MET:HE3	2.14	0.47
1:B:230:LEU:CD2	1:B:230:LEU:H	2.27	0.47
1:B:231:SER:H	1:B:234:GLU:CD	2.23	0.47
1:D:220:LEU:HD12	1:D:224:LEU:HB2	1.97	0.47
1:A:81:GLN:H	1:A:238:ILE:HG21	1.80	0.47
1:B:47:LEU:HD13	1:D:43:LEU:HD11	1.96	0.47
1:A:34:LYS:O	1:A:35:TRP:O	2.33	0.47
1:B:80:ASN:HD22	1:B:80:ASN:HA	1.56	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85:LEU:HD23	1:B:89:PHE:HE1	1.80	0.46
1:B:243:TYR:CZ	1:B:245:PRO:HB3	2.50	0.46
1:C:46:ALA:HB1	1:C:174:LEU:HD12	1.97	0.46
1:A:123:ASN:OD1	1:A:125:ASN:HB2	2.15	0.46
1:B:201:PRO:O	1:B:202:ALA:HB3	2.15	0.46
1:C:80:ASN:HD21	1:C:235:ARG:CB	2.26	0.46
1:C:186:PRO:HB3	1:C:207:TRP:CD2	2.50	0.46
1:B:202:ALA:O	1:B:205:VAL:HG23	2.15	0.46
1:A:149:SER:HB3	1:A:161:PRO:HB3	1.96	0.46
1:A:185:ASN:ND2	1:A:187:ALA:HB3	2.30	0.46
1:D:2:LYS:H	1:D:2:LYS:HG3	1.54	0.46
1:B:165:ILE:O	1:B:168:SER:HB3	2.15	0.46
2:D:266:PS6:H2A	2:D:266:PS6:C14	2.46	0.46
1:B:224:LEU:HD12	1:B:224:LEU:HA	1.73	0.46
1:D:106:TYR:CE1	1:D:114:ARG:HD3	2.50	0.46
1:B:122:LEU:CD1	1:B:126:THR:HG22	2.40	0.46
1:A:70:PRO:HG3	1:A:94:GLN:HE21	1.81	0.46
1:A:81:GLN:O	1:A:81:GLN:NE2	2.49	0.45
1:A:82:ILE:HA	3:A:280:HOH:O	2.16	0.45
1:A:115:GLY:HA3	1:A:116:ASN:HB3	1.98	0.45
1:A:233:SER:O	1:A:233:SER:OG	2.30	0.45
1:B:125:ASN:ND2	3:B:279:HOH:O	2.49	0.45
1:A:232:LEU:O	1:A:234:GLU:N	2.49	0.45
1:C:110:PRO:HG2	1:C:113:ALA:HB2	1.98	0.45
1:D:85:LEU:HD22	1:D:85:LEU:N	2.31	0.45
1:B:68:ILE:O	1:B:69:ASN:HB2	2.17	0.45
1:D:90:TYR:O	1:D:94:GLN:HG3	2.17	0.45
1:D:127:THR:HG22	1:D:130:GLN:OE1	2.16	0.45
1:D:237:ALA:HB1	1:D:243:TYR:CA	2.46	0.45
1:B:128:GLN:O	1:B:131:ALA:N	2.50	0.45
1:D:122:LEU:HD23	1:D:122:LEU:HA	1.74	0.45
1:D:194:VAL:O	1:D:196:MET:N	2.50	0.45
1:A:241:GLY:HA2	3:A:281:HOH:O	2.16	0.45
1:B:196:MET:O	1:B:198:ARG:N	2.50	0.45
1:C:4:GLU:O	1:C:6:CYS:N	2.50	0.45
1:C:70:PRO:O	1:C:74:LEU:HG	2.16	0.45
1:B:236:VAL:HA	1:B:239:ILE:HG13	1.98	0.45
1:C:101:GLY:HA2	1:C:104:ILE:HD12	1.99	0.45
1:D:36:PRO:HG3	1:D:112:ASN:ND2	2.32	0.45
1:B:34:LYS:NZ	3:B:281:HOH:O	2.49	0.45
1:B:120:ASN:ND2	1:B:189:SER:OG	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:116:ASN:OD1	1:C:116:ASN:N	2.50	0.45
1:B:22:LEU:HD22	1:B:100:ALA:HB3	1.99	0.44
1:B:5:VAL:HG21	1:B:85:LEU:CD2	2.39	0.44
1:B:126:THR:HA	1:B:130:GLN:NE2	2.33	0.44
1:B:144:LEU:HD13	1:C:60:LEU:HD21	2.00	0.44
1:C:230:LEU:HD22	1:C:230:LEU:HA	1.67	0.44
1:A:24:PHE:HB2	1:A:56:LEU:HD12	1.99	0.44
1:B:67:HIS:ND1	1:B:90:TYR:HD2	2.15	0.44
1:B:117:LEU:HD23	1:B:192:PRO:HB3	1.99	0.44
1:A:104:ILE:O	1:A:108:VAL:HB	2.17	0.44
1:B:83:SER:HA	1:B:238:ILE:O	2.17	0.44
1:A:85:LEU:HD12	1:A:85:LEU:HA	1.88	0.44
1:A:231:SER:HA	1:A:232:LEU:HA	1.72	0.44
1:D:183:SER:O	1:D:184:MET:HB2	2.17	0.44
1:A:119:VAL:HG22	1:A:192:PRO:HB2	1.99	0.44
1:B:119:VAL:HG12	1:B:192:PRO:HB2	2.00	0.44
1:C:16:ALA:HB3	1:C:64:SER:HB3	1.99	0.44
1:D:123:ASN:O	1:D:123:ASN:OD1	2.36	0.44
1:D:134:VAL:HG21	1:D:179:PHE:CD1	2.53	0.44
1:B:17:GLU:OE2	1:B:64:SER:OG	2.29	0.43
1:C:116:ASN:CB	1:C:118:ALA:H	2.29	0.43
1:C:123:ASN:O	1:C:124:ASN:HB3	2.17	0.43
1:D:4:GLU:O	1:D:7:SER:OG	2.32	0.43
1:A:224:LEU:HD12	1:A:224:LEU:N	2.33	0.43
1:B:231:SER:N	1:B:234:GLU:OE2	2.50	0.43
1:A:230:LEU:HD13	1:A:234:GLU:CD	2.44	0.43
1:B:42:ILE:HD13	1:B:178:TYR:OH	2.18	0.43
1:C:230:LEU:HB3	1:C:235:ARG:CG	2.48	0.43
1:D:46:ALA:O	1:D:170:THR:HG22	2.18	0.43
1:B:20:ALA:O	1:B:23:ILE:N	2.49	0.43
1:B:198:ARG:NH1	1:B:200:SER:HA	2.33	0.43
1:B:120:ASN:HD21	1:B:188:ARG:CD	2.30	0.43
1:B:142:LEU:O	1:B:146:ILE:HG13	2.19	0.43
1:B:203:HIS:HA	1:B:206:PHE:CD2	2.54	0.43
1:B:238:ILE:HG23	1:B:243:TYR:CB	2.48	0.43
1:C:3:LYS:HB3	1:C:3:LYS:HE2	1.41	0.43
1:C:111:LEU:H	1:C:111:LEU:CD1	2.22	0.43
1:C:35:TRP:O	1:C:39:LEU:HD13	2.19	0.43
1:C:117:LEU:HB2	1:C:196:MET:SD	2.58	0.43
1:A:165:ILE:N	1:A:165:ILE:HD13	2.34	0.43
1:B:132:MET:SD	1:B:208:VAL:HG22	2.59	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:VAL:O	1:A:240:LYS:HG3	2.18	0.43
1:B:31:SER:HB2	3:B:274:HOH:O	2.19	0.42
1:D:80:ASN:ND2	1:D:235:ARG:HH11	2.13	0.42
1:D:132:MET:HA	1:D:205:VAL:HG13	2.00	0.42
1:B:135:GLU:CD	1:B:183:SER:HB3	2.44	0.42
1:B:156:SER:HA	1:B:157:PRO:HD3	1.91	0.42
1:A:47:LEU:HD23	1:C:43:LEU:HD11	2.01	0.42
1:B:132:MET:HA	1:B:205:VAL:HG13	2.01	0.42
1:B:238:ILE:H	1:B:238:ILE:HG12	1.72	0.42
1:C:208:VAL:O	1:C:212:VAL:HG23	2.19	0.42
1:B:2:LYS:HD2	1:B:3:LYS:NZ	2.34	0.42
1:C:24:PHE:HB2	1:C:56:LEU:CD1	2.50	0.42
1:C:72:ILE:HD12	1:C:146:ILE:HD12	2.00	0.42
1:C:132:MET:HG3	1:C:204:TRP:CH2	2.54	0.42
1:D:25:VAL:O	1:D:29:LEU:HG	2.19	0.42
1:D:238:ILE:HG12	1:D:238:ILE:H	1.65	0.42
1:C:122:LEU:HD22	1:C:126:THR:HG22	2.01	0.42
1:D:127:THR:HG22	1:D:130:GLN:HB2	2.01	0.42
1:A:138:LEU:O	1:A:184:MET:HE3	2.20	0.42
1:B:80:ASN:ND2	1:B:238:ILE:HG13	2.34	0.42
1:B:23:ILE:O	1:B:26:PHE:HB3	2.19	0.42
1:B:44:GLN:NE2	1:D:42:ILE:HG21	2.35	0.42
1:B:136:LEU:HD21	1:B:216:LEU:CD2	2.49	0.42
1:A:243:TYR:CE2	1:A:244:GLU:HG2	2.53	0.42
1:C:123:ASN:O	1:C:125:ASN:N	2.50	0.42
1:C:158:VAL:HG22	1:C:158:VAL:O	2.19	0.42
1:B:74:LEU:O	1:B:78:VAL:HG23	2.19	0.42
1:C:70:PRO:HB3	1:C:91:VAL:HG13	2.02	0.42
1:D:170:THR:O	1:D:174:LEU:HD22	2.20	0.42
1:A:163:LEU:HD23	1:A:163:LEU:HA	1.84	0.42
1:C:12:LYS:HA	1:C:12:LYS:HD2	1.71	0.42
1:C:91:VAL:HA	1:C:94:GLN:HE21	1.84	0.42
1:C:136:LEU:HA	1:C:209:GLY:O	2.20	0.42
1:C:46:ALA:CB	1:C:174:LEU:HD12	2.50	0.41
1:A:50:GLY:HA3	1:A:170:THR:OG1	2.19	0.41
1:A:127:THR:HG22	1:A:130:GLN:N	2.08	0.41
1:B:129:GLY:O	1:B:133:VAL:HG23	2.20	0.41
1:A:186:PRO:HB3	1:A:207:TRP:CE3	2.55	0.41
1:B:198:ARG:HA	1:B:198:ARG:HD2	1.59	0.41
1:C:122:LEU:HD12	1:C:202:ALA:CB	2.50	0.41
1:D:80:ASN:ND2	1:D:235:ARG:HG2	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:190:PHE:CD1	1:A:207:TRP:HZ2	2.37	0.41
1:C:190:PHE:CD1	1:C:207:TRP:HZ2	2.38	0.41
1:B:230:LEU:HD12	1:B:238:ILE:CD1	2.50	0.41
1:C:128:GLN:O	1:C:132:MET:HB2	2.19	0.41
1:C:220:LEU:O	1:C:224:LEU:HB2	2.21	0.41
1:D:46:ALA:CB	1:D:174:LEU:HD13	2.50	0.41
1:C:116:ASN:CG	1:C:118:ALA:H	2.28	0.41
1:C:122:LEU:HD12	1:C:202:ALA:HB2	2.03	0.41
1:D:204:TRP:CE3	1:D:208:VAL:HG21	2.55	0.41
1:C:7:SER:OG	1:C:9:ALA:HB3	2.19	0.41
1:D:102:ALA:HB3	1:D:195:VAL:HG21	2.02	0.41
1:A:94:GLN:HE21	1:A:94:GLN:HB2	1.57	0.41
1:A:186:PRO:HG3	1:A:207:TRP:HA	2.03	0.41
1:B:183:SER:O	1:B:184:MET:HB2	2.20	0.41
1:C:70:PRO:HA	1:C:91:VAL:HG22	2.02	0.41
1:C:188:ARG:O	1:C:192:PRO:HD2	2.20	0.41
1:C:243:TYR:CZ	1:C:245:PRO:HB3	2.55	0.41
1:D:47:LEU:HD11	2:D:266:PS6:H25A	2.03	0.41
1:A:236:VAL:HG23	1:A:240:LYS:HD2	2.03	0.41
1:B:114:ARG:NE	3:B:266:HOH:O	2.50	0.41
1:B:144:LEU:HD23	1:B:144:LEU:HA	1.92	0.41
1:B:50:GLY:HA3	1:B:170:THR:OG1	2.21	0.40
1:C:132:MET:HE3	1:C:132:MET:HB3	1.81	0.40
1:D:80:ASN:HD21	1:D:235:ARG:HG2	1.86	0.40
1:B:198:ARG:CZ	1:B:200:SER:HA	2.51	0.40
1:D:8:VAL:O	1:D:12:LYS:HG3	2.21	0.40
1:D:144:LEU:HD12	1:D:144:LEU:HA	1.97	0.40
1:D:184:MET:HE3	3:D:288:HOH:O	2.20	0.40
1:A:70:PRO:CA	1:A:94:GLN:HE22	2.34	0.40
1:A:101:GLY:HA2	1:A:104:ILE:HD12	2.03	0.40
1:B:70:PRO:HB2	1:B:186:PRO:HG2	2.04	0.40
1:B:127:THR:HG23	1:B:130:GLN:HG3	2.02	0.40
1:B:238:ILE:HD12	1:B:238:ILE:HG21	1.90	0.40
1:B:244:GLU:HA	1:B:245:PRO:HD3	1.84	0.40
1:D:49:PHE:O	1:D:52:ALA:HB3	2.21	0.40

There are no symmetry-related clashes.

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	243/266 (91%)	222 (91%)	16 (7%)	5 (2%)	5	2
1	B	243/266 (91%)	221 (91%)	19 (8%)	3 (1%)	10	6
1	C	241/266 (91%)	213 (88%)	17 (7%)	11 (5%)	2	0
1	D	243/266 (91%)	223 (92%)	16 (7%)	4 (2%)	7	3
All	All	970/1064 (91%)	879 (91%)	68 (7%)	23 (2%)	4	2

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	116	ASN
1	A	231	SER
1	A	233	SER
1	C	5	VAL
1	C	116	ASN
1	C	198	ARG
1	C	230	LEU
1	C	231	SER
1	A	81	GLN
1	B	182	CYS
1	C	123	ASN
1	D	37	SER
1	B	121	ALA
1	B	154	ARG
1	C	197	ASN
1	D	195	VAL
1	A	35	TRP
1	C	4	GLU
1	D	244	GLU
1	C	182	CYS
1	C	244	GLU
1	C	115	GLY
1	D	194	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	188/209 (90%)	159 (85%)	29 (15%)	2	1
1	B	188/209 (90%)	151 (80%)	37 (20%)	1	0
1	C	186/209 (89%)	160 (86%)	26 (14%)	3	2
1	D	188/209 (90%)	168 (89%)	20 (11%)	6	4
All	All	750/836 (90%)	638 (85%)	112 (15%)	3	1

All (112) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	3	LYS
1	A	4	GLU
1	A	7	SER
1	A	37	SER
1	A	39	LEU
1	A	72	ILE
1	A	80	ASN
1	A	81	GLN
1	A	84	LEU
1	A	85	LEU
1	A	108	VAL
1	A	114	ARG
1	A	122	LEU
1	A	124	ASN
1	A	127	THR
1	A	132	MET
1	A	155	THR
1	A	165	ILE
1	A	169	VAL
1	A	197	ASN
1	A	200	SER
1	A	203	HIS
1	A	220	LEU

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Mol	Chain	Res	Type
1	A	229	SER
1	A	231	SER
1	A	233	SER
1	A	236	VAL
1	A	244	GLU
1	B	2	LYS
1	B	7	SER
1	B	11	LEU
1	B	22	LEU
1	B	34	LYS
1	B	37	SER
1	B	42	ILE
1	B	58	GLN
1	B	68	ILE
1	B	80	ASN
1	B	81	GLN
1	B	83	SER
1	B	85	LEU
1	B	108	VAL
1	B	111	LEU
1	B	114	ARG
1	B	117	LEU
1	B	122	LEU
1	B	123	ASN
1	B	126	THR
1	B	127	THR
1	B	137	ILE
1	B	138	LEU
1	B	156	SER
1	B	164	SER
1	B	196	MET
1	B	197	ASN
1	B	198	ARG
1	B	208	VAL
1	B	224	LEU
1	B	230	LEU
1	B	232	LEU
1	B	233	SER
1	B	238	ILE
1	B	239	ILE
1	B	240	LYS
1	B	244	GLU

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Mol	Chain	Res	Type
1	C	3	LYS
1	C	7	SER
1	C	8	VAL
1	C	11	LEU
1	C	12	LYS
1	C	34	LYS
1	C	37	SER
1	C	39	LEU
1	C	60	LEU
1	C	68	ILE
1	C	116	ASN
1	C	123	ASN
1	C	126	THR
1	C	132	MET
1	C	164	SER
1	C	180	THR
1	C	196	MET
1	C	197	ASN
1	C	211	ILE
1	C	216	LEU
1	C	224	LEU
1	C	229	SER
1	C	230	LEU
1	C	231	SER
1	C	236	VAL
1	C	244	GLU
1	D	2	LYS
1	D	7	SER
1	D	14	VAL
1	D	34	LYS
1	D	37	SER
1	D	39	LEU
1	D	62	PRO
1	D	111	LEU
1	D	114	ARG
1	D	155	THR
1	D	169	VAL
1	D	174	LEU
1	D	197	ASN
1	D	224	LEU
1	D	229	SER
1	D	230	LEU

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Mol	Chain	Res	Type
1	D	233	SER
1	D	238	ILE
1	D	242	THR
1	D	244	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	94	GLN
1	A	125	ASN
1	A	173	HIS
1	A	185	ASN
1	A	197	ASN
1	B	44	GLN
1	B	81	GLN
1	B	120	ASN
1	B	124	ASN
1	B	128	GLN
1	B	130	GLN
1	B	173	HIS
1	C	67	HIS
1	C	94	GLN
1	C	173	HIS
1	D	58	GLN
1	D	80	ASN
1	D	81	GLN
1	D	120	ASN
1	D	128	GLN
1	D	173	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 oligosaccharides in this entry.

5.6 Ligand geometry

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PS6	D	266	-	36,37,37	0.63	0	38,44,44	1.36	5 (13%)

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	PS6	D	266	-	-	27/43/43/43	-

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	266	PS6	O1-CB-CA	5.37	112.75	108.06
2	D	266	PS6	O52-C5-C6	2.66	119.95	111.83
2	D	266	PS6	O11-C1-C13	2.48	116.85	111.48
2	D	266	PS6	O52-C5-O51	-2.38	117.67	123.63
2	D	266	PS6	C3-O11-C1	-2.16	112.64	117.80

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	266	PS6	C2-O2-P-O1

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Mol	Chain	Res	Type	Atoms
2	D	266	PS6	CB-O1-P-O2
2	D	266	PS6	CB-O1-P-O4
2	D	266	PS6	C13-C1-O11-C3
2	D	266	PS6	N-CA-CB-O1
2	D	266	PS6	C-CA-CB-O1
2	D	266	PS6	O-C-CA-CB
2	D	266	PS6	OXT-C-CA-CB
2	D	266	PS6	O12-C1-O11-C3
2	D	266	PS6	C1-C13-C14-C15
2	D	266	PS6	C18-C19-C20-C21
2	D	266	PS6	C7-C8-C9-C10
2	D	266	PS6	C9-C10-C18-C19
2	D	266	PS6	O2-C2-C3-C4
2	D	266	PS6	C2-C3-O11-C1
2	D	266	PS6	C21-C22-C23-C24
2	D	266	PS6	O11-C1-C13-C14
2	D	266	PS6	O2-C2-C3-O11
2	D	266	PS6	C2-O2-P-O4
2	D	266	PS6	CB-O1-P-O3
2	D	266	PS6	C22-C23-C24-C25
2	D	266	PS6	C13-C14-C15-C16
2	D	266	PS6	O-C-CA-N
2	D	266	PS6	CA-CB-O1-P
2	D	266	PS6	O11-C3-C4-O52
2	D	266	PS6	O12-C1-C13-C14
2	D	266	PS6	OXT-C-CA-N

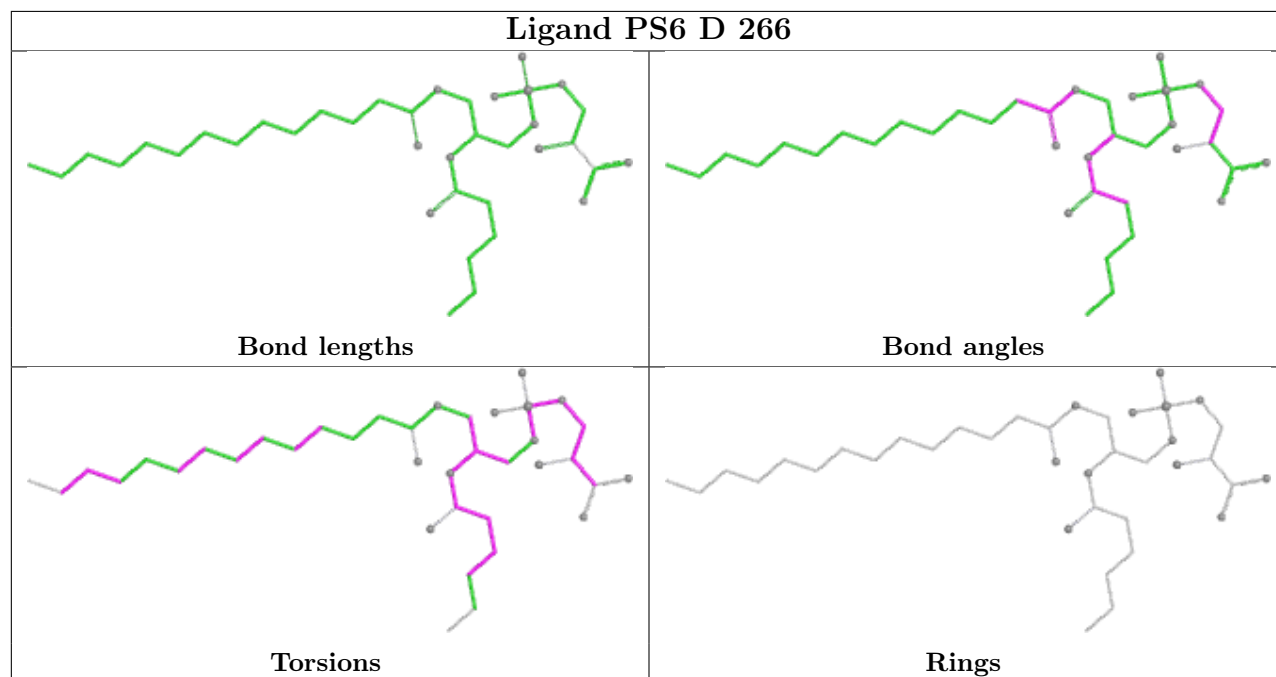
There are no ring outliers.

1 monomer is involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	266	PS6	16	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and

any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	245/266 (92%)	-1.00	0 100 100	18, 31, 68, 90	1 (0%)
1	B	245/266 (92%)	-0.95	0 100 100	19, 36, 72, 103	1 (0%)
1	C	243/266 (91%)	-0.94	0 100 100	18, 34, 57, 118	0
1	D	245/266 (92%)	-1.12	0 100 100	16, 29, 59, 104	1 (0%)
All	All	978/1064 (91%)	-1.00	0 100 100	16, 33, 66, 118	3 (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 oligosaccharides in this entry.

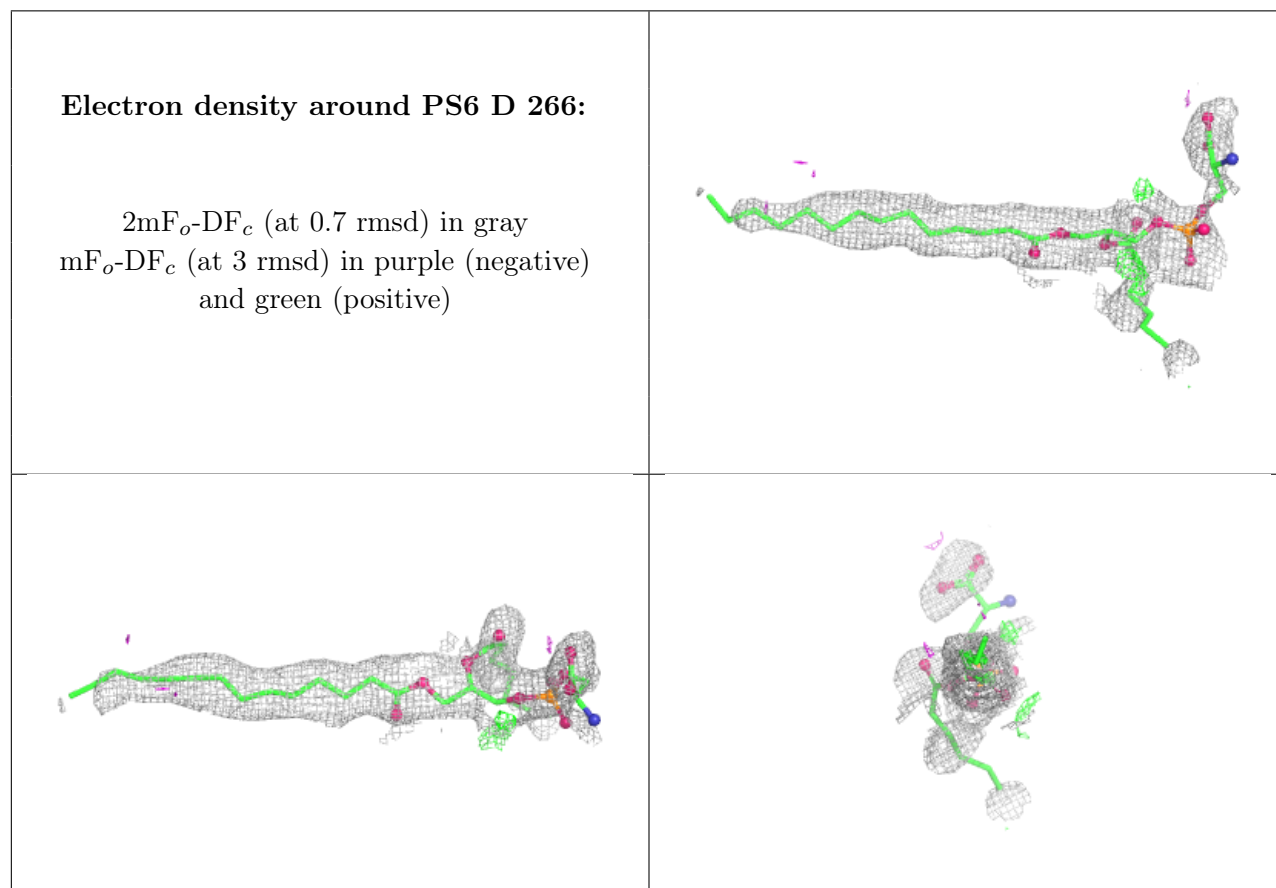
6.4 Ligands [i](#)

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
2	PS6	D	266	38/38	0.98	0.08	26,70,125,154	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.