



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 12:34 AM UTC

PDB ID : 1QI9 / pdb_00001qi9
Title : X-RAY SIRAS STRUCTURE DETERMINATION OF A VANADIUM-DEPENDENT HALOPEROXIDASE FROM ASCOPHYLLUM NODOSUM AT 2.0 Å RESOLUTION
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Deposited on : 1999-06-10
Resolution : 2.05 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

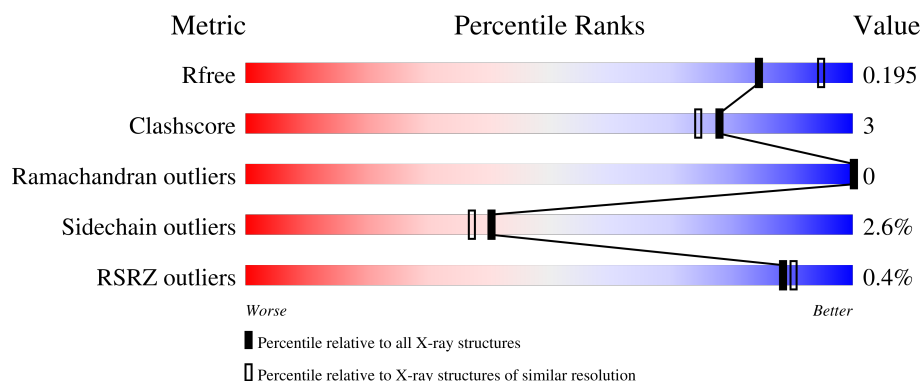
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	556	 86% 12% .
2	B	556	 86% 12% .

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	VO4	A	600	-	-	X	-
3	VO4	B	600	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 10264 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.

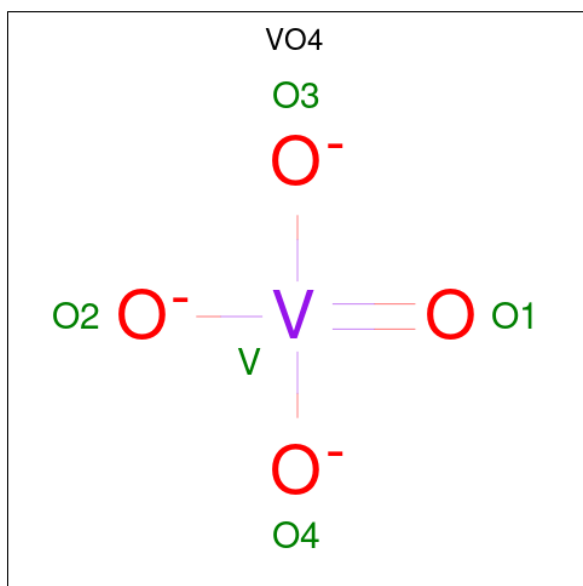
- Molecule 1 is a protein called Vanadium-dependent bromoperoxidase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	555	Total	C	I	N	O	S	0	0	0
			4244	2682	3	709	836	14			

- Molecule 2 is a protein called Vanadium-dependent bromoperoxidase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	556	Total	C	I	N	O	S	0	0	0
			4251	2687	2	710	838	14			

- Molecule 3 is VANADATE ION (CCD ID: VO4) (formula: O₄V).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	V	0	0
			5	4	1		
3	B	1	Total	O	V	0	0
			5	4	1		

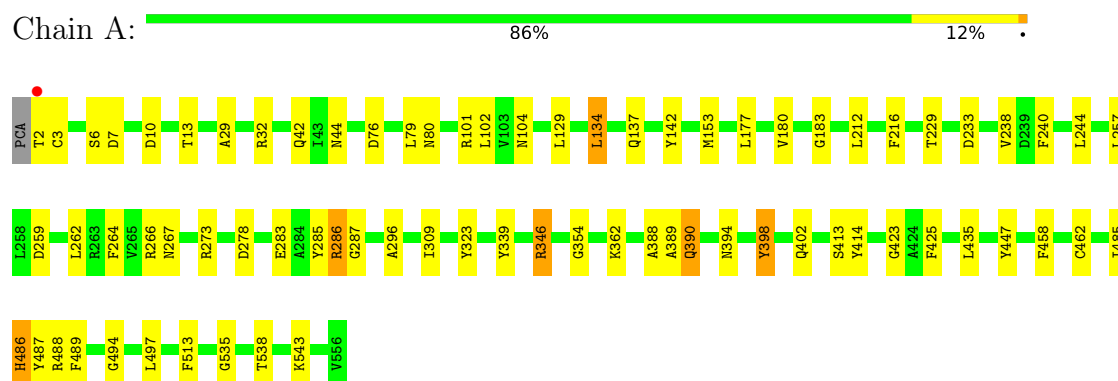
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	877	Total 877	O 877	0	0
4	B	882	Total 882	O 882	0	0

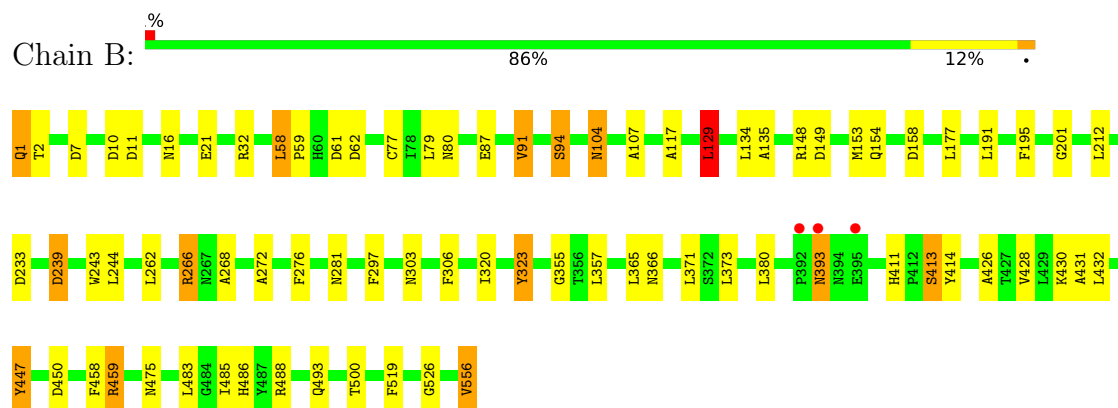
3 Residue-property plots

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: Vanadium-dependent bromoperoxidase



- Molecule 2: Vanadium-dependent bromoperoxidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	113.19Å 113.19Å 272.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.00 – 2.05 31.00 – 2.05	Depositor EDS
% Data completeness (in resolution range)	95.0 (31.00-2.05) 94.7 (31.00-2.05)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.57 (at 2.05Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.165 , 0.219 0.151 , 0.195	Depositor DCC
R_{free} test set	5320 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	7.0	Xtriage
Anisotropy	0.978	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 75.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10264	wwPDB-VP
Average B, all atoms (Å ²)	9.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.66% 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: IYR, PCA, VO4, TYI

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	0.89	0/4308	1.65	54/5869 (0.9%)
2	B	0.89	1/4323 (0.0%)	1.65	62/5893 (1.1%)
All	All	0.89	1/8631 (0.0%)	1.65	116/11762 (1.0%)

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
2	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	233	ASP	C-O	5.71	1.31	1.24

All (116) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	THR	N-CA-CB	-11.54	91.89	111.50
1	A	323	TYR	CA-CB-CG	9.70	131.36	113.90
1	A	286	ARG	NE-CZ-NH2	-9.34	110.80	119.20
2	B	104	ASN	O-C-N	-9.32	117.24	121.53
2	B	485	ILE	N-CA-C	9.18	120.89	112.17
2	B	323	TYR	CA-CB-CG	8.26	128.77	113.90
1	A	390	GLN	CA-CB-CG	8.04	130.18	114.10
1	A	44	ASN	CA-CB-CG	7.78	120.38	112.60
2	B	475	ASN	CA-CB-CG	7.61	120.21	112.60
1	A	296	ALA	N-CA-C	-7.42	102.56	111.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	101	ARG	CD-NE-CZ	7.37	134.72	124.40
2	B	306	PHE	CA-CB-CG	-7.26	106.54	113.80
1	A	489	PHE	CA-CB-CG	7.22	121.02	113.80
2	B	413	SER	N-CA-C	7.20	119.20	111.36
1	A	264	PHE	CA-CB-CG	7.13	120.93	113.80
2	B	80	ASN	N-CA-C	7.10	122.11	113.17
2	B	32	ARG	NE-CZ-NH1	-7.09	114.41	121.50
1	A	497	LEU	CA-C-N	6.98	127.69	119.94
1	A	497	LEU	C-N-CA	6.98	127.69	119.94
2	B	104	ASN	N-CA-CB	-6.91	103.99	111.66
1	A	285	TYR	CA-C-O	6.71	127.53	120.42
2	B	2	THR	CA-CB-CG2	6.67	121.84	110.50
2	B	500	THR	CA-C-N	6.66	128.96	120.56
2	B	500	THR	C-N-CA	6.66	128.96	120.56
2	B	80	ASN	CA-CB-CG	-6.65	105.95	112.60
1	A	29	ALA	CA-C-N	6.55	129.06	120.28
1	A	29	ALA	C-N-CA	6.55	129.06	120.28
2	B	32	ARG	NH1-CZ-NH2	6.42	127.65	119.30
2	B	366	ASN	CA-CB-CG	-6.36	106.24	112.60
2	B	10	ASP	CA-CB-CG	6.32	118.92	112.60
1	A	287	GLY	CA-C-N	6.31	128.64	120.44
1	A	287	GLY	C-N-CA	6.31	128.64	120.44
1	A	402	GLN	O-C-N	-6.31	115.81	123.19
2	B	556	VAL	N-CA-CB	-6.30	100.78	111.50
1	A	486	HIS	CE1-NE2-CD2	-6.28	102.72	109.00
1	A	323	TYR	N-CA-CB	-6.23	100.82	109.91
1	A	80	ASN	N-CA-C	6.18	120.83	113.16
2	B	526	GLY	N-CA-C	6.18	122.16	114.37
2	B	380	LEU	CA-C-O	-6.16	114.02	120.55
2	B	7	ASP	CA-C-O	-6.16	114.05	121.32
2	B	117	ALA	CA-C-N	6.12	129.09	120.28
2	B	117	ALA	C-N-CA	6.12	129.09	120.28
1	A	286	ARG	CB-CG-CD	-6.11	97.24	111.30
1	A	494	GLY	N-CA-C	-6.10	104.96	112.77
1	A	339	TYR	O-C-N	6.10	128.59	122.12
2	B	519	PHE	CA-CB-CG	6.07	119.87	113.80
2	B	459	ARG	CD-NE-CZ	6.02	132.83	124.40
2	B	428	VAL	O-C-N	5.98	127.77	121.91
1	A	413	SER	N-CA-C	5.96	117.85	111.36
2	B	195	PHE	CA-CB-CG	5.93	119.73	113.80
2	B	281	ASN	CA-CB-CG	5.93	118.53	112.60
2	B	129	LEU	CA-CB-CG	5.89	136.90	116.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	13	THR	CA-C-N	5.87	123.94	119.66
1	A	13	THR	C-N-CA	5.87	123.94	119.66
2	B	94	SER	N-CA-C	5.87	120.43	113.16
1	A	240	PHE	CA-CB-CG	-5.84	107.96	113.80
1	A	134	LEU	N-CA-CB	5.83	118.69	110.12
1	A	423	GLY	N-CA-C	-5.82	105.32	112.77
2	B	426	ALA	CA-C-O	5.80	126.69	120.55
1	A	414	TYR	CB-CG-CD1	-5.78	112.12	120.80
1	A	267	ASN	CA-CB-CG	5.77	118.37	112.60
1	A	513	PHE	CA-CB-CG	5.72	119.52	113.80
1	A	414	TYR	N-CA-C	5.72	118.57	109.64
1	A	462	CYS	CA-C-O	-5.71	114.73	120.96
2	B	366	ASN	CB-CG-ND2	-5.71	107.84	116.40
2	B	91	VAL	CB-CA-C	-5.71	102.12	110.62
1	A	354	GLY	CA-C-N	5.70	126.15	119.94
1	A	354	GLY	C-N-CA	5.70	126.15	119.94
2	B	79	LEU	N-CA-C	5.63	117.50	111.36
1	A	259	ASP	CA-CB-CG	5.58	118.18	112.60
2	B	11	ASP	CA-C-N	5.57	125.41	119.28
2	B	11	ASP	C-N-CA	5.57	125.41	119.28
2	B	107	ALA	N-CA-C	5.57	118.06	111.33
1	A	425	PHE	N-CA-CB	5.56	118.39	110.16
2	B	414	TYR	N-CA-C	5.55	118.30	109.64
1	A	102	LEU	CA-C-O	-5.54	114.61	120.92
1	A	104	ASN	CA-CB-CG	5.51	118.11	112.60
2	B	148	ARG	NE-CZ-NH2	-5.49	114.26	119.20
1	A	487	TYR	CA-CB-CG	5.48	123.76	113.90
2	B	21	GLU	CA-C-O	-5.47	114.75	120.55
1	A	389	ALA	CA-C-N	5.44	129.88	120.58
1	A	389	ALA	C-N-CA	5.44	129.88	120.58
2	B	268	ALA	CA-C-O	-5.43	114.79	120.55
1	A	42	GLN	OE1-CD-NE2	-5.41	117.19	122.60
2	B	266	ARG	CD-NE-CZ	5.38	131.94	124.40
2	B	519	PHE	N-CA-CB	5.38	120.18	110.83
2	B	320	ILE	N-CA-C	5.37	117.76	111.05
2	B	493	GLN	OE1-CD-NE2	-5.35	117.25	122.60
1	A	10	ASP	CA-CB-CG	5.34	117.94	112.60
1	A	233	ASP	N-CA-C	5.32	122.14	110.80
2	B	16	ASN	OD1-CG-ND2	-5.32	117.28	122.60
2	B	94	SER	CA-CB-OG	-5.31	100.48	111.10
2	B	149	ASP	CA-CB-CG	-5.29	107.31	112.60
2	B	158	ASP	CA-CB-CG	-5.29	107.31	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	7	ASP	CA-CB-CG	5.27	117.87	112.60
2	B	303	ASN	CA-CB-CG	5.26	117.86	112.60
1	A	180	VAL	O-C-N	-5.26	117.31	122.67
1	A	346	ARG	CB-CA-C	-5.22	104.17	111.80
2	B	149	ASP	CA-C-N	5.22	126.35	122.59
2	B	149	ASP	C-N-CA	5.22	126.35	122.59
1	A	435	LEU	N-CA-C	5.21	116.64	111.07
2	B	201	GLY	CA-C-O	5.18	124.53	119.04
2	B	276	PHE	CA-CB-CG	5.17	118.97	113.80
1	A	104	ASN	CB-CG-ND2	5.15	124.13	116.40
1	A	79	LEU	N-CA-C	5.15	116.70	111.14
1	A	142	TYR	O-C-N	-5.11	116.81	122.07
2	B	365	LEU	CA-C-O	-5.10	115.66	121.58
2	B	239	ASP	CA-CB-CG	5.09	117.69	112.60
2	B	117	ALA	CA-C-O	-5.05	116.31	121.87
1	A	485	ILE	N-CA-CB	-5.05	105.74	112.26
2	B	411	HIS	CA-CB-CG	5.05	118.85	113.80
2	B	431	ALA	O-C-N	5.04	127.46	122.12
2	B	297	PHE	CA-CB-CG	-5.04	108.76	113.80
2	B	430	LYS	N-CA-C	5.03	116.77	111.28
2	B	393	ASN	CA-CB-CG	-5.03	107.57	112.60
1	A	183	GLY	N-CA-C	-5.02	106.12	112.54

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	1	PCA	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	4244	0	4068	22	0
2	B	4251	0	4077	20	0
3	A	5	0	0	4	0
3	B	5	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	877	0	0	1	0
4	B	882	0	0	4	0
All	All	10264	0	8145	47	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 3.

All (47) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:362:LYS:NZ	1:A:398:IYR:IE	2.86	0.75
1:A:447:TYI:I2	1:A:458:PHE:HA	2.63	0.69
2:B:459:ARG:HD3	4:B:1318:HOH:O	1.91	0.69
1:A:153:MET:HG3	1:A:447:TYI:I1	2.65	0.67
3:A:600:VO4:V	3:A:600:VO4:O3	1.52	0.66
3:B:600:VO4:O1	3:B:600:VO4:V	1.54	0.63
1:A:216:PHE:CE1	1:A:286:ARG:HD3	2.35	0.62
3:B:600:VO4:V	3:B:600:VO4:O3	1.57	0.61
3:A:600:VO4:V	3:A:600:VO4:O1	1.59	0.58
2:B:134:LEU:O	2:B:134:LEU:HG	1.99	0.58
2:B:153:MET:HG3	2:B:447:TYI:I1	2.73	0.58
2:B:129:LEU:HA	2:B:134:LEU:HD22	1.84	0.58
2:B:58:LEU:HB3	2:B:59:PRO:CD	2.34	0.57
2:B:62:ASP:HB2	4:B:1458:HOH:O	2.05	0.56
2:B:153:MET:HE3	4:B:1002:HOH:O	2.06	0.55
1:A:278:ASP:OD2	1:A:283:GLU:OE2	2.25	0.54
1:A:134:LEU:HG	1:A:134:LEU:O	2.07	0.54
1:A:129:LEU:HD23	1:A:129:LEU:C	2.34	0.53
1:A:362:LYS:HG2	1:A:398:IYR:IE	2.79	0.53
1:A:2:THR:HA	4:A:1269:HOH:O	2.08	0.53
2:B:243:TRP:CZ2	2:B:355:GLY:HA3	2.44	0.52
2:B:486:HIS:NE2	3:B:600:VO4:O1	2.42	0.52
1:A:3:CYS:HB2	1:A:6:SER:O	2.11	0.51
1:A:538:THR:HG22	1:A:543:LYS:HA	1.93	0.50
1:A:177:LEU:HB3	1:A:488:ARG:HD2	1.93	0.50
1:A:346:ARG:HH11	1:A:346:ARG:HG3	1.76	0.50
3:A:600:VO4:O3	3:A:600:VO4:O4	2.30	0.49
2:B:154:GLN:HG2	4:B:1388:HOH:O	2.12	0.49
1:A:486:HIS:NE2	3:A:600:VO4:O3	2.47	0.48
1:A:238:VAL:HG21	1:A:390:GLN:HG2	1.96	0.48
2:B:447:TYI:I2	2:B:458:PHE:HA	2.85	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:THR:HA	1:A:273:ARG:HG2	1.97	0.46
2:B:486:HIS:NE2	3:B:600:VO4:O3	2.49	0.45
1:A:137:GLN:HG2	1:A:177:LEU:HD22	1.99	0.45
1:A:388:ALA:O	1:A:394:ASN:HA	2.18	0.44
2:B:323:TYR:HA	2:B:432:LEU:HD13	2.01	0.42
1:A:32:ARG:NH2	1:A:535:GLY:HA3	2.34	0.42
1:A:346:ARG:HG3	1:A:346:ARG:NH1	2.34	0.42
2:B:135:ALA:HB1	2:B:373:LEU:HB2	2.02	0.42
2:B:272:ALA:HB1	2:B:483:LEU:HD23	2.03	0.41
2:B:154:GLN:NE2	2:B:450:ASP:OD2	2.47	0.41
3:B:600:VO4:O3	3:B:600:VO4:O4	2.38	0.41
1:A:238:VAL:CG2	1:A:390:GLN:HG2	2.51	0.41
2:B:58:LEU:HB3	2:B:59:PRO:HD2	2.02	0.41
2:B:177:LEU:HB3	2:B:488:ARG:HD2	2.02	0.41
1:A:309:ILE:HB	2:B:87:GLU:CD	2.46	0.40
2:B:61:ASP:OD1	2:B:61:ASP:C	2.65	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	551/556 (99%)	536 (97%)	15 (3%)	0	100	100
2	B	553/556 (100%)	540 (98%)	13 (2%)	0	100	100
All	All	1104/1112 (99%)	1076 (98%)	28 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	449/449 (100%)	443 (99%)	6 (1%)	61	63
2	B	450/450 (100%)	433 (96%)	17 (4%)	29	24
All	All	899/899 (100%)	876 (97%)	23 (3%)	40	37

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	ASP
1	A	212	LEU
1	A	244	LEU
1	A	257	LEU
1	A	262	LEU
1	A	266	ARG
2	B	58	LEU
2	B	77	CYS
2	B	91	VAL
2	B	94	SER
2	B	104	ASN
2	B	129	LEU
2	B	191	LEU
2	B	212	LEU
2	B	239	ASP
2	B	244	LEU
2	B	262	LEU
2	B	266	ARG
2	B	357	LEU
2	B	371	LEU
2	B	393	ASN
2	B	413	SER
2	B	556	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	GLN
1	A	366	ASN
1	A	402	GLN
2	B	16	ASN
2	B	402	GLN
2	B	493	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
2	PCA	B	1	2	7,8,9	1.23	1 (14%)	9,10,12	1.87	1 (11%)
1	TYI	A	447	1	13,14,15	0.85	0	14,19,21	0.77	0
2	TYI	B	447	2	13,14,15	1.10	1 (7%)	14,19,21	0.94	1 (7%)
1	IYR	A	398	1	12,13,14	0.83	0	12,17,19	1.82	3 (25%)

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	PCA	B	1	2	-	0/0/11/13	0/1/1/1
1	TYI	A	447	1	-	0/5/6/8	0/1/1/1
2	TYI	B	447	2	-	1/5/6/8	0/1/1/1
1	IYR	A	398	1	-	1/5/6/8	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	447	TYI	CE2-I2	-2.04	2.05	2.10
2	B	1	PCA	O-C	2.03	1.27	1.20

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	PCA	O-C-CA	-4.72	112.64	124.77
1	A	398	IYR	CH-CG-CF	-4.00	116.50	120.50
1	A	398	IYR	CG-CF-CE	2.85	121.96	119.27
1	A	398	IYR	CB-CC-CD	-2.55	116.06	120.43
2	B	447	TYI	CD2-CG-CD1	-2.08	116.19	119.01

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	398	IYR	O-C-CA-CB
2	B	447	TYI	O-C-CA-CB

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	447	TYI	2	0
2	B	447	TYI	2	0
1	A	398	IYR	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	VO4	A	600	1	0,4,4	-	-	-		
3	VO4	B	600	2	0,4,4	-	-	-		

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	600	VO4	4	0
3	B	600	VO4	5	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	553/556 (99%)	-0.83	1 (0%) 91 93	2, 6, 16, 38	0
2	B	554/556 (99%)	-0.83	3 (0%) 87 89	2, 6, 14, 36	0
All	All	1107/1112 (99%)	-0.83	4 (0%) 88 90	2, 6, 15, 38	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	395	GLU	3.9
2	B	392	PRO	3.5
2	B	393	ASN	2.6
1	A	2	THR	2.5

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	TYI	A	447	14/15	0.96	0.08	6,16,20,24	2
1	IYR	A	398	13/14	0.97	0.05	3,7,11,12	1
2	PCA	B	1	8/9	0.97	0.06	5,7,8,10	0
2	TYI	B	447	14/15	0.97	0.08	6,16,19,21	2

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
3	VO4	A	600	5/5	0.99	0.05	2,2,3,3	0
3	VO4	B	600	5/5	0.99	0.05	2,2,5,6	0

6.5 Other polymers [i](#)

There are no such residues in this entry.