



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:53 AM UTC

PDB ID : 2AEL / pdb_00002ael
Title : R304K Trichodiene Synthase: Complex With Mg, Pyrophosphate, and (4R)-7-Azabisabolene
Authors : Vedula, L.S.; Cane, D.E.; Christianson, D.W.
Deposited on : 2005-07-22
Resolution : 2.50 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

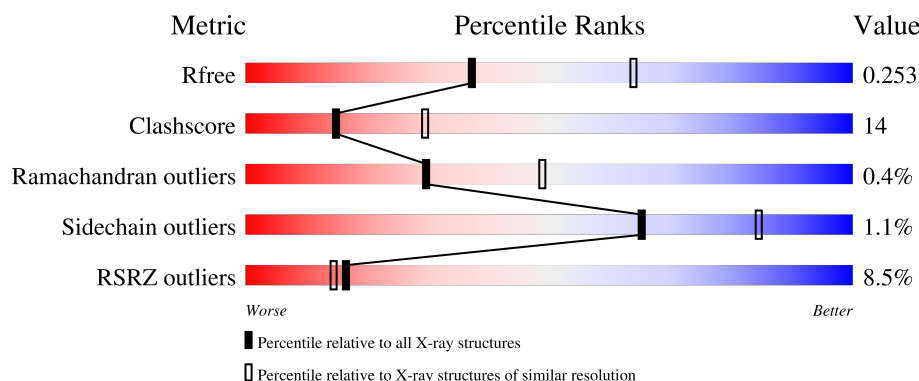
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	374	 2% 65% 28% • 6%
1	B	374	 13% 70% 23% • 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
2	EDO	A	750	-	X	-	-
4	SAZ	B	704	X	-	-	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6073 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 Trichodiene synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	352	Total	C	N	O	S	54	0	0
			2921	1872	489	543	17			
1	B	353	Total	C	N	O	S	150	0	0
			2930	1877	490	546	17			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	304	LYS	ARG	engineered mutation	UNP P13513
B	304	LYS	ARG	engineered mutation	UNP P13513

- Molecule 2 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

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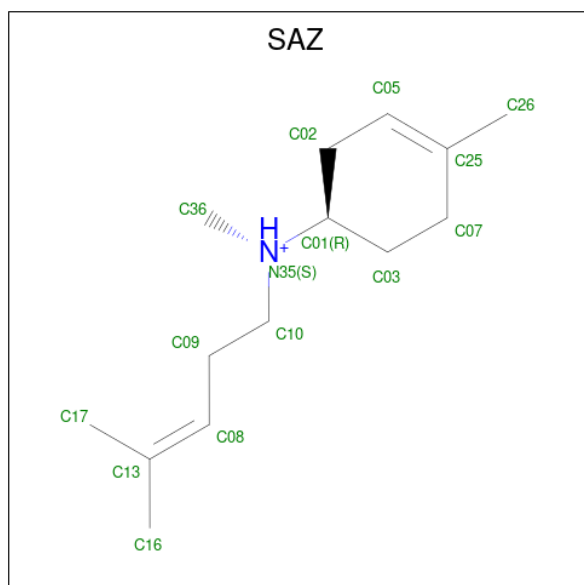
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

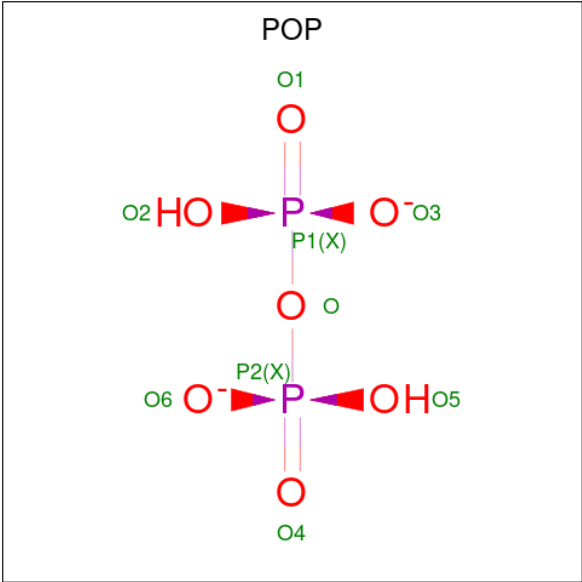
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	3	Total	Mg	0	0
			3	3		

- Molecule 4 is (1S)-N,4-DIMETHYL-N-(4-METHYLPENT-3-ENYL)CYCLOHEX-3-ENAMINIUM (CCD ID: SAZ) (formula: C₁₄H₂₆N).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	N	0	0
			15	14	1		

- Molecule 5 is PYROPHOSPHATE 2- (CCD ID: POP) (formula: H₂O₇P₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	O	P	0	0
			9	7	2		

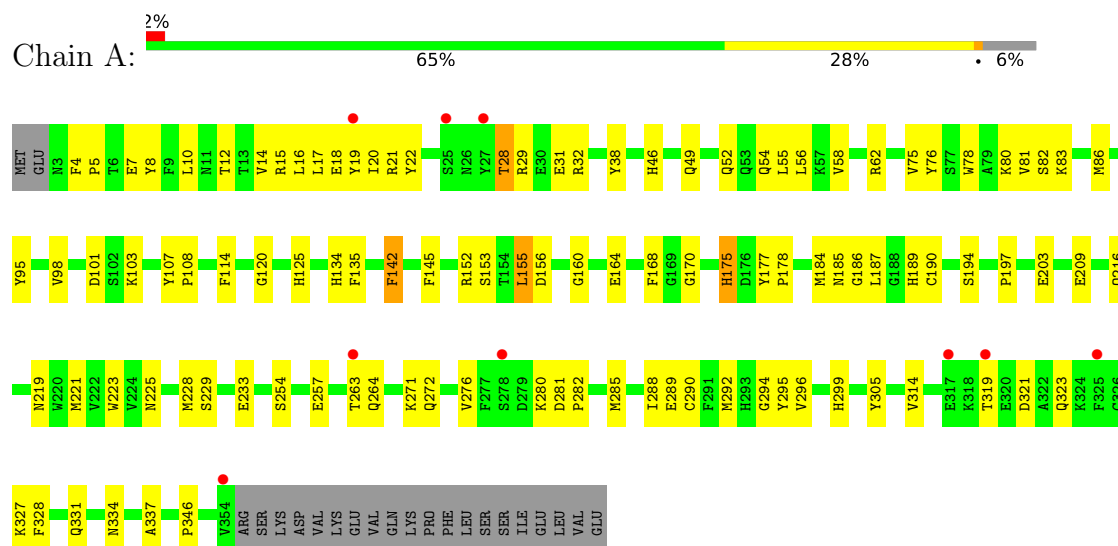
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	107	Total	O	0	0
			107	107		
6	B	76	Total	O	0	0
			76	76		

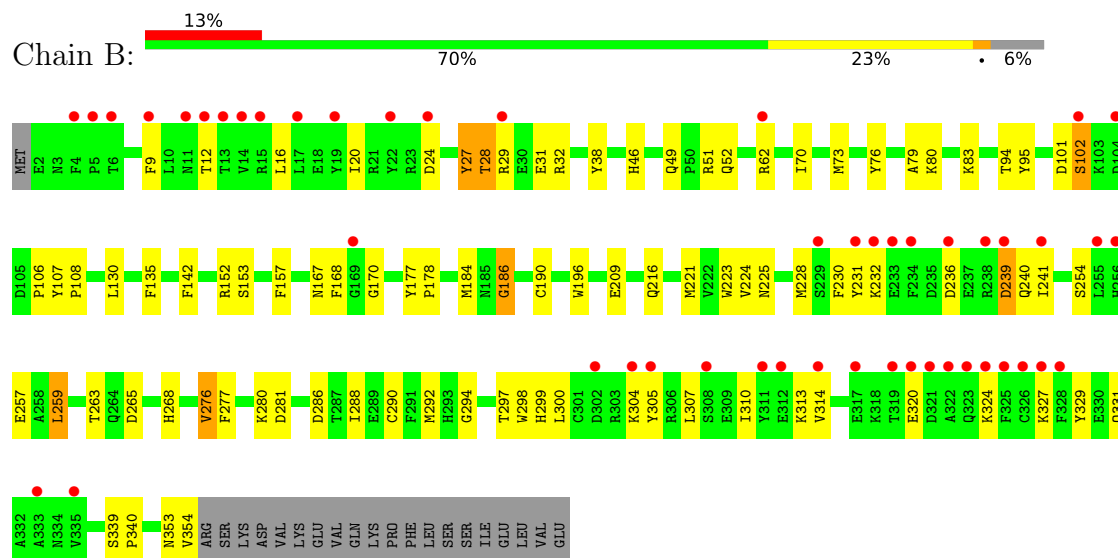
3 Residue-property plots [i](#)

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

• Molecule 1: Trichodiene synthase



• Molecule 1: Trichodiene synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	122.71Å 122.71Å 151.04Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	80.00 – 2.50 80.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	92.3 (80.00-2.50) 92.3 (80.00-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.34 (at 2.48Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.216 , 0.251 0.222 , 0.253	Depositor DCC
R_{free} test set	2130 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	50.1	Xtriage
Anisotropy	0.250	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.019 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6073	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.45% 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: POP, EDO, SAZ, MG

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	0.43	0/3006	0.92	12/4081 (0.3%)
1	B	0.41	0/3015	0.88	9/4093 (0.2%)
All	All	0.42	0/6021	0.90	21/8174 (0.3%)

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	B	0	1

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	264	GLN	OE1-CD-NE2	-9.83	112.77	122.60
1	B	240	GLN	OE1-CD-NE2	-9.81	112.79	122.60
1	B	28	THR	N-CA-C	-8.03	97.24	110.17
1	A	103	LYS	N-CA-C	-7.11	105.14	113.88
1	A	264	GLN	CG-CD-NE2	6.71	126.47	116.40
1	B	240	GLN	CG-CD-NE2	6.69	126.44	116.40
1	B	276	VAL	N-CA-C	5.51	116.16	110.82
1	A	142	PHE	N-CA-C	5.50	118.05	109.52
1	A	346	PRO	CA-C-N	-5.43	114.36	119.85
1	A	346	PRO	C-N-CA	-5.43	114.36	119.85
1	B	277	PHE	N-CA-C	5.39	119.84	113.16
1	A	155	LEU	N-CA-C	-5.31	105.57	111.36
1	B	79	ALA	N-CA-C	5.30	118.84	112.38
1	B	196	TRP	N-CA-C	5.23	118.14	108.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	114	PHE	N-CA-C	5.18	116.62	111.07
1	B	186	GLY	N-CA-C	5.13	120.24	113.37
1	A	28	THR	N-CA-C	5.12	117.13	110.53
1	A	125	HIS	CA-C-N	5.10	125.13	119.32
1	A	125	HIS	C-N-CA	5.10	125.13	119.32
1	A	175	HIS	N-CA-C	5.04	118.20	111.75
1	B	142	PHE	N-CA-C	5.03	117.31	109.52

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	27	TYR	Sidechain

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2921	0	2775	82	0
1	B	2930	0	2781	77	0
2	A	8	0	12	0	0
2	B	4	0	6	0	0
3	B	3	0	0	0	0
4	B	15	0	26	7	0
5	B	9	0	0	0	0
6	A	107	0	0	1	0
6	B	76	0	0	0	0
All	All	6073	0	5600	153	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 14.

All (153) 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:184:MET:HE1	1:B:152:ARG:NH2	1.73	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:LEU:HD21	1:A:296:VAL:HG11	1.42	1.00
1:A:152:ARG:NH2	1:B:184:MET:HE1	1.84	0.93
1:A:75:VAL:HA	1:A:86:MET:HE2	1.56	0.86
1:B:12:THR:HB	1:B:263:THR:HG21	1.72	0.71
1:B:12:THR:HG21	1:B:259:LEU:HD12	1.70	0.71
1:B:225:ASN:HD22	1:B:299:HIS:HE1	1.39	0.68
1:A:107:TYR:HB3	1:A:108:PRO:HD3	1.76	0.67
1:A:54:GLN:HG3	1:A:55:LEU:HD23	1.78	0.66
1:A:184:MET:HE1	1:B:152:ARG:HH22	1.60	0.65
1:A:38:TYR:CD2	1:A:83:LYS:HB3	2.33	0.62
1:A:323:GLN:O	1:A:327:LYS:HG3	2.00	0.61
1:B:73:MET:SD	4:B:704:SAZ:H163	2.40	0.61
1:B:27:TYR:HB2	1:B:32:ARG:HG3	1.82	0.61
1:B:167:ASN:HA	1:B:241:ILE:HG13	1.84	0.60
1:B:80:LYS:HE2	1:B:290:CYS:SG	2.42	0.60
1:A:292:MET:O	1:A:296:VAL:HG23	2.02	0.59
1:B:231:TYR:CD1	1:B:310:ILE:HD11	2.38	0.59
1:A:19:TYR:CE2	1:A:271:LYS:HG3	2.38	0.59
1:A:254:SER:OG	1:A:257:GLU:HG3	2.03	0.59
1:A:19:TYR:CD2	1:A:271:LYS:HG3	2.38	0.58
1:A:276:VAL:O	1:A:280:LYS:HD3	2.03	0.58
1:B:254:SER:OG	1:B:257:GLU:HG3	2.03	0.58
1:B:16:LEU:O	1:B:20:ILE:HG12	2.03	0.58
1:B:300:LEU:HD23	1:B:307:LEU:CD1	2.34	0.58
1:B:353:ASN:O	1:B:354:VAL:C	2.47	0.58
1:A:184:MET:HE1	1:B:152:ARG:CZ	2.34	0.57
1:A:62:ARG:NH1	1:A:101:ASP:OD2	2.38	0.57
1:A:76:TYR:O	1:A:294:GLY:HA3	2.04	0.56
1:B:12:THR:CG2	1:B:263:THR:HG21	2.36	0.56
1:A:46:HIS:HA	1:A:49:GLN:HE21	1.71	0.56
1:B:12:THR:CB	1:B:263:THR:HG21	2.36	0.55
1:A:14:VAL:O	1:A:18:GLU:HG3	2.06	0.55
1:A:80:LYS:HE2	1:A:290:CYS:SG	2.46	0.55
1:B:310:ILE:O	1:B:314:VAL:HG22	2.06	0.54
1:B:12:THR:HB	1:B:263:THR:CG2	2.37	0.54
1:B:80:LYS:HE3	1:B:286:ASP:OD2	2.06	0.54
1:A:285:MET:HE3	1:A:289:GLU:HB2	1.89	0.54
1:A:12:THR:HG21	1:A:263:THR:HG22	1.89	0.54
1:B:225:ASN:HD22	1:B:299:HIS:CE1	2.24	0.54
1:A:62:ARG:HB3	1:A:62:ARG:HH11	1.72	0.54
1:A:280:LYS:O	1:A:281:ASP:C	2.51	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:ARG:NH1	1:A:62:ARG:HB3	2.23	0.53
1:A:82:SER:O	1:A:86:MET:HG3	2.08	0.53
1:B:225:ASN:ND2	1:B:299:HIS:HE1	2.07	0.52
1:B:153:SER:HB3	1:B:190:CYS:HB2	1.92	0.52
1:B:46:HIS:HE1	1:B:95:TYR:OH	1.92	0.52
1:A:177:TYR:HB3	1:A:178:PRO:HD3	1.93	0.51
1:B:276:VAL:O	1:B:280:LYS:HD3	2.10	0.51
1:B:231:TYR:CG	1:B:310:ILE:HD11	2.45	0.51
1:B:305:TYR:OH	4:B:704:SAZ:H172	2.11	0.51
1:A:17:LEU:HD23	1:A:20:ILE:HD11	1.92	0.50
1:B:62:ARG:HH12	1:B:101:ASP:HB3	1.75	0.50
1:A:19:TYR:CE2	1:A:271:LYS:HE3	2.46	0.50
1:A:229:SER:HG	1:A:305:TYR:HE1	1.58	0.50
1:A:288:ILE:HG22	1:A:292:MET:HE2	1.92	0.50
1:A:17:LEU:HD22	1:A:22:TYR:HB2	1.94	0.50
1:A:285:MET:CE	1:A:289:GLU:HB2	2.42	0.50
1:B:51:ARG:NH1	1:B:130:LEU:HD23	2.27	0.50
1:B:46:HIS:O	1:B:52:GLN:HG3	2.12	0.50
1:B:223:TRP:CZ3	1:B:265:ASP:HB3	2.47	0.50
1:A:228:MET:HE1	1:A:296:VAL:HG13	1.94	0.49
1:A:80:LYS:CE	1:A:290:CYS:SG	3.00	0.49
1:B:236:ASP:HB3	1:B:239:ASP:HB2	1.94	0.49
1:A:46:HIS:HE1	1:A:95:TYR:OH	1.95	0.49
1:A:17:LEU:HD22	1:A:22:TYR:CG	2.48	0.49
1:B:76:TYR:O	1:B:294:GLY:HA3	2.11	0.49
1:B:288:ILE:HG22	1:B:292:MET:HE2	1.94	0.49
1:B:62:ARG:HH12	1:B:101:ASP:CB	2.26	0.48
1:B:9:PHE:HE1	1:B:230:PHE:HB3	1.79	0.48
1:B:70:ILE:HG13	1:B:94:THR:HA	1.95	0.48
1:A:10:LEU:HD11	1:A:314:VAL:HG11	1.94	0.48
1:A:15:ARG:O	1:A:19:TYR:CD1	2.67	0.48
1:B:186:GLY:O	4:B:704:SAZ:H362	2.14	0.48
1:B:46:HIS:HA	1:B:49:GLN:HE21	1.79	0.48
1:A:28:THR:OG1	1:A:31:GLU:HG3	2.14	0.47
1:A:186:GLY:O	1:A:187:LEU:HB2	2.14	0.47
1:A:331:GLN:O	1:A:334:ASN:HB3	2.13	0.47
1:B:186:GLY:O	1:B:221:MET:HE1	2.14	0.47
1:B:300:LEU:HD23	1:B:307:LEU:HD12	1.96	0.47
1:B:329:TYR:CD1	1:B:329:TYR:O	2.68	0.47
1:B:29:ARG:HG3	1:B:29:ARG:HH11	1.79	0.47
1:A:152:ARG:HH22	1:B:184:MET:HE1	1.73	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:ASP:OD1	1:B:152:ARG:HD2	2.14	0.47
1:A:120:GLY:HA2	6:A:784:HOH:O	2.15	0.47
1:A:285:MET:HE3	1:A:285:MET:O	2.15	0.47
1:B:101:ASP:O	1:B:102:SER:O	2.34	0.46
1:B:228:MET:HE3	1:B:299:HIS:HB3	1.97	0.46
1:A:228:MET:HE3	1:A:299:HIS:HB2	1.96	0.46
1:B:28:THR:OG1	1:B:31:GLU:HG3	2.16	0.46
1:A:46:HIS:O	1:A:52:GLN:HG3	2.14	0.46
1:A:225:ASN:HD22	1:A:299:HIS:HE1	1.62	0.46
1:B:38:TYR:CD2	1:B:83:LYS:HB3	2.50	0.46
1:A:7:GLU:HG3	1:A:8:TYR:N	2.31	0.46
1:B:101:ASP:O	1:B:102:SER:C	2.59	0.46
1:A:328:PHE:O	1:A:331:GLN:HB2	2.16	0.45
1:B:221:MET:SD	4:B:704:SAZ:H361	2.57	0.45
1:A:4:PHE:HA	1:A:5:PRO:HD3	1.84	0.45
1:A:229:SER:HB3	1:A:233:GLU:OE2	2.16	0.45
1:B:300:LEU:HA	1:B:307:LEU:HD12	1.97	0.45
1:A:272:GLN:O	1:A:276:VAL:HG23	2.15	0.45
1:B:288:ILE:O	1:B:292:MET:HG3	2.17	0.44
1:A:78:TRP:O	1:A:81:VAL:HG22	2.17	0.44
1:B:27:TYR:CB	1:B:32:ARG:HG3	2.46	0.44
1:B:313:LYS:O	1:B:313:LYS:HG2	2.18	0.44
1:B:327:LYS:O	1:B:331:GLN:HG3	2.18	0.44
1:A:219:ASN:HB3	1:A:223:TRP:CZ3	2.53	0.44
1:A:56:LEU:HG	1:A:58:VAL:HG23	2.00	0.44
1:A:16:LEU:O	1:A:20:ILE:HG12	2.18	0.44
1:A:281:ASP:C	1:A:281:ASP:OD1	2.60	0.43
1:A:29:ARG:NH2	1:A:337:ALA:O	2.48	0.43
1:A:168:PHE:CE2	1:A:170:GLY:HA2	2.53	0.43
1:A:58:VAL:HG21	1:A:98:VAL:CG1	2.49	0.43
1:B:329:TYR:CD1	1:B:329:TYR:C	2.96	0.43
4:B:704:SAZ:H072	4:B:704:SAZ:H162	2.00	0.43
1:B:216:GLN:HE22	1:B:276:VAL:HG21	1.84	0.43
1:B:320:GLU:O	1:B:324:LYS:HG3	2.19	0.42
1:A:81:VAL:HG12	1:A:197:PRO:HG2	2.00	0.42
1:B:228:MET:HE3	1:B:299:HIS:CB	2.50	0.42
1:A:58:VAL:HG21	1:A:98:VAL:HG11	2.00	0.42
1:A:319:THR:HG22	1:A:321:ASP:H	1.85	0.42
1:A:142:PHE:CE2	1:A:194:SER:HA	2.55	0.42
1:A:295:TYR:O	1:A:299:HIS:HD2	2.02	0.42
1:B:280:LYS:O	1:B:281:ASP:C	2.63	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:228:MET:CE	1:A:296:VAL:HG13	2.50	0.41
1:B:52:GLN:OE1	1:B:95:TYR:OH	2.38	0.41
1:A:153:SER:HB3	1:A:190:CYS:HB2	2.03	0.41
1:B:168:PHE:CZ	1:B:170:GLY:HA2	2.56	0.41
1:B:224:VAL:O	1:B:228:MET:HG2	2.20	0.41
1:B:106:PRO:O	1:B:107:TYR:C	2.63	0.41
1:A:221:MET:HE3	1:A:221:MET:HB3	2.00	0.41
1:B:107:TYR:HB3	1:B:108:PRO:CD	2.50	0.41
1:A:160:GLY:O	1:A:164:GLU:HG3	2.21	0.41
1:A:175:HIS:HD2	1:B:268:HIS:NE2	2.19	0.41
1:B:73:MET:HB2	1:B:298:TRP:CE2	2.55	0.41
1:B:24:ASP:OD2	1:B:297:THR:OG1	2.38	0.41
1:A:134:HIS:O	1:A:134:HIS:CG	2.74	0.41
1:A:145:PHE:HB3	1:A:203:GLU:CD	2.46	0.41
1:A:155:LEU:HD23	1:B:152:ARG:HG2	2.02	0.41
1:A:185:ASN:OD1	1:A:185:ASN:C	2.64	0.41
1:A:281:ASP:HA	1:A:282:PRO:HD3	1.90	0.41
1:B:107:TYR:C	1:B:107:TYR:CD1	2.99	0.41
1:A:216:GLN:HE22	1:A:276:VAL:HG11	1.86	0.40
1:A:12:THR:HG21	1:A:263:THR:CG2	2.51	0.40
1:B:177:TYR:N	1:B:178:PRO:CD	2.85	0.40
4:B:704:SAZ:H01	4:B:704:SAZ:C08	2.51	0.40
1:A:20:ILE:O	1:A:21:ARG:HB2	2.21	0.40
1:B:46:HIS:HA	1:B:49:GLN:NE2	2.37	0.40
1:A:152:ARG:NH2	1:A:189:HIS:ND1	2.65	0.40
1:A:17:LEU:HD22	1:A:22:TYR:CB	2.51	0.40
1:B:157:PHE:CZ	4:B:704:SAZ:H263	2.57	0.40
1:B:300:LEU:HD23	1:B:307:LEU:HD13	2.02	0.40
1:B:339:SER:HA	1:B:340:PRO:HD3	1.89	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	350/374 (94%)	339 (97%)	11 (3%)	0	100	100
1	B	351/374 (94%)	323 (92%)	25 (7%)	3 (1%)	14	27
All	All	701/748 (94%)	662 (94%)	36 (5%)	3 (0%)	30	49

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	102	SER
1	B	304	LYS
1	B	239	ASP

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	318/340 (94%)	315 (99%)	3 (1%)	70	87
1	B	319/340 (94%)	315 (99%)	4 (1%)	61	82
All	All	637/680 (94%)	630 (99%)	7 (1%)	65	84

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	ARG
1	A	135	PHE
1	A	209	GLU
1	B	135	PHE
1	B	209	GLU
1	B	232	LYS
1	B	259	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN

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Mol	Chain	Res	Type
1	A	46	HIS
1	A	49	GLN
1	A	137	ASN
1	A	175	HIS
1	A	200	GLN
1	A	216	GLN
1	A	272	GLN
1	A	283	GLN
1	A	299	HIS
1	A	323	GLN
1	A	350	GLN
1	A	353	ASN
1	B	46	HIS
1	B	49	GLN
1	B	137	ASN
1	B	175	HIS
1	B	179	GLN
1	B	216	GLN
1	B	225	ASN
1	B	264	GLN
1	B	272	GLN
1	B	299	HIS
1	B	353	ASN

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 ⓘ

Of 8 ligands modelled in this entry, 3 are monoatomic - leaving 5 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	$\# Z > 2$
5	POP	B	700	3	6,8,8	1.54	1 (16%)	12,13,13	1.60	1 (8%)
2	EDO	B	751	-	3,3,3	2.29	2 (66%)	2,2,2	0.48	0
2	EDO	A	752	-	3,3,3	2.41	2 (66%)	2,2,2	0.43	0
4	SAZ	B	704	-	14,15,15	2.75	5 (35%)	18,19,19	1.85	5 (27%)
2	EDO	A	750	-	3,3,3	2.26	2 (66%)	2,2,2	0.48	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
5	POP	B	700	3	-	0/6/6/6	-
2	EDO	B	751	-	-	0/1/1/1	-
2	EDO	A	752	-	-	0/1/1/1	-
4	SAZ	B	704	-	1/1/4/6	5/10/20/20	0/1/1/1
2	EDO	A	750	-	-	1/1/1/1	-

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	704	SAZ	C05-C25	5.78	1.55	1.33
4	B	704	SAZ	C01-N35	5.52	1.54	1.47
4	B	704	SAZ	C02-C01	3.46	1.58	1.52
4	B	704	SAZ	C02-C05	3.23	1.56	1.50
2	A	752	EDO	O2-C2	3.19	1.58	1.42
4	B	704	SAZ	C07-C25	-3.14	1.38	1.49
2	B	751	EDO	O2-C2	2.90	1.56	1.42
2	A	750	EDO	O2-C2	2.88	1.56	1.42
2	B	751	EDO	O1-C1	2.58	1.55	1.42
2	A	752	EDO	O1-C1	2.55	1.55	1.42
2	A	750	EDO	O1-C1	2.53	1.55	1.42
5	B	700	POP	P2-O6	-2.45	1.45	1.54

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	700	POP	O2-P1-O	4.71	120.43	104.64
4	B	704	SAZ	C36-N35-C10	-4.01	104.64	110.51
4	B	704	SAZ	C03-C07-C25	2.91	123.29	113.51
4	B	704	SAZ	C02-C05-C25	-2.82	112.63	123.96
4	B	704	SAZ	C01-C02-C05	-2.71	107.38	111.32
4	B	704	SAZ	C10-C09-C08	2.38	114.56	111.14

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	B	704	SAZ	C01

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	704	SAZ	C02-C01-N35-C10
4	B	704	SAZ	C02-C01-N35-C36
4	B	704	SAZ	C03-C01-N35-C10
4	B	704	SAZ	C08-C09-C10-N35
4	B	704	SAZ	C03-C01-N35-C36
2	A	750	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	704	SAZ	7	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

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	346/374 (92%)	0.20	9 (2%) 57 52	31, 48, 76, 102	0
1	B	337/374 (90%)	0.67	49 (14%) 6 5	33, 53, 100, 123	0
All	All	683/748 (91%)	0.43	58 (8%) 16 14	31, 51, 92, 123	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	320	GLU	8.7
1	B	22	TYR	8.5
1	B	239	ASP	5.6
1	B	314	VAL	5.4
1	B	236	ASP	5.3
1	A	25	SER	4.6
1	B	319	THR	4.6
1	B	317	GLU	4.5
1	B	9	PHE	4.5
1	A	354	VAL	4.4
1	B	241	ILE	4.2
1	B	238	ARG	4.0
1	B	4	PHE	4.0
1	B	6	THR	4.0
1	B	19	TYR	4.0
1	B	232	LYS	3.9
1	B	11	ASN	3.9
1	B	321	ASP	3.7
1	B	24	ASP	3.7
1	B	102	SER	3.6
1	B	17	LEU	3.6
1	B	327	LYS	3.6
1	B	324	LYS	3.5
1	B	14	VAL	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	328	PHE	3.5
1	B	322	ALA	3.4
1	B	325	PHE	3.3
1	A	263	THR	3.3
1	A	319	THR	3.3
1	B	5	PRO	3.3
1	A	325	PHE	3.2
1	B	311	TYR	3.1
1	B	323	GLN	3.0
1	B	326	CYS	2.9
1	B	304	LYS	2.8
1	B	233	GLU	2.8
1	B	312	GLU	2.7
1	A	19	TYR	2.7
1	B	231	TYR	2.6
1	B	62	ARG	2.5
1	B	229	SER	2.5
1	A	317	GLU	2.5
1	B	335	VAL	2.5
1	A	27	TYR	2.4
1	A	278	SER	2.4
1	B	305	TYR	2.4
1	B	234	PHE	2.3
1	B	302	ASP	2.3
1	B	255	LEU	2.2
1	B	29	ARG	2.2
1	B	104	ASP	2.2
1	B	13	THR	2.2
1	B	12	THR	2.2
1	B	15	ARG	2.1
1	B	333	ALA	2.1
1	B	169	GLY	2.1
1	B	308	SER	2.1
1	B	256	HIS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MG	B	702	1/1	0.86	0.18	60,60,60,60	0
2	EDO	A	752	4/4	0.87	0.14	53,57,59,62	0
3	MG	B	703	1/1	0.87	0.19	70,70,70,70	0
4	SAZ	B	704	15/15	0.87	0.26	95,97,98,99	0
5	POP	B	700	9/9	0.90	0.12	90,91,93,94	0
3	MG	B	701	1/1	0.95	0.07	65,65,65,65	0
2	EDO	B	751	4/4	0.96	0.09	37,39,42,44	0
2	EDO	A	750	4/4	0.97	0.06	40,44,45,46	0

6.5 Other polymers [i](#)

There are no such residues in this entry.