



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 29, 2026 – 06:55 AM UTC

PDB ID : 1WMP / pdb_00001wmp
Title : Crystal structure of amine oxidase complexed with cobalt ion
Authors : Okajima, T.; Kishishita, S.; Chiu, Y.C.; Murakawa, T.; Kim, M.; Yamaguchi, H.; Hirota, S.; Kuroda, S.; Tanizawa, K.
Deposited on : 2004-07-13
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
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

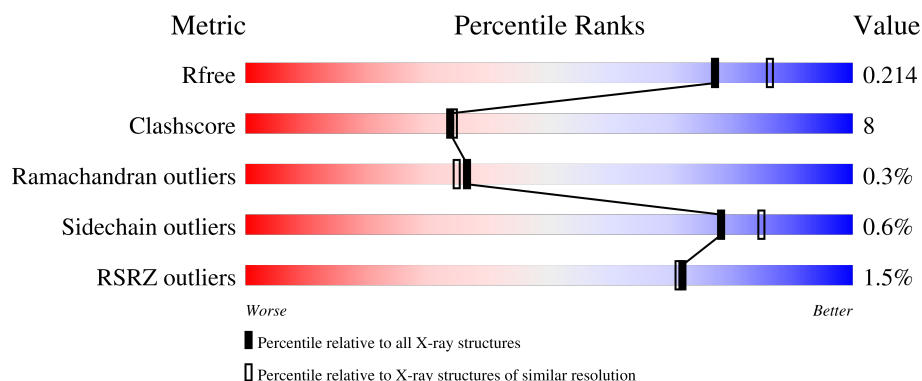
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	638	
1	B	638	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10709 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 Phenylethylamine oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	620	Total	C	N	O	S	0	0	0
			4866	3074	855	928	9			
1	B	620	Total	C	N	O	S	0	0	0
			4866	3074	855	928	9			

- Molecule 2 is COBALT (II) ION (CCD ID: CO) (formula: Co).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Co	0	0
			1	1		
2	B	1	Total	Co	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	442	Total	O	0	0
			442	442		
3	B	533	Total	O	0	0
			533	533		

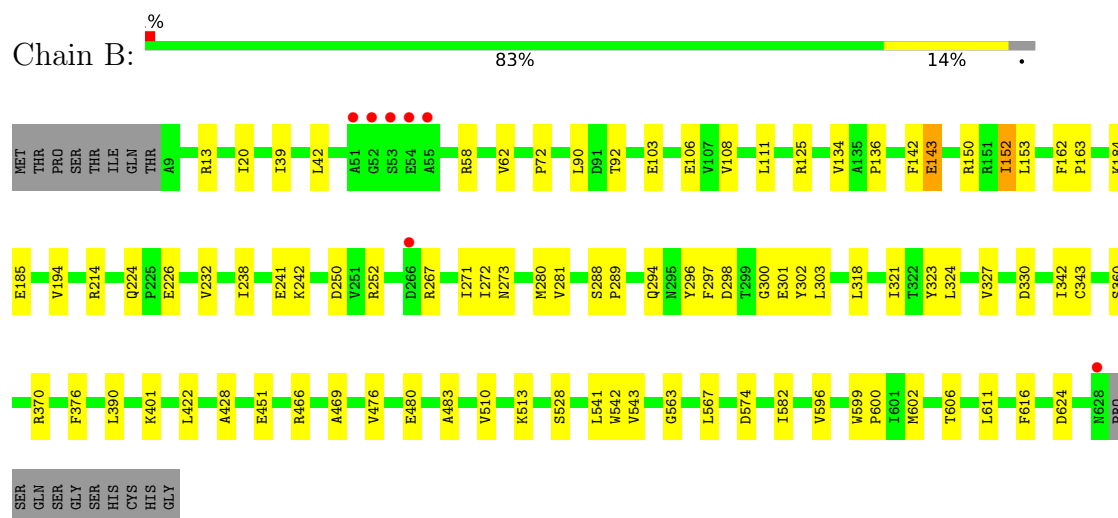
3 Residue-property plots

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: Phenylethylamine oxidase



• Molecule 1: Phenylethylamine oxidase



4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	157.58Å 63.38Å 184.20Å 90.00° 112.47° 90.00°	Depositor
Resolution (Å)	10.00 – 2.00 10.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.00) 97.7 (10.00-2.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 1.99Å)	Xtrriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.188 , 0.220 0.181 , 0.214	Depositor DCC
R_{free} test set	5637 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	21.6	Xtrriage
Anisotropy	0.689	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.53 , 92.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10709	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 36.17 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.2107e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.39	0/4989	0.88	8/6795 (0.1%)
1	B	0.41	0/4989	0.90	12/6795 (0.2%)
All	All	0.40	0/9978	0.89	20/13590 (0.1%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	602	MET	N-CA-C	6.74	118.56	110.07
1	B	602	MET	N-CA-C	6.58	118.08	109.93
1	B	108	VAL	N-CA-C	6.41	116.89	110.23
1	A	582	ILE	N-CA-C	5.82	117.23	108.96
1	B	294	GLN	N-CA-C	5.63	117.26	109.15
1	A	592	HIS	N-CA-C	5.53	117.91	108.90
1	B	321	ILE	N-CA-C	5.48	116.48	108.53
1	A	540	ASP	N-CA-C	5.33	117.17	111.36
1	B	582	ILE	N-CA-C	5.31	117.28	108.99
1	B	300	GLY	N-CA-C	5.26	119.50	112.77
1	A	408	THR	N-CA-C	5.25	117.86	110.14
1	B	224	GLN	CA-C-N	5.20	126.34	119.84
1	B	224	GLN	C-N-CA	5.20	126.34	119.84
1	A	152	ILE	N-CA-C	5.19	116.84	108.85
1	B	152	ILE	N-CA-C	5.13	117.15	109.20
1	A	269	ARG	N-CA-C	5.12	116.35	109.24
1	B	92	THR	N-CA-C	5.12	116.86	111.28
1	A	300	GLY	N-CA-C	5.09	119.29	112.77
1	B	469	ALA	N-CA-C	5.09	117.81	110.28
1	B	563	GLY	N-CA-C	-5.08	104.74	112.51

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	4866	0	4688	102	0
1	B	4866	0	4688	53	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	442	0	0	5	0
3	B	533	0	0	4	0
All	All	10709	0	9376	147	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 8.

All (147) 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:72:PRO:HG2	1:A:90:LEU:HB2	1.59	0.83
1:A:143:GLU:HG3	1:A:144:TYR:H	1.49	0.77
1:A:104:GLU:O	1:A:107:VAL:HG22	1.85	0.77
1:A:108:VAL:HG22	1:A:179:VAL:HG21	1.67	0.76
1:A:422:LEU:HD11	1:A:428:ALA:HB2	1.70	0.74
1:A:297:PHE:HB2	1:A:301:GLU:HG3	1.68	0.74
1:A:451:GLU:HG3	3:A:1154:HOH:O	1.92	0.70
1:A:247:VAL:HG11	1:A:344:MET:HE3	1.74	0.69
1:A:105:PHE:HD1	1:A:136:PRO:HB2	1.57	0.68
1:A:324:LEU:HB2	1:A:342:ILE:HB	1.74	0.68
1:A:247:VAL:HG11	1:A:344:MET:CE	2.24	0.68
1:B:324:LEU:HB2	1:B:342:ILE:HB	1.74	0.68
1:A:66:ASP:OD1	1:A:68:SER:HB3	1.94	0.67
1:A:162:PHE:HB2	1:A:163:PRO:HD2	1.77	0.66
1:A:483:ALA:HB1	1:A:543:VAL:HB	1.77	0.66
1:A:238:ILE:HG13	1:A:344:MET:HE1	1.77	0.65
1:A:170:HIS:HD2	1:A:198:PRO:O	1.80	0.65
1:A:206:ASP:OD1	1:A:208:GLU:HG2	1.95	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247:VAL:HG21	1:A:324:LEU:HD11	1.80	0.64
1:A:624:ASP:CG	1:B:214:ARG:HD2	2.22	0.64
1:A:505:ARG:HD3	1:A:618:ASP:HB3	1.79	0.64
1:B:72:PRO:HG2	1:B:90:LEU:HB2	1.81	0.63
1:A:280:MET:HE1	1:A:384:TYR:CE2	2.33	0.63
1:A:105:PHE:CD1	1:A:136:PRO:HB2	2.37	0.60
1:B:599:TRP:CD2	1:B:600:PRO:HA	2.37	0.59
1:B:451:GLU:HG3	3:B:1346:HOH:O	2.01	0.59
1:A:347:GLU:OE1	1:A:370:ARG:HD3	2.04	0.58
1:A:16:SER:OG	1:A:19:GLU:HG3	2.04	0.58
1:B:143:GLU:H	1:B:143:GLU:CD	2.12	0.58
1:A:599:TRP:CD2	1:A:600:PRO:HA	2.39	0.57
1:A:78:SER:HB2	1:A:85:ILE:HD11	1.86	0.57
1:A:241:GLU:O	1:A:242:LYS:HB2	2.04	0.56
1:B:232:VAL:HG22	1:B:238:ILE:CD1	2.36	0.56
1:B:241:GLU:O	1:B:242:LYS:HB2	2.05	0.56
1:A:134:VAL:O	1:A:136:PRO:HD3	2.08	0.54
1:A:247:VAL:O	1:A:247:VAL:HG23	2.07	0.54
1:A:280:MET:HG3	1:A:435:PHE:CD2	2.42	0.54
1:B:111:LEU:HD11	1:B:184:LYS:HA	1.90	0.53
1:A:297:PHE:HB2	1:A:301:GLU:CG	2.38	0.53
1:A:370:ARG:O	1:A:370:ARG:HG2	2.09	0.53
1:A:624:ASP:OD2	1:B:214:ARG:HD2	2.09	0.53
1:B:150:ARG:NH2	1:B:185:GLU:OE1	2.42	0.53
1:B:106:GLU:HB3	3:B:1503:HOH:O	2.08	0.52
1:A:218:LYS:HD2	1:B:226:GLU:HB2	1.92	0.52
1:B:318:LEU:HD12	3:B:1511:HOH:O	2.09	0.52
1:A:154:ARG:HD2	1:A:293:TRP:CE3	2.45	0.52
1:B:297:PHE:HB2	1:B:301:GLU:HG3	1.91	0.52
1:A:170:HIS:HE1	3:A:1234:HOH:O	1.93	0.51
1:B:20:ILE:HD12	1:B:327:VAL:HG12	1.93	0.50
1:A:214:ARG:HE	1:B:624:ASP:CG	2.18	0.50
1:A:232:VAL:HG22	1:A:238:ILE:CD1	2.42	0.50
1:A:206:ASP:CG	1:A:208:GLU:HG2	2.36	0.50
1:B:152:ILE:HG22	1:B:153:LEU:N	2.26	0.49
1:A:180:ASP:OD1	1:A:182:VAL:HG22	2.12	0.49
1:A:143:GLU:HG3	1:A:144:TYR:N	2.25	0.49
1:A:95:THR:O	1:A:557:PHE:HB2	2.13	0.48
1:A:292:SER:OG	1:B:466:ARG:HD2	2.13	0.48
1:A:619:ARG:HA	1:B:596:VAL:HG12	1.95	0.48
1:A:319:GLY:C	1:A:321:ILE:HD12	2.39	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:504:ASN:HD21	1:A:508:GLU:HB3	1.78	0.48
1:A:26:ILE:HD11	1:A:82:GLY:HA2	1.94	0.47
1:B:142:PHE:HB3	1:B:143:GLU:OE1	2.15	0.47
1:A:36:GLU:O	1:A:67:VAL:HG22	2.14	0.47
1:A:154:ARG:HD2	1:A:293:TRP:CD2	2.50	0.47
1:B:42:LEU:C	1:B:42:LEU:HD23	2.39	0.47
1:A:143:GLU:CG	1:A:144:TYR:H	2.23	0.47
1:A:182:VAL:HG22	3:A:1352:HOH:O	2.14	0.47
1:A:28:ARG:HG3	1:A:33:LEU:HD23	1.97	0.46
1:B:125:ARG:HG2	1:B:194:VAL:HG23	1.96	0.46
1:A:66:ASP:CG	1:A:70:ALA:HB3	2.41	0.46
1:A:182:VAL:HG23	1:A:183:SER:N	2.31	0.46
1:A:238:ILE:CG1	1:A:344:MET:HE1	2.44	0.46
1:A:152:ILE:HG22	1:A:153:LEU:N	2.31	0.46
1:A:20:ILE:HD12	1:A:327:VAL:HG12	1.97	0.46
1:A:323:TYR:HE1	1:A:343:CYS:SG	2.39	0.46
1:A:443:ILE:H	1:A:448:ASN:HD21	1.64	0.46
1:A:321:ILE:HD12	1:A:321:ILE:N	2.29	0.46
1:A:15:ALA:HA	1:A:19:GLU:OE1	2.16	0.46
1:B:323:TYR:CE1	1:B:343:CYS:SG	3.09	0.46
1:A:246:ASP:HB2	1:A:258:HIS:HB2	1.98	0.46
1:A:63:PHE:HB3	1:A:98:LEU:HD13	1.98	0.45
1:B:280:MET:CG	1:B:298:ASP:HB2	2.46	0.45
1:B:451:GLU:HG2	1:B:476:VAL:HG22	1.98	0.45
1:A:140:GLY:HA3	1:A:293:TRP:O	2.16	0.45
1:A:150:ARG:NH2	1:A:185:GLU:OE1	2.50	0.45
1:B:143:GLU:CD	1:B:143:GLU:N	2.74	0.45
1:B:152:ILE:CG2	1:B:153:LEU:N	2.80	0.45
1:B:483:ALA:HB1	1:B:543:VAL:HB	1.98	0.45
1:A:183:SER:OG	1:A:185:GLU:HG2	2.17	0.44
1:B:13:ARG:NH2	1:B:58:ARG:HG3	2.32	0.44
1:B:323:TYR:HE1	1:B:343:CYS:SG	2.41	0.44
1:A:302:TYR:O	1:A:303:LEU:HB2	2.18	0.44
1:A:550:GLU:HG2	1:A:570:TYR:CE1	2.53	0.44
1:A:49:ARG:HB2	3:A:1379:HOH:O	2.17	0.44
1:A:232:VAL:HG22	1:A:238:ILE:HD12	2.00	0.44
1:A:146:GLU:C	1:A:148:ARG:H	2.25	0.44
1:A:152:ILE:CG2	1:A:153:LEU:N	2.80	0.44
1:A:247:VAL:HG21	1:A:324:LEU:CD1	2.46	0.44
1:A:317:CYS:HB3	1:A:321:ILE:HG12	2.00	0.44
1:A:318:LEU:C	1:A:321:ILE:HD11	2.42	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:542:TRP:CG	1:B:567:LEU:HD13	2.53	0.43
1:A:133:ARG:NE	1:A:133:ARG:HA	2.33	0.43
1:A:460:MET:HE2	1:A:466:ARG:C	2.43	0.43
1:A:247:VAL:CG2	1:A:324:LEU:HD11	2.46	0.43
1:A:328:ILE:HD11	1:A:336:ARG:NH2	2.34	0.43
1:A:484:VAL:CG1	1:A:539:LYS:HG3	2.48	0.43
1:A:22:GLU:O	1:A:26:ILE:HG13	2.19	0.43
1:A:42:LEU:C	1:A:42:LEU:HD23	2.43	0.43
1:A:590:LEU:HD21	1:A:607:VAL:HB	2.00	0.43
1:B:267:ARG:HG2	1:B:267:ARG:HH11	1.83	0.43
1:A:250:ASP:OD2	1:A:252:ARG:HB3	2.19	0.43
1:A:147:GLU:C	1:A:150:ARG:HG3	2.44	0.43
1:B:39:ILE:HD12	1:B:62:VAL:HG11	2.00	0.43
1:B:480:GLU:HB2	1:B:574:ASP:HA	2.00	0.42
1:A:267:ARG:NH2	3:A:1137:HOH:O	2.52	0.42
1:B:250:ASP:OD2	1:B:252:ARG:HB3	2.20	0.42
1:A:486:GLU:OE1	1:A:539:LYS:NZ	2.52	0.42
1:B:302:TYR:O	1:B:303:LEU:HB2	2.19	0.42
1:A:590:LEU:CD2	1:A:607:VAL:HB	2.50	0.42
1:A:245:LEU:HD23	1:A:245:LEU:N	2.34	0.42
1:B:271:ILE:HG22	1:B:272:ILE:HG13	2.02	0.42
1:A:460:MET:HE1	1:B:528:SER:HA	2.01	0.41
1:B:103:GLU:O	1:B:106:GLU:HG2	2.20	0.41
1:A:247:VAL:HG11	1:A:344:MET:HE1	2.00	0.41
1:B:513:LYS:O	1:B:611:LEU:HA	2.20	0.41
1:A:61:ARG:HH22	1:A:556:ASP:CG	2.29	0.41
1:A:72:PRO:HB2	1:A:98:LEU:HD11	2.02	0.41
1:A:152:ILE:HG23	1:A:179:VAL:O	2.19	0.41
1:B:273:ASN:ND2	3:B:1079:HOH:O	2.40	0.41
1:B:288:SER:HA	1:B:289:PRO:HD3	1.93	0.41
1:B:370:ARG:HA	1:B:390:LEU:O	2.21	0.41
1:A:163:PRO:O	1:B:360:SER:HA	2.20	0.41
1:B:542:TRP:CD2	1:B:567:LEU:HD13	2.55	0.41
1:A:13:ARG:NH2	1:A:58:ARG:HG3	2.35	0.41
1:A:323:TYR:CE1	1:A:343:CYS:SG	3.14	0.41
1:A:371:MET:HB3	1:A:390:LEU:HB2	2.03	0.41
1:B:162:PHE:HB2	1:B:163:PRO:HD2	2.02	0.41
1:B:281:VAL:HA	1:B:296:TYR:O	2.21	0.41
1:B:510:VAL:HB	1:B:616:PHE:HA	2.03	0.41
1:B:401:LYS:HG2	1:B:606:THR:HG22	2.02	0.40
1:A:507:ASN:HD22	1:A:507:ASN:HA	1.70	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:134:VAL:O	1:B:136:PRO:HD3	2.20	0.40
1:A:115:ASP:OD1	1:A:117:ARG:HB2	2.22	0.40
1:A:261:ALA:CB	1:A:268:LEU:HD11	2.51	0.40
1:A:241:GLU:O	1:A:242:LYS:CB	2.69	0.40
1:B:330:ASP:OD2	1:B:330:ASP:C	2.64	0.40
1:B:422:LEU:HD11	1:B:428:ALA:HB2	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	618/638 (97%)	582 (94%)	32 (5%)	4 (1%)	21	17
1	B	618/638 (97%)	590 (96%)	28 (4%)	0	100	100
All	All	1236/1276 (97%)	1172 (95%)	60 (5%)	4 (0%)	36	35

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	148	ARG
1	A	55	ALA
1	A	143	GLU
1	A	144	TYR

5.3.2 Protein sidechains [i](#)

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	514/530 (97%)	511 (99%)	3 (1%)	78	85
1	B	514/530 (97%)	511 (99%)	3 (1%)	78	85
All	All	1028/1060 (97%)	1022 (99%)	6 (1%)	78	85

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	296	TYR
1	A	370	ARG
1	A	376	PHE
1	B	143	GLU
1	B	376	PHE
1	B	541	LEU

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

Mol	Chain	Res	Type
1	A	170	HIS
1	A	306	GLN
1	A	309	ASN
1	A	340	ASN
1	A	418	ASN
1	A	432	GLN
1	A	458	GLN
1	A	507	ASN
1	A	515	HIS
1	A	573	GLN
1	B	224	GLN
1	B	306	GLN
1	B	340	ASN
1	B	418	ASN
1	B	421	GLN
1	B	458	GLN
1	B	468	ASN
1	B	507	ASN
1	B	519	GLN
1	B	628	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

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	620/638 (97%)	-0.34	12 (1%) 66 66	16, 28, 48, 93	0
1	B	620/638 (97%)	-0.63	7 (1%) 78 77	13, 22, 40, 80	0
All	All	1240/1276 (97%)	-0.49	19 (1%) 72 71	13, 24, 45, 93	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	52	GLY	5.7
1	A	144	TYR	4.7
1	B	51	ALA	4.6
1	A	53	SER	4.4
1	B	52	GLY	4.1
1	A	51	ALA	3.9
1	A	145	ALA	3.8
1	B	55	ALA	3.6
1	A	148	ARG	3.5
1	B	54	GLU	3.4
1	B	628	ASN	3.3
1	A	54	GLU	2.8
1	B	53	SER	2.7
1	B	266	ASP	2.6
1	A	628	ASN	2.4
1	A	141	VAL	2.3
1	A	142	PHE	2.2
1	A	140	GLY	2.1
1	A	343	CYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CO	A	1001	1/1	0.99	0.07	45,45,45,45	0
2	CO	B	1002	1/1	0.99	0.07	49,49,49,49	0

6.5 Other polymers [i](#)

There are no such residues in this entry.