



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 9, 2017 – 07:41 AM EDT

PDB ID : 2PIN
Title : Thyroid receptor beta in complex with inhibitor
Authors : Estebanez-Perpina, E.; Jouravel, N.; Baxter, J.D.; Guy, L.R.; Webb, P.; Fletcherick, R.J.
Deposited on : unknown
Resolution : 2.30 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030345
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030345

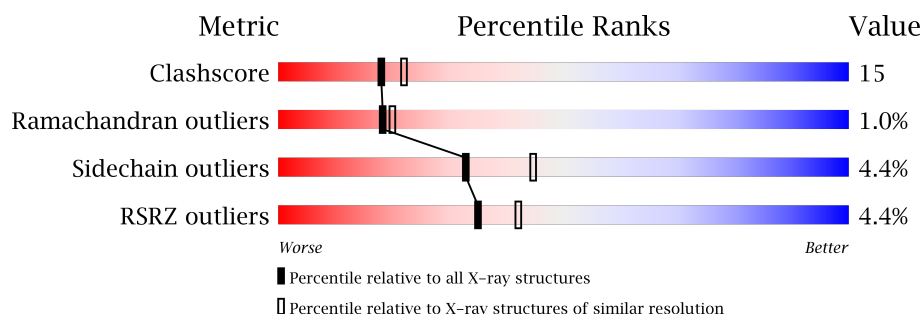
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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))
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	253	<div> <div>4%</div> <div>65%</div> <div>29%</div> <div>..</div> </div>
1	B	253	<div> <div>5%</div> <div>64%</div> <div>32%</div> <div>..</div> </div>

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	LEG	A	501	-	-	-	X
3	LEG	B	502	-	-	-	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4326 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.

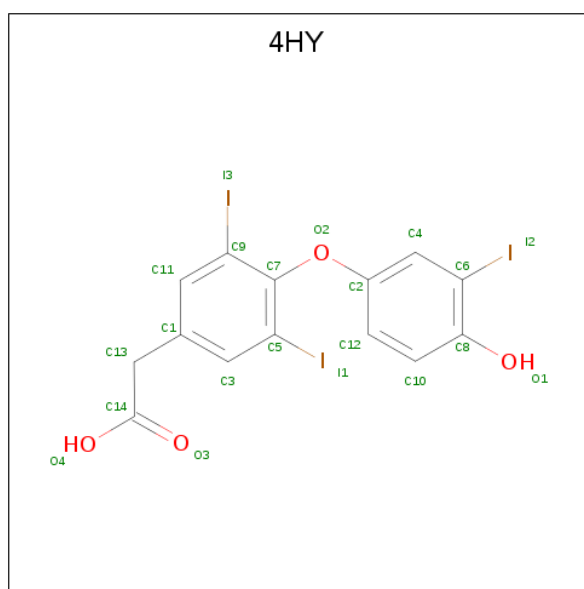
- Molecule 1 is a protein called Thyroid hormone receptor beta-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	245	Total	C	N	O	S	0	0	0
			1949	1251	329	353	16			
1	B	250	Total	C	N	O	S	0	0	0
			1978	1270	332	360	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	355	ARG	ASP	ENGINEERED	UNP P10828
B	355	ARG	ASP	ENGINEERED	UNP P10828

- Molecule 2 is [4-(4-HYDROXY-3-iodo-PHENOXY)-3,5-DIIODO-PHENYL]-ACETIC ACID (three-letter code: 4HY) (formula: C₁₄H₉I₃O₄).



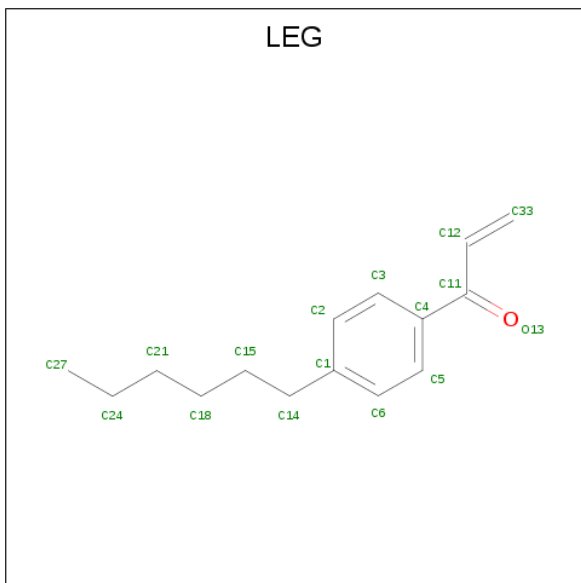
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	I	O	0	0
			21	14	3	4		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	I	O	0	0
			21	14	3	4		

- Molecule 3 is 1-(4-HEXYLPHENYL)PROP-2-EN-1-ONE (three-letter code: LEG) (formula: C₁₅H₂₀O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			16	15	1		
3	B	1	Total	C	O	0	0
			16	15	1		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		

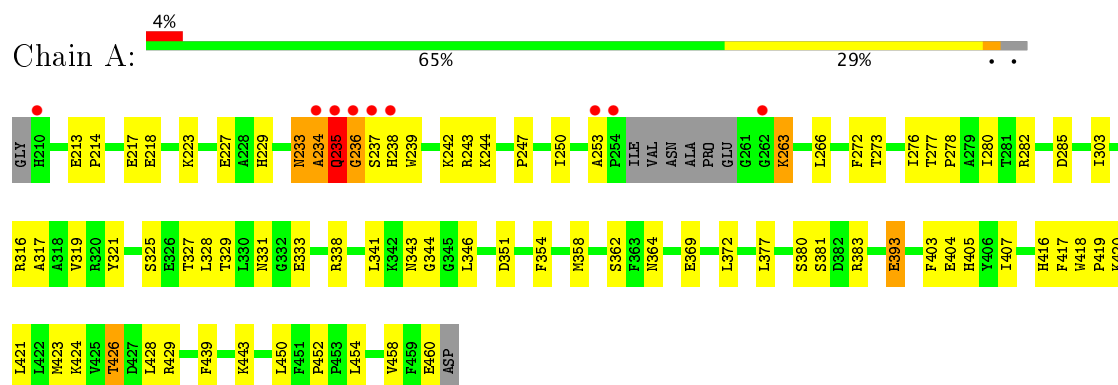
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	166	Total	O	0	0
			166	166		
5	B	154	Total	O	0	0
			154	154		

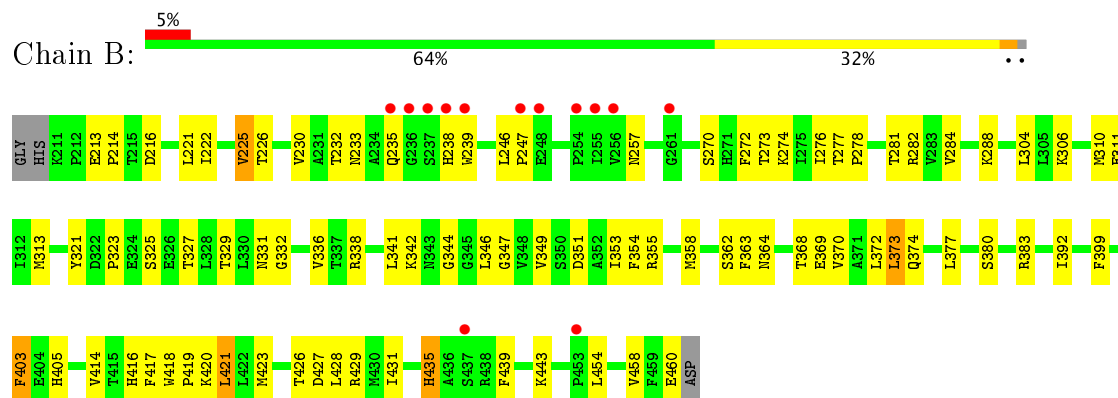
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Thyroid hormone receptor beta-1



• Molecule 1: Thyroid hormone receptor beta-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	55.13Å 92.87Å 58.35Å 90.00° 109.65° 90.00°	Depositor
Resolution (Å)	25.00 – 2.30 54.95 – 2.29	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-2.30) 99.9 (54.95-2.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 2.29Å)	Xtriage
Refinement program	ELVES, CNS	Depositor
R, R_{free}	0.216 , 0.256 0.218 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	26.4	Xtriage
Anisotropy	0.433	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 46.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4326	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.76% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, 4HY, LEG

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	0.32	1/1994 (0.1%)	0.57	3/2695 (0.1%)
1	B	0.27	0/2024	0.43	0/2739
All	All	0.29	1/4018 (0.0%)	0.50	3/5434 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	234	ALA	C-N	6.80	1.49	1.34

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	234	ALA	O-C-N	-12.87	102.11	122.70
1	A	234	ALA	C-N-CA	10.88	148.89	121.70
1	A	234	ALA	CA-C-N	8.91	136.81	117.20

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1949	0	1955	63	0
1	B	1978	0	1982	60	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	21	0	8	2	0
2	B	21	0	8	2	0
3	A	16	0	20	2	0
3	B	16	0	20	2	0
4	B	5	0	0	0	0
5	A	166	0	0	3	0
5	B	154	0	0	1	0
All	All	4326	0	3993	121	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 15.

All (121) 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:454:LEU:HD11	3:A:501:LEG:H273	1.60	0.82
1:B:418:TRP:HB3	1:B:419:PRO:HD3	1.63	0.81
1:A:235:GLN:HG2	1:A:239:TRP:HD1	1.47	0.80
1:A:418:TRP:HB3	1:A:419:PRO:HD3	1.67	0.75
1:A:423:MET:O	1:A:426:THR:HG22	1.86	0.75
1:B:377:LEU:HD11	1:B:428:LEU:HD13	1.70	0.73
1:B:272:PHE:HE1	1:B:344:GLY:HA3	1.57	0.69
1:A:303:ILE:HG23	5:A:659:HOH:O	1.93	0.68
1:B:277:THR:HB	1:B:278:PRO:HD3	1.76	0.67
1:A:223:LYS:O	1:A:227:GLU:HG3	1.96	0.66
1:B:310:MET:SD	1:B:313:MET:HE3	2.36	0.66
1:B:213:GLU:HB3	1:B:214:PRO:HD2	1.78	0.66
1:A:362:SER:HB2	1:B:364:ASN:HD21	1.62	0.65
1:A:277:THR:HB	1:A:278:PRO:HD3	1.79	0.64
1:B:342:LYS:HG3	1:B:347:GLY:HA2	1.81	0.63
1:A:273:THR:O	1:A:276:ILE:HG22	1.99	0.63
1:A:253:ALA:HB2	1:A:343:ASN:HD21	1.62	0.63
1:A:381:SER:OG	1:A:393:GLU:HG2	1.99	0.62
1:A:213:GLU:HB3	1:A:214:PRO:HD2	1.81	0.62
1:B:272:PHE:CE1	1:B:344:GLY:HA3	2.35	0.61
1:B:246:LEU:HD12	1:B:247:PRO:HD2	1.83	0.61
1:A:214:PRO:N	1:A:405:HIS:HB3	2.15	0.60
1:B:354:PHE:O	1:B:358:MET:HG3	2.02	0.60
1:A:416:HIS:HB2	1:A:420:LYS:HE3	1.83	0.59
1:A:377:LEU:HD11	1:A:428:LEU:HD13	1.84	0.59
1:A:235:GLN:HA	5:A:635:HOH:O	2.03	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:364:ASN:HD21	1:B:362:SER:HB2	1.68	0.58
1:A:235:GLN:HG2	1:A:239:TRP:CD1	2.33	0.57
1:A:235:GLN:O	1:A:237:SER:N	2.38	0.57
1:A:372:LEU:HD13	1:A:403:PHE:HA	1.86	0.57
1:B:273:THR:O	1:B:276:ILE:HG22	2.04	0.57
1:B:372:LEU:HD13	1:B:403:PHE:HA	1.87	0.57
1:A:280:ILE:HG21	1:A:458:VAL:HG21	1.87	0.56
1:B:221:LEU:O	1:B:225:VAL:HG22	2.06	0.56
1:A:263:LYS:HD3	1:A:263:LYS:H	1.71	0.56
1:A:266:LEU:H	1:A:266:LEU:HD22	1.72	0.55
1:B:272:PHE:HE2	1:B:336:VAL:HG11	1.72	0.54
1:A:253:ALA:HB2	1:A:343:ASN:ND2	2.21	0.54
1:A:454:LEU:CD1	3:A:501:LEG:H273	2.34	0.54
1:A:380:SER:O	1:A:383:ARG:HG2	2.08	0.53
1:A:404:GLU:HG2	5:A:632:HOH:O	2.07	0.53
1:B:235:GLN:HG2	1:B:239:TRP:HA	1.90	0.53
1:A:234:ALA:HB2	1:A:282:ARG:HH12	1.74	0.53
1:A:272:PHE:HE1	1:A:344:GLY:HA3	1.73	0.53
1:B:338:ARG:HD3	1:B:351:ASP:OD1	2.09	0.53
1:A:235:GLN:C	1:A:237:SER:N	2.62	0.53
1:B:439:PHE:CE2	1:B:443:LYS:HD2	2.44	0.52
1:A:341:LEU:HG	1:A:346:LEU:HD12	1.92	0.52
1:A:273:THR:HA	1:A:276:ILE:HB	1.91	0.52
1:A:280:ILE:CG2	1:A:458:VAL:HG21	2.41	0.51
1:B:454:LEU:O	1:B:458:VAL:HG23	2.10	0.51
1:B:311:GLU:OE2	1:B:429:ARG:HD3	2.11	0.51
1:B:427:ASP:O	1:B:431:ILE:HG13	2.11	0.51
1:B:351:ASP:O	1:B:355:ARG:HG3	2.11	0.51
1:A:235:GLN:C	1:A:237:SER:H	2.14	0.51
1:A:243:ARG:O	1:A:244:LYS:HD2	2.11	0.51
1:A:331:ASN:H	2:A:500:4HY:C14	2.25	0.50
1:A:217:GLU:CD	1:A:217:GLU:H	2.14	0.50
1:B:321:TYR:CE2	1:B:323:PRO:HG3	2.46	0.50
1:B:284:VAL:O	1:B:288:LYS:HG2	2.12	0.49
1:A:407:ILE:HD12	1:A:417:PHE:CE2	2.47	0.49
1:B:277:THR:O	1:B:281:THR:HG23	2.12	0.49
1:B:221:LEU:HD11	1:B:399:PHE:CE2	2.48	0.49
1:B:331:ASN:H	2:B:500:4HY:C14	2.25	0.49
1:A:403:PHE:O	1:A:407:ILE:HG12	2.13	0.49
1:B:454:LEU:CD2	3:B:502:LEG:H2	2.43	0.49
1:B:310:MET:HB3	1:B:435:HIS:ND1	2.28	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:423:MET:O	1:B:426:THR:HB	2.14	0.48
1:B:355:ARG:HA	1:B:358:MET:CE	2.44	0.48
1:B:417:PHE:CE2	1:B:421:LEU:HD12	2.48	0.48
1:B:270:SER:O	1:B:274:LYS:HG3	2.14	0.47
1:B:214:PRO:N	1:B:405:HIS:HB3	2.30	0.47
1:B:327:THR:HB	1:B:336:VAL:O	2.15	0.47
1:B:380:SER:O	1:B:383:ARG:HG2	2.14	0.47
1:B:370:VAL:O	1:B:374:GLN:HG3	2.15	0.46
1:B:214:PRO:HD3	1:B:405:HIS:O	2.16	0.46
1:A:321:TYR:HA	1:A:328:LEU:HD23	1.96	0.46
1:A:239:TRP:O	1:A:243:ARG:HB3	2.15	0.45
1:B:369:GLU:O	1:B:373:LEU:HB2	2.16	0.45
1:B:341:LEU:HG	1:B:346:LEU:HD12	1.99	0.45
1:A:276:ILE:HG23	1:A:277:THR:N	2.32	0.45
1:B:346:LEU:HB3	1:B:349:VAL:HB	1.97	0.45
1:B:325:SER:O	1:B:327:THR:HG23	2.16	0.45
1:A:229:HIS:O	1:A:233:ASN:HB2	2.17	0.44
1:A:247:PRO:HD2	1:A:250:ILE:HD12	1.99	0.44
1:B:214:PRO:HD3	1:B:405:HIS:C	2.38	0.44
1:A:317:ALA:HB2	2:A:500:4HY:H111	1.99	0.44
1:B:226:THR:O	1:B:230:VAL:HG23	2.18	0.44
1:B:232:THR:O	1:B:282:ARG:HD3	2.18	0.44
1:B:222:ILE:HG23	1:B:368:THR:HG23	1.98	0.44
1:A:277:THR:HG23	1:A:454:LEU:HD12	1.98	0.44
1:A:354:PHE:O	1:A:358:MET:HG3	2.18	0.43
1:A:316:ARG:O	1:A:319:VAL:HG22	2.18	0.43
1:A:338:ARG:HH11	1:A:351:ASP:HB3	1.83	0.43
1:B:429:ARG:HD2	5:B:646:HOH:O	2.17	0.43
1:A:235:GLN:NE2	1:A:236:GLY:H	2.17	0.43
1:B:276:ILE:HG23	1:B:277:THR:N	2.34	0.43
1:B:416:HIS:O	1:B:420:LYS:HG3	2.19	0.43
1:B:304:LEU:HD11	1:B:392:ILE:HD13	2.01	0.43
1:B:353:ILE:HD13	2:B:500:4HY:I3	2.89	0.42
1:A:325:SER:O	1:A:327:THR:HG23	2.19	0.42
1:A:341:LEU:CG	1:A:346:LEU:HD12	2.49	0.42
1:A:403:PHE:CE1	1:A:407:ILE:HD11	2.54	0.42
1:A:424:LYS:HD3	1:A:424:LYS:HA	1.91	0.42
1:A:450:LEU:O	1:A:452:PRO:HD3	2.19	0.41
1:A:416:HIS:C	1:A:419:PRO:HD2	2.40	0.41
1:A:407:ILE:HD12	1:A:417:PHE:HE2	1.83	0.41
1:B:306:LYS:HG2	3:B:502:LEG:H332	2.01	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:214:PRO:HD3	1:A:405:HIS:C	2.41	0.41
1:B:363:PHE:HD2	1:B:417:PHE:CE1	2.39	0.41
1:A:439:PHE:CE2	1:A:443:LYS:HD2	2.55	0.41
1:A:214:PRO:HA	1:A:218:GLU:OE1	2.21	0.41
1:A:239:TRP:HZ3	1:A:329:THR:HG1	1.69	0.41
1:B:278:PRO:HB3	1:B:282:ARG:HH12	1.86	0.41
1:B:338:ARG:HA	1:B:354:PHE:CE1	2.56	0.41
1:B:341:LEU:CG	1:B:346:LEU:HD12	2.51	0.41
1:B:418:TRP:HB3	1:B:419:PRO:CD	2.43	0.41
1:A:243:ARG:HA	1:A:333:GLU:HA	2.03	0.40
1:B:281:THR:HG22	1:B:454:LEU:CD2	2.52	0.40
1:B:329:THR:HG22	1:B:332:GLY:HA2	2.04	0.40
1:A:369:GLU:HG2	1:A:403:PHE:HE1	1.87	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	241/253 (95%)	227 (94%)	11 (5%)	3 (1%)	15	16
1	B	248/253 (98%)	235 (95%)	11 (4%)	2 (1%)	22	26
All	All	489/506 (97%)	462 (94%)	22 (4%)	5 (1%)	18	20

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	235	GLN
1	B	238	HIS
1	B	257	ASN
1	A	236	GLY
1	A	238	HIS

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	215/221 (97%)	205 (95%)	10 (5%)	30	41
1	B	217/221 (98%)	208 (96%)	9 (4%)	35	48
All	All	432/442 (98%)	413 (96%)	19 (4%)	33	45

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	233	ASN
1	A	235	GLN
1	A	242	LYS
1	A	263	LYS
1	A	285	ASP
1	A	393	GLU
1	A	421	LEU
1	A	426	THR
1	A	429	ARG
1	A	460	GLU
1	B	216	ASP
1	B	225	VAL
1	B	233	ASN
1	B	373	LEU
1	B	403	PHE
1	B	414	VAL
1	B	421	LEU
1	B	435	HIS
1	B	460	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	233	ASN
1	A	241	GLN
1	A	271	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	343	ASN
1	A	364	ASN
1	A	396	GLN
1	B	233	ASN
1	B	252	GLN
1	B	364	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

5 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	$\# Z > 2$
2	4HY	A	500	-	19,22,22	1.61	4 (21%)	28,31,31	1.09	1 (3%)
3	LEG	A	501	-	15,16,16	0.81	0	18,19,19	0.57	0
2	4HY	B	500	-	19,22,22	1.64	4 (21%)	28,31,31	1.10	1 (3%)
4	SO4	B	501	-	4,4,4	0.30	0	6,6,6	0.06	0
3	LEG	B	502	-	15,16,16	0.81	0	18,19,19	0.58	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
2	4HY	A	500	-	-	0/6/8/8	0/2/2/2
3	LEG	A	501	-	-	0/10/12/12	0/1/1/1
2	4HY	B	500	-	-	0/6/8/8	0/2/2/2
4	SO4	B	501	-	-	0/0/0/0	0/0/0/0
3	LEG	B	502	-	-	0/10/12/12	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	4HY	C12-C2	2.05	1.42	1.38
2	A	500	4HY	C12-C2	2.15	1.43	1.38
2	B	500	4HY	O1-C8	2.16	1.40	1.36
2	A	500	4HY	O1-C8	2.18	1.40	1.36
2	A	500	4HY	C11-C9	2.55	1.45	1.39
2	B	500	4HY	C11-C9	2.71	1.45	1.39
2	A	500	4HY	C3-C1	3.55	1.45	1.39
2	B	500	4HY	C3-C1	3.81	1.46	1.39

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	4HY	C2-O2-C7	3.81	124.98	118.53
2	B	500	4HY	C2-O2-C7	3.91	125.16	118.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	4HY	2	0
3	A	501	LEG	2	0
2	B	500	4HY	2	0
3	B	502	LEG	2	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	245/253 (96%)	0.45	9 (3%) 42 49	15, 30, 57, 78	0
1	B	250/253 (98%)	0.47	13 (5%) 28 35	14, 29, 66, 80	0
All	All	495/506 (97%)	0.46	22 (4%) 35 42	14, 29, 61, 80	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	237	SER	9.4
1	B	235	GLN	7.8
1	B	256	VAL	7.3
1	A	236	GLY	6.6
1	A	238	HIS	6.5
1	B	237	SER	6.2
1	B	255	ILE	5.3
1	A	254	PRO	4.8
1	A	262	GLY	4.3
1	B	236	GLY	3.5
1	B	238	HIS	3.4
1	A	210	HIS	2.9
1	B	239	TRP	2.6
1	B	261	GLY	2.6
1	A	253	ALA	2.6
1	B	254	PRO	2.5
1	A	234	ALA	2.3
1	B	437	SER	2.2
1	A	235	GLN	2.2
1	B	453	PRO	2.2
1	B	248	GLU	2.2
1	B	247	PRO	2.2

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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	LEG	B	502	16/16	0.76	0.28	4.62	45,47,50,53	0
3	LEG	A	501	16/16	0.76	0.27	3.79	45,47,50,53	0
4	SO4	B	501	5/5	0.99	0.18	0.95	34,36,37,37	0
2	4HY	B	500	21/21	0.99	0.14	-0.64	23,25,30,34	0
2	4HY	A	500	21/21	0.99	0.14	-0.69	20,23,31,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.