



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:49 AM UTC

PDB ID : 1JMY / pdb_00001jmy
Title : Truncated Recombinant Human Bile Salt Stimulated Lipase
Authors : Moore, S.A.; Kingston, R.L.; Loomes, K.M.; Hernell, O.; Blackberg, L.; Baker, H.M.; Baker, E.N.
Deposited on : 2001-07-20
Resolution : 2.60 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

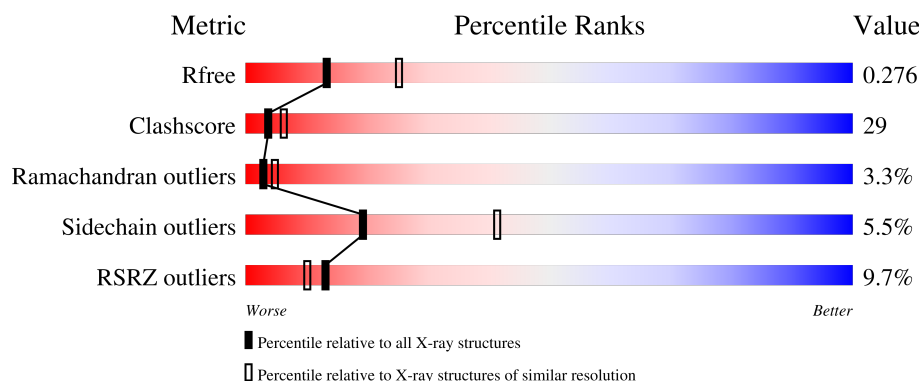
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	522	10% 52% 39% 5% ..

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4058 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.

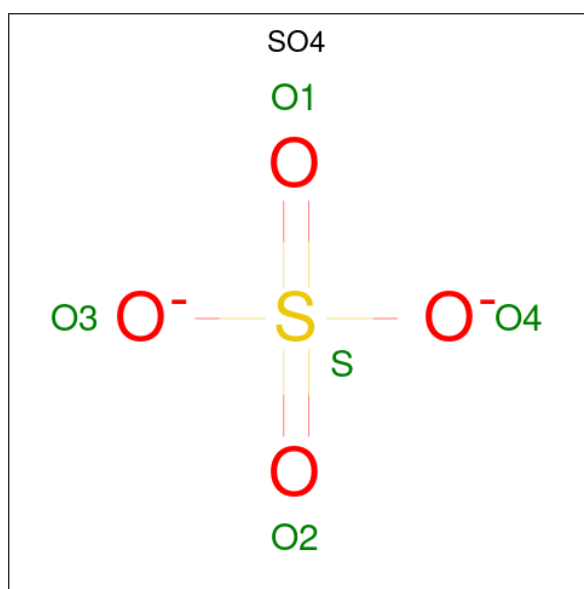
- Molecule 1 is a protein called BILE-SALT-ACTIVATED LIPASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	515	3990	2556	681	737	16	0	0	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	331	ARG	ALA	conflict	UNP P19835
A	519	PRO	-	cloning artifact	UNP P19835
A	520	GLY	-	cloning artifact	UNP P19835
A	521	ILE	-	cloning artifact	UNP P19835
A	522	HIS	-	cloning artifact	UNP P19835

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	63	Total 63	O 63	0	0

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: BILE-SALT-ACTIVATED LIPASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.64Å 90.08Å 103.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	100.00 – 2.60 67.88 – 2.60	Depositor EDS
% Data completeness (in resolution range)	89.6 (100.00-2.60) 89.6 (67.88-2.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.58 (at 2.61Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.236 , 0.276 0.238 , 0.276	Depositor DCC
R_{free} test set	1570 reflections (10.06%)	wwPDB-VP
Wilson B-factor (Å ²)	34.9	Xtriage
Anisotropy	1.013	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 55.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4058	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.41% 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: SO4

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	0.42	0/4096	1.15	42/5566 (0.8%)

There are no bond length outliers.

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	430	TRP	N-CA-C	-13.59	96.01	112.59
1	A	284	TYR	N-CA-C	11.12	125.95	112.38
1	A	341	GLU	N-CA-C	-9.27	102.10	113.97
1	A	351	PHE	N-CA-C	-9.12	102.26	112.57
1	A	22	LEU	N-CA-C	-8.96	102.12	112.87
1	A	394	LEU	N-CA-C	8.46	120.50	111.28
1	A	449	THR	CA-C-N	8.10	128.86	119.47
1	A	449	THR	C-N-CA	8.10	128.86	119.47
1	A	427	TYR	CA-C-N	8.00	128.00	119.76
1	A	427	TYR	C-N-CA	8.00	128.00	119.76
1	A	353	ILE	N-CA-C	-7.89	105.32	112.90
1	A	484	VAL	CA-C-N	7.82	127.37	118.85
1	A	484	VAL	C-N-CA	7.82	127.37	118.85
1	A	521	ILE	N-CA-C	7.55	125.05	109.34
1	A	296	ILE	CA-C-N	7.41	127.43	119.28
1	A	296	ILE	C-N-CA	7.41	127.43	119.28
1	A	23	LEU	N-CA-C	7.11	125.95	110.80
1	A	22	LEU	CA-C-N	6.95	134.81	121.54
1	A	22	LEU	C-N-CA	6.95	134.81	121.54
1	A	484	VAL	N-CA-C	6.77	114.75	109.19
1	A	121	ASN	N-CA-C	6.71	125.08	110.80
1	A	428	PRO	N-CA-C	6.43	121.18	111.14
1	A	23	LEU	CA-CB-CG	-6.36	94.03	116.30
1	A	424	MET	N-CA-C	-6.36	100.40	109.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	371	TRP	N-CA-C	-6.04	104.05	112.45
1	A	439	ILE	N-CA-C	5.99	116.64	110.36
1	A	178	ILE	N-CA-C	5.77	119.13	111.17
1	A	25	ASP	N-CA-C	5.76	118.33	110.55
1	A	411	ALA	N-CA-C	5.75	118.34	110.24
1	A	166	ASP	N-CA-C	-5.69	104.99	111.14
1	A	21	GLY	N-CA-C	5.51	126.23	113.18
1	A	473	LYS	N-CA-C	5.44	117.02	111.14
1	A	329	MET	CA-C-N	5.37	126.55	119.84
1	A	329	MET	C-N-CA	5.37	126.55	119.84
1	A	146	GLY	CA-C-N	5.32	125.39	119.32
1	A	146	GLY	C-N-CA	5.32	125.39	119.32
1	A	410	SER	N-CA-C	-5.24	106.57	113.12
1	A	519	PRO	N-CA-C	-5.20	102.43	111.32
1	A	194	SER	CB-CA-C	-5.13	110.25	117.23
1	A	271	LYS	N-CA-C	5.10	121.66	110.80
1	A	518	PHE	N-CA-C	5.08	120.22	108.93
1	A	316	THR	N-CA-CB	-5.04	103.10	111.57

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	3990	0	3927	229	0
2	A	5	0	0	0	0
3	A	63	0	0	6	0
All	All	4058	0	3927	229	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 29.

All (229) 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:316:THR:HG23	3:A:867:HOH:O	1.35	1.23
1:A:316:THR:CG2	3:A:867:HOH:O	1.90	1.19
1:A:382:LYS:HE2	1:A:430:TRP:HE1	1.16	1.10
1:A:280:PRO:HD2	1:A:327:ILE:HG21	1.31	1.09
1:A:358:ARG:HG3	1:A:359:GLY:H	1.17	1.04
1:A:329:MET:HE2	1:A:331:ARG:HE	1.24	1.01
1:A:119:PHE:CE2	1:A:449:THR:HG21	2.02	0.94
1:A:482:SER:HB2	3:A:833:HOH:O	1.72	0.89
1:A:74:THR:CG2	1:A:268:LEU:HD11	2.02	0.89
1:A:74:THR:HG22	1:A:268:LEU:HD11	1.60	0.84
1:A:92:LYS:O	1:A:93:GLN:HG3	1.79	0.82
1:A:280:PRO:HD2	1:A:327:ILE:CG2	2.09	0.82
1:A:315:GLY:HA3	1:A:416:TYR:CE1	2.16	0.81
1:A:280:PRO:CB	1:A:324:PHE:HD1	1.95	0.80
1:A:362:THR:O	1:A:366:VAL:HG23	1.81	0.80
1:A:68:THR:O	1:A:71:GLN:HG2	1.81	0.80
1:A:21:GLY:C	1:A:23:LEU:H	1.88	0.79
1:A:271:LYS:HG2	1:A:272:VAL:N	1.96	0.79
1:A:329:MET:HE2	1:A:331:ARG:NE	1.98	0.78
1:A:280:PRO:HB3	1:A:324:PHE:HD1	1.48	0.78
1:A:21:GLY:C	1:A:23:LEU:N	2.40	0.78
1:A:20:LEU:CD1	1:A:132:THR:HG21	2.15	0.77
1:A:134:GLY:HA3	1:A:469:THR:HG21	1.65	0.77
1:A:358:ARG:HG3	1:A:359:GLY:N	1.96	0.75
1:A:314:ALA:HB3	1:A:401:LEU:HD21	1.68	0.75
1:A:2:LYS:HD2	1:A:2:LYS:N	2.01	0.74
1:A:382:LYS:HE2	1:A:430:TRP:NE1	1.99	0.73
1:A:326:SER:HA	1:A:332:ILE:HD11	1.69	0.72
1:A:342:GLU:HB3	1:A:346:LYS:HE2	1.70	0.72
1:A:291:ILE:HD11	1:A:300:PRO:HD3	1.70	0.72
1:A:281:MET:HB2	1:A:284:TYR:HD1	1.55	0.71
1:A:122:ASN:OD1	1:A:445:LYS:NZ	2.24	0.71
1:A:91:ARG:N	3:A:862:HOH:O	1.97	0.70
1:A:252:ALA:O	1:A:256:GLN:HG2	1.94	0.68
1:A:512:ARG:O	1:A:513:SER:HB2	1.94	0.67
1:A:20:LEU:HD12	1:A:132:THR:HG21	1.75	0.67
1:A:315:GLY:HA3	1:A:416:TYR:CD1	2.30	0.67
1:A:110:LEU:CD2	1:A:272:VAL:HG11	2.27	0.65
1:A:248:VAL:HA	1:A:254:MET:HE2	1.77	0.65
1:A:119:PHE:O	1:A:120:LEU:O	2.15	0.64
1:A:281:MET:CB	1:A:284:TYR:HD1	2.11	0.63
1:A:429:LYS:HA	1:A:431:VAL:HG23	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:498:TYR:HD2	1:A:514:LEU:HG	1.64	0.63
1:A:42:LEU:HB3	1:A:158:LEU:HD11	1.79	0.63
1:A:280:PRO:HB3	1:A:324:PHE:CD1	2.33	0.62
1:A:281:MET:HE3	1:A:283:HIS:HB2	1.82	0.62
1:A:445:LYS:HB2	1:A:446:PRO:HD3	1.81	0.62
1:A:280:PRO:CG	1:A:327:ILE:HB	2.30	0.61
1:A:373:GLN:O	1:A:373:GLN:HG2	2.00	0.61
1:A:281:MET:HB2	1:A:284:TYR:CD1	2.36	0.61
1:A:326:SER:CA	1:A:332:ILE:HD11	2.31	0.60
1:A:250:ASP:HB3	1:A:253:ARG:HB2	1.84	0.60
1:A:119:PHE:O	1:A:120:LEU:C	2.45	0.59
1:A:521:ILE:HG22	1:A:522:HIS:O	2.02	0.59
1:A:74:THR:HG23	1:A:268:LEU:HD11	1.82	0.59
1:A:342:GLU:O	1:A:346:LYS:HG3	2.03	0.59
1:A:271:LYS:O	1:A:272:VAL:HB	2.03	0.59
1:A:66:GLN:OE1	1:A:111:MET:HB2	2.02	0.58
1:A:74:THR:CG2	1:A:268:LEU:CD1	2.77	0.58
1:A:332:ILE:HA	1:A:381:LYS:HD3	1.85	0.58
1:A:9:GLU:HG2	1:A:177:ASN:HD21	1.68	0.58
1:A:291:ILE:CD1	1:A:300:PRO:HD3	2.32	0.58
1:A:110:LEU:HD22	1:A:272:VAL:HG11	1.86	0.58
1:A:387:PHE:O	1:A:391:VAL:HG23	2.03	0.58
1:A:230:GLN:HB3	1:A:290:VAL:HG12	1.85	0.58
1:A:1:ALA:HB3	1:A:90:GLY:O	2.04	0.57
1:A:79:ASP:HA	3:A:840:HOH:O	2.04	0.57
1:A:200:VAL:O	1:A:204:THR:HG23	2.03	0.57
1:A:174:VAL:O	1:A:178:ILE:HB	2.04	0.57
1:A:217:ILE:HG12	1:A:313:ILE:HB	1.86	0.57
1:A:95:SER:HB3	1:A:98:LEU:HD21	1.87	0.57
1:A:77:ASP:OD2	1:A:78:GLU:N	2.38	0.56
1:A:282:LEU:HD22	1:A:351:PHE:CD2	2.40	0.56
1:A:338:LYS:HB2	1:A:377:GLN:CD	2.31	0.56
1:A:349:SER:O	1:A:352:THR:HG23	2.05	0.56
1:A:500:GLU:O	1:A:505:MET:HE3	2.04	0.56
1:A:280:PRO:HB2	1:A:324:PHE:HD1	1.68	0.56
1:A:368:THR:HB	1:A:371:TRP:HB2	1.88	0.56
1:A:123:TYR:O	1:A:445:LYS:NZ	2.32	0.56
1:A:117:ALA:O	1:A:124:LEU:HB2	2.06	0.55
1:A:382:LYS:CE	1:A:430:TRP:HE1	2.02	0.55
1:A:297:PRO:HG2	1:A:303:LEU:HD21	1.88	0.55
1:A:363:THR:O	1:A:366:VAL:HB	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:395:VAL:HB	1:A:396:PRO:HD3	1.88	0.55
1:A:495:ASN:HD22	1:A:512:ARG:CZ	2.19	0.55
1:A:498:TYR:CD2	1:A:514:LEU:HG	2.41	0.55
1:A:328:ASP:C	1:A:330:PRO:HD3	2.32	0.54
1:A:119:PHE:CE2	1:A:449:THR:CG2	2.86	0.54
1:A:323:ILE:HG23	1:A:324:PHE:N	2.23	0.54
1:A:357:LEU:O	1:A:361:LYS:HG2	2.07	0.54
1:A:34:PRO:HB3	1:A:82:TYR:CE1	2.43	0.54
1:A:241:ALA:HB1	1:A:246:CYS:HB2	1.90	0.54
1:A:329:MET:HE3	1:A:329:MET:HA	1.89	0.54
1:A:391:VAL:HG22	1:A:521:ILE:HD11	1.91	0.53
1:A:336:ASN:OD1	1:A:337:LYS:HG3	2.08	0.53
1:A:345:TYR:OH	1:A:361:LYS:HE3	2.09	0.53
1:A:496:SER:O	1:A:514:LEU:HB2	2.09	0.52
1:A:161:ASN:ND2	1:A:164:LEU:HD12	2.24	0.52
1:A:345:TYR:CE1	1:A:361:LYS:HD3	2.44	0.52
1:A:378:GLU:O	1:A:382:LYS:HB2	2.08	0.52
1:A:339:VAL:HG13	1:A:339:VAL:O	2.08	0.52
1:A:329:MET:O	1:A:332:ILE:HG12	2.10	0.52
1:A:119:PHE:O	1:A:119:PHE:CD2	2.63	0.52
1:A:398:GLU:HG2	1:A:514:LEU:HD22	1.92	0.51
1:A:147:PRO:O	1:A:151:LEU:HB2	2.10	0.51
1:A:2:LYS:HG3	1:A:17:ASN:ND2	2.25	0.51
1:A:323:ILE:HG23	1:A:324:PHE:H	1.76	0.51
1:A:297:PRO:HG2	1:A:303:LEU:CD2	2.41	0.51
1:A:325:ALA:HB2	1:A:388:GLU:HG3	1.92	0.50
1:A:119:PHE:CE2	1:A:122:ASN:HB3	2.46	0.50
1:A:271:LYS:HG2	1:A:272:VAL:H	1.73	0.50
1:A:314:ALA:O	1:A:415:ALA:HA	2.11	0.50
1:A:318:ASN:OD1	1:A:419:SER:HA	2.11	0.50
1:A:391:VAL:HG22	1:A:521:ILE:CD1	2.40	0.50
1:A:323:ILE:HG23	1:A:324:PHE:CD2	2.47	0.50
1:A:435:HIS:O	1:A:436:ALA:HB3	2.12	0.50
1:A:454:ARG:CZ	1:A:456:GLN:HE21	2.24	0.50
1:A:37:ALA:O	1:A:38:PRO:C	2.55	0.50
1:A:70:THR:HB	1:A:72:ASP:OD2	2.12	0.50
1:A:344:PHE:O	1:A:348:VAL:HG13	2.11	0.49
1:A:124:LEU:HG	1:A:125:TYR:CE1	2.46	0.49
1:A:258:LEU:O	1:A:266:LEU:HD11	2.12	0.49
1:A:280:PRO:HG2	1:A:327:ILE:HB	1.94	0.49
1:A:155:ASP:C	1:A:155:ASP:OD1	2.56	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:280:PRO:HG3	1:A:324:PHE:HA	1.95	0.49
1:A:56:LYS:HD2	1:A:58:LYS:NZ	2.27	0.49
1:A:326:SER:N	1:A:332:ILE:HD11	2.28	0.49
1:A:65:LEU:HD23	1:A:148:LEU:HD11	1.94	0.49
1:A:358:ARG:CG	1:A:359:GLY:H	2.04	0.49
1:A:342:GLU:CB	1:A:346:LYS:HE2	2.42	0.49
1:A:368:THR:O	1:A:371:TRP:HB2	2.12	0.49
1:A:456:GLN:OE1	1:A:503:LYS:HD3	2.11	0.49
1:A:494:GLU:CD	1:A:494:GLU:H	2.05	0.49
1:A:20:LEU:H	1:A:26:SER:HA	1.78	0.48
1:A:95:SER:C	1:A:96:ARG:HG2	2.37	0.48
1:A:349:SER:O	1:A:352:THR:CG2	2.62	0.48
1:A:454:ARG:NE	1:A:456:GLN:HE21	2.11	0.48
1:A:152:SER:OG	1:A:153:THR:N	2.45	0.48
1:A:165:ARG:O	1:A:169:MET:HG2	2.13	0.48
1:A:9:GLU:OE2	1:A:176:ARG:NH1	2.47	0.48
1:A:69:ILE:O	1:A:71:GLN:HG3	2.13	0.48
1:A:103:TRP:CZ3	1:A:192:GLY:HA2	2.49	0.47
1:A:366:VAL:O	1:A:369:GLU:HG3	2.13	0.47
1:A:128:GLU:O	1:A:132:THR:HG23	2.15	0.47
1:A:374:ASP:OD1	1:A:374:ASP:O	2.33	0.47
1:A:378:GLU:O	1:A:382:LYS:HE3	2.15	0.47
1:A:512:ARG:O	1:A:513:SER:CB	2.63	0.47
1:A:313:ILE:HG13	1:A:471:PHE:CZ	2.49	0.47
1:A:308:ALA:HB1	1:A:409:LYS:HG3	1.96	0.46
1:A:378:GLU:HG2	1:A:382:LYS:NZ	2.31	0.46
1:A:378:GLU:OE2	1:A:382:LYS:NZ	2.47	0.46
1:A:496:SER:HB3	1:A:514:LEU:HB2	1.98	0.46
1:A:160:GLY:O	1:A:161:ASN:HB2	2.16	0.46
1:A:401:LEU:HD13	1:A:415:ALA:HB2	1.95	0.46
1:A:24:GLY:CA	1:A:91:ARG:HD2	2.46	0.46
1:A:315:GLY:HA3	1:A:416:TYR:CZ	2.50	0.46
1:A:350:GLU:C	1:A:352:THR:H	2.24	0.46
1:A:64:CYS:O	1:A:66:GLN:HG2	2.16	0.46
1:A:85:ILE:HG22	1:A:86:TRP:N	2.31	0.46
1:A:134:GLY:HA3	1:A:469:THR:CG2	2.41	0.46
1:A:280:PRO:CD	1:A:327:ILE:HD12	2.45	0.46
1:A:30:PHE:N	1:A:30:PHE:CD1	2.84	0.46
1:A:261:THR:HG22	1:A:262:ASP:N	2.31	0.46
1:A:299:ASP:OD1	1:A:301:ILE:HG22	2.15	0.46
1:A:349:SER:HA	1:A:352:THR:HG23	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:280:PRO:CD	1:A:327:ILE:CD1	2.93	0.45
1:A:117:ALA:O	1:A:124:LEU:CB	2.64	0.45
1:A:119:PHE:CD2	1:A:119:PHE:C	2.92	0.45
1:A:316:THR:HG21	1:A:394:LEU:HD13	1.99	0.45
1:A:486:THR:HG22	1:A:487:HIS:N	2.31	0.45
1:A:42:LEU:HD13	1:A:158:LEU:HD21	1.98	0.45
1:A:238:LYS:HG2	1:A:254:MET:HE1	1.99	0.45
1:A:416:TYR:HB3	1:A:488:TRP:CH2	2.51	0.45
1:A:193:GLU:HA	1:A:219:GLN:O	2.16	0.45
1:A:489:GLU:HA	1:A:490:PRO:HD3	1.86	0.45
1:A:288:VAL:HB	1:A:289:PRO:CD	2.46	0.45
1:A:313:ILE:HD11	1:A:471:PHE:CG	2.52	0.45
1:A:281:MET:CB	1:A:284:TYR:CD1	2.96	0.44
1:A:83:LEU:HD21	1:A:173:TRP:CE3	2.52	0.44
1:A:32:GLY:O	1:A:82:TYR:HD1	2.00	0.44
1:A:390:ASP:CG	1:A:515:ARG:HH22	2.25	0.44
1:A:77:ASP:OD2	1:A:77:ASP:C	2.61	0.44
1:A:16:VAL:HG13	1:A:58:LYS:HG2	1.99	0.44
1:A:376:SER:C	1:A:378:GLU:N	2.74	0.44
1:A:416:TYR:HB3	1:A:488:TRP:CZ2	2.52	0.44
1:A:241:ALA:HA	1:A:258:LEU:HD21	2.00	0.44
1:A:378:GLU:HG2	1:A:382:LYS:CE	2.48	0.43
1:A:119:PHE:CD2	1:A:122:ASN:HB3	2.54	0.43
1:A:345:TYR:O	1:A:348:VAL:HG22	2.18	0.43
1:A:418:PHE:CE1	1:A:464:MET:HE2	2.53	0.43
1:A:28:ASP:OD2	1:A:89:GLN:HG3	2.18	0.43
1:A:454:ARG:CZ	1:A:456:GLN:NE2	2.82	0.43
1:A:2:LYS:N	1:A:2:LYS:CD	2.78	0.43
1:A:271:LYS:CG	1:A:272:VAL:H	2.29	0.43
1:A:23:LEU:HA	1:A:23:LEU:HD12	1.43	0.43
1:A:120:LEU:HD12	1:A:120:LEU:HA	1.86	0.43
1:A:280:PRO:HB2	1:A:324:PHE:CD1	2.52	0.43
1:A:448:ALA:C	1:A:449:THR:CG2	2.92	0.43
1:A:448:ALA:C	1:A:449:THR:HG23	2.43	0.43
1:A:449:THR:HA	1:A:450:PRO:HD2	1.83	0.43
1:A:317:ASN:HB2	1:A:320:ASP:CG	2.44	0.42
1:A:74:THR:HG22	1:A:268:LEU:CD1	2.40	0.42
1:A:280:PRO:CB	1:A:324:PHE:CD1	2.87	0.42
1:A:153:THR:OG1	1:A:155:ASP:OD1	2.38	0.42
1:A:32:GLY:O	1:A:82:TYR:CD1	2.74	0.41
1:A:283:HIS:ND1	1:A:350:GLU:OE1	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:329:MET:O	1:A:331:ARG:N	2.53	0.41
1:A:235:PHE:CD1	1:A:235:PHE:C	2.98	0.41
1:A:244:VAL:HB	1:A:261:THR:HG21	2.01	0.41
1:A:329:MET:HG3	1:A:331:ARG:CG	2.51	0.41
1:A:375:PRO:HB2	1:A:380:LYS:NZ	2.35	0.41
1:A:492:THR:OG1	1:A:494:GLU:HG2	2.19	0.41
1:A:288:VAL:HB	1:A:289:PRO:HD2	2.03	0.41
1:A:329:MET:HA	1:A:329:MET:CE	2.51	0.41
1:A:199:SER:O	1:A:203:GLN:HG2	2.21	0.41
1:A:314:ALA:HB3	1:A:401:LEU:CD2	2.44	0.41
1:A:338:LYS:HB2	1:A:377:GLN:OE1	2.20	0.41
1:A:42:LEU:HD23	1:A:42:LEU:N	2.36	0.40
1:A:16:VAL:HG12	1:A:58:LYS:HE3	2.03	0.40
1:A:21:GLY:O	1:A:23:LEU:N	2.53	0.40
1:A:92:LYS:O	1:A:93:GLN:CG	2.60	0.40
1:A:193:GLU:CD	1:A:194:SER:HB2	2.45	0.40
1:A:147:PRO:HD2	3:A:827:HOH:O	2.21	0.40
1:A:521:ILE:HG22	1:A:522:HIS:N	2.36	0.40
1:A:151:LEU:HD13	1:A:287:PHE:CE1	2.56	0.40
1:A:349:SER:O	1:A:352:THR:OG1	2.24	0.40
1:A:380:LYS:O	1:A:383:THR:HB	2.22	0.40
1:A:134:GLY:O	1:A:136:VAL:HG13	2.21	0.40
1:A:282:LEU:HD22	1:A:351:PHE:HD2	1.85	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	511/522 (98%)	451 (88%)	43 (8%)	17 (3%)	3 5

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	21	GLY
1	A	120	LEU
1	A	121	ASN
1	A	271	LYS
1	A	521	ILE
1	A	23	LEU
1	A	116	GLY
1	A	340	THR
1	A	358	ARG
1	A	513	SER
1	A	92	LYS
1	A	429	LYS
1	A	93	GLN
1	A	62	LYS
1	A	330	PRO
1	A	431	VAL
1	A	374	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	419/426 (98%)	396 (94%)	23 (6%)	19	41

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

Mol	Chain	Res	Type
1	A	72	ASP
1	A	75	TYR
1	A	91	ARG
1	A	111	MET
1	A	119	PHE
1	A	120	LEU
1	A	152	SER
1	A	178	ILE
1	A	191	PHE
1	A	194	SER

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Mol	Chain	Res	Type
1	A	225	SER
1	A	253	ARG
1	A	283	HIS
1	A	316	THR
1	A	320	ASP
1	A	329	MET
1	A	331	ARG
1	A	348	VAL
1	A	352	THR
1	A	371	TRP
1	A	383	THR
1	A	494	GLU
1	A	521	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	GLN
1	A	59	ASN
1	A	93	GLN
1	A	121	ASN
1	A	186	ASN
1	A	373	GLN
1	A	495	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](#)

1 ligand is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	600	-	4,4,4	0.43	0	6,6,6	0.06	0

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 ⓘ

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	515/522 (98%)	0.53	50 (9%) 13 10	17, 42, 81, 95	7 (1%)

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	120	LEU	6.0
1	A	124	LEU	5.5
1	A	280	PRO	5.5
1	A	119	PHE	5.4
1	A	123	TYR	5.2
1	A	284	TYR	4.1
1	A	521	ILE	4.1
1	A	353	ILE	4.0
1	A	356	GLY	3.9
1	A	121	ASN	3.7
1	A	22	LEU	3.6
1	A	111	MET	3.4
1	A	522	HIS	3.3
1	A	352	THR	3.3
1	A	117	ALA	3.3
1	A	122	ASN	3.1
1	A	269	ALA	3.1
1	A	1	ALA	3.1
1	A	428	PRO	3.0
1	A	272	VAL	3.0
1	A	114	GLY	2.9
1	A	372	ALA	2.9
1	A	345	TYR	2.8
1	A	23	LEU	2.7
1	A	358	ARG	2.7
1	A	371	TRP	2.7
1	A	493	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	270	TYR	2.6
1	A	244	VAL	2.6
1	A	247	PRO	2.6
1	A	118	ASN	2.5
1	A	113	SER	2.5
1	A	281	MET	2.5
1	A	355	LYS	2.4
1	A	426	VAL	2.4
1	A	258	LEU	2.3
1	A	368	THR	2.3
1	A	427	TYR	2.3
1	A	370	SER	2.2
1	A	3	LEU	2.2
1	A	282	LEU	2.2
1	A	248	VAL	2.2
1	A	268	LEU	2.1
1	A	75	TYR	2.1
1	A	24	GLY	2.1
1	A	286	GLY	2.1
1	A	240	VAL	2.0
1	A	351	PHE	2.0
1	A	357	LEU	2.0
1	A	348	VAL	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.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	A	600	5/5	0.76	0.20	52,52,53,53	5

6.5 Other polymers [i](#)

There are no such residues in this entry.