



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 07:56 PM UTC

PDB ID : 1ZLP / pdb_00001zlp
Title : Petal death protein PSR132 with cysteine-linked glutaraldehyde forming a thiohemiacetal adduct
Authors : Teplyakov, A.; Liu, S.; Lu, Z.; Howard, A.; Dunaway-Mariano, D.; Herzberg, O.
Deposited on : 2005-05-08
Resolution : 2.70 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

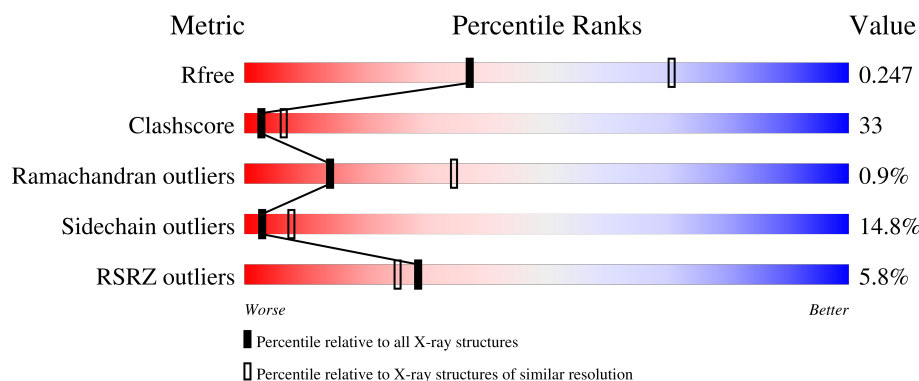
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	318	4% 46% 36% 7% 11%
1	B	318	6% 47% 33% 9% 10%

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	GAQ	A	402	-	-	X	-
3	GAQ	B	402	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4411 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.

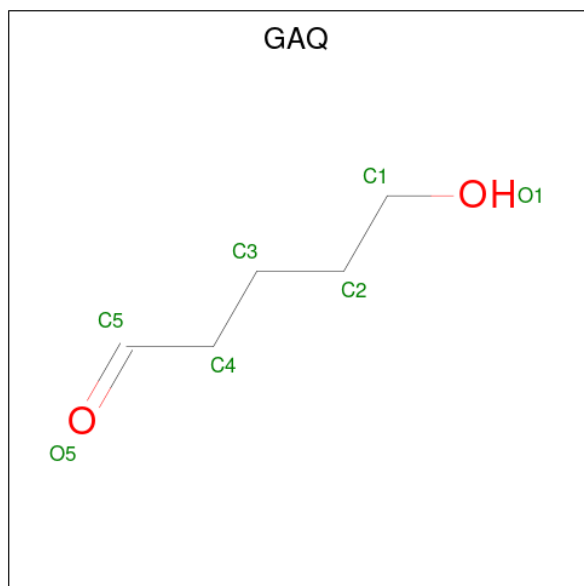
- Molecule 1 is a protein called petal death protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	284	Total	C	N	O	S	0	0	0
			2159	1364	375	409	11			
1	B	285	Total	C	N	O	S	0	0	0
			2167	1368	377	411	11			

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 3 is 5-HYDROXYPENTANAL (CCD ID: GAQ) (formula: C₅H₁₀O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	5	2		
3	B	1	Total	C	O	0	0
			7	5	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	38	Total	O	0	0
			38	38		
4	B	31	Total	O	0	0
			31	31		

4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	156.19Å 156.19Å 75.05Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 2.70 10.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	100.0 (10.00-2.70) 88.0 (10.00-2.70)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.88 (at 2.68Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.195 , 0.250 0.205 , 0.247	Depositor DCC
R_{free} test set	2045 reflections (7.77%)	wwPDB-VP
Wilson B-factor (Å ²)	57.0	Xtriage
Anisotropy	1.144	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 94.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.024 for -h,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4411	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% 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: GAQ, 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	0.82	1/2198 (0.0%)	1.18	11/2969 (0.4%)
1	B	0.58	0/2206	0.98	3/2980 (0.1%)
All	All	0.71	1/4404 (0.0%)	1.09	14/5949 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	298	ILE	CA-CB	7.89	1.63	1.54

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	299	SER	N-CA-C	11.73	127.58	111.74
1	A	257	SER	N-CA-C	8.87	120.71	111.14
1	A	66	SER	N-CA-C	8.03	120.71	109.15
1	A	148	ARG	N-CA-C	7.12	118.69	111.07
1	A	299	SER	N-CA-CB	-6.98	101.59	112.13
1	A	136	GLU	N-CA-C	6.38	120.10	109.95
1	B	148	ARG	N-CA-C	6.25	118.09	111.28
1	A	82	LEU	N-CA-C	-6.17	105.66	113.43
1	A	298	ILE	CB-CA-C	-6.01	104.31	111.65
1	A	106	VAL	N-CA-C	5.56	116.31	108.36
1	B	257	SER	N-CA-C	5.39	117.58	111.11
1	A	261	VAL	CB-CA-C	-5.35	104.91	112.14
1	A	137	ASP	N-CA-C	5.35	118.90	112.38
1	B	254	ILE	N-CA-C	5.34	114.43	106.85

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	2159	0	2164	142	0
1	B	2167	0	2170	147	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	7	0	9	6	0
3	B	7	0	9	5	0
4	A	38	0	0	4	0
4	B	31	0	0	1	0
All	All	4411	0	4352	287	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 33.

All (287) 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:29:THR:HG22	1:A:32:HIS:CD2	1.66	1.28
1:A:43:MET:CE	1:A:105:VAL:HB	1.65	1.26
1:A:43:MET:HE1	1:A:105:VAL:HB	1.23	1.08
1:A:29:THR:CG2	1:A:32:HIS:CD2	2.43	1.01
1:A:295:ASN:N	1:A:295:ASN:HD22	1.58	1.00
1:A:213:ASN:C	1:A:213:ASN:HD22	1.72	0.97
1:A:142:LYS:HE3	1:A:143:LYS:O	1.65	0.96
1:B:280:THR:HG22	1:B:282:ARG:H	1.30	0.94
1:A:47:GLN:HB3	1:A:69:SER:HB3	1.52	0.88
1:A:29:THR:CG2	1:A:32:HIS:HD2	1.82	0.88
1:A:43:MET:HE1	1:A:105:VAL:CB	2.03	0.88
1:B:69:SER:OG	1:B:261:VAL:HG21	1.74	0.88
1:B:254:ILE:HD12	1:B:254:ILE:N	1.89	0.87
1:B:280:THR:CG2	1:B:282:ARG:HB3	2.06	0.86
1:B:281:THR:HG21	1:B:288:MET:SD	2.16	0.86
1:A:295:ASN:HD22	1:A:295:ASN:H	1.19	0.85
1:A:145:GLY:H	3:A:402:GAQ:H31	1.43	0.83
1:A:228:ARG:H	1:A:252:HIS:CD2	1.96	0.83
1:B:133:VAL:HG22	1:B:174:PHE:CE2	2.13	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:ASP:HB3	1:A:152:VAL:HG22	1.59	0.83
1:B:40:SER:HB3	1:B:254:ILE:HD11	1.60	0.82
1:B:265:ALA:O	1:B:269:VAL:HG23	1.78	0.81
1:A:43:MET:HE3	1:A:105:VAL:HB	1.61	0.81
1:B:295:ASN:HD22	1:B:300:LEU:HD22	1.46	0.80
1:A:227:LEU:HA	1:A:252:HIS:CD2	2.17	0.80
1:B:233:ILE:HD11	3:B:402:GAQ:H32	1.63	0.80
1:B:256:HIS:HD2	1:B:259:THR:OG1	1.65	0.79
1:B:30:THR:O	1:B:34:LEU:HB2	1.82	0.79
1:A:29:THR:HG22	1:A:32:HIS:HD2	1.06	0.79
1:B:133:VAL:HG22	1:B:174:PHE:HE2	1.49	0.78
1:B:227:LEU:HA	1:B:252:HIS:CD2	2.21	0.76
1:A:295:ASN:N	1:A:295:ASN:ND2	2.31	0.76
1:B:292:SER:O	1:B:296:GLU:CB	2.35	0.75
1:B:280:THR:HG21	1:B:282:ARG:HB3	1.69	0.74
1:B:229:ILE:CG2	1:B:230:ALA:N	2.49	0.73
1:A:144:CYS:SG	1:A:145:GLY:N	2.60	0.73
1:B:133:VAL:CG2	1:B:174:PHE:HE2	2.01	0.73
1:B:280:THR:HG22	1:B:282:ARG:N	2.04	0.73
1:A:295:ASN:H	1:A:295:ASN:ND2	1.82	0.72
1:B:299:SER:HB3	1:B:302:SER:HB3	1.71	0.72
1:B:152:VAL:HB	1:B:194:ARG:NH2	2.06	0.71
1:B:184:ALA:HB3	1:B:185:PRO:HD3	1.73	0.71
1:B:227:LEU:HD23	1:B:252:HIS:CD2	2.26	0.71
1:B:190:GLU:OE1	1:B:190:GLU:HA	1.91	0.71
1:B:190:GLU:OE2	1:B:194:ARG:HD2	1.91	0.71
1:B:280:THR:HG22	1:B:282:ARG:HB3	1.71	0.70
1:B:229:ILE:HG13	1:B:253:LEU:HB3	1.72	0.70
1:B:295:ASN:HB3	1:B:300:LEU:HB2	1.72	0.70
1:A:301:GLU:OE1	1:A:301:GLU:HA	1.90	0.70
1:A:29:THR:HG23	1:A:32:HIS:H	1.57	0.70
1:A:29:THR:HG22	1:A:32:HIS:CG	2.26	0.69
1:B:47:GLN:HB3	1:B:69:SER:HB3	1.74	0.69
1:B:56:GLU:HB2	1:B:102:LEU:HG	1.75	0.69
1:B:69:SER:OG	1:B:261:VAL:CG2	2.41	0.68
1:A:158:HIS:O	1:A:162:ILE:HG12	1.95	0.67
1:A:213:ASN:C	1:A:213:ASN:ND2	2.46	0.67
1:A:145:GLY:N	3:A:402:GAQ:H31	2.10	0.66
1:A:133:VAL:HG13	1:A:174:PHE:CZ	2.30	0.66
1:B:288:MET:HE3	1:B:288:MET:H	1.59	0.66
1:A:71:SER:HB3	1:A:77:LEU:O	1.95	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:294:PHE:O	1:B:297:LEU:HB2	1.96	0.65
1:B:292:SER:O	1:B:296:GLU:HB2	1.95	0.65
1:A:233:ILE:CD1	3:A:402:GAQ:H22	2.27	0.64
1:A:144:CYS:HB3	1:A:147:MET:HG3	1.78	0.64
1:B:43:MET:HE1	1:B:105:VAL:HG21	1.78	0.64
1:B:179:ARG:HD2	1:B:207:PHE:CE2	2.33	0.64
1:A:106:VAL:HG21	1:A:125:LEU:HD22	1.80	0.64
1:A:281:THR:HG22	1:A:285:LEU:HD23	1.80	0.64
1:B:96:THR:HG22	1:B:97:ALA:N	2.13	0.64
1:B:291:PHE:N	1:B:291:PHE:HD1	1.97	0.63
1:B:253:LEU:C	1:B:254:ILE:HD12	2.23	0.63
1:B:298:ILE:O	1:B:299:SER:HB2	1.96	0.63
1:A:290:THR:HB	1:A:293:GLU:HB2	1.80	0.63
1:A:133:VAL:HG13	1:A:174:PHE:HZ	1.62	0.62
1:A:213:ASN:ND2	1:A:216:GLU:H	1.97	0.62
1:A:306:MET:HE1	1:B:167:GLU:OE1	1.99	0.62
1:B:292:SER:O	1:B:296:GLU:HB3	1.98	0.62
1:B:177:VAL:HG13	1:B:229:ILE:HD13	1.82	0.62
1:B:227:LEU:HA	1:B:252:HIS:HD2	1.62	0.61
1:B:291:PHE:N	1:B:291:PHE:CD1	2.65	0.61
1:A:43:MET:HE2	1:A:64:PHE:HA	1.83	0.61
1:A:142:LYS:CE	1:A:143:LYS:O	2.46	0.61
1:B:58:THR:HG22	1:B:58:THR:O	2.00	0.61
1:B:256:HIS:CD2	1:B:259:THR:OG1	2.52	0.61
1:A:153:VAL:O	1:A:194:ARG:NH2	2.34	0.61
1:A:84:THR:O	1:A:88:VAL:HG23	2.00	0.60
1:B:69:SER:CB	1:B:261:VAL:HG21	2.32	0.60
1:A:192:ILE:HD13	1:A:223:LYS:HB2	1.83	0.60
1:A:260:ALA:HB3	4:A:405:HOH:O	2.01	0.60
1:A:228:ARG:H	1:A:252:HIS:HD2	1.46	0.59
1:B:213:ASN:HD22	1:B:213:ASN:C	2.10	0.59
1:A:142:LYS:HD3	1:A:142:LYS:C	2.27	0.59
1:B:152:VAL:HB	1:B:194:ARG:HH21	1.67	0.59
1:B:295:ASN:O	1:B:299:SER:N	2.34	0.59
1:A:144:CYS:HB3	1:A:147:MET:CG	2.33	0.58
1:B:182:ALA:HB3	1:B:191:GLY:HA2	1.85	0.58
1:B:288:MET:HE3	1:B:288:MET:N	2.17	0.58
1:B:71:SER:HB3	1:B:77:LEU:O	2.02	0.58
1:A:152:VAL:CG1	1:A:194:ARG:NH2	2.67	0.58
1:B:301:GLU:HA	1:B:301:GLU:OE1	2.04	0.58
1:B:213:ASN:ND2	1:B:216:GLU:H	2.01	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:229:ILE:HG22	1:B:230:ALA:N	2.18	0.58
1:A:295:ASN:O	1:A:299:SER:N	2.37	0.58
1:A:74:MET:HE2	1:A:94:ARG:HH12	1.68	0.58
1:A:280:THR:HG23	1:A:282:ARG:HB3	1.86	0.57
1:A:43:MET:HE2	1:A:64:PHE:CA	2.34	0.57
1:B:229:ILE:HG23	1:B:230:ALA:H	1.69	0.57
1:A:184:ALA:HB3	1:A:185:PRO:HD3	1.87	0.57
1:A:64:PHE:CZ	1:A:107:ASP:HB2	2.40	0.57
1:B:110:THR:HB	4:B:405:HOH:O	2.03	0.57
1:A:184:ALA:HB3	1:A:185:PRO:CD	2.34	0.57
1:B:238:THR:CG2	1:B:239:PRO:HD2	2.34	0.57
1:B:254:ILE:N	1:B:254:ILE:CD1	2.61	0.57
1:A:259:THR:HB	4:A:404:HOH:O	2.04	0.57
1:B:238:THR:HG23	1:B:239:PRO:HD2	1.87	0.57
1:A:117:ASN:H	1:A:117:ASN:HD22	1.51	0.57
1:B:133:VAL:CG2	1:B:174:PHE:CE2	2.83	0.57
1:A:58:THR:O	1:A:58:THR:HG22	2.05	0.56
1:A:143:LYS:HD2	1:A:149:GLY:O	2.05	0.56
1:B:117:ASN:HD22	1:B:117:ASN:H	1.53	0.56
1:B:229:ILE:HG23	1:B:230:ALA:N	2.20	0.56
1:A:43:MET:CE	1:A:105:VAL:CB	2.60	0.56
1:B:66:SER:O	1:B:70:VAL:HG23	2.05	0.56
1:B:290:THR:O	1:B:294:PHE:N	2.34	0.56
1:B:295:ASN:HB3	1:B:300:LEU:CB	2.36	0.55
1:B:29:THR:HG22	1:B:31:MET:H	1.70	0.55
1:A:125:LEU:HD13	1:A:133:VAL:CG1	2.36	0.55
1:B:284:ASP:OD1	1:B:284:ASP:N	2.33	0.55
1:A:56:GLU:HB2	1:A:102:LEU:HG	1.89	0.55
1:B:29:THR:HG22	1:B:30:THR:N	2.20	0.55
1:B:47:GLN:CB	1:B:69:SER:HB3	2.37	0.55
1:A:179:ARG:HD2	1:A:207:PHE:CE2	2.42	0.54
1:A:43:MET:HE1	1:A:105:VAL:CG1	2.38	0.53
1:A:213:ASN:HD21	1:A:216:GLU:H	1.54	0.53
1:B:140:TRP:CD2	1:B:141:PRO:HA	2.42	0.53
1:A:69:SER:HA	1:A:261:VAL:HG21	1.89	0.53
1:A:176:LEU:HB3	1:A:204:ASP:OD1	2.09	0.53
1:B:166:ARG:O	1:B:166:ARG:HD3	2.08	0.53
1:A:244:GLU:CD	1:A:244:GLU:H	2.17	0.53
1:A:47:GLN:HG3	1:A:73:ALA:HB2	1.91	0.53
1:B:40:SER:CB	1:B:254:ILE:HD11	2.36	0.53
1:A:183:ARG:NH1	1:A:216:GLU:OE1	2.41	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:70:VAL:O	1:B:74:MET:HB3	2.09	0.52
1:B:295:ASN:OD1	1:B:295:ASN:N	2.42	0.52
1:B:179:ARG:NH1	3:B:402:GAQ:O5	2.43	0.52
1:B:258:LEU:HD11	3:B:402:GAQ:O1	2.10	0.51
1:A:162:ILE:O	1:A:165:ALA:HB3	2.11	0.51
1:B:85:THR:O	1:B:89:VAL:HG23	2.09	0.51
1:A:29:THR:HG21	1:A:175:PHE:HD1	1.76	0.51
1:A:101:ASN:O	1:A:102:LEU:C	2.52	0.51
1:B:83:LEU:HD21	1:B:88:VAL:HG12	1.92	0.51
1:B:210:ALA:N	1:B:211:PRO:HD3	2.26	0.51
1:B:271:ILE:HD11	1:B:288:MET:HE2	1.91	0.51
1:A:214:VAL:HG13	1:A:249:MET:HE2	1.92	0.51
1:B:156:GLU:CD	1:B:156:GLU:H	2.18	0.51
1:B:227:LEU:CD2	1:B:252:HIS:CD2	2.93	0.51
1:A:299:SER:HB3	1:A:302:SER:HB3	1.93	0.50
1:A:96:THR:HG21	1:A:129:GLY:HA3	1.93	0.50
1:A:136:GLU:HG3	4:A:412:HOH:O	2.11	0.50
1:A:228:ARG:N	1:A:252:HIS:CD2	2.75	0.50
1:B:78:PRO:O	1:B:80:PHE:N	2.43	0.50
1:A:106:VAL:HG21	1:A:125:LEU:CD2	2.41	0.50
1:A:227:LEU:HA	1:A:252:HIS:HD2	1.73	0.50
1:A:57:LYS:HD3	1:A:57:LYS:C	2.37	0.50
1:A:229:ILE:HG13	1:A:253:LEU:HB3	1.93	0.50
1:B:117:ASN:H	1:B:117:ASN:ND2	2.09	0.50
1:A:47:GLN:HB3	1:A:69:SER:CB	2.35	0.50
1:B:242:THR:N	1:B:245:GLU:OE1	2.38	0.50
1:A:108:GLY:O	1:A:109:ASP:C	2.54	0.49
1:B:43:MET:SD	1:B:64:PHE:HB2	2.52	0.49
1:B:177:VAL:CG1	1:B:229:ILE:HD13	2.42	0.49
1:A:30:THR:HA	1:A:33:ARG:NH1	2.27	0.49
1:B:152:VAL:CG2	1:B:194:ARG:NH2	2.75	0.49
1:B:228:ARG:H	1:B:252:HIS:CD2	2.30	0.49
1:A:238:THR:HG23	1:A:239:PRO:HD2	1.95	0.49
1:B:282:ARG:HG2	1:B:283:ASP:N	2.25	0.49
1:B:299:SER:O	1:B:300:LEU:C	2.55	0.49
1:A:85:THR:O	1:A:89:VAL:HG23	2.13	0.49
1:A:47:GLN:CB	1:A:69:SER:HB3	2.34	0.49
1:A:68:TYR:CD1	1:A:258:LEU:HD22	2.47	0.49
1:B:33:ARG:O	1:B:37:GLU:HG3	2.13	0.48
1:B:288:MET:N	1:B:288:MET:CE	2.76	0.48
1:B:179:ARG:HH12	3:B:402:GAQ:H42	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:171:ASP:O	1:A:172:SER:C	2.56	0.48
1:B:43:MET:CE	1:B:105:VAL:HG21	2.43	0.48
1:B:227:LEU:CA	1:B:252:HIS:HD2	2.26	0.48
1:A:290:THR:CG2	1:A:291:PHE:N	2.77	0.48
1:B:47:GLN:O	1:B:48:ASP:HB3	2.13	0.48
1:A:306:MET:HE3	1:B:119:GLN:HE22	1.78	0.48
1:B:290:THR:HG22	1:B:291:PHE:H	1.78	0.48
1:A:233:ILE:CD1	3:A:402:GAQ:C2	2.92	0.48
1:B:68:TYR:CD2	1:B:79:ASP:HB2	2.49	0.48
1:B:144:CYS:HB3	1:B:147:MET:SD	2.53	0.47
1:B:295:ASN:ND2	1:B:300:LEU:HD22	2.23	0.47
1:B:297:LEU:HD23	1:B:297:LEU:HA	1.77	0.47
1:B:121:PHE:CE1	1:B:125:LEU:HD21	2.50	0.47
1:A:185:PRO:C	1:A:186:HIS:ND1	2.73	0.47
1:A:256:HIS:HD2	1:A:259:THR:OG1	1.98	0.47
1:A:46:VAL:HG23	1:A:55:VAL:HG21	1.97	0.46
1:A:280:THR:CG2	1:A:282:ARG:HB3	2.44	0.46
1:A:307:GLU:OE1	1:A:311:LYS:HE3	2.15	0.46
1:A:173:ASP:C	1:A:173:ASP:OD1	2.58	0.46
1:B:131:LYS:O	1:B:174:PHE:HA	2.15	0.46
1:B:279:GLY:O	1:B:280:THR:OG1	2.31	0.46
1:A:233:ILE:HD13	3:A:402:GAQ:H22	1.97	0.46
1:A:29:THR:HG21	1:A:32:HIS:CD2	2.45	0.46
1:A:74:MET:HE1	1:A:91:ALA:HB2	1.98	0.46
1:A:290:THR:HG22	1:A:291:PHE:N	2.31	0.46
1:B:39:GLY:HA3	1:B:247:LYS:HD2	1.96	0.46
1:A:227:LEU:CA	1:A:252:HIS:HD2	2.29	0.46
1:A:125:LEU:HD13	1:A:133:VAL:HG12	1.98	0.46
1:A:117:ASN:H	1:A:117:ASN:ND2	2.15	0.45
1:B:47:GLN:HG3	1:B:73:ALA:CB	2.46	0.45
1:A:118:VAL:O	1:A:122:ILE:HG12	2.17	0.45
1:A:300:LEU:O	1:A:304:TYR:HD2	1.99	0.45
1:A:301:GLU:OE1	1:A:301:GLU:CA	2.58	0.45
1:B:31:MET:CE	1:B:253:LEU:HD21	2.47	0.45
1:B:155:ALA:O	1:B:156:GLU:C	2.59	0.45
1:A:281:THR:HG21	1:A:288:MET:CE	2.46	0.45
1:B:232:MET:HE2	1:B:232:MET:HA	1.99	0.45
1:A:137:ASP:HB2	1:A:181:ASP:H	1.81	0.45
1:A:152:VAL:HG13	1:A:194:ARG:NH2	2.31	0.45
1:A:242:THR:HB	1:A:243:PRO:HD2	1.99	0.45
1:A:245:GLU:O	1:A:248:GLU:HB2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:134:PHE:CE1	1:B:177:VAL:HG21	2.52	0.44
1:A:110:THR:O	1:A:110:THR:CG2	2.63	0.44
1:A:291:PHE:O	1:A:295:ASN:ND2	2.51	0.44
1:B:114:GLY:O	1:B:115:PRO:C	2.61	0.44
1:B:137:ASP:HB2	1:B:181:ASP:H	1.82	0.44
1:B:234:GLU:H	1:B:234:GLU:HG2	1.44	0.44
1:A:253:LEU:HD23	1:A:253:LEU:HA	1.85	0.44
1:A:109:ASP:HB3	1:A:142:LYS:HG2	2.00	0.44
1:A:228:ARG:N	1:A:252:HIS:HD2	2.12	0.43
1:B:114:GLY:H	1:B:117:ASN:ND2	2.16	0.43
1:A:82:LEU:HD23	1:A:82:LEU:HA	1.74	0.43
1:B:152:VAL:CB	1:B:194:ARG:NH2	2.78	0.43
1:B:245:GLU:O	1:B:248:GLU:HB2	2.18	0.43
1:A:258:LEU:CB	1:A:262:TYR:CE2	3.01	0.43
1:A:52:ALA:HB2	1:A:95:ILE:HG23	2.00	0.43
1:A:137:ASP:HB3	1:A:152:VAL:CG2	2.39	0.43
1:B:169:ILE:CD1	1:B:174:PHE:HD2	2.31	0.43
1:B:253:LEU:HD23	1:B:253:LEU:HA	1.80	0.43
1:A:50:LEU:O	1:A:54:VAL:HG23	2.19	0.43
1:A:69:SER:OG	1:A:261:VAL:CG2	2.67	0.43
1:A:47:GLN:NE2	4:A:407:HOH:O	2.51	0.43
1:A:184:ALA:N	1:A:185:PRO:HD2	2.34	0.43
1:B:47:GLN:HG3	1:B:73:ALA:HB2	2.01	0.43
1:B:271:ILE:HD11	1:B:288:MET:CE	2.48	0.43
1:B:290:THR:HG22	1:B:291:PHE:N	2.34	0.43
1:B:242:THR:HB	1:B:243:PRO:HD2	2.01	0.42
1:A:233:ILE:HD11	3:A:402:GAQ:C2	2.49	0.42
1:A:303:TRP:O	1:A:304:TYR:C	2.62	0.42
1:B:134:PHE:CD1	1:B:177:VAL:HB	2.55	0.42
1:B:214:VAL:O	1:B:215:ASP:C	2.63	0.42
1:B:179:ARG:HH12	3:B:402:GAQ:C4	2.32	0.42
1:B:290:THR:C	1:B:291:PHE:HD1	2.27	0.42
1:A:90:GLU:O	1:A:91:ALA:C	2.61	0.42
1:A:69:SER:CB	1:A:261:VAL:HG21	2.50	0.42
1:A:227:LEU:HA	1:A:252:HIS:NE2	2.34	0.42
1:A:85:THR:CG2	1:A:86:THR:N	2.82	0.42
1:B:213:ASN:HA	1:B:239:PRO:HD3	2.02	0.42
1:B:300:LEU:O	1:B:301:GLU:C	2.63	0.42
1:A:281:THR:OG1	1:A:288:MET:HE1	2.20	0.41
1:B:190:GLU:O	1:B:191:GLY:C	2.62	0.41
1:B:291:PHE:O	1:B:292:SER:C	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:SER:CB	1:B:261:VAL:CG2	2.98	0.41
1:B:74:MET:HE3	1:B:74:MET:HB2	1.71	0.41
1:B:308:SER:C	1:B:310:PHE:N	2.77	0.41
1:A:143:LYS:C	1:A:144:CYS:O	2.63	0.41
1:A:183:ARG:NH1	1:A:216:GLU:OE2	2.53	0.41
1:A:242:THR:HB	1:A:244:GLU:OE1	2.20	0.41
1:B:98:ALA:C	1:B:100:PRO:HD3	2.46	0.41
1:A:35:ILE:O	1:A:38:HIS:O	2.38	0.41
1:A:281:THR:CG2	1:A:288:MET:HE1	2.51	0.41
1:B:101:ASN:O	1:B:102:LEU:C	2.64	0.41
1:B:279:GLY:C	1:B:280:THR:OG1	2.63	0.41
1:A:42:LEU:HB2	1:A:243:PRO:HG3	2.03	0.40
1:B:117:ASN:ND2	1:B:117:ASN:N	2.69	0.40
1:A:258:LEU:HB3	1:A:262:TYR:CE2	2.57	0.40
1:B:166:ARG:HD3	1:B:166:ARG:C	2.46	0.40
1:A:47:GLN:HG3	1:A:73:ALA:CB	2.51	0.40
1:B:150:LYS:HE3	1:B:150:LYS:HB2	1.75	0.40
1:B:158:HIS:O	1:B:159:ALA:C	2.62	0.40
1:A:41:VAL:HB	1:A:253:LEU:HD23	2.03	0.40
1:A:169:ILE:O	1:A:172:SER:HB2	2.21	0.40
1:B:140:TRP:CG	1:B:141:PRO:HA	2.57	0.40
1:B:273:LYS:CD	1:B:277:GLU:OE2	2.69	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	282/318 (89%)	251 (89%)	27 (10%)	4 (1%)	9	23
1	B	283/318 (89%)	259 (92%)	23 (8%)	1 (0%)	30	54
All	All	565/636 (89%)	510 (90%)	50 (9%)	5 (1%)	14	35

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	282	ARG
1	A	103	CYS
1	A	144	CYS
1	A	304	TYR
1	B	48	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	223/249 (90%)	195 (87%)	28 (13%)	4	11
1	B	224/249 (90%)	186 (83%)	38 (17%)	2	6
All	All	447/498 (90%)	381 (85%)	66 (15%)	3	8

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LYS
1	A	40	SER
1	A	46	VAL
1	A	75	LEU
1	A	96	THR
1	A	110	THR
1	A	117	ASN
1	A	142	LYS
1	A	152	VAL
1	A	183	ARG
1	A	188	LEU
1	A	206	THR
1	A	213	ASN
1	A	234	GLU
1	A	253	LEU
1	A	254	ILE
1	A	256	HIS
1	A	257	SER

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Mol	Chain	Res	Type
1	A	271	ILE
1	A	275	LEU
1	A	281	THR
1	A	287	GLN
1	A	295	ASN
1	A	296	GLU
1	A	299	SER
1	A	301	GLU
1	A	302	SER
1	A	308	SER
1	B	28	LYS
1	B	31	MET
1	B	34	LEU
1	B	40	SER
1	B	46	VAL
1	B	66	SER
1	B	71	SER
1	B	88	VAL
1	B	96	THR
1	B	110	THR
1	B	117	ASN
1	B	120	ARG
1	B	131	LYS
1	B	133	VAL
1	B	142	LYS
1	B	183	ARG
1	B	190	GLU
1	B	206	THR
1	B	213	ASN
1	B	214	VAL
1	B	223	LYS
1	B	229	ILE
1	B	233	ILE
1	B	234	GLU
1	B	253	LEU
1	B	254	ILE
1	B	256	HIS
1	B	261	VAL
1	B	271	ILE
1	B	275	LEU
1	B	282	ARG
1	B	284	ASP

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Mol	Chain	Res	Type
1	B	286	ASP
1	B	288	MET
1	B	292	SER
1	B	295	ASN
1	B	296	GLU
1	B	302	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	HIS
1	A	117	ASN
1	A	158	HIS
1	A	213	ASN
1	A	252	HIS
1	A	256	HIS
1	A	287	GLN
1	A	295	ASN
1	B	38	HIS
1	B	117	ASN
1	B	119	GLN
1	B	186	HIS
1	B	213	ASN
1	B	252	HIS
1	B	256	HIS

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 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	$\# Z > 2$
3	GAQ	B	402	1	5,6,6	1.95	1 (20%)	4,5,5	1.58	2 (50%)
3	GAQ	A	402	1	5,6,6	1.95	1 (20%)	4,5,5	1.57	2 (50%)

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
3	GAQ	B	402	1	-	1/4/4/4	-
3	GAQ	A	402	1	-	2/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	GAQ	O5-C5	4.30	1.44	1.20
3	A	402	GAQ	O5-C5	4.29	1.44	1.20

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	GAQ	C2-C3-C4	-2.28	102.91	113.86
3	B	402	GAQ	O5-C5-C4	-2.25	107.45	126.30
3	B	402	GAQ	C2-C3-C4	-2.13	103.64	113.86
3	A	402	GAQ	O5-C5-C4	-2.12	108.55	126.30

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	402	GAQ	C3-C4-C5-O5
3	B	402	GAQ	C1-C2-C3-C4
3	A	402	GAQ	C2-C3-C4-C5

There are no ring outliers.

2 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	GAQ	5	0
3	A	402	GAQ	6	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	284/318 (89%)	0.27	13 (4%) 37 34	64, 79, 101, 130	0
1	B	285/318 (89%)	0.54	20 (7%) 22 19	68, 80, 100, 125	0
All	All	569/636 (89%)	0.40	33 (5%) 29 25	64, 80, 101, 130	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	292	SER	9.0
1	B	289	ALA	4.2
1	B	296	GLU	3.6
1	A	300	LEU	3.3
1	B	145	GLY	3.1
1	B	141	PRO	2.9
1	B	283	ASP	2.9
1	A	141	PRO	2.9
1	A	302	SER	2.8
1	A	145	GLY	2.7
1	B	312	ASN	2.7
1	A	299	SER	2.5
1	A	291	PHE	2.5
1	B	290	THR	2.4
1	A	282	ARG	2.4
1	B	285	LEU	2.3
1	B	277	GLU	2.3
1	B	40	SER	2.3
1	A	86	THR	2.3
1	A	146	HIS	2.3
1	A	147	MET	2.3
1	A	303	TRP	2.3
1	B	86	THR	2.2
1	A	221	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	73	ALA	2.2
1	B	244	GLU	2.2
1	B	278	LYS	2.1
1	B	310	PHE	2.1
1	B	282	ARG	2.1
1	B	212	ALA	2.1
1	A	283	ASP	2.1
1	B	250	GLY	2.0
1	B	279	GLY	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
3	GAQ	B	402	7/7	0.38	0.26	68,74,81,84	0
3	GAQ	A	402	7/7	0.66	0.19	66,75,80,83	0
2	MG	B	401	1/1	0.79	0.08	98,98,98,98	0
2	MG	A	401	1/1	0.88	0.07	94,94,94,94	0

6.5 Other polymers [i](#)

There are no such residues in this entry.