



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 11:01 PM UTC

PDB ID : 2VT6 / pdb_00002vt6
Title : Native Torpedo californica acetylcholinesterase collected with a cumulated dose of 9400000 Gy
Authors : Colletier, J.P.; Bourgeois, D.; Sanson, B.; Fournier, D.; Sussman, J.L.; Silman, I.; Weik, M.
Deposited on : 2008-05-09
Resolution : 2.40 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

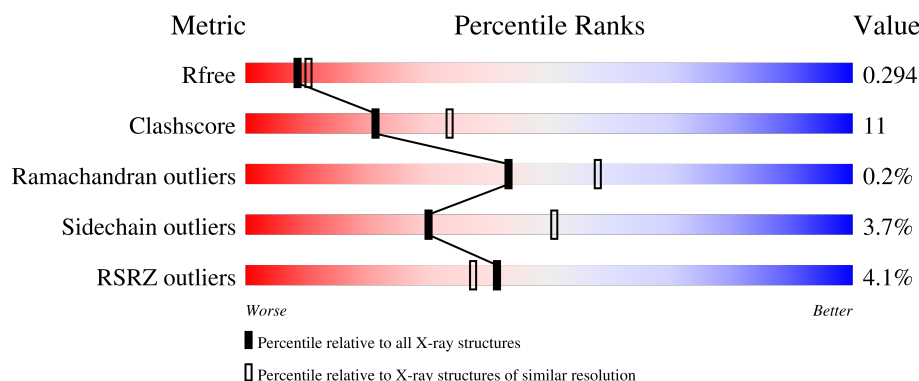
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	537	5% 75% 20% ..
1	B	537	3% 74% 22% ..

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
4	PGE	B	1540	-	X	-	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 9558 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 ACETYLCHOLINESTERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	529	Total	C	N	O	S	0	0	0
			4215	2705	714	774	22			
1	B	532	Total	C	N	O	S	0	0	0
			4242	2720	720	780	22			

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

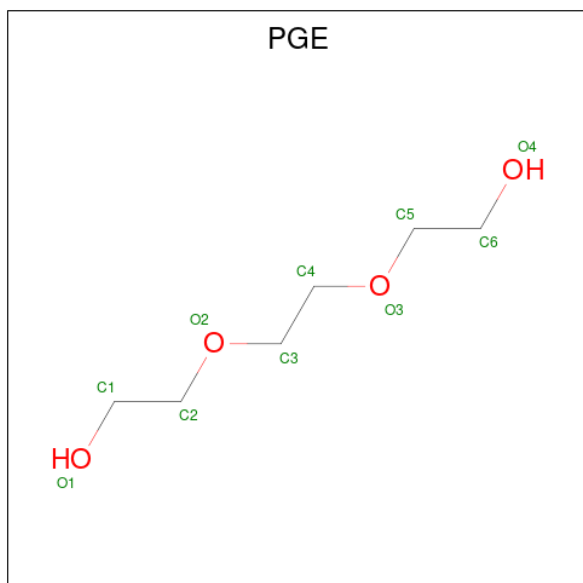


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		
3	B	1	Total	Cl	0	0
			1	1		

- Molecule 4 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			10	6	4		
4	A	1	Total	C	O	0	0
			10	6	4		
4	A	1	Total	C	O	0	0
			10	6	4		
4	A	1	Total	C	O	0	0
			10	6	4		
4	B	1	Total	C	O	0	0
			10	6	4		
4	B	1	Total	C	O	0	0
			10	6	4		
4	B	1	Total	C	O	0	0
			10	6	4		

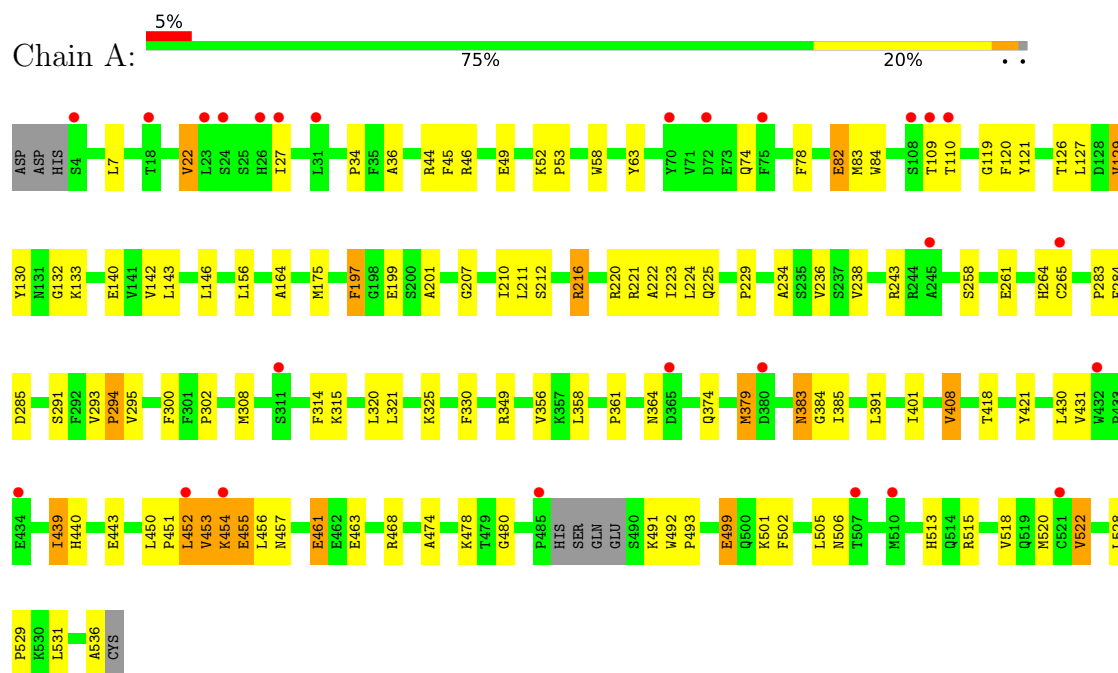
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	468	Total 468	O 468	0	0
5	B	495	Total 495	O 495	0	0

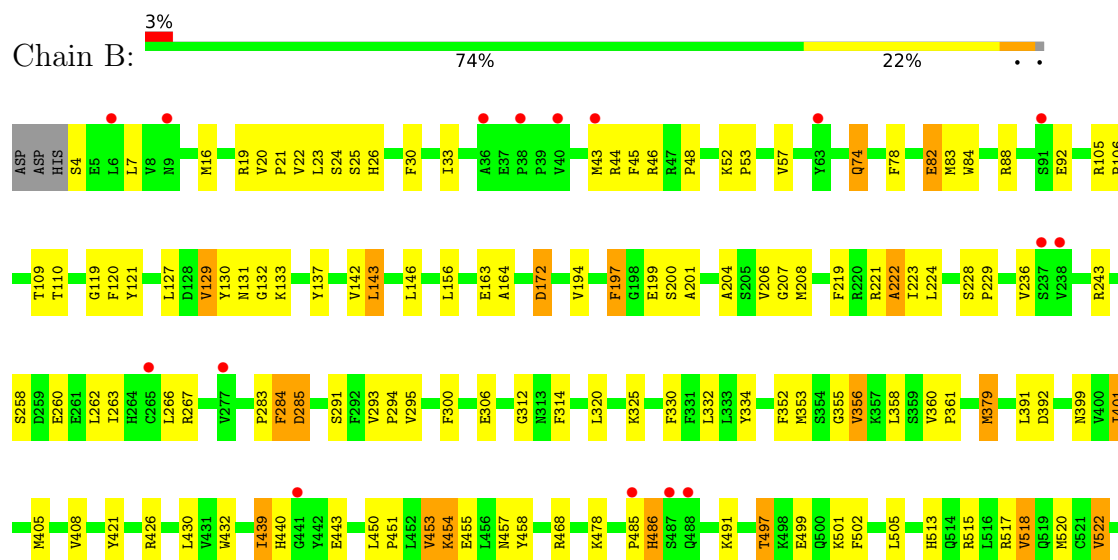
3 Residue-property plots [i](#)

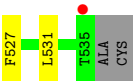
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: ACETYLCHOLINESTERASE



• Molecule 1: ACETYLCHOLINESTERASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	91.66Å 106.08Å 150.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.96 – 2.40 19.96 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.4 (19.96-2.40) 98.2 (19.96-2.40)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.47 (at 2.41Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.193 , 0.234 0.281 , 0.294	Depositor DCC
R_{free} test set	2664 reflections (4.67%)	wwPDB-VP
Wilson B-factor (Å ²)	15.8	Xtriage
Anisotropy	1.040	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 63.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.27$, $\langle L^2 \rangle = 0.11$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	9558	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 22.97 % 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 5.1222e-03.*

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: PGE, CL, NAG

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	0.40	0/4336	0.94	16/5886 (0.3%)
1	B	0.39	0/4365	0.95	22/5926 (0.4%)
All	All	0.39	0/8701	0.95	38/11812 (0.3%)

There are no bond length outliers.

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	291	SER	N-CA-C	8.00	119.69	110.97
1	A	285	ASP	N-CA-C	-7.62	98.39	109.59
1	B	291	SER	N-CA-C	7.44	119.08	110.97
1	A	216	ARG	N-CA-C	7.43	120.32	111.33
1	A	401	ILE	N-CA-C	7.31	117.83	110.23
1	B	401	ILE	N-CA-C	6.93	117.07	110.42
1	B	443	GLU	N-CA-C	-6.87	105.06	113.50
1	B	454	LYS	N-CA-C	6.80	119.56	111.33
1	A	439	ILE	N-CA-C	6.78	118.82	109.80
1	A	443	GLU	N-CA-C	-6.73	104.54	112.89
1	B	527	PHE	N-CA-C	6.72	118.30	110.97
1	B	164	ALA	CA-C-N	6.19	125.75	119.19
1	B	164	ALA	C-N-CA	6.19	125.75	119.19
1	A	379	MET	N-CA-C	-6.06	105.89	113.28
1	A	201	ALA	N-CA-C	-6.02	104.83	111.82
1	B	439	ILE	N-CA-C	5.86	119.04	109.78
1	B	285	ASP	N-CA-C	-5.83	101.21	109.96
1	A	156	LEU	N-CA-C	-5.80	101.46	110.10
1	B	222	ALA	N-CA-C	5.70	117.94	108.99
1	B	156	LEU	N-CA-C	-5.68	101.64	110.10
1	B	522	VAL	N-CA-C	-5.66	105.00	110.72
1	B	300	PHE	N-CA-C	-5.66	105.21	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	164	ALA	CA-C-N	5.65	125.27	119.56
1	A	164	ALA	C-N-CA	5.65	125.27	119.56
1	B	399	ASN	N-CA-C	5.60	120.23	112.90
1	A	454	LYS	N-CA-C	5.58	118.08	111.33
1	A	522	VAL	N-CA-C	-5.51	105.16	110.72
1	B	457	ASN	N-CA-C	5.45	119.10	111.74
1	B	379	MET	N-CA-C	-5.45	106.59	113.18
1	B	172	ASP	N-CA-C	-5.39	105.30	111.07
1	B	201	ALA	N-CA-C	-5.38	105.58	111.82
1	A	300	PHE	N-CA-C	-5.27	105.70	111.82
1	B	360	VAL	CA-C-N	5.22	125.52	119.47
1	B	360	VAL	C-N-CA	5.22	125.52	119.47
1	A	234	ALA	N-CA-C	5.17	119.44	113.18
1	B	491	LYS	N-CA-C	5.17	117.40	109.07
1	A	457	ASN	N-CA-C	5.07	118.23	111.24
1	B	92	GLU	N-CA-C	-5.02	106.67	112.89

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	4215	0	4074	97	0
1	B	4242	0	4091	83	0
2	A	28	0	26	0	0
2	B	28	0	26	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	40	0	56	7	0
4	B	40	0	56	9	0
5	A	468	0	0	14	0
5	B	495	0	0	9	0
All	All	9558	0	8329	184	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 11.

All (184) 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:499:GLU:HG2	1:A:501:LYS:HE3	1.39	1.03
1:A:454:LYS:HA	1:A:454:LYS:HE2	1.56	0.88
1:B:194:VAL:HG12	5:B:2228:HOH:O	1.80	0.80
1:B:194:VAL:HG13	1:B:219:PHE:HA	1.61	0.80
1:B:48:PRO:HG3	1:B:172:ASP:OD1	1.80	0.79
4:B:1542:PGE:O4	4:B:1542:PGE:H32	1.83	0.79
1:B:194:VAL:CG1	1:B:219:PHE:HA	2.15	0.76
1:A:364:ASN:ND2	1:A:536:ALA:HB2	2.04	0.73
1:A:383:ASN:C	1:A:383:ASN:HD22	1.96	0.71
1:A:364:ASN:HD22	1:A:536:ALA:HB2	1.54	0.71
1:A:243:ARG:HH11	1:A:283:PRO:HB3	1.55	0.69
1:A:265:CYS:HB2	5:A:2267:HOH:O	1.91	0.69
1:B:355:GLY:HA3	1:B:391:LEU:HD21	1.75	0.69
1:A:36:ALA:HB2	1:A:175:MET:HE2	1.74	0.68
1:A:430:LEU:HD11	5:A:2403:HOH:O	1.94	0.66
1:A:452:LEU:HD22	1:A:463:GLU:HG3	1.77	0.65
4:B:1539:PGE:H62	4:B:1540:PGE:O1	1.97	0.64
1:B:284:PHE:CD2	1:B:361:PRO:HB2	2.32	0.64
1:A:520:MET:HE2	1:A:520:MET:HA	1.81	0.63
1:B:426:ARG:CZ	1:B:430:LEU:HD23	2.28	0.63
1:B:515:ARG:HB3	1:B:518:VAL:HG13	1.79	0.63
1:B:515:ARG:HB3	1:B:518:VAL:CG1	2.28	0.63
1:B:468:ARG:HD3	5:B:2424:HOH:O	1.99	0.61
1:A:34:PRO:HG3	1:A:58:TRP:CZ2	2.37	0.60
1:A:506:ASN:HB2	5:A:2395:HOH:O	2.03	0.59
1:A:374:GLN:OE1	1:A:520:MET:HE1	2.03	0.58
1:A:132:GLY:HA3	1:A:143:LEU:CD2	2.32	0.58
1:B:353:MET:HG2	5:B:2351:HOH:O	2.04	0.58
1:A:110:THR:OG1	1:A:478:LYS:HG2	2.04	0.57
1:A:384:GLY:H	4:A:1543:PGE:H42	1.69	0.57
1:A:207:GLY:HA3	1:A:229:PRO:HD3	1.86	0.57
1:A:49:GLU:O	1:A:175:MET:HE1	2.05	0.57
1:B:499:GLU:HG2	1:B:501:LYS:HE3	1.87	0.56
1:A:474:ALA:O	1:A:478:LYS:HG3	2.05	0.56
1:A:236:VAL:HG23	1:A:295:VAL:HG12	1.88	0.56
1:A:518:VAL:O	1:A:522:VAL:HG23	2.06	0.55
1:B:43:MET:HA	1:B:46:ARG:HD2	1.88	0.55
1:A:109:THR:CG2	1:A:142:VAL:HG23	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:453:VAL:HG13	1:A:456:LEU:HD12	1.88	0.55
1:A:197:PHE:HB3	1:A:223:ILE:HB	1.88	0.55
1:B:236:VAL:HG23	1:B:295:VAL:HG12	1.88	0.55
1:B:531:LEU:C	1:B:531:LEU:HD23	2.31	0.55
1:B:19:ARG:NH2	1:B:26:HIS:HB2	2.22	0.55
1:B:332:LEU:HD11	1:B:392:ASP:HA	1.89	0.54
1:A:49:GLU:C	1:A:175:MET:HE1	2.32	0.54
1:B:207:GLY:HA3	1:B:229:PRO:HD3	1.88	0.54
1:A:461:GLU:H	1:A:461:GLU:CD	2.14	0.54
1:A:212:SER:O	1:A:216:ARG:HG3	2.08	0.54
1:A:379:MET:HE3	5:A:2361:HOH:O	2.07	0.54
1:A:515:ARG:HB3	1:A:518:VAL:HB	1.90	0.54
1:B:163:GLU:HB3	1:B:267:ARG:HH22	1.73	0.54
1:B:223:ILE:HA	1:B:320:LEU:O	2.09	0.53
4:B:1541:PGE:H42	5:B:2274:HOH:O	2.08	0.53
1:B:30:PHE:HB3	1:B:33:ILE:HD11	1.90	0.53
1:A:450:LEU:N	1:A:451:PRO:CD	2.72	0.52
1:B:197:PHE:HB3	1:B:223:ILE:HB	1.90	0.52
1:B:197:PHE:CB	1:B:223:ILE:HB	2.39	0.52
1:B:204:ALA:O	1:B:208:MET:HG3	2.08	0.52
1:B:199:GLU:OE1	4:B:1540:PGE:H62	2.08	0.52
1:A:383:ASN:ND2	1:A:385:ILE:H	2.08	0.52
1:A:264:HIS:HB2	5:A:2273:HOH:O	2.09	0.52
1:B:163:GLU:HB3	1:B:267:ARG:NH2	2.25	0.52
1:A:384:GLY:N	4:A:1543:PGE:H42	2.25	0.52
1:A:224:LEU:HD12	1:A:224:LEU:N	2.26	0.51
1:A:431:VAL:HB	5:A:2399:HOH:O	2.09	0.51
1:B:121:TYR:OH	4:B:1540:PGE:H1	2.10	0.51
1:A:121:TYR:OH	4:A:1541:PGE:H22	2.11	0.51
1:A:450:LEU:O	1:A:453:VAL:HG13	2.11	0.51
1:B:267:ARG:HH11	4:B:1541:PGE:C1	2.24	0.51
1:B:132:GLY:HA3	1:B:143:LEU:HD22	1.93	0.51
1:A:468:ARG:HD3	5:A:2203:HOH:O	2.10	0.51
1:B:405:MET:HA	1:B:408:VAL:HG12	1.93	0.50
1:B:224:LEU:N	1:B:224:LEU:HD12	2.27	0.50
1:B:110:THR:OG1	1:B:478:LYS:HG2	2.12	0.50
1:B:453:VAL:HA	5:B:2408:HOH:O	2.12	0.49
1:A:220:ARG:HG3	1:A:221:ARG:HG3	1.95	0.49
1:A:83:MET:HE2	1:A:84:TRP:CZ2	2.47	0.49
1:B:44:ARG:O	1:B:45:PHE:HB2	2.13	0.49
1:A:34:PRO:HG3	1:A:58:TRP:CH2	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:452:LEU:HD22	1:A:463:GLU:CG	2.42	0.48
1:B:146:LEU:C	1:B:146:LEU:HD12	2.37	0.48
1:A:197:PHE:CB	1:A:223:ILE:HB	2.43	0.48
1:A:520:MET:HE3	5:A:2168:HOH:O	2.13	0.48
1:B:426:ARG:NH1	1:B:432:TRP:O	2.46	0.48
1:B:405:MET:O	1:B:408:VAL:HG12	2.13	0.48
1:A:211:LEU:HD23	1:A:314:PHE:HB3	1.96	0.47
1:B:515:ARG:HD3	1:B:518:VAL:HG11	1.95	0.47
1:A:453:VAL:HG22	1:A:456:LEU:HG	1.96	0.47
1:A:27:ILE:HD11	1:A:133:LYS:HB2	1.95	0.47
1:A:78:PHE:O	1:A:82:GLU:HB2	2.15	0.47
1:A:454:LYS:HE2	1:A:454:LYS:CA	2.36	0.47
1:B:243:ARG:HH11	1:B:283:PRO:HG3	1.79	0.47
1:A:82:GLU:HB3	5:A:2101:HOH:O	2.15	0.47
1:A:528:LEU:HB3	1:A:529:PRO:HD3	1.97	0.47
1:B:200:SER:HB2	1:B:440:HIS:CE1	2.50	0.46
1:B:74:GLN:HG3	1:B:334:TYR:CE2	2.50	0.46
1:A:132:GLY:HA3	1:A:143:LEU:HD23	1.97	0.46
1:B:78:PHE:O	1:B:82:GLU:HB2	2.16	0.46
1:A:502:PHE:CZ	1:A:513:HIS:HB2	2.50	0.46
1:B:16:MET:HB2	1:B:57:VAL:CG1	2.45	0.46
1:B:439:ILE:HG22	1:B:440:HIS:N	2.31	0.46
1:B:23:LEU:O	1:B:24:SER:HB2	2.15	0.45
1:A:210:ILE:HD11	1:A:222:ALA:CB	2.47	0.45
1:B:258:SER:HB2	1:B:260:GLU:OE2	2.16	0.45
1:B:520:MET:HG2	5:B:2473:HOH:O	2.15	0.45
1:A:119:GLY:O	1:A:120:PHE:HB2	2.17	0.45
1:B:25:SER:HB3	1:B:137:TYR:CE2	2.52	0.45
1:A:258:SER:OG	1:A:261:GLU:HG3	2.15	0.45
1:B:497:THR:HG22	4:B:1542:PGE:O1	2.16	0.45
1:B:16:MET:HB2	1:B:57:VAL:HG11	1.98	0.45
1:A:210:ILE:HD11	1:A:222:ALA:HB3	1.99	0.45
1:A:325:LYS:NZ	5:A:2317:HOH:O	2.50	0.45
1:A:520:MET:HE2	1:A:520:MET:CA	2.46	0.45
1:B:306:GLU:HA	1:B:306:GLU:OE1	2.17	0.45
1:A:408:VAL:HG22	1:A:418:THR:HG21	1.98	0.44
1:A:453:VAL:HG22	1:A:453:VAL:O	2.17	0.44
1:B:502:PHE:CZ	1:B:513:HIS:HB2	2.52	0.44
1:A:383:ASN:HD22	1:A:384:GLY:N	2.13	0.44
1:B:206:VAL:CG1	1:B:222:ALA:HB1	2.48	0.44
1:A:315:LYS:NZ	5:A:2311:HOH:O	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:383:ASN:HA	4:A:1543:PGE:H52	1.99	0.44
1:A:302:PRO:HD2	1:A:308:MET:CE	2.48	0.44
1:A:384:GLY:H	4:A:1543:PGE:H52	1.83	0.44
1:B:83:MET:HE2	1:B:84:TRP:CZ2	2.53	0.44
1:A:531:LEU:C	1:A:531:LEU:HD23	2.43	0.44
1:B:228:SER:HB2	1:B:229:PRO:HD2	1.99	0.44
1:B:485:PRO:O	1:B:486:HIS:C	2.61	0.44
1:A:44:ARG:O	1:A:45:PHE:HB2	2.18	0.43
1:A:293:VAL:HB	1:A:294:PRO:HD2	2.00	0.43
1:B:497:THR:HG21	5:B:2446:HOH:O	2.18	0.43
1:B:450:LEU:N	1:B:451:PRO:CD	2.81	0.43
1:A:238:VAL:HG21	4:A:1542:PGE:H2	2.00	0.43
1:B:260:GLU:H	1:B:260:GLU:CD	2.24	0.43
1:B:293:VAL:HB	1:B:294:PRO:HD2	2.00	0.43
1:A:132:GLY:HA3	1:A:143:LEU:HD22	1.99	0.43
1:B:163:GLU:HB2	1:B:263:ILE:HD13	2.01	0.43
1:B:352:PHE:O	1:B:356:VAL:HG13	2.17	0.43
1:A:27:ILE:HG22	5:A:2137:HOH:O	2.17	0.43
1:A:321:LEU:HD23	1:A:321:LEU:N	2.34	0.43
1:B:109:THR:CG2	1:B:142:VAL:HG23	2.49	0.43
1:A:320:LEU:C	1:A:320:LEU:HD23	2.44	0.43
1:A:491:LYS:O	1:A:493:PRO:HD3	2.18	0.43
4:B:1540:PGE:H4	4:B:1540:PGE:H22	1.81	0.43
1:A:349:ARG:HA	1:A:349:ARG:HD2	1.83	0.43
1:A:127:LEU:HD12	1:A:130:TYR:CE2	2.54	0.43
1:B:194:VAL:HG13	1:B:219:PHE:CA	2.40	0.43
1:B:119:GLY:O	1:B:120:PHE:HB2	2.19	0.42
1:B:131:ASN:OD1	1:B:133:LYS:HG2	2.19	0.42
1:A:46:ARG:HD2	5:A:2040:HOH:O	2.19	0.42
1:A:439:ILE:HG22	1:A:440:HIS:N	2.34	0.42
1:A:22:VAL:O	1:A:22:VAL:HG13	2.19	0.42
1:B:52:LYS:HA	1:B:53:PRO:HD3	1.94	0.42
1:B:22:VAL:O	1:B:22:VAL:HG13	2.19	0.42
1:A:7:LEU:C	1:A:7:LEU:HD13	2.44	0.42
1:A:284:PHE:CD2	1:A:361:PRO:HB2	2.55	0.42
1:B:88:ARG:HG2	1:B:88:ARG:HH21	1.85	0.42
1:A:421:TYR:HB3	1:A:492:TRP:CZ2	2.54	0.42
1:B:4:SER:HB3	1:B:7:LEU:HB3	2.02	0.42
1:B:421:TYR:C	1:B:421:TYR:CD1	2.97	0.42
1:A:27:ILE:HD11	1:A:133:LYS:CB	2.50	0.42
1:B:451:PRO:HA	1:B:458:TYR:CD1	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:461:GLU:CD	1:A:461:GLU:N	2.78	0.41
1:B:401:ILE:HG21	1:B:517:ARG:HD3	2.03	0.41
1:B:454:LYS:HA	1:B:454:LYS:HD2	1.78	0.41
1:A:211:LEU:O	1:A:308:MET:HE1	2.20	0.41
1:B:20:VAL:HA	1:B:21:PRO:HD3	1.84	0.41
1:A:520:MET:HG2	5:A:2450:HOH:O	2.20	0.41
1:B:84:TRP:CE3	4:B:1540:PGE:H5	2.55	0.41
1:B:129:VAL:HG22	5:B:2109:HOH:O	2.20	0.41
1:B:379:MET:HE3	5:B:2357:HOH:O	2.20	0.41
1:B:518:VAL:O	1:B:522:VAL:HG23	2.20	0.41
1:A:199:GLU:HA	1:A:225:GLN:O	2.21	0.41
1:A:221:ARG:HD3	1:A:480:GLY:HA2	2.01	0.41
1:A:391:LEU:HD12	1:A:391:LEU:HA	1.89	0.41
1:B:127:LEU:HD12	1:B:130:TYR:CE2	2.56	0.41
1:B:262:LEU:O	1:B:266:LEU:HG	2.21	0.41
1:A:384:GLY:H	4:A:1543:PGE:C4	2.33	0.41
1:A:129:VAL:HA	1:A:450:LEU:HD11	2.03	0.40
1:A:146:LEU:C	1:A:146:LEU:HD12	2.46	0.40
1:B:105:ARG:HA	1:B:106:PRO:HD3	1.87	0.40
1:A:63:TYR:CD1	1:A:126:THR:HG22	2.56	0.40
1:A:52:LYS:HA	1:A:53:PRO:HD3	1.88	0.40
1:A:453:VAL:CG2	1:A:455:GLU:HG2	2.51	0.40
1:B:312:GLY:HA2	1:B:314:PHE:CE2	2.57	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	525/537 (98%)	504 (96%)	20 (4%)	1 (0%)	43 58
1	B	530/537 (99%)	506 (96%)	23 (4%)	1 (0%)	43 58

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1055/1074 (98%)	1010 (96%)	43 (4%)	2 (0%)	43 58

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	486	HIS
1	A	22	VAL

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	461/469 (98%)	444 (96%)	17 (4%)	30 51
1	B	464/469 (99%)	447 (96%)	17 (4%)	30 51
All	All	925/938 (99%)	891 (96%)	34 (4%)	30 51

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	GLN
1	A	82	GLU
1	A	129	VAL
1	A	140	GLU
1	A	197	PHE
1	A	294	PRO
1	A	330	PHE
1	A	356	VAL
1	A	358	LEU
1	A	383	ASN
1	A	408	VAL
1	A	452	LEU
1	A	453	VAL
1	A	455	GLU
1	A	461	GLU
1	A	499	GLU

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Mol	Chain	Res	Type
1	A	505	LEU
1	B	74	GLN
1	B	82	GLU
1	B	129	VAL
1	B	143	LEU
1	B	197	PHE
1	B	221	ARG
1	B	284	PHE
1	B	285	ASP
1	B	325	LYS
1	B	330	PHE
1	B	356	VAL
1	B	358	LEU
1	B	453	VAL
1	B	455	GLU
1	B	497	THR
1	B	505	LEU
1	B	518	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	ASN
1	A	68	GLN
1	A	74	GLN
1	A	383	ASN
1	A	514	GLN
1	B	66	ASN
1	B	68	GLN
1	B	74	GLN
1	B	162	GLN
1	B	257	ASN
1	B	374	GLN
1	B	457	ASN
1	B	514	GLN

5.3.3 RNA ⓘ

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 14 ligands modelled in this entry, 2 are monoatomic - leaving 12 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$
4	PGE	B	1540	-	9,9,9	2.25	5 (55%)	8,8,8	2.01	4 (50%)
4	PGE	A	1543	-	9,9,9	1.42	2 (22%)	8,8,8	1.14	1 (12%)
2	NAG	A	1544	1	14,14,15	0.64	0	17,19,21	0.78	1 (5%)
4	PGE	A	1541	-	9,9,9	1.42	2 (22%)	8,8,8	1.36	1 (12%)
4	PGE	B	1539	-	9,9,9	1.38	2 (22%)	8,8,8	0.98	0
2	NAG	A	1538	1	14,14,15	0.55	0	17,19,21	0.94	1 (5%)
2	NAG	B	1536	1	14,14,15	0.55	0	17,19,21	0.79	1 (5%)
4	PGE	A	1540	-	9,9,9	1.64	2 (22%)	8,8,8	1.80	1 (12%)
2	NAG	B	1537	1	14,14,15	0.72	0	17,19,21	0.59	0
4	PGE	B	1542	-	9,9,9	1.40	1 (11%)	8,8,8	1.93	4 (50%)
4	PGE	A	1542	-	9,9,9	1.68	3 (33%)	8,8,8	0.94	0
4	PGE	B	1541	-	9,9,9	0.89	0	8,8,8	1.24	1 (12%)

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
4	PGE	B	1540	-	-	4/7/7/7	-
4	PGE	A	1543	-	-	3/7/7/7	-
2	NAG	A	1544	1	-	0/6/23/26	0/1/1/1
4	PGE	A	1541	-	-	1/7/7/7	-
4	PGE	B	1539	-	-	2/7/7/7	-
2	NAG	A	1538	1	-	2/6/23/26	0/1/1/1
2	NAG	B	1536	1	-	0/6/23/26	0/1/1/1
4	PGE	A	1540	-	-	3/7/7/7	-
2	NAG	B	1537	1	-	2/6/23/26	0/1/1/1
4	PGE	B	1542	-	-	4/7/7/7	-
4	PGE	A	1542	-	-	3/7/7/7	-
4	PGE	B	1541	-	-	2/7/7/7	-

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1540	PGE	O2-C2	3.70	1.58	1.42
4	B	1540	PGE	C4-C3	3.26	1.65	1.49
4	B	1540	PGE	O3-C5	2.93	1.54	1.42
4	B	1542	PGE	O3-C5	2.85	1.54	1.42
4	B	1540	PGE	O2-C3	2.50	1.52	1.42
4	A	1542	PGE	O2-C2	2.46	1.52	1.42
4	A	1540	PGE	O2-C3	2.41	1.52	1.42
4	A	1541	PGE	O2-C3	2.37	1.52	1.42
4	B	1540	PGE	C5-C6	2.33	1.61	1.49
4	A	1541	PGE	O3-C4	2.32	1.52	1.42
4	A	1542	PGE	O3-C4	2.31	1.52	1.42
4	B	1539	PGE	O3-C4	2.23	1.51	1.42
4	A	1542	PGE	C4-C3	2.15	1.59	1.49
4	B	1539	PGE	C4-C3	2.09	1.59	1.49
4	A	1540	PGE	O2-C2	2.09	1.51	1.42
4	A	1543	PGE	O2-C2	2.08	1.51	1.42
4	A	1543	PGE	O2-C3	2.06	1.51	1.42

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1540	PGE	C5-O3-C4	4.01	130.81	113.26
4	B	1540	PGE	O2-C2-C1	3.61	126.04	110.11
4	B	1540	PGE	O3-C4-C3	2.86	123.39	110.35
4	B	1541	PGE	O3-C4-C3	2.77	122.99	110.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1541	PGE	O2-C3-C4	-2.67	98.20	110.35
2	A	1538	NAG	C2-N2-C7	-2.62	119.39	122.90
4	A	1543	PGE	O3-C5-C6	2.59	121.52	110.11
4	B	1542	PGE	O3-C5-C6	2.54	121.28	110.11
4	B	1542	PGE	O3-C4-C3	2.42	121.37	110.35
2	B	1536	NAG	C2-N2-C7	-2.42	119.66	122.90
4	B	1540	PGE	O3-C5-C6	2.34	120.42	110.11
2	A	1544	NAG	C2-N2-C7	-2.33	119.78	122.90
4	B	1542	PGE	O2-C3-C4	-2.30	99.86	110.35
4	B	1542	PGE	C3-O2-C2	2.26	123.16	113.26
4	B	1540	PGE	C3-O2-C2	-2.23	103.49	113.26

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1538	NAG	C8-C7-N2-C2
2	B	1537	NAG	O5-C5-C6-O6
2	A	1538	NAG	O7-C7-N2-C2
2	B	1537	NAG	C4-C5-C6-O6
4	B	1542	PGE	C6-C5-O3-C4
4	A	1543	PGE	C6-C5-O3-C4
4	B	1539	PGE	C3-C4-O3-C5
4	B	1540	PGE	C6-C5-O3-C4
4	A	1542	PGE	C3-C4-O3-C5
4	A	1542	PGE	C1-C2-O2-C3
4	A	1543	PGE	C1-C2-O2-C3
4	A	1540	PGE	C6-C5-O3-C4
4	A	1540	PGE	C3-C4-O3-C5
4	B	1541	PGE	C3-C4-O3-C5
4	A	1541	PGE	C6-C5-O3-C4
4	B	1540	PGE	C3-C4-O3-C5
4	B	1542	PGE	C3-C4-O3-C5
4	B	1542	PGE	C1-C2-O2-C3
4	B	1540	PGE	C4-C3-O2-C2
4	B	1541	PGE	O2-C3-C4-O3
4	B	1539	PGE	O2-C3-C4-O3
4	A	1542	PGE	O2-C3-C4-O3
4	A	1543	PGE	O2-C3-C4-O3
4	B	1542	PGE	O2-C3-C4-O3
4	B	1540	PGE	O2-C3-C4-O3
4	A	1540	PGE	O2-C3-C4-O3

There are no ring outliers.

7 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1540	PGE	5	0
4	A	1543	PGE	5	0
4	A	1541	PGE	1	0
4	B	1539	PGE	1	0
4	B	1542	PGE	2	0
4	A	1542	PGE	1	0
4	B	1541	PGE	2	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	529/537 (98%)	0.77	26 (4%) 35 31	29, 46, 65, 78	0
1	B	532/537 (99%)	0.64	17 (3%) 50 46	29, 46, 68, 91	0
All	All	1061/1074 (98%)	0.71	43 (4%) 41 37	29, 46, 66, 91	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	485	PRO	4.8
1	B	485	PRO	4.0
1	B	38	PRO	3.8
1	B	535	THR	3.4
1	B	40	VAL	3.4
1	A	23	LEU	3.2
1	B	237	SER	3.1
1	A	108	SER	2.9
1	A	18	THR	2.8
1	A	110	THR	2.8
1	A	245	ALA	2.8
1	B	441	GLY	2.7
1	B	277	VAL	2.6
1	B	63	TYR	2.6
1	B	265	CYS	2.6
1	A	4	SER	2.6
1	B	36	ALA	2.5
1	A	507	THR	2.5
1	A	510	MET	2.5
1	A	27	ILE	2.4
1	A	365	ASP	2.4
1	A	109	THR	2.3
1	A	265	CYS	2.3
1	A	434	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	70	TYR	2.3
1	B	91	SER	2.2
1	A	72	ASP	2.2
1	A	26	HIS	2.2
1	A	454	LYS	2.2
1	A	380	ASP	2.2
1	A	75	PHE	2.2
1	B	6	LEU	2.2
1	A	311	SER	2.2
1	A	432	TRP	2.1
1	B	9	ASN	2.1
1	B	43	MET	2.1
1	A	24	SER	2.1
1	A	521	CYS	2.1
1	B	238	VAL	2.0
1	A	31	LEU	2.0
1	A	452	LEU	2.0
1	B	488	GLN	2.0
1	B	487	SER	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
4	PGE	B	1540	10/10	0.74	0.27	76,79,80,82	0
2	NAG	A	1538	14/15	0.75	0.20	77,82,85,85	0
4	PGE	A	1543	10/10	0.77	0.21	72,76,77,77	0
2	NAG	A	1544	14/15	0.80	0.15	61,67,70,71	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	PGE	A	1540	10/10	0.80	0.25	81,81,82,83	0
4	PGE	B	1539	10/10	0.81	0.18	72,77,81,82	0
4	PGE	A	1541	10/10	0.82	0.20	72,79,80,81	0
4	PGE	B	1542	10/10	0.82	0.18	66,72,73,75	0
4	PGE	B	1541	10/10	0.85	0.19	69,74,78,78	0
4	PGE	A	1542	10/10	0.87	0.19	77,81,84,85	0
2	NAG	B	1536	14/15	0.87	0.14	72,75,76,76	0
2	NAG	B	1537	14/15	0.88	0.11	50,56,60,62	0
3	CL	B	1538	1/1	0.88	0.14	78,78,78,78	0
3	CL	A	1539	1/1	0.92	0.16	75,75,75,75	0

6.5 Other polymers [i](#)

There are no such residues in this entry.