



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 10:43 PM UTC

PDB ID : 7ABE / pdb_00007abe
Title : Crystal structure of human phosphodiesterase 4D2 catalytic domain with inhibitor NPD-769
Authors : Singh, A.K.; Blaazer, A.R.; Zara, L.; de Esch, I.J.P.; Leurs, R.; Brown, D.G.
Deposited on : 2020-09-07
Resolution : 1.83 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

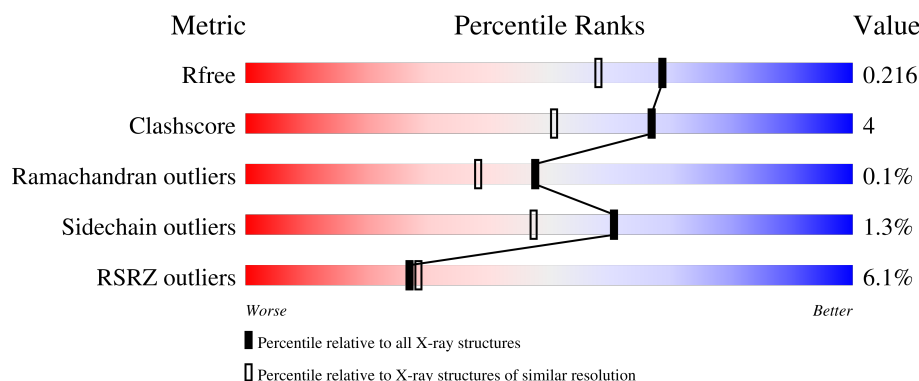
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.83 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	1296 (1.84-1.84)
Clashscore	190562	1329 (1.84-1.84)
Ramachandran outliers	187476	1318 (1.84-1.84)
Sidechain outliers	187428	1318 (1.84-1.84)
RSRZ outliers	180081	1296 (1.84-1.84)

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	364	9% 83% 8% 9%
1	B	364	4% 80% 7% 11%
1	C	364	7% 84% 5% 10%
1	D	364	2% 83% 5% 11%

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	EDO	B	509	-	-	X	-
6	EPE	A	522	-	-	X	-
7	DMS	B	511	-	-	X	-
8	PEG	D	511	-	-	X	-

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 11744 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 cAMP-specific 3',5'-cyclic phosphodiesterase 4D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total	C	N	O	S	0	1	0
			2674	1694	455	510	15			
1	B	324	Total	C	N	O	S	0	0	0
			2622	1659	448	501	14			
1	C	326	Total	C	N	O	S	0	0	0
			2638	1668	450	506	14			
1	D	324	Total	C	N	O	S	0	1	0
			2627	1663	448	501	15			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	75	GLY	-	expression tag	UNP Q08499
A	76	SER	-	expression tag	UNP Q08499
A	77	HIS	-	expression tag	UNP Q08499
A	78	MET	-	expression tag	UNP Q08499
B	75	GLY	-	expression tag	UNP Q08499
B	76	SER	-	expression tag	UNP Q08499
B	77	HIS	-	expression tag	UNP Q08499
B	78	MET	-	expression tag	UNP Q08499
C	75	GLY	-	expression tag	UNP Q08499
C	76	SER	-	expression tag	UNP Q08499
C	77	HIS	-	expression tag	UNP Q08499
C	78	MET	-	expression tag	UNP Q08499
D	75	GLY	-	expression tag	UNP Q08499
D	76	SER	-	expression tag	UNP Q08499
D	77	HIS	-	expression tag	UNP Q08499
D	78	MET	-	expression tag	UNP Q08499

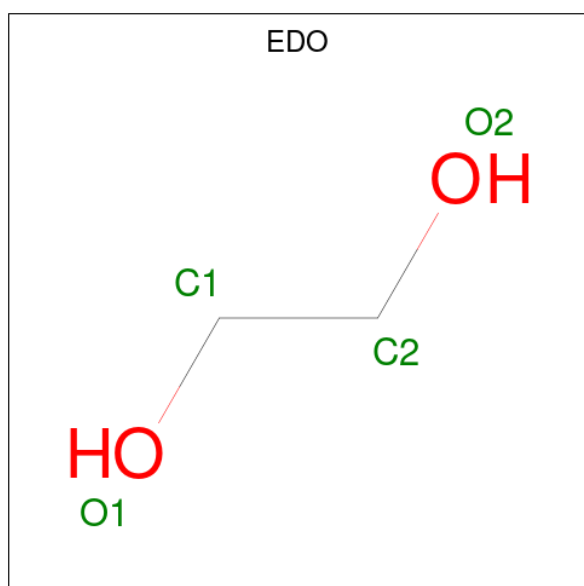
- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		

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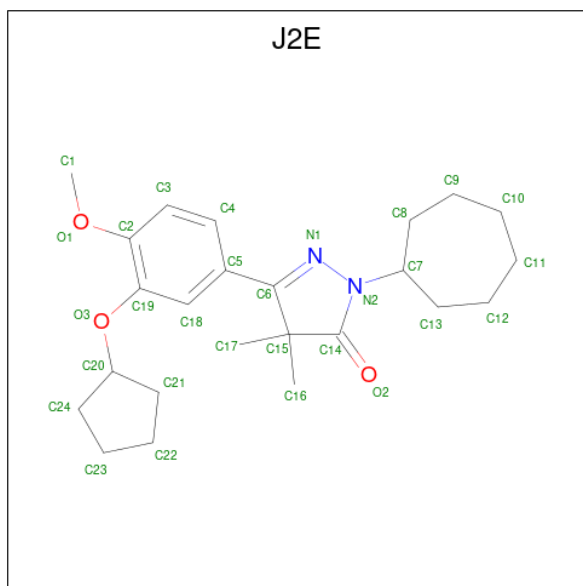
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		

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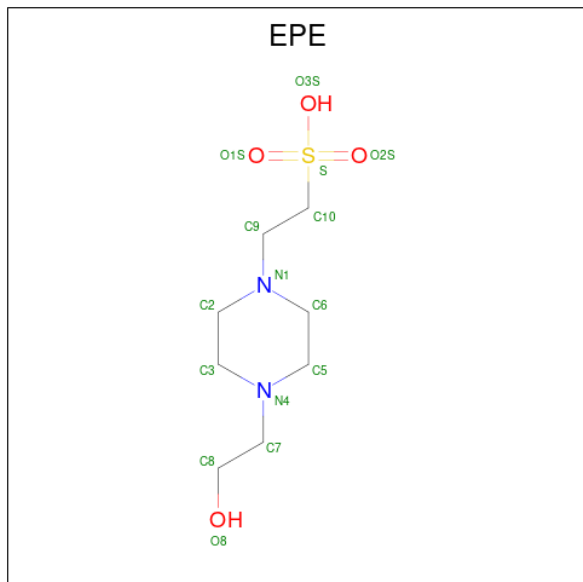
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is 1-cycloheptyl-3-[3-(cyclopentyloxy)-4-methoxyphenyl]-4,4-dimethyl-4,5-dihydro-1H-pyrazol-5-one (CCD ID: J2E) (formula: C₂₄H₃₄N₂O₃) (labeled as "Ligand of Interest" by depositor).



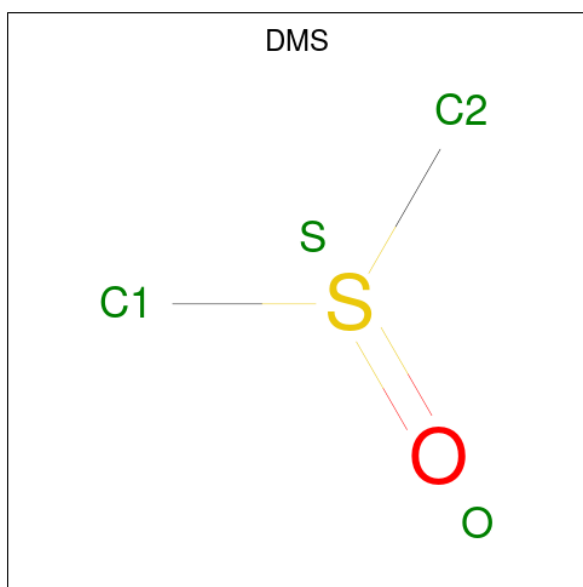
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			29	24	2	3		
5	B	1	Total	C	N	O	0	0
			29	24	2	3		
5	C	1	Total	C	N	O	0	0
			29	24	2	3		
5	D	1	Total	C	N	O	0	0
			29	24	2	3		

- Molecule 6 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).



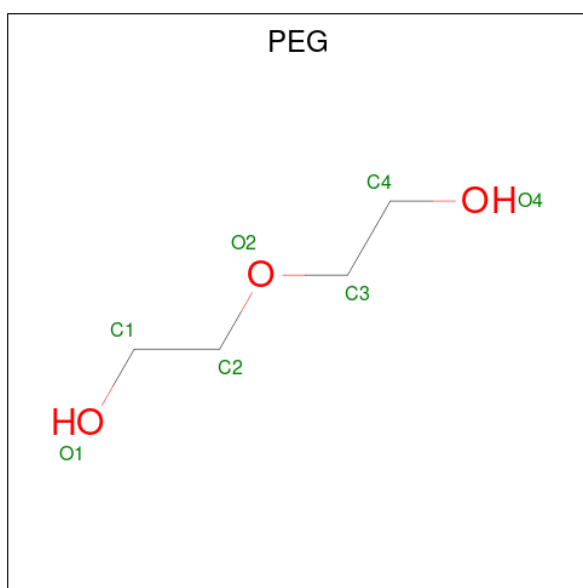
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
6	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
6	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
6	C	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
6	D	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 7 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C_2H_6OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	B	1	Total	C	O	S	0	0
			4	2	1	1		
7	B	1	Total	C	O	S	0	0
			4	2	1	1		

- Molecule 8 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	D	1	Total	C	O	0	0
			7	4	3		
8	D	1	Total	C	O	0	0
			7	4	3		

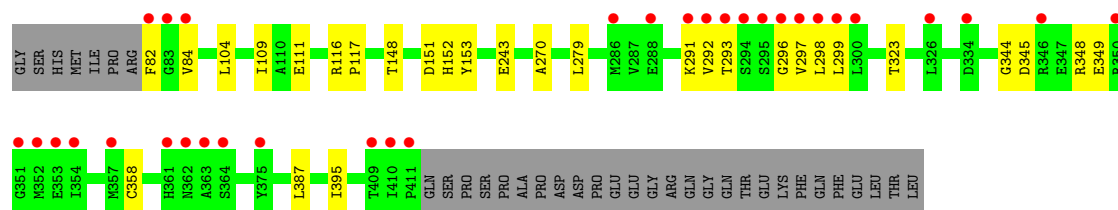
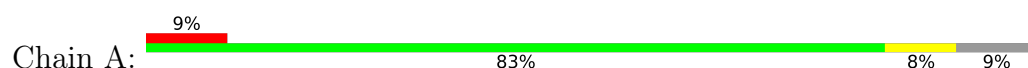
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	182	Total 182	O 182	0	0
9	B	203	Total 203	O 203	0	0
9	C	147	Total 147	O 147	0	0
9	D	230	Total 230	O 230	0	0

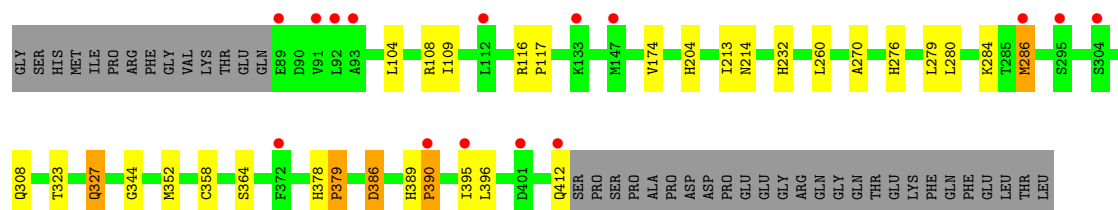
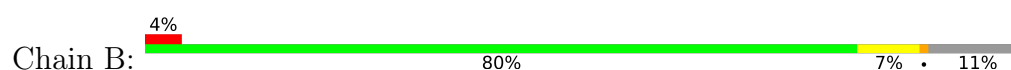
3 Residue-property plots [i](#)

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.

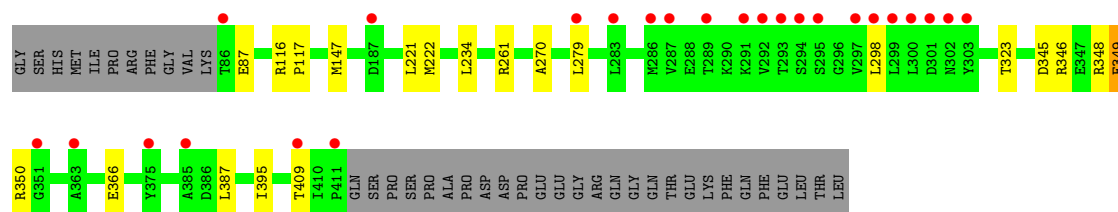
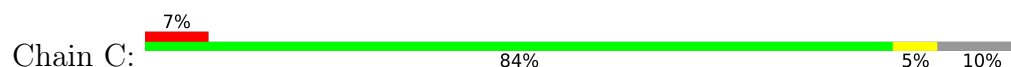
- Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4D



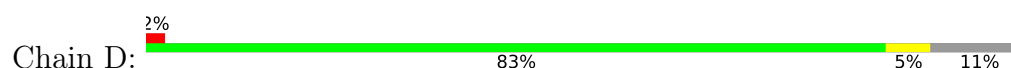
- Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4D

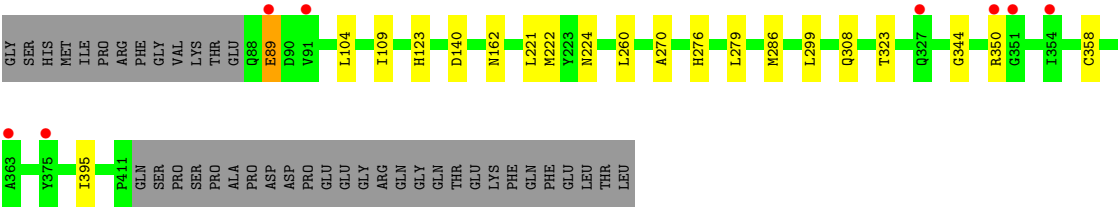


- Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4D



- Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4D





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	98.70Å 111.25Å 160.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	84.16 – 1.83 84.16 – 1.83	Depositor EDS
% Data completeness (in resolution range)	99.9 (84.16-1.83) 99.9 (84.16-1.83)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 1.83Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.179 , 0.208 0.188 , 0.216	Depositor DCC
R_{free} test set	7830 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	31.7	Xtriage
Anisotropy	0.170	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 38.8	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	0.96	EDS
Total number of atoms	11744	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.20% 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: J2E, PEG, DMS, EPE, ZN, EDO, 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	1.05	0/2732	1.26	0/3710
1	B	1.16	5/2676 (0.2%)	1.32	6/3636 (0.2%)
1	C	1.05	0/2692	1.24	1/3658 (0.0%)
1	D	1.06	1/2684 (0.0%)	1.20	1/3646 (0.0%)
All	All	1.08	6/10784 (0.1%)	1.26	8/14650 (0.1%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	379	PRO	N-CA	16.01	1.69	1.47
1	B	378	HIS	C-N	7.07	1.43	1.33
1	D	276	HIS	CE1-NE2	6.25	1.38	1.32
1	B	204	HIS	CE1-NE2	6.22	1.38	1.32
1	B	232	HIS	CE1-NE2	5.58	1.38	1.32
1	B	276	HIS	CE1-NE2	5.08	1.37	1.32

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	378	HIS	CA-C-N	14.63	134.45	119.24
1	B	378	HIS	C-N-CA	14.63	134.45	119.24
1	B	378	HIS	O-C-N	-8.70	113.51	120.38
1	B	379	PRO	N-CA-C	-8.38	100.21	113.78
1	B	379	PRO	CA-N-CD	-6.06	103.52	112.00
1	C	349	GLU	CB-CA-C	-5.70	101.27	110.74
1	B	390	PRO	N-CA-CB	-5.24	96.83	102.60
1	D	89	GLU	CB-CG-CD	5.18	121.41	112.60

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	2674	0	2634	23	0
1	B	2622	0	2578	23	0
1	C	2638	0	2591	16	0
1	D	2627	0	2587	19	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	76	0	114	6	0
4	B	32	0	48	8	0
4	C	40	0	60	3	0
4	D	52	0	78	1	0
5	A	29	0	0	0	0
5	B	29	0	0	0	0
5	C	29	0	0	1	0
5	D	29	0	0	0	0
6	A	15	0	17	7	0
6	B	30	0	36	1	0
6	C	15	0	18	0	0
6	D	15	0	18	0	0
7	B	8	0	12	5	0
8	D	14	0	20	6	0
9	A	182	0	0	0	0
9	B	203	0	0	0	0
9	C	147	0	0	0	0
9	D	230	0	0	0	0
All	All	11744	0	10811	81	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 4.

All (81) 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:B:379:PRO:N	1:B:379:PRO:CA	1.69	1.39
1:A:111:GLU:OE2	6:A:522:EPE:H52	1.44	1.17
1:A:151:ASP:O	4:A:520:EDO:H21	1.62	0.98
1:A:153:TYR:O	4:A:520:EDO:H11	1.62	0.98
1:B:352:MET:HE2	4:B:509:EDO:C2	1.97	0.94
1:A:111:GLU:OE2	6:A:522:EPE:C5	2.16	0.94
1:D:350:ARG:HH11	1:D:350:ARG:HG3	1.37	0.87
1:A:111:GLU:OE2	6:A:522:EPE:H72	1.75	0.86
1:A:111:GLU:CD	6:A:522:EPE:H52	2.03	0.84
1:D:123:HIS:HD1	8:D:518:PEG:H22	1.43	0.84
1:B:352:MET:HE2	4:B:509:EDO:H22	1.64	0.77
1:B:379:PRO:N	1:B:379:PRO:C	2.48	0.71
1:B:286:MET:HE3	1:B:308:GLN:NE2	2.08	0.69
1:C:234:LEU:HB3	8:D:511:PEG:H22	1.75	0.67
1:A:292:VAL:HG23	1:A:292:VAL:O	1.98	0.64
1:B:352:MET:HE2	4:B:509:EDO:C1	2.29	0.62
1:C:346:ARG:O	1:C:350:ARG:HD2	1.99	0.62
1:A:298:LEU:HD11	1:A:387:LEU:HG	1.83	0.60
1:D:350:ARG:HG3	1:D:350:ARG:NH1	2.12	0.59
1:D:123:HIS:ND1	8:D:518:PEG:H22	2.16	0.59
1:C:366:GLU:HG2	1:C:409:THR:HG23	1.82	0.59
1:A:148:THR:HG23	4:A:513:EDO:H12	1.85	0.59
1:C:298:LEU:HD11	1:C:387:LEU:HG	1.84	0.59
1:A:111:GLU:OE2	6:A:522:EPE:C7	2.51	0.58
1:D:299:LEU:C	1:D:299:LEU:HD13	2.28	0.58
1:B:213:ILE:CD1	7:B:511:DMS:H11	2.34	0.57
6:A:522:EPE:O8	6:A:522:EPE:H51	2.04	0.57
1:B:327:GLN:NE2	6:B:515:EPE:O1S	2.38	0.57
1:B:270:ALA:HB1	1:B:279:LEU:HD11	1.88	0.55
1:B:352:MET:HE2	4:B:509:EDO:H21	1.86	0.55
1:A:270:ALA:HB1	1:A:279:LEU:HD11	1.89	0.55
4:C:510:EDO:H21	5:C:513:J2E:O2	2.07	0.54
7:B:510:DMS:H21	1:D:140:ASP:HA	1.89	0.54
1:A:293:THR:OG1	1:A:299:LEU:HD22	2.07	0.53
1:D:350:ARG:HH11	1:D:350:ARG:CG	2.11	0.52
1:D:270:ALA:HB1	1:D:279:LEU:HD11	1.91	0.52
1:C:222:MET:HE3	1:D:222[B]:MET:HE1	1.92	0.52
1:B:213:ILE:CD1	7:B:511:DMS:C1	2.89	0.51
1:B:214:ASN:HD22	4:B:509:EDO:H21	1.76	0.51
1:B:213:ILE:HD12	7:B:511:DMS:C1	2.42	0.50
1:C:221:LEU:HD23	1:C:221:LEU:O	2.13	0.49
1:D:286:MET:HE3	1:D:308:GLN:OE1	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:104:LEU:HD11	1:D:109:ILE:HD11	1.95	0.48
1:A:243:GLU:OE1	4:A:513:EDO:H12	2.13	0.48
1:A:152:HIS:HE1	4:A:513:EDO:H11	1.79	0.48
1:C:270:ALA:HB1	1:C:279:LEU:HD11	1.96	0.48
1:B:323:THR:HB	1:B:395:ILE:HG23	1.96	0.47
1:D:350:ARG:NH1	1:D:350:ARG:CG	2.75	0.47
1:B:213:ILE:HD12	7:B:511:DMS:H13	1.96	0.47
1:B:174:VAL:CG1	4:B:513:EDO:H11	2.46	0.46
1:B:280:LEU:HG	1:B:284:LYS:HE2	1.98	0.45
1:C:261:ARG:HB2	4:C:504:EDO:C1	2.47	0.44
1:D:104:LEU:HD11	1:D:109:ILE:CD1	2.47	0.44
1:A:349:GLU:HG3	1:C:147:MET:HE1	2.00	0.44
1:B:386:ASP:O	1:B:389:HIS:HD2	2.01	0.44
1:B:214:ASN:HD22	4:B:509:EDO:C2	2.31	0.44
1:C:234:LEU:CB	8:D:511:PEG:H22	2.43	0.43
1:A:323:THR:HB	1:A:395:ILE:HG23	2.01	0.42
1:A:104:LEU:HD11	1:A:109:ILE:HD11	2.00	0.42
1:B:104:LEU:HD11	1:B:109:ILE:HD11	2.01	0.42
1:D:323:THR:HB	1:D:395:ILE:HG23	2.02	0.42
1:C:261:ARG:HB2	4:C:504:EDO:H12	2.01	0.42
1:D:221:LEU:O	8:D:511:PEG:H41	2.20	0.42
1:B:344:GLY:HA3	1:B:358:CYS:O	2.20	0.41
1:D:224:ASN:OD1	8:D:511:PEG:H21	2.19	0.41
1:C:323:THR:HB	1:C:395:ILE:HG23	2.01	0.41
1:D:162:ASN:OD1	4:D:508:EDO:H11	2.21	0.41
1:C:345:ASP:O	1:C:349:GLU:HG3	2.20	0.41
1:C:116:ARG:N	1:C:117:PRO:CD	2.84	0.41
1:A:345:ASP:OD1	1:A:348:ARG:NH2	2.54	0.41
1:A:344:GLY:HA3	1:A:358:CYS:O	2.21	0.40
1:C:221:LEU:HD23	1:C:221:LEU:C	2.46	0.40
1:C:345:ASP:OD1	1:C:348:ARG:NH2	2.54	0.40
1:A:104:LEU:HD11	1:A:109:ILE:CD1	2.51	0.40
1:A:111:GLU:CD	6:A:522:EPE:C5	2.80	0.40
1:B:214:ASN:HD22	4:B:509:EDO:C1	2.35	0.40
1:D:344:GLY:HA3	1:D:358:CYS:O	2.21	0.40
1:B:116:ARG:N	1:B:117:PRO:CD	2.85	0.40
1:D:286:MET:CE	1:D:308:GLN:OE1	2.70	0.40
1:A:116:ARG:N	1:A:117:PRO:CD	2.85	0.40
1:A:153:TYR:O	4:A:520:EDO:C1	2.51	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	329/364 (90%)	320 (97%)	8 (2%)	1 (0%)	36	25
1	B	322/364 (88%)	317 (98%)	5 (2%)	0	100	100
1	C	324/364 (89%)	318 (98%)	6 (2%)	0	100	100
1	D	323/364 (89%)	317 (98%)	6 (2%)	0	100	100
All	All	1298/1456 (89%)	1272 (98%)	25 (2%)	1 (0%)	48	38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	296	GLY

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	302/331 (91%)	298 (99%)	4 (1%)	61	47
1	B	296/331 (89%)	287 (97%)	9 (3%)	36	19
1	C	298/331 (90%)	297 (100%)	1 (0%)	86	81
1	D	297/331 (90%)	295 (99%)	2 (1%)	76	68
All	All	1193/1324 (90%)	1177 (99%)	16 (1%)	61	47

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	82	PHE
1	A	84	VAL
1	A	291	LYS
1	A	297	VAL
1	B	108	ARG
1	B	260	LEU
1	B	286	MET
1	B	327	GLN
1	B	364	SER
1	B	386	ASP
1	B	390	PRO
1	B	396	LEU
1	B	412	GLN
1	C	87	GLU
1	D	89	GLU
1	D	260	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	210	GLN
1	A	258	GLN
1	B	210	GLN
1	B	214	ASN
1	B	258	GLN
1	B	389	HIS
1	C	210	GLN
1	C	250	GLN
1	C	321	ASN
1	C	331	GLN
1	D	88	GLN
1	D	127	GLN
1	D	210	GLN
1	D	258	GLN
1	D	308	GLN
1	D	312	ASN
1	D	321	ASN
1	D	407	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 71 ligands modelled in this entry, 8 are monoatomic - leaving 63 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	EDO	D	507	-	3,3,3	0.35	0	2,2,2	0.21	0
5	J2E	B	514	-	31,32,32	0.68	1 (3%)	38,46,46	1.53	5 (13%)
8	PEG	D	511	-	6,6,6	0.93	0	5,5,5	0.57	0
4	EDO	A	508	-	3,3,3	0.24	0	2,2,2	0.21	0
4	EDO	A	510	-	3,3,3	0.26	0	2,2,2	0.23	0
4	EDO	A	515	-	3,3,3	0.42	0	2,2,2	0.44	0
5	J2E	C	513	-	31,32,32	0.74	1 (3%)	38,46,46	1.39	7 (18%)
6	EPE	D	519	-	15,15,15	1.75	1 (6%)	19,20,20	2.03	4 (21%)
7	DMS	B	510	-	3,3,3	0.22	0	3,3,3	0.20	0
4	EDO	A	511	-	3,3,3	0.19	0	2,2,2	0.38	0
6	EPE	C	509	-	15,15,15	2.12	2 (13%)	19,20,20	1.85	4 (21%)
4	EDO	C	506	-	3,3,3	0.15	0	2,2,2	0.37	0
4	EDO	C	504	-	3,3,3	0.14	0	2,2,2	0.04	0
4	EDO	B	504	-	3,3,3	0.04	0	2,2,2	0.13	0
4	EDO	A	503	-	3,3,3	0.09	0	2,2,2	0.09	0
4	EDO	D	501	-	3,3,3	0.37	0	2,2,2	0.53	0
4	EDO	D	506	-	3,3,3	0.37	0	2,2,2	0.42	0
4	EDO	A	507	-	3,3,3	0.10	0	2,2,2	0.04	0
4	EDO	A	518	-	3,3,3	0.28	0	2,2,2	0.33	0
6	EPE	B	515	-	15,15,15	1.76	1 (6%)	19,20,20	2.07	7 (36%)
4	EDO	B	505	-	3,3,3	0.14	0	2,2,2	0.06	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	EDO	A	517	-	3,3,3	0.18	0	2,2,2	0.15	0
4	EDO	B	508	-	3,3,3	0.03	0	2,2,2	0.11	0
4	EDO	A	514	-	3,3,3	0.09	0	2,2,2	0.11	0
4	EDO	B	507	-	3,3,3	0.20	0	2,2,2	0.20	0
4	EDO	C	507	-	3,3,3	0.04	0	2,2,2	0.05	0
4	EDO	D	516	-	3,3,3	0.59	0	2,2,2	0.57	0
4	EDO	C	505	-	3,3,3	0.09	0	2,2,2	0.25	0
4	EDO	D	513	-	3,3,3	0.56	0	2,2,2	0.51	0
4	EDO	D	509	-	3,3,3	0.16	0	2,2,2	0.67	0
4	EDO	D	505	-	3,3,3	0.03	0	2,2,2	0.20	0
4	EDO	D	508	-	3,3,3	0.36	0	2,2,2	0.47	0
4	EDO	D	515	-	3,3,3	0.28	0	2,2,2	0.44	0
4	EDO	C	503	-	3,3,3	0.16	0	2,2,2	0.25	0
4	EDO	A	504	-	3,3,3	0.26	0	2,2,2	0.25	0
4	EDO	B	503	-	3,3,3	0.32	0	2,2,2	0.53	0
4	EDO	C	514	-	3,3,3	0.12	0	2,2,2	0.12	0
4	EDO	D	512	-	3,3,3	1.07	0	2,2,2	0.60	0
4	EDO	B	506	-	3,3,3	0.11	0	2,2,2	0.20	0
6	EPE	A	522	-	15,15,15	3.19	1 (6%)	19,20,20	1.96	5 (26%)
4	EDO	A	506	-	3,3,3	0.07	0	2,2,2	0.16	0
8	PEG	D	518	-	6,6,6	0.66	0	5,5,5	0.23	0
4	EDO	A	513	-	3,3,3	0.71	0	2,2,2	0.15	0
5	J2E	A	519	-	31,32,32	0.61	1 (3%)	38,46,46	1.49	6 (15%)
4	EDO	A	509	-	3,3,3	0.08	0	2,2,2	0.58	0
5	J2E	D	517	-	31,32,32	0.81	2 (6%)	38,46,46	1.58	5 (13%)
4	EDO	C	512	-	3,3,3	0.35	0	2,2,2	0.23	0
4	EDO	A	523	-	3,3,3	0.15	0	2,2,2	0.24	0
4	EDO	A	520	-	3,3,3	0.29	0	2,2,2	0.61	0
6	EPE	B	512	-	15,15,15	1.81	1 (6%)	19,20,20	1.80	2 (10%)
4	EDO	C	510	-	3,3,3	0.19	0	2,2,2	0.27	0
4	EDO	D	504	-	3,3,3	0.38	0	2,2,2	0.70	0
4	EDO	A	505	-	3,3,3	0.26	0	2,2,2	0.41	0
4	EDO	A	516	-	3,3,3	0.44	0	2,2,2	0.26	0
4	EDO	B	509	-	3,3,3	0.55	0	2,2,2	0.82	0
7	DMS	B	511	-	3,3,3	0.33	0	3,3,3	0.23	0
4	EDO	D	510	-	3,3,3	0.61	0	2,2,2	0.63	0
4	EDO	A	512	-	3,3,3	0.52	0	2,2,2	0.69	0
4	EDO	D	514	-	3,3,3	0.23	0	2,2,2	0.16	0
4	EDO	C	508	-	3,3,3	0.11	0	2,2,2	0.29	0
4	EDO	A	521	-	3,3,3	0.11	0	2,2,2	0.21	0
4	EDO	C	511	-	3,3,3	0.22	0	2,2,2	0.27	0
4	EDO	B	513	-	3,3,3	0.11	0	2,2,2	0.25	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
4	EDO	D	507	-	-	1/1/1/1	-
5	J2E	B	514	-	-	3/12/50/50	0/4/4/4
8	PEG	D	511	-	-	3/4/4/4	-
4	EDO	A	508	-	-	1/1/1/1	-
4	EDO	A	510	-	-	0/1/1/1	-
4	EDO	A	515	-	-	0/1/1/1	-
5	J2E	C	513	-	-	3/12/50/50	0/4/4/4
6	EPE	D	519	-	-	2/9/19/19	0/1/1/1
4	EDO	A	511	-	-	1/1/1/1	-
6	EPE	C	509	-	-	2/9/19/19	0/1/1/1
4	EDO	C	506	-	-	0/1/1/1	-
4	EDO	C	504	-	-	1/1/1/1	-
4	EDO	B	504	-	-	1/1/1/1	-
4	EDO	A	503	-	-	1/1/1/1	-
4	EDO	D	501	-	-	0/1/1/1	-
4	EDO	D	506	-	-	0/1/1/1	-
4	EDO	A	507	-	-	0/1/1/1	-
4	EDO	A	518	-	-	1/1/1/1	-
6	EPE	B	515	-	-	1/9/19/19	0/1/1/1
4	EDO	B	505	-	-	0/1/1/1	-
4	EDO	A	517	-	-	1/1/1/1	-
4	EDO	B	508	-	-	1/1/1/1	-
4	EDO	A	514	-	-	1/1/1/1	-
4	EDO	B	507	-	-	0/1/1/1	-
4	EDO	C	507	-	-	1/1/1/1	-
4	EDO	D	516	-	-	1/1/1/1	-
4	EDO	C	505	-	-	0/1/1/1	-
4	EDO	D	513	-	-	0/1/1/1	-
4	EDO	D	509	-	-	1/1/1/1	-
4	EDO	D	505	-	-	0/1/1/1	-
4	EDO	D	508	-	-	1/1/1/1	-
4	EDO	D	515	-	-	1/1/1/1	-
4	EDO	C	503	-	-	0/1/1/1	-
4	EDO	A	504	-	-	1/1/1/1	-
4	EDO	B	503	-	-	0/1/1/1	-
4	EDO	C	514	-	-	1/1/1/1	-
4	EDO	D	512	-	-	0/1/1/1	-
4	EDO	B	506	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	EPE	A	522	-	-	3/9/19/19	0/1/1/1
4	EDO	A	506	-	-	1/1/1/1	-
8	PEG	D	518	-	-	1/4/4/4	-
4	EDO	A	513	-	-	0/1/1/1	-
5	J2E	A	519	-	-	8/12/50/50	0/4/4/4
4	EDO	A	509	-	-	1/1/1/1	-
5	J2E	D	517	-	-	5/12/50/50	0/4/4/4
4	EDO	C	512	-	-	1/1/1/1	-
4	EDO	A	523	-	-	1/1/1/1	-
4	EDO	A	520	-	-	0/1/1/1	-
6	EPE	B	512	-	-	5/9/19/19	0/1/1/1
4	EDO	C	510	-	-	0/1/1/1	-
4	EDO	D	504	-	-	0/1/1/1	-
4	EDO	A	505	-	-	1/1/1/1	-
4	EDO	A	516	-	-	1/1/1/1	-
4	EDO	B	509	-	-	1/1/1/1	-
4	EDO	D	510	-	-	0/1/1/1	-
4	EDO	A	512	-	-	1/1/1/1	-
4	EDO	D	514	-	-	1/1/1/1	-
4	EDO	C	508	-	-	1/1/1/1	-
4	EDO	A	521	-	-	1/1/1/1	-
4	EDO	C	511	-	-	1/1/1/1	-
4	EDO	B	513	-	-	0/1/1/1	-

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	522	EPE	C10-S	-12.01	1.60	1.77
6	C	509	EPE	C10-S	-7.37	1.67	1.77
6	D	519	EPE	C10-S	-6.35	1.68	1.77
6	B	515	EPE	C10-S	-6.32	1.68	1.77
6	B	512	EPE	C10-S	-6.13	1.69	1.77
5	C	513	J2E	C7-N2	3.21	1.51	1.47
6	C	509	EPE	O1S-S	2.79	1.53	1.45
5	B	514	J2E	C7-N2	2.78	1.50	1.47
5	D	517	J2E	C7-N2	2.74	1.50	1.47
5	D	517	J2E	C14-N2	2.72	1.37	1.35
5	A	519	J2E	C7-N2	2.66	1.50	1.47

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	519	EPE	O2S-S-C10	6.22	116.12	106.73
5	A	519	J2E	C14-N2-N1	-5.81	109.77	113.43
6	B	512	EPE	O1S-S-C10	5.39	114.87	106.73
5	D	517	J2E	O2-C14-N2	5.26	131.43	127.01
6	C	509	EPE	O2S-S-C10	5.13	114.48	106.73
6	A	522	EPE	O3S-S-C10	4.81	115.41	106.00
5	D	517	J2E	C14-N2-N1	-4.56	110.55	113.43
5	B	514	J2E	C14-N2-N1	-4.48	110.60	113.43
6	B	515	EPE	C5-N4-C3	4.20	117.89	108.84
5	B	514	J2E	O3-C19-C18	-4.19	114.47	123.87
6	D	519	EPE	O2S-S-O1S	-4.18	100.24	113.82
6	B	515	EPE	O3S-S-C10	3.91	113.65	106.00
6	A	522	EPE	O1S-S-C10	3.80	112.47	106.73
5	C	513	J2E	C14-N2-N1	-3.56	111.18	113.43
6	A	522	EPE	O3S-S-O2S	-3.47	102.72	111.40
5	D	517	J2E	O2-C14-C15	-3.31	123.74	126.93
5	D	517	J2E	C6-N1-N2	-3.29	104.33	108.30
5	B	514	J2E	C19-O3-C20	-3.20	109.86	120.14
6	B	515	EPE	C2-C3-N4	3.19	117.08	110.65
6	B	512	EPE	C6-N1-C2	3.15	115.62	108.84
6	C	509	EPE	O3S-S-O1S	-3.12	103.58	111.40
5	A	519	J2E	C6-N1-N2	-3.09	104.57	108.30
5	B	514	J2E	C4-C5-C6	3.01	126.49	120.67
5	B	514	J2E	O3-C19-C2	2.90	125.30	116.70
5	C	513	J2E	O2-C14-C15	-2.83	124.20	126.93
5	C	513	J2E	O2-C14-N2	2.77	129.34	127.01
6	B	515	EPE	O1S-S-C10	-2.77	102.55	106.73
5	C	513	J2E	O3-C19-C18	-2.58	118.09	123.87
6	B	515	EPE	O2S-S-O1S	-2.57	105.46	113.82
5	C	513	J2E	C19-O3-C20	-2.56	111.91	120.14
6	C	509	EPE	C2-C3-N4	2.55	115.78	110.65
6	C	509	EPE	O3S-S-O2S	-2.45	105.27	111.40
5	C	513	J2E	C4-C5-C6	2.39	125.30	120.67
5	C	513	J2E	O1-C2-C3	2.30	128.18	124.30
5	D	517	J2E	O3-C19-C18	-2.27	118.78	123.87
6	B	515	EPE	O3S-S-O1S	2.24	117.00	111.40
5	A	519	J2E	O1-C2-C3	2.24	128.07	124.30
5	A	519	J2E	C19-O3-C20	-2.23	112.97	120.14
6	B	515	EPE	C5-C6-N1	-2.20	106.20	110.65
5	A	519	J2E	O3-C19-C18	-2.15	119.04	123.87
6	A	522	EPE	C2-C3-N4	-2.09	106.43	110.65
6	A	522	EPE	C5-C6-N1	-2.08	106.45	110.65
5	A	519	J2E	C1-O1-C2	2.07	120.55	117.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	519	EPE	C10-C9-N1	2.02	119.99	112.36
6	D	519	EPE	C3-C2-N1	-2.00	106.61	110.65

There are no chirality outliers.

All (65) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	519	J2E	C8-C7-N2-C14
5	A	519	J2E	C21-C20-O3-C19
5	A	519	J2E	C24-C20-O3-C19
5	B	514	J2E	C21-C20-O3-C19
5	C	513	J2E	C8-C7-N2-C14
5	D	517	J2E	C13-C7-N2-C14
5	D	517	J2E	C21-C20-O3-C19
6	A	522	EPE	C10-C9-N1-C2
6	B	512	EPE	C8-C7-N4-C3
6	D	519	EPE	C9-C10-S-O2S
8	D	511	PEG	O2-C3-C4-O4
8	D	511	PEG	O1-C1-C2-O2
8	D	518	PEG	O1-C1-C2-O2
4	A	509	EDO	O1-C1-C2-O2
4	A	512	EDO	O1-C1-C2-O2
4	C	507	EDO	O1-C1-C2-O2
5	B	514	J2E	C24-C20-O3-C19
5	D	517	J2E	C24-C20-O3-C19
4	A	505	EDO	O1-C1-C2-O2
4	A	508	EDO	O1-C1-C2-O2
4	A	511	EDO	O1-C1-C2-O2
4	A	517	EDO	O1-C1-C2-O2
4	B	509	EDO	O1-C1-C2-O2
6	B	512	EPE	C10-C9-N1-C2
4	D	508	EDO	O1-C1-C2-O2
4	D	509	EDO	O1-C1-C2-O2
6	C	509	EPE	C8-C7-N4-C5
6	B	512	EPE	C9-C10-S-O2S
6	B	515	EPE	N4-C7-C8-O8
5	C	513	J2E	C21-C20-O3-C19
5	C	513	J2E	C24-C20-O3-C19
5	A	519	J2E	C4-C5-C6-C15
6	B	512	EPE	C8-C7-N4-C5
6	C	509	EPE	C8-C7-N4-C3
6	A	522	EPE	S-C10-C9-N1

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Mol	Chain	Res	Type	Atoms
4	A	523	EDO	O1-C1-C2-O2
5	A	519	J2E	C4-C5-C6-N1
4	A	506	EDO	O1-C1-C2-O2
5	A	519	J2E	C18-C5-C6-N1
6	B	512	EPE	C9-C10-S-O3S
6	D	519	EPE	C9-C10-S-O3S
5	A	519	J2E	C18-C5-C6-C15
4	D	515	EDO	O1-C1-C2-O2
5	A	519	J2E	C2-C19-O3-C20
5	B	514	J2E	C2-C19-O3-C20
8	D	511	PEG	C4-C3-O2-C2
6	A	522	EPE	N4-C7-C8-O8
4	A	503	EDO	O1-C1-C2-O2
4	A	516	EDO	O1-C1-C2-O2
4	A	521	EDO	O1-C1-C2-O2
4	B	508	EDO	O1-C1-C2-O2
4	C	511	EDO	O1-C1-C2-O2
4	C	514	EDO	O1-C1-C2-O2
4	D	516	EDO	O1-C1-C2-O2
4	A	518	EDO	O1-C1-C2-O2
4	B	504	EDO	O1-C1-C2-O2
4	C	504	EDO	O1-C1-C2-O2
4	C	508	EDO	O1-C1-C2-O2
4	A	514	EDO	O1-C1-C2-O2
4	C	512	EDO	O1-C1-C2-O2
4	D	507	EDO	O1-C1-C2-O2
5	D	517	J2E	C4-C5-C6-C15
5	D	517	J2E	C2-C19-O3-C20
4	A	504	EDO	O1-C1-C2-O2
4	D	514	EDO	O1-C1-C2-O2

There are no ring outliers.

14 monomers are involved in 37 short contacts:

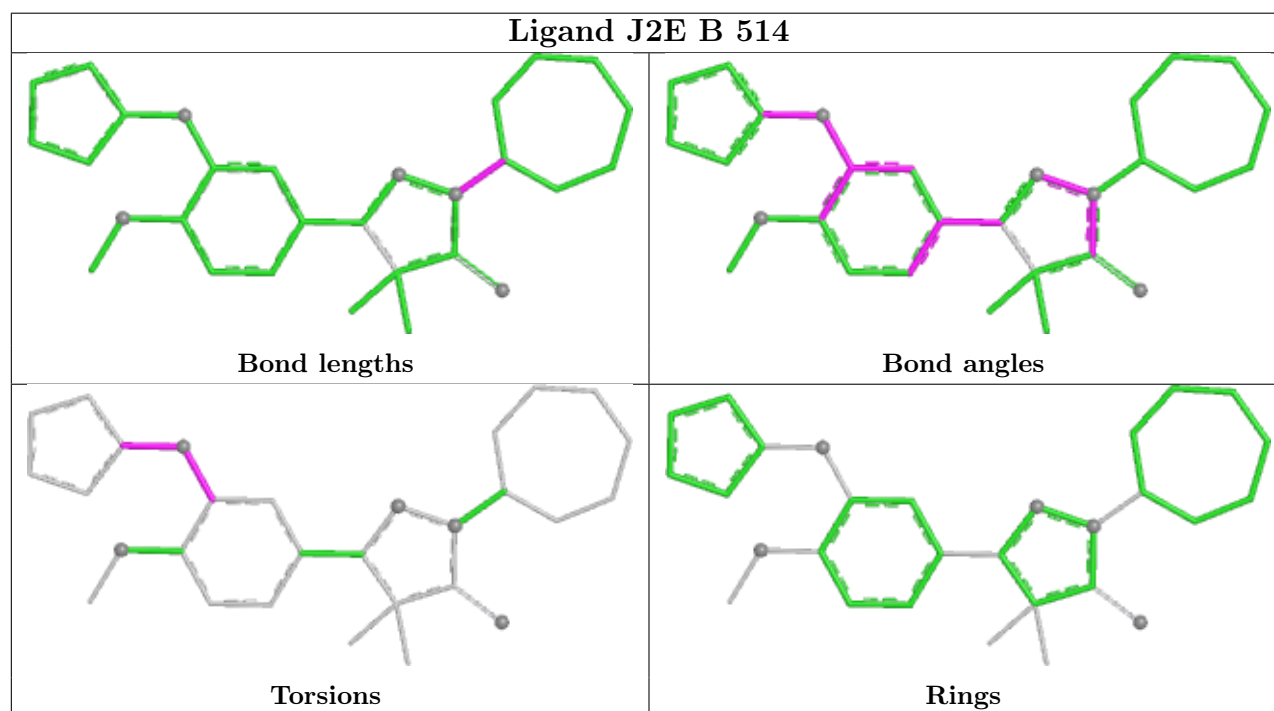
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	D	511	PEG	4	0
5	C	513	J2E	1	0
7	B	510	DMS	1	0
4	C	504	EDO	2	0
6	B	515	EPE	1	0
4	D	508	EDO	1	0
6	A	522	EPE	7	0

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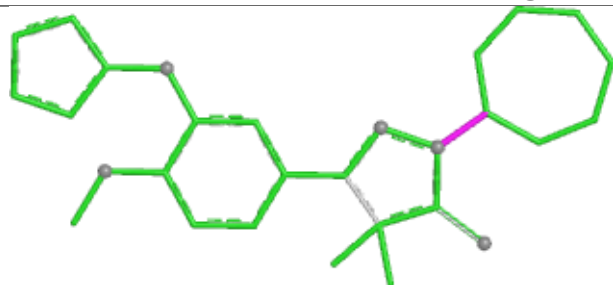
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	D	518	PEG	2	0
4	A	513	EDO	3	0
4	A	520	EDO	3	0
4	C	510	EDO	1	0
4	B	509	EDO	7	0
7	B	511	DMS	4	0
4	B	513	EDO	1	0

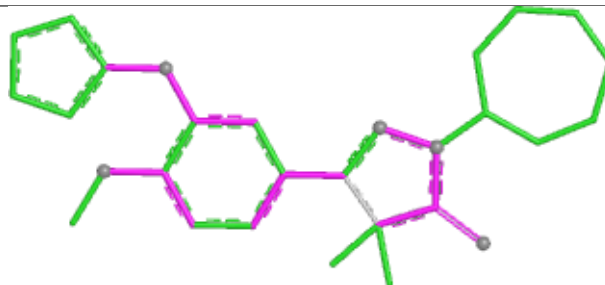
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



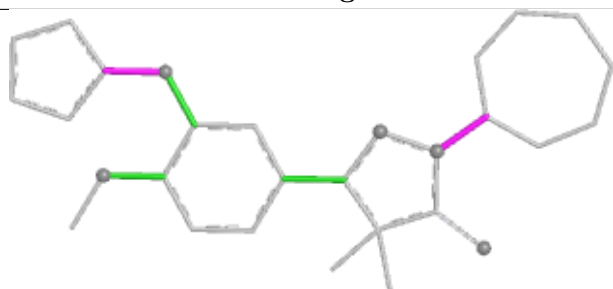
Ligand J2E C 513



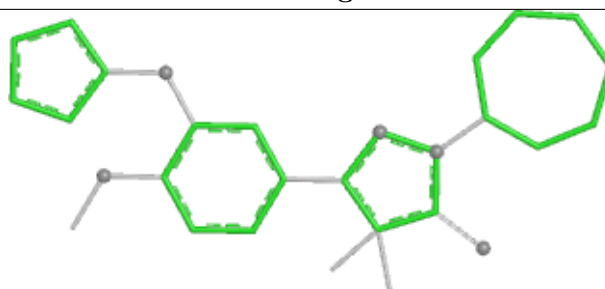
Bond lengths



Bond angles

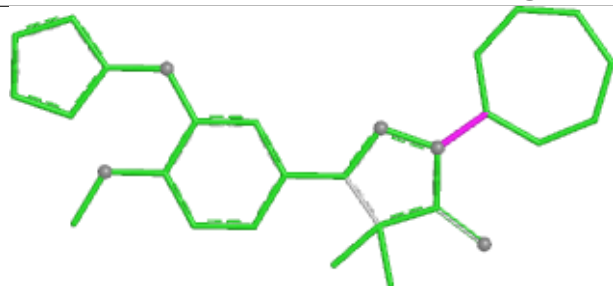


Torsions

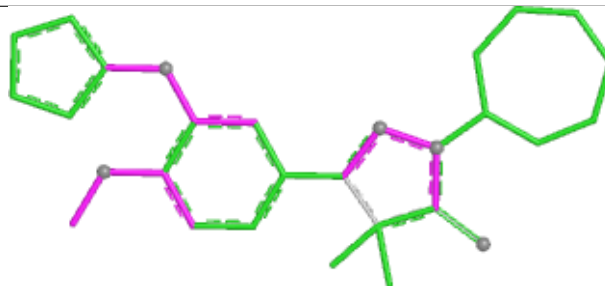


Rings

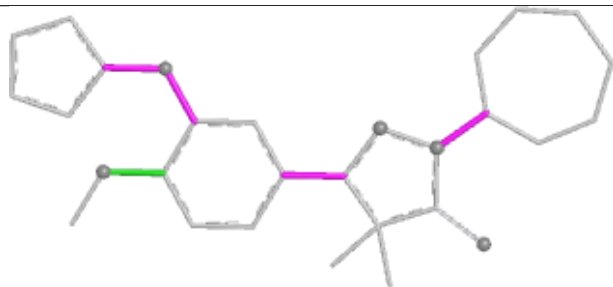
Ligand J2E A 519



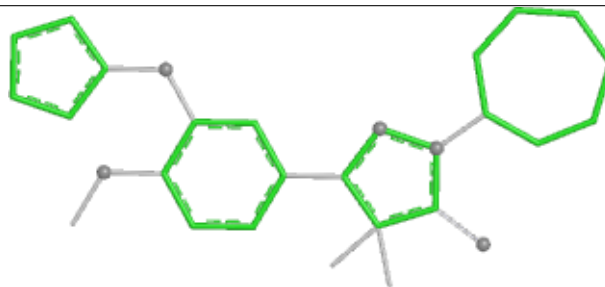
Bond lengths



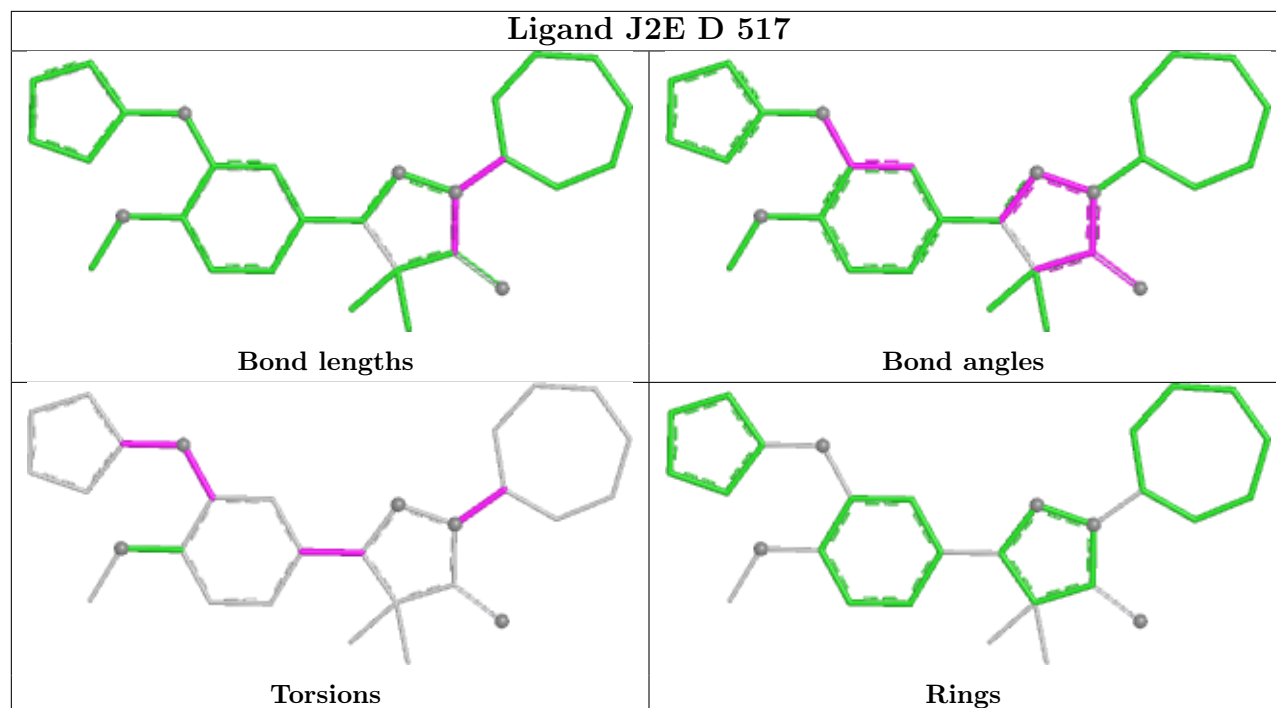
Bond angles



Torsions



Rings



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	330/364 (90%)	0.56	32 (9%) 13 14	21, 36, 71, 123	1 (0%)
1	B	324/364 (89%)	0.63	15 (4%) 37 39	23, 38, 56, 85	0
1	C	326/364 (89%)	0.66	25 (7%) 19 20	23, 37, 74, 118	0
1	D	324/364 (89%)	0.08	8 (2%) 58 64	19, 29, 53, 78	1 (0%)
All	All	1304/1456 (89%)	0.48	80 (6%) 27 28	19, 36, 65, 123	2 (0%)

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	82	PHE	7.2
1	A	293	THR	4.8
1	A	84	VAL	4.2
1	A	83	GLY	4.1
1	A	299	LEU	3.9
1	C	299	LEU	3.8
1	C	303	TYR	3.5
1	C	292	VAL	3.5
1	A	411	PRO	3.4
1	B	91	VAL	3.4
1	A	298	LEU	3.2
1	D	354	ILE	3.1
1	C	298	LEU	3.1
1	C	300	LEU	3.1
1	C	411	PRO	3.1
1	A	297	VAL	3.1
1	D	351	GLY	3.0
1	C	295	SER	3.0
1	C	86	THR	2.9
1	A	351	GLY	2.9
1	A	292	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	301	ASP	2.8
1	C	302	ASN	2.8
1	B	147	MET	2.8
1	B	286	MET	2.8
1	C	294	SER	2.7
1	A	410	ILE	2.7
1	A	363	ALA	2.6
1	C	293	THR	2.6
1	A	350	ARG	2.6
1	C	187	ASP	2.6
1	C	363	ALA	2.6
1	D	363	ALA	2.6
1	A	362	ASN	2.6
1	B	89	GLU	2.6
1	A	357	MET	2.5
1	A	291	LYS	2.5
1	A	295	SER	2.5
1	A	364	SER	2.5
1	A	409	THR	2.5
1	A	334	ASP	2.5
1	D	350	ARG	2.4
1	C	289	THR	2.4
1	D	91	VAL	2.4
1	A	346	ARG	2.4
1	A	375	TYR	2.4
1	B	92	LEU	2.4
1	C	297	VAL	2.4
1	C	351	GLY	2.4
1	B	93	ALA	2.4
1	C	385	ALA	2.4
1	A	296	GLY	2.4
1	C	286	MET	2.4
1	A	352	MET	2.3
1	B	295	SER	2.3
1	C	279	LEU	2.3
1	C	283	LEU	2.3
1	D	327	GLN	2.3
1	A	300	LEU	2.3
1	C	375	TYR	2.3
1	A	294	SER	2.3
1	B	401	ASP	2.3
1	B	395	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	412	GLN	2.2
1	D	375	TYR	2.1
1	C	287	VAL	2.1
1	C	409	THR	2.1
1	B	304	SER	2.1
1	A	353	GLU	2.1
1	A	361	HIS	2.1
1	B	372	PHE	2.1
1	A	326	LEU	2.1
1	B	112	LEU	2.1
1	C	291	LYS	2.1
1	D	89	GLU	2.1
1	B	390	PRO	2.1
1	A	354	ILE	2.0
1	B	133	LYS	2.0
1	A	288	GLU	2.0
1	A	286	MET	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(Å ²)	Q<0.9
4	EDO	A	520	4/4	0.77	0.17	45,48,49,55	0
4	EDO	D	513	4/4	0.77	0.18	52,53,55,60	0
7	DMS	B	511	4/4	0.77	0.30	75,86,90,91	0
5	J2E	A	519	29/29	0.78	0.33	61,79,102,107	9
4	EDO	D	508	4/4	0.79	0.26	44,51,58,65	0
4	EDO	D	516	4/4	0.79	0.24	54,56,60,61	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	EDO	A	512	4/4	0.80	0.20	41,53,54,57	0
4	EDO	A	517	4/4	0.80	0.17	57,59,60,74	0
4	EDO	C	507	4/4	0.81	0.21	61,68,72,73	0
4	EDO	A	504	4/4	0.81	0.18	47,54,59,60	0
4	EDO	A	518	4/4	0.82	0.17	41,52,53,56	0
4	EDO	D	510	4/4	0.82	0.18	36,43,48,50	0
4	EDO	B	505	4/4	0.83	0.19	53,59,60,68	0
4	EDO	B	503	4/4	0.83	0.21	52,53,55,56	0
5	J2E	C	513	29/29	0.83	0.23	41,66,110,115	0
4	EDO	C	512	4/4	0.83	0.20	54,59,64,67	0
8	PEG	D	511	7/7	0.83	0.17	37,41,59,61	0
8	PEG	D	518	7/7	0.83	0.18	51,53,60,63	0
4	EDO	B	508	4/4	0.84	0.17	53,58,58,63	0
4	EDO	A	513	4/4	0.84	0.21	38,49,54,56	0
4	EDO	B	506	4/4	0.84	0.17	49,55,55,57	0
5	J2E	D	517	29/29	0.85	0.23	39,60,103,104	0
4	EDO	A	523	4/4	0.85	0.16	41,57,58,62	0
4	EDO	A	516	4/4	0.86	0.17	39,48,52,53	0
4	EDO	A	505	4/4	0.86	0.16	38,52,55,56	0
7	DMS	B	510	4/4	0.86	0.23	69,74,77,79	0
4	EDO	A	506	4/4	0.86	0.16	42,55,63,76	0
4	EDO	A	514	4/4	0.86	0.15	59,71,73,74	0
5	J2E	B	514	29/29	0.86	0.22	41,66,113,117	0
4	EDO	A	521	4/4	0.87	0.18	35,56,58,73	0
4	EDO	C	508	4/4	0.87	0.15	55,60,60,66	0
4	EDO	D	515	4/4	0.87	0.16	33,42,47,48	0
4	EDO	C	510	4/4	0.87	0.14	50,52,62,65	0
4	EDO	B	513	4/4	0.87	0.17	57,64,69,71	0
4	EDO	C	503	4/4	0.87	0.17	59,61,63,66	0
6	EPE	A	522	15/15	0.88	0.18	45,87,101,101	0
4	EDO	B	509	4/4	0.88	0.15	48,52,53,58	0
4	EDO	A	508	4/4	0.88	0.16	57,57,65,67	0
4	EDO	A	511	4/4	0.88	0.19	49,52,53,56	0
4	EDO	B	504	4/4	0.88	0.14	50,54,54,55	0
4	EDO	C	514	4/4	0.89	0.14	66,68,71,74	0
4	EDO	C	505	4/4	0.89	0.13	47,48,51,52	0
4	EDO	C	506	4/4	0.89	0.19	54,60,62,67	0
4	EDO	C	511	4/4	0.89	0.18	56,64,73,74	0
4	EDO	A	509	4/4	0.89	0.17	44,52,59,65	0
4	EDO	B	507	4/4	0.90	0.13	48,53,53,55	0
4	EDO	D	514	4/4	0.90	0.16	61,62,63,70	0
4	EDO	D	512	4/4	0.90	0.14	40,43,45,45	0

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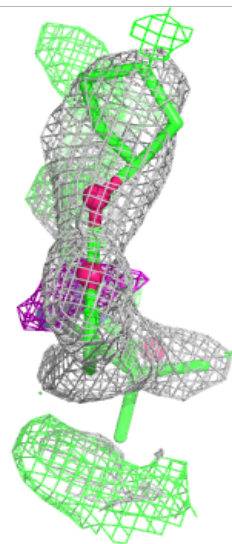
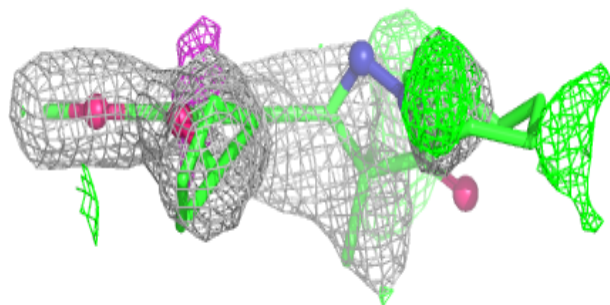
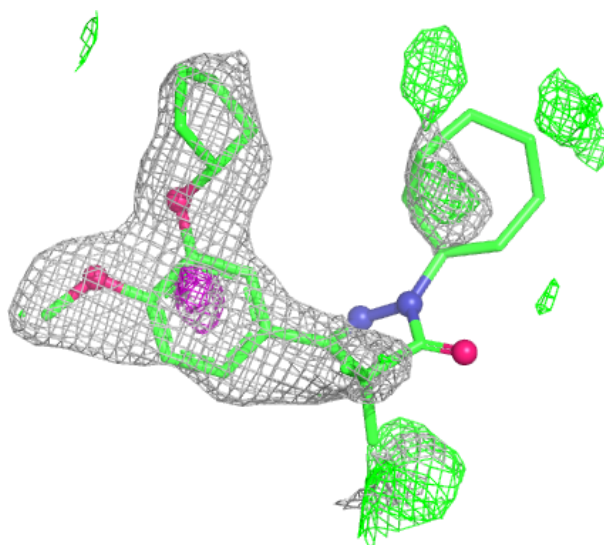
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	EPE	B	515	15/15	0.90	0.16	42,93,107,108	0
4	EDO	A	503	4/4	0.91	0.13	58,60,62,63	0
4	EDO	A	507	4/4	0.91	0.14	52,57,60,61	0
6	EPE	B	512	15/15	0.91	0.13	50,60,63,67	0
4	EDO	A	515	4/4	0.91	0.14	34,36,42,51	0
4	EDO	D	509	4/4	0.92	0.13	35,39,45,49	0
4	EDO	D	506	4/4	0.92	0.13	40,45,46,54	0
4	EDO	D	501	4/4	0.92	0.12	42,43,44,52	0
4	EDO	D	505	4/4	0.93	0.11	53,55,57,58	0
6	EPE	D	519	15/15	0.93	0.15	34,70,83,85	0
4	EDO	C	504	4/4	0.94	0.11	57,58,60,60	0
4	EDO	D	504	4/4	0.94	0.09	33,34,36,37	0
4	EDO	A	510	4/4	0.95	0.09	32,35,36,39	0
4	EDO	D	507	4/4	0.95	0.08	33,35,35,38	0
6	EPE	C	509	15/15	0.97	0.08	37,45,65,76	0
3	MG	C	502	1/1	0.98	0.06	20,20,20,20	0
3	MG	B	502	1/1	0.99	0.07	22,22,22,22	0
2	ZN	A	501	1/1	0.99	0.02	29,29,29,29	0
3	MG	D	503	1/1	0.99	0.06	18,18,18,18	0
2	ZN	D	502	1/1	1.00	0.01	26,26,26,26	0
3	MG	A	502	1/1	1.00	0.04	21,21,21,21	0
2	ZN	B	501	1/1	1.00	0.04	30,30,30,30	0
2	ZN	C	501	1/1	1.00	0.02	30,30,30,30	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

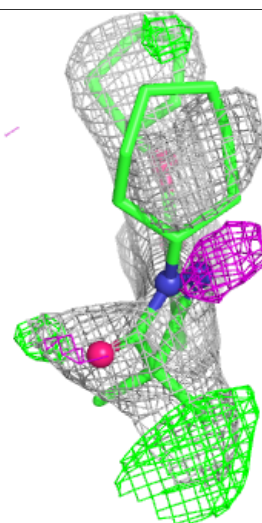
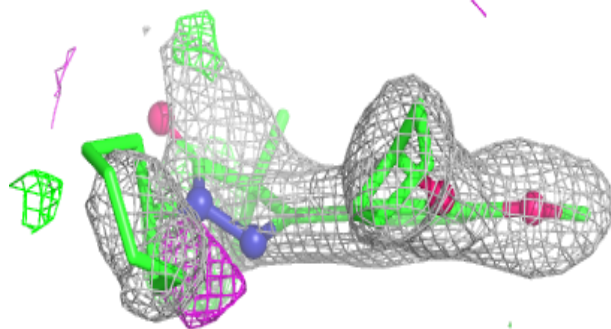
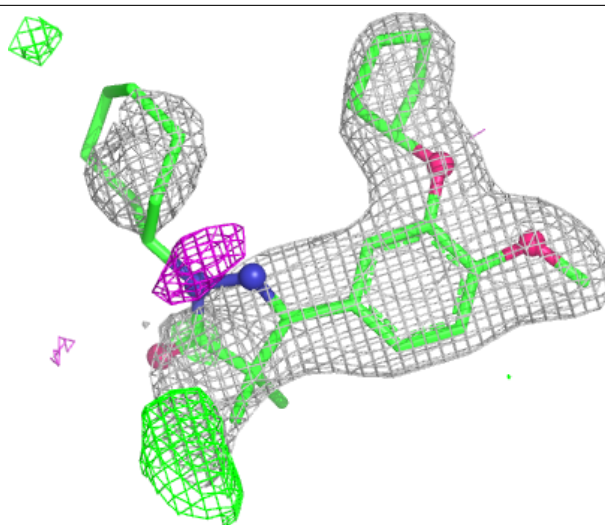
Electron density around J2E A 519:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



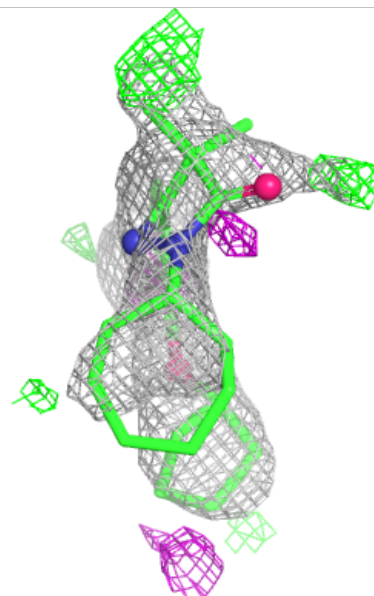
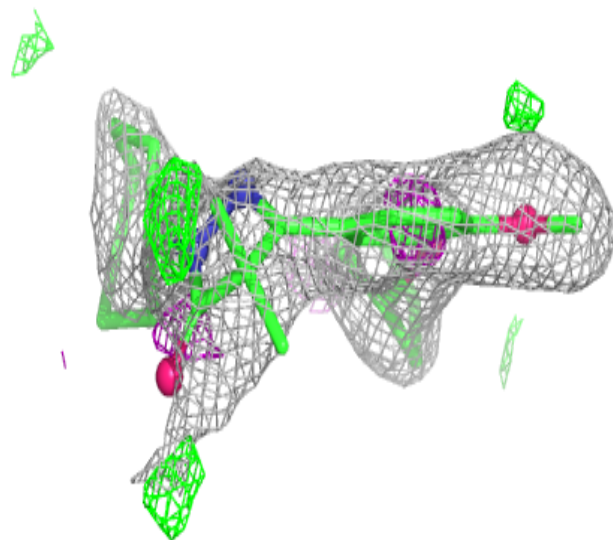
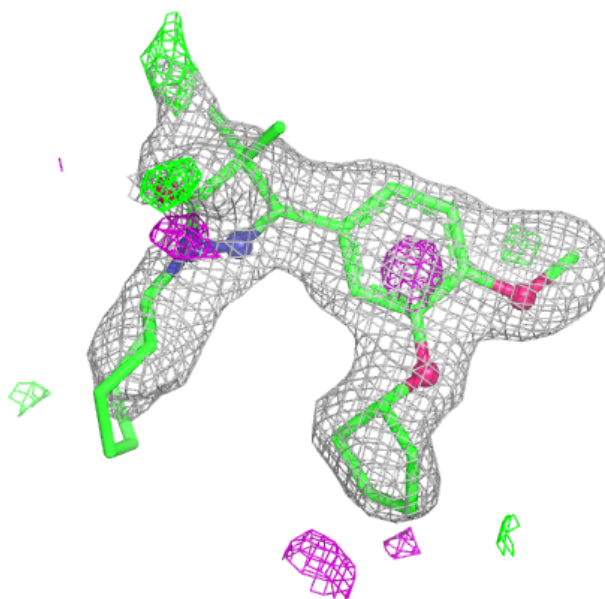
Electron density around J2E C 513:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



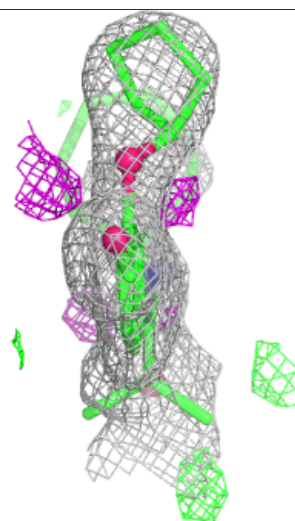
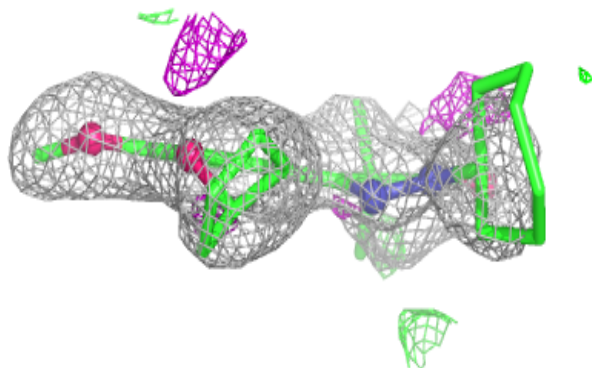
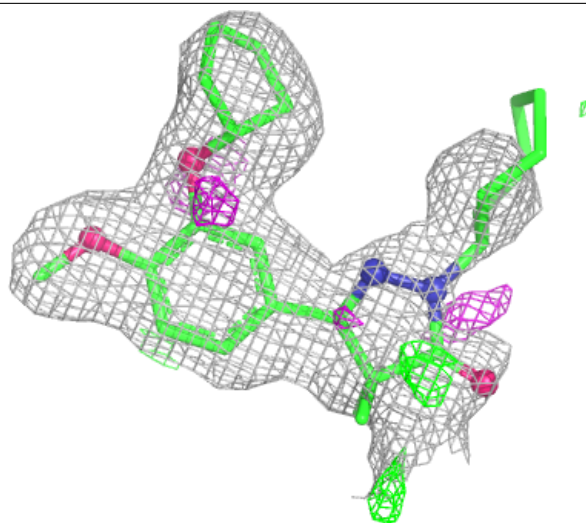
Electron density around J2E D 517:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around J2E B 514:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.